

Concise Machine Learning

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Abstract

This report contains lecture notes for UC Berkeley's introductory class on Machine Learning. It covers many methods for classification and regression, and several methods for clustering and dimensionality reduction. It is concise because not a word is included that cannot be written or spoken in a single semester's lectures (with whiteboard lectures and almost no slides!) and because the choice of topics is limited to a small selection of particularly useful, popular algorithms.

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Keywords: machine learning, classification, regression, density estimation, dimensionality reduction, clustering, perceptrons, support vector machines (SVMs), Gaussian discriminant analysis, linear discriminant analysis (LDA), quadratic discriminant analysis (QDA), logistic regression, decision trees, neural networks, convolutional neural networks (CNNs, ConvNets), nearest neighbor search, least-squares linear regression, logistic regression, polynomial regression, ridge regression, Lasso, maximum likelihood estimation (MLE), principal components analysis (PCA), singular value decomposition (SVD), latent factor analysis, latent semantic indexing, k -means clustering, hierarchical clustering, spectral graph clustering

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About this Report

This report compiles my lectures notes for UC Berkeley’s class CS 189/289A, *Machine Learning*, which is both an undergraduate and introductory graduate course. I hope it will serve as a fast introduction to the subject for readers who are already comfortable with vector calculus, linear algebra, probability, and statistics. Please consult my CS 189/289A web page¹ as an addendum to this report; it includes an extended description of each lecture and additional web links and reading assignments related to the lectures. Consider this report and the web page to be living documents; both will be refined a bit every time I teach the class.

The term “lecture notes” has shifted to include long textbook-style treatments written by professors as supplements to their classes. Not so here. This report compiles the actual notes that I lecture from. I call it *Concise Machine Learning* because I do not include a single word that I do not have time to write or speak during one fourteen-week semester of twice-weekly 80-minute lectures. Words that appear [in brackets] are spoken; everything else is written on the “whiteboard”—in my class, a tablet computer. My whiteboard software permits me to incorporate (and write on) figures, included here. However, I am largely anti-Powerpoint and I resort to prepared slides for just three or four brief segments during the semester.

These notes might be ideal for mathematically sophisticated readers who want to learn the basics of machine learning as quickly as possible. But they’re not ideal for everybody. The time limitation necessitates that many details are omitted. I believe that the smartest, most mathematically well-prepared readers will be able to fill in those details themselves. But many readers, including most students who take the class, will need additional readings or discussion sections for greater detail. My class web page lists additional readings for most of the lectures, many of them from two textbooks that have been kindly made available for free on the web: *An Introduction to Statistical Learning with Applications in R*,² by Gareth James, Daniela Witten, Trevor Hastie, and Robert Tibshirani, Springer, New York, 2013, ISBN # 978-1-4614-7137-0; and *The Elements of Statistical Learning: Data Mining, Inference, and Prediction*,³ second edition, by Trevor Hastie, Robert Tibshirani, and Jerome Friedman, Springer, New York, 2008. Wikipedia also has excellent introductions to many machine learning algorithms. Readers wanting the verbose kind of “lecture notes” should consider the fine ones written by Stanford University’s Andrew Ng.⁴ That web site also includes good primers on linear algebra and probability that can serve as useful reviews before reading this report. I have no interest in duplicating these efforts; instead, I’m aiming for the neglected niche of “shortest introduction.” (And perhaps also “best stolen illustrations.”)

The other thing that makes this report concise is the choice of topics. CS 189/289A was introduced at UC Berkeley in the spring of 2013 by Prof. Jitendra Malik, and most of his topic choices remain intact here. Jitendra told me that he only taught a machine learning algorithm if he or his collaborators had used it successfully for some application. He said, “the machine learning course is too important to leave to the machine learning experts”—that is, users of machine learning algorithms often have a more clear-eyed view of their usefulness than inventors of machine learning algorithms.

I thank Peter Bartlett, Alyosha Efros, Isabelle Guyon, and Jitendra Malik—the previous teachers of CS 189/289A—for their lectures and lecture notes, from which I learned the topic myself. While I’ve given the lectures my own twist and rearranged the material a lot, I am ultimately making incremental improvements (and perhaps incremental worsenings) to a structure they handed down to me. I also thank Carlos Flores for sharing screenshots from my lectures.

¹<http://www.cs.berkeley.edu/~jrs/189/>

²<http://www-bcf.usc.edu/~gareth/ISL/>

³<http://statweb.stanford.edu/~tibs/ElemStatLearn/>

⁴<http://cs229.stanford.edu/materials.html>

1 Introduction

CS 189 / 289A [Spring 2019]

Machine Learning

Jonathan Shewchuk

<http://www.cs.berkeley.edu/~jrs/189>

TAs: Marc Khouri, Carlos Campo, Patrick Chao, Melih Elibol, Panna Felsen, Yang Gao, Noah Golmant, Michael McDonald, Fanyu Meng, Alex Krentsel, Yunqi “Evelyn” Li, Ivon Liu, Vasilis Oikonomou, Keiran Paster, Zipeng Qin, Alan Rosenthal, Faraz Tavakoli, Andy Zhang, Tijana Zrnic.

Questions: Please use Piazza, not email. [Piazza has an option for private questions, but please use public for most questions so other people can benefit.]

For personal matters only, jrs@cory.eecs.berkeley.edu

Discussion sections (22, Tue/Wed):

Attend any section. If room is too full, please go to another one.

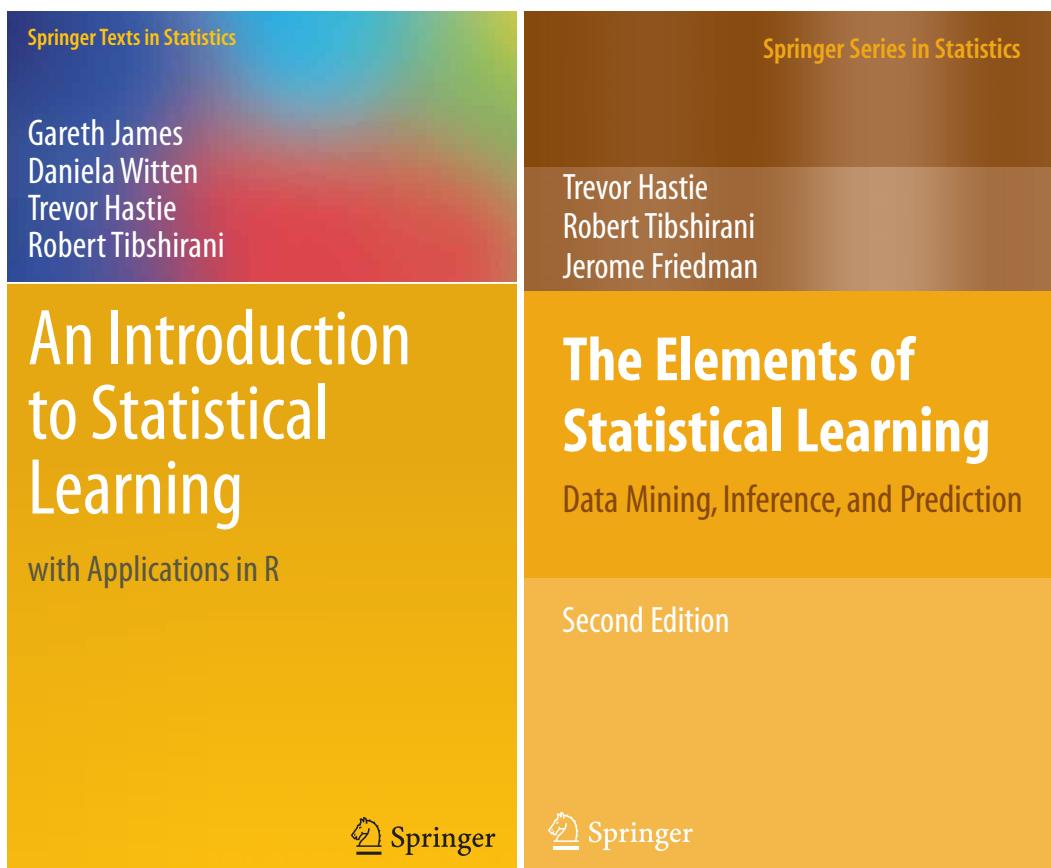
[We hope to have a few advanced sections, including research discussion or exam problem preparation.]

Sections start Tuesday. [Next week.]

Homework 1 (see Piazza) due next Wednesday. HW party this Friday, 5–7 PM, Cory 540AB.

[Enrollment: 744 students max. 383 waitlisted. Expecting many drops. EECS grads have highest priority; undergrads second; non-EECS grads third; little hope for concurrent enrollment students.]

[Textbooks: Available free online. Linked from class web page.]



Prerequisites

Math 53 (vector calculus)

Math 54, Math 110, or EE 16A+16B (linear algebra)

CS 70 or Stat 134 (probability)

NOT CS 188 [Might still be listed as a prerequisite, but we're having it removed.]

Grading: 189

40% 7 Homeworks. Late policy: 5 slip days total

20% Midterm: *Tentatively* Monday, March 18, in class (6:30–8 pm)

40% Final Exam: Friday, May 17, 3–6 PM (Exam group 19)

Grading: 289A

40% HW

20% Midterm

20% Final

20% Project

Cheating

- Discussion of HW problems is encouraged.
- All homeworks, including programming, must be written individually.
- We will actively check for plagiarism.
- Typical penalty is a large NEGATIVE score, but I reserve right to give an instant F for even one violation, and will always give an F for two.

[Last time I taught CS 61B, we had to punish roughly 100 people for cheating. It was very painful. Please don't put me through that again.]

CORE MATERIAL

- Finding patterns in data; using them to make predictions.
- Models and statistics help us understand patterns.
- Optimization algorithms “learn” the patterns.

[The most important part of this is the data. Data drives everything else.

You cannot learn much if you don't have enough data.

You cannot learn much if your data sucks.

But it's amazing what you can do if you have lots of good data.

Machine learning has changed a lot in the last decade because the internet has made truly vast quantities of data available. For instance, with a little patience you can download tens of millions of photographs. Then you can build a 3D model of Paris.

Some techniques that had fallen out of favor, like neural nets, have come back big in recent years because researchers found that they work so much better when you have vast quantities of data.]

CLASSIFICATION

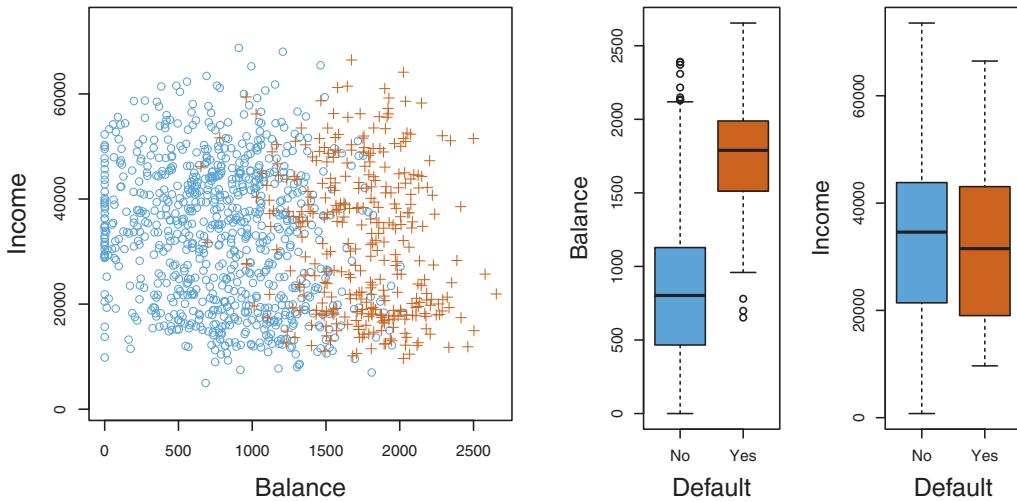
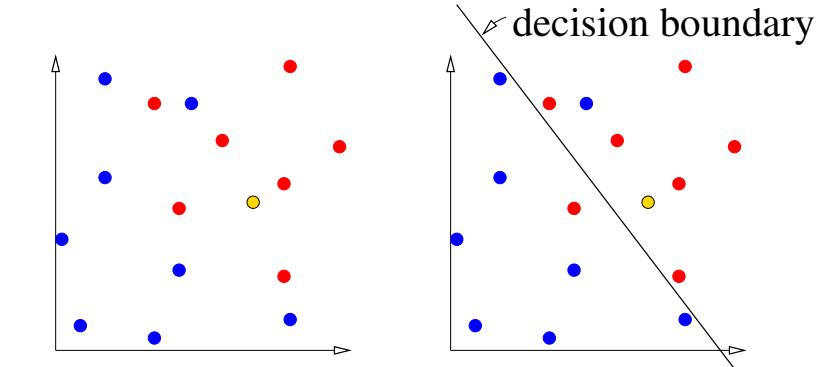


FIGURE 4.1. The `Default` data set. Left: The annual incomes and monthly credit card balances of a number of individuals. The individuals who defaulted on their credit card payments are shown in orange, and those who did not are shown in blue. Center: Boxplots of `balance` as a function of `default` status. Right: Boxplots of `income` as a function of `default` status.

`creditcards.pdf` (ISL, Figure 4.1) [The problem of classification. We are given data points, each belonging to one of two classes. Then we are given additional points whose class is unknown, and we are asked to predict what class each new point is in. Given the credit card balance and annual income of a cardholder, predict whether they will default on their debt.]

- Collect training data: reliable debtors & defaulted debtors
- Evaluate new applicants (prediction)



[Draw this figure by hand. `classify.pdf`]

[Draw 2 colors of dots, almost but not quite linearly separable.]

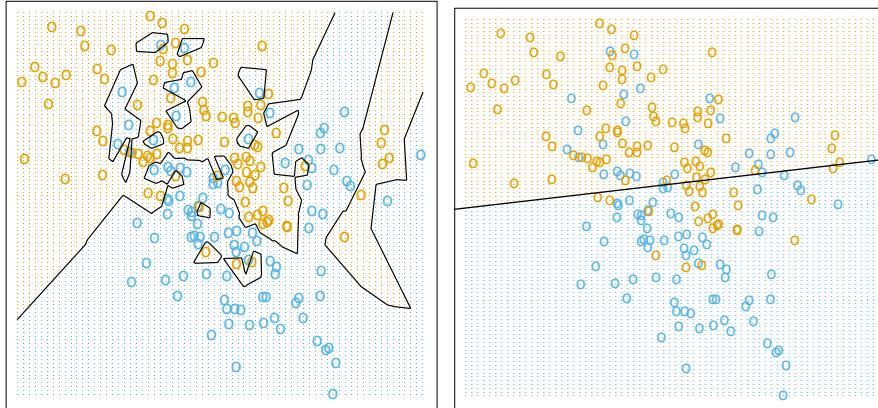
[“How do we classify a new point?” Draw a point in a third color.]

[One possibility: look at its nearest neighbor.]

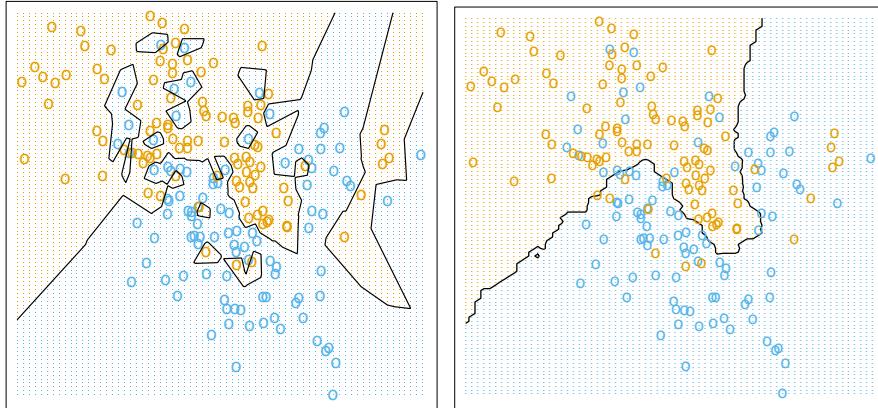
[Another possibility: draw a linear decision boundary; label it.]

[Those are two different *models* for the data.]

[We'll learn some ways to draw these linear decision boundaries in the next several lectures. But for now, let's compare these two methods.]



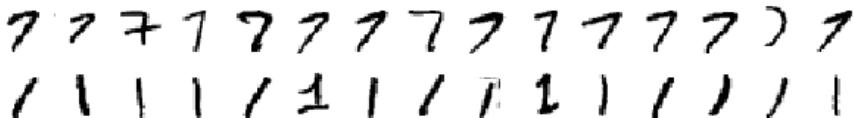
`classnear.pdf, classlinear.pdf` (ESL, Figures 2.3 & 2.1) [Here are two examples of classifiers for the same data. At left we have a nearest neighbor classifier, which classifies a point by finding the nearest point in the input data, and assigning it the same class. At right we have a linear classifier, which guesses that everything above the line is brown, and everything below the line is blue. The decision boundaries are in black.]



`classnear.pdf, classnear15.pdf` (ESL, Figures 2.3 & 2.2) [At right we have a 15-nearest neighbor classifier. Instead of looking at the nearest neighbor of a new point, it looks at the 15 nearest neighbors and lets them vote for the correct class. The 1-nearest neighbor classifier at left has a big advantage: it classifies all the training data correctly, whereas the 15-nearest neighbor classifier at right figure does not. But the right figure has an advantage too. Somebody please tell me what.]

[The left figure is an example of what's called overfitting. In the left figure, observe how intricate the decision boundary is that separates the positive examples from the negative examples. It's a bit too intricate to reflect reality. In the right figure, the decision boundary is smoother. Intuitively, that smoothness is probably more likely to correspond to reality.]

Classifying Digits



`sevensones.pdf` [In this simplified digit recognition problem, we are given handwritten 7's and 1's, and we are asked to learn to distinguish the 7's from the 1's.]

Express these images as vectors

$$\begin{array}{|c|c|c|c|} \hline
 3 & 3 & 3 & 3 \\ \hline
 0 & 0 & 2 & 3 \\ \hline
 0 & 0 & 1 & 3 \\ \hline
 3 & 3 & 3 & 3 \\ \hline
 \end{array} \rightarrow \begin{bmatrix} 3 \\ 3 \\ 3 \\ 3 \\ 3 \\ 0 \\ 0 \\ 2 \\ 3 \\ 0 \\ 0 \\ 1 \\ 3 \\ 3 \\ 3 \\ 3 \\ 3 \\ 3 \end{bmatrix}$$

Images are points in 16-dimensional space. Linear decision boundary is a hyperplane.

Validation

- Train a classifier: it learns to distinguish 7 from not 7
- Test the classifier on NEW images

2 kinds of error:

- Training set error: fraction of training images not classified correctly
[This is zero with the 1-nearest neighbor classifier, but nonzero with the 15-nearest neighbor and linear classifiers we've just seen.]
- Test set error: fraction of misclassified NEW images, not seen during training.

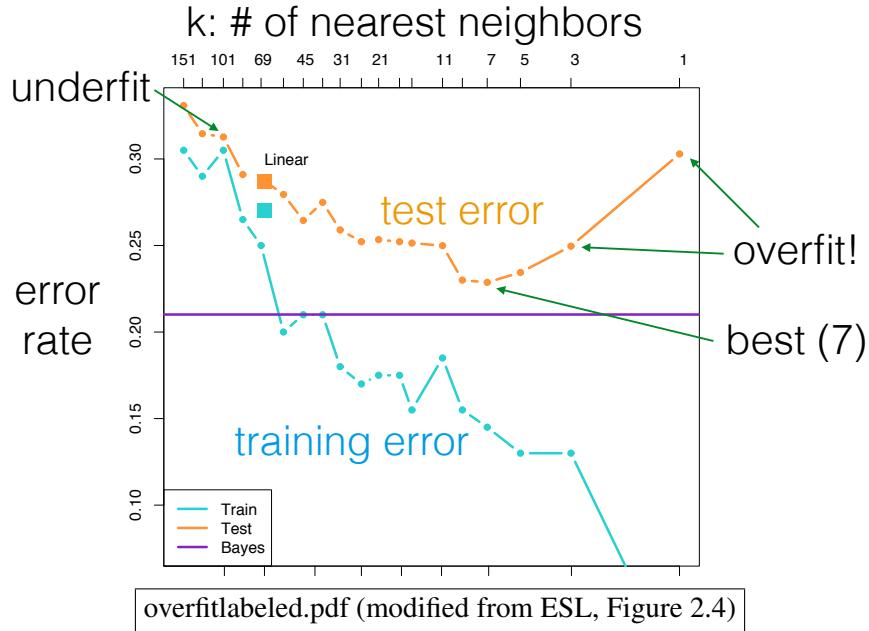
[When I underline a word or phrase, that usually means it's a definition. If you want to do well in this course, my advice to you is to memorize the definitions I cover in class.]

outliers: points whose labels are atypical (e.g. solvent borrower who defaulted anyway).

overfitting: when the test error deteriorates because the classifier becomes too sensitive to outliers or other spurious patterns.

[In machine learning, the goal is to create a classifier that generalizes to new examples we haven't seen yet. Overfitting is counterproductive to that goal. So we're always seeking a compromise: we want decision boundaries that make fine distinctions without being downright superstitious.]

Most ML algorithms have a few hyperparameters that control over/underfitting, e.g. k in k -nearest neighbors.



We select them by validation:

- Hold back a subset of the labeled data, called the validation set.
- Train the classifier multiple times with different hyperparameter settings.
- Choose the settings that work best on validation set.

Now we have 3 sets:

training set used to learn model weights

validation set used to tune hyperparameters, choose among different models

test set used as FINAL evaluation of model. Keep in a vault. Run ONCE, at the very end.

[It's very bad when researchers in medicine or pharmaceuticals peek into the test set prematurely!]

Kaggle.com:

- Runs ML competitions, including our HWS
 - We use 2 data sets:
 - “public” set results available during competition
 - “private” set revealed only after due date
- [If your public results are a lot better than your private results, we will know that you overfitted.]

Techniques [taught in this class, NOT a complete list]

Supervised learning:

- Classification: is this email spam?
- Regression: how likely does this patient have cancer?

Unsupervised learning:

- Clustering: which DNA sequences are similar to each other?
- Dimensionality reduction: what are common features of faces? common differences?

2 Linear Classifiers and Perceptrons

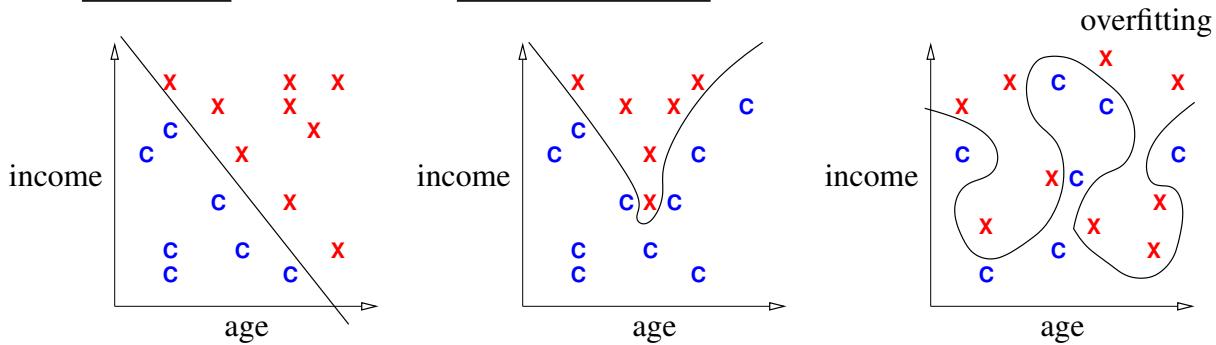
CLASSIFIERS

You are given sample of n observations, each with d features [aka predictors]. Some observations belong to class C; some do not.

Example: Observations are bank loans
 Features are income & age ($d = 2$)
 Some are in class “defaulted”, some are not

Goal: Predict whether future borrowers will default,
 based on their income & age.

Represent each observation as a point in d -dimensional space, called a sample point / a feature vector / independent variables.



[Draw this by hand; decision boundaries last. [classify3.pdf](#)]

[We draw these lines/curves separating C's from X's. Then we use these curves to predict which future borrowers will default. In the last example, though, we're probably overfitting, which could hurt our predictions.]

decision boundary: the boundary chosen by our classifier to separate items in the class from those not.

overfitting: When sinuous decision boundary fits sample points so well that it doesn't classify future points well.

[A reminder that underlined phrase are definitions, worth memorizing.]

Some (not all) classifiers work by computing a

decision function: A function $f(x)$ that maps a sample point x to a scalar such that

$$\begin{array}{ll} f(x) > 0 & \text{if } x \in \text{class C;} \\ f(x) \leq 0 & \text{if } x \notin \text{class C.} \end{array}$$

Aka predictor function or discriminant function.

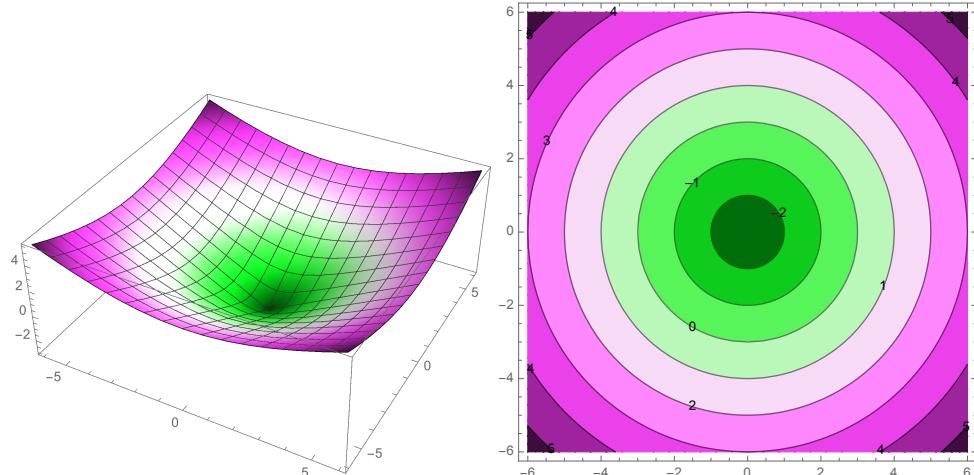
For these classifiers, the decision boundary is $\{x \in \mathbb{R}^d : f(x) = 0\}$

[That is, the set of all points where the decision function is zero.]

Usually, this set is a $(d - 1)$ -dimensional surface in \mathbb{R}^d .

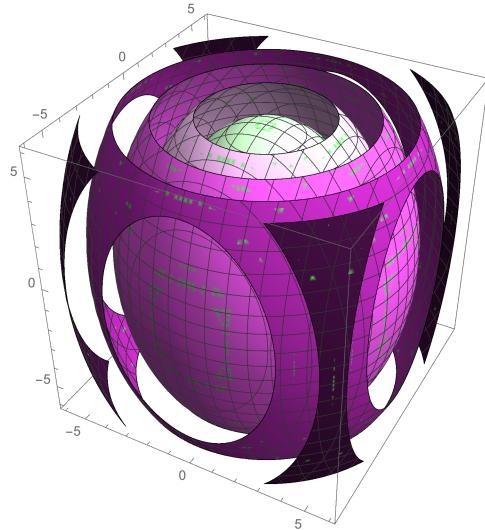
$\{x : f(x) = 0\}$ is also called an isosurface of f for the isovalue 0.

f has other isosurfaces for other isovalues, e.g. $\{x : f(x) = 1\}$.



`radiusplot.pdf, radiusiso.pdf` [Plot and isocontours of the cone] $f(x, y) = \sqrt{x^2 + y^2} - 3$.

[Imagine a decision function in \mathbb{R}^d , and imagine its $(d - 1)$ -dimensional isosurfaces.]



`radiusiso3d.pdf`

[One of these spheres could be the decision boundary.]

linear classifier: The decision boundary is a line/plane.

Usually uses a linear decision function. [Sometimes no decision fn.]

Math Review

[I will write vectors in matrix notation.]

$$\text{Vectors: } \mathbf{x} = \begin{bmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \\ x_5 \end{bmatrix} = [x_1 \ x_2 \ x_3 \ x_4 \ x_5]^T$$

Think of x as a point in 5-dimensional space.

Conventions (often, but not always):

uppercase roman	= matrix, random variable, set	X
lowercase roman	= vector	x
Greek	= scalar	α
Other scalars:		$n = \# \text{ of sample points}$ $d = \# \text{ of features (per point)}$ $= \text{dimension of sample points}$
		$i \ j \ k = \text{indices}$
function (often scalar)		$f(\), s(\), \dots$

inner product (aka dot product): $x \cdot y = x_1y_1 + x_2y_2 + \dots + x_dy_d$

also written $x^T y$

Clearly, $f(x) = w \cdot x + \alpha$ is a linear function in x .

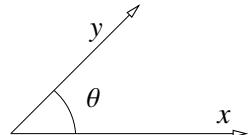
Euclidean norm: $|x| = \sqrt{x \cdot x} = \sqrt{x_1^2 + x_2^2 + \dots + x_d^2}$

$|x|$ is the length (aka Euclidean length) of a vector x .

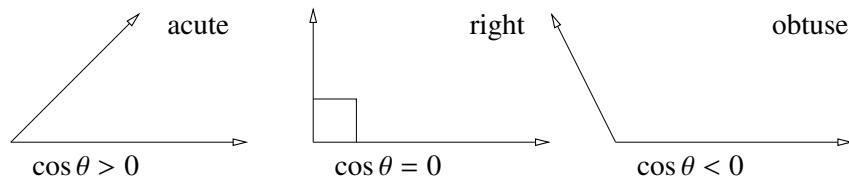
Given a vector x , $\frac{x}{|x|}$ is a unit vector (length 1).

“Normalize a vector x ”: replace x with $\frac{x}{|x|}$.

Use dot products to compute angles:



$$\cos \theta = \frac{x \cdot y}{|x||y|} = \underbrace{\frac{x}{|x|}}_{\text{length 1}} \cdot \underbrace{\frac{y}{|y|}}_{\text{length 1}}$$



Given a linear decision function $f(x) = w \cdot x + \alpha$, the decision boundary is

$$H = \{x : w \cdot x = -\alpha\}.$$

The set H is called a hyperplane. (A line in 2D, a plane in 3D.)

[I want you to understand what a hyperplane is. In 2D, it's a line. In 3D, it's a plane. Now take that concept and generalize it to higher dimensions. In d dimensions, a hyperplane is a flat, infinite thing with dimension $d - 1$. A hyperplane divides the d -dimensional space into two halves.]

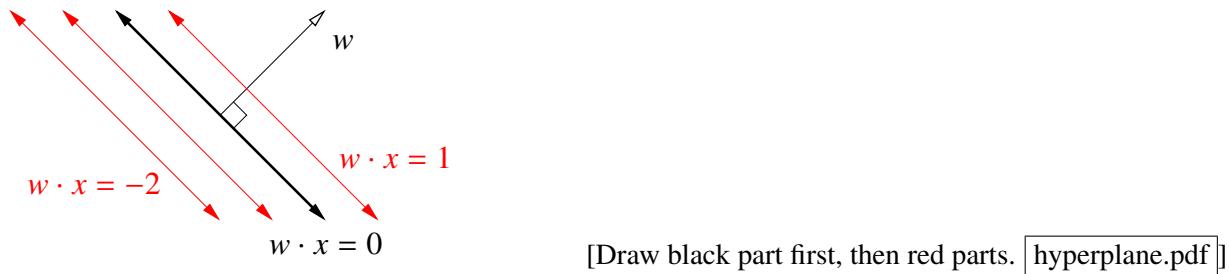
Theorem: Let x, y be 2 points that lie on H . Then $w \cdot (y - x) = 0$.

Proof: $w \cdot (y - x) = -\alpha - (-\alpha) = 0$.

w is called the normal vector of H ,

because (as the theorem shows) w is normal (perpendicular) to H .

[I.e. w is perpendicular to every line through any pair of points in H .]



If w is a unit vector, then $w \cdot x + \alpha$ is the signed distance from x to H .

I.e. positive on one side of H ; negative on other side.

Moreover, the distance from H to the origin is α . [How do we know that?]

Hence $\alpha = 0$ if and only if H passes through origin.

[w does not have to be a unit vector for the classifier to work.]

If w is not a unit vector, $w \cdot x + \alpha$ is the signed distance times some real.

If you want to fix that, you can rescale the equation by computing $|w|$ and dividing both w and α by $|w|$.]

The coefficients in w , plus α , are called weights or sometimes regression coefficients.

[That's why I named the vector w ; "w" stands for "weights."]

The input data is linearly separable if there exists a hyperplane that separates all the sample points in class C from all the points NOT in class C.

[At the beginning of this lecture, I showed you one plot that's linearly separable and two that are not.]

[We will investigate some linear classifiers that only work for linearly separable data, then we'll move on to more sophisticated linear classifiers that do a decent job with non-separable data. Obviously, if your data is not linearly separable, a linear classifier cannot do a *perfect* job. But we're still happy if we can find a classifier that usually predicts correctly.]

A Simple Classifier

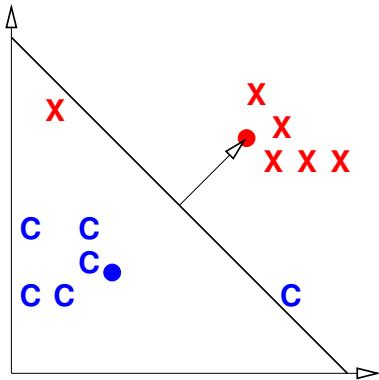
Centroid method: compute mean μ_C of all vectors in class C and mean μ_X of all vectors NOT in C.

We use the decision function

$$f(x) = \underbrace{(\mu_C - \mu_X)}_{\text{normal vector}} \cdot x - (\mu_C - \mu_X) \cdot \underbrace{\frac{\mu_C + \mu_X}{2}}_{\text{midpoint between } \mu_C, \mu_X}$$

so the decision boundary is the hyperplane that bisects line segment w/Endpoints μ_C, μ_X .

[Better yet, we can adjust the scalar term to minimize the number of misclassified points. Then the hyperplane has the same normal vector, but a different position.]



[Draw data, then μ_C, μ_X , then line & normal. [centroid.pdf](#)]

[In this example, there's clearly a better linear classifier that classifies every sample point correctly. Note that this is hardly the worst example I could have given. If you're in the mood for an easy puzzle, pull out a sheet of paper and think of an example, with lots of sample points, where the centroid method misclassifies every sample point but one.]

[Nevertheless, there are cases where this method works well, like when all your positive examples come from one Gaussian distribution, and all your negative examples come from another.]

Perceptron Algorithm (Frank Rosenblatt, 1957)

Slow, but correct for linearly separable points.

Uses a numerical optimization algorithm, namely, gradient descent.

[Poll:

How many of you know what gradient descent is?

How many of you know what a linear program is?

How many of you know what the simplex algorithm for linear programming is?

How many of you know what a quadratic program is?

We're going to learn what most of these things are. As machine learning people, we will be heavy users of optimization methods. Unfortunately, I won't have time to teach you *algorithms* for many optimization problems, but we'll learn a few. To learn more, take EECS 127.]

Consider n sample points X_1, X_2, \dots, X_n .

[The reason I'm using capital X here is because we typically store these vectors as rows of a matrix X . So the subscript picks out a row of X , representing a specific sample point.]

For each sample point, the label $y_i = \begin{cases} 1 & \text{if } X_i \in \text{class C, and} \\ -1 & \text{if } X_i \notin \text{C.} \end{cases}$

For simplicity, consider only decision boundaries that pass through the origin. (We'll fix this later.)

Goal: find weights w such that

$$\begin{aligned} X_i \cdot w &\geq 0 && \text{if } y_i = 1, \text{ and} \\ X_i \cdot w &\leq 0 && \text{if } y_i = -1. \end{aligned} \quad [\text{remember, } X_i \cdot w \text{ is the signed distance}]$$

Equivalently: $y_i X_i \cdot w \geq 0$. \leftarrow inequality called a constraint.

Idea: We define a risk function R that is positive if some constraints are violated. Then we use optimization to choose w that minimizes R .

Define the loss function

$$L(z, y_i) = \begin{cases} 0 & \text{if } y_i z \geq 0, \text{ and} \\ -y_i z & \text{otherwise.} \end{cases}$$

[Here, z is the classifier's prediction, and y_i is the correct answer.]

Idea: if z has the same sign as y_i , the loss function is zero (happiness).

But if z has the wrong sign, the loss function is positive.

[For each sample point, you want to get the loss function down to zero, or as close to zero as possible. It's called the "loss function" because the bigger it is, the bigger a loser you are.]

Define risk function (aka objective function or cost function)

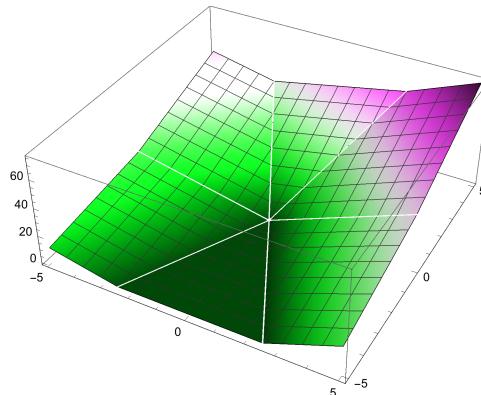
$$\begin{aligned} R(w) &= \sum_{i=1}^n L(X_i \cdot w, y_i), \\ &= \sum_{i \in V} -y_i X_i \cdot w \end{aligned} \quad \text{where } V \text{ is the set of indices } i \text{ for which } y_i X_i \cdot w < 0.$$

If w classifies all X_1, \dots, X_n correctly, then $R(w) = 0$.

Otherwise, $R(w)$ is positive, and we want to find a better value of w .

Goal: Solve this optimization problem:

Find w that minimizes $R(w)$.



riskplot.pdf [Plot of risk $R(w)$. Every point in the dark green flat spot is a minimum. We'll look at this more next lecture.]

3 Perceptron Learning; Maximum Margin Classifiers

Perceptron Algorithm (cont'd)

Recall:

- linear decision fn $f(x) = w \cdot x$ (for simplicity, no α)
- decision boundary $\{x : f(x) = 0\}$ (a hyperplane through the origin)
- sample points $X_1, X_2, \dots, X_n \in \mathbb{R}^d$; classifications $y_1, \dots, y_n = \pm 1$
- goal: find weights w such that $y_i X_i \cdot w \geq 0$
- goal, rewritten: find w that minimizes $R(w) = \sum_{i \in V} -y_i X_i \cdot w$ [risk function]
where V is the set of indices i for which $y_i X_i \cdot w < 0$.

[Our original problem was to find a separating hyperplane in one space, which I'll call x -space. But we've transformed this into a problem of finding an optimal point in a different space, which I'll call w -space. It's important to understand transformations like this, where a geometric structure in one space becomes a point in another space.]

Objects in x -space transform to objects in w -space:

x -space	w -space
hyperplane: $\{z : w \cdot z = 0\}$	point: w
point: x	hyperplane: $\{z : x \cdot z = 0\}$

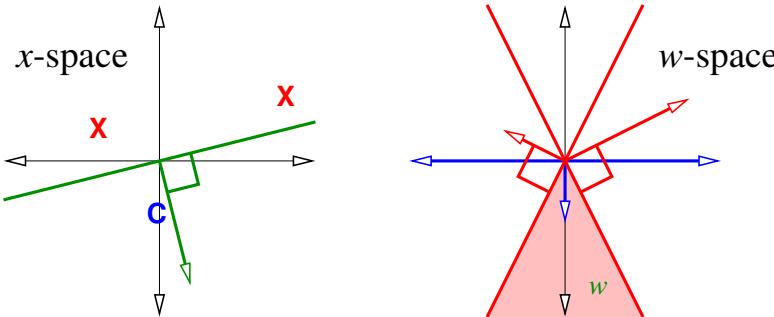
Point x lies on hyperplane $\{z : w \cdot z = 0\} \Leftrightarrow w \cdot x = 0 \Leftrightarrow$ point w lies on hyperplane $\{z : x \cdot z = 0\}$ in w -space.

[So a hyperplane transforms to its normal vector. And a sample point transforms to the hyperplane whose normal vector is the sample point.]

[In this case, the transformations happen to be symmetric: a hyperplane in x -space transforms to a point in w -space the same way that a hyperplane in w -space transforms to a point in x -space. That won't always be true for the weight spaces we use this semester.]

If we want to enforce inequality $x \cdot w \geq 0$, that means

- in x -space, x should be on the same side of $\{z : z \cdot w = 0\}$ as w
- in w -space, w " " " " " " " " $\{z : x \cdot z = 0\}$ as x

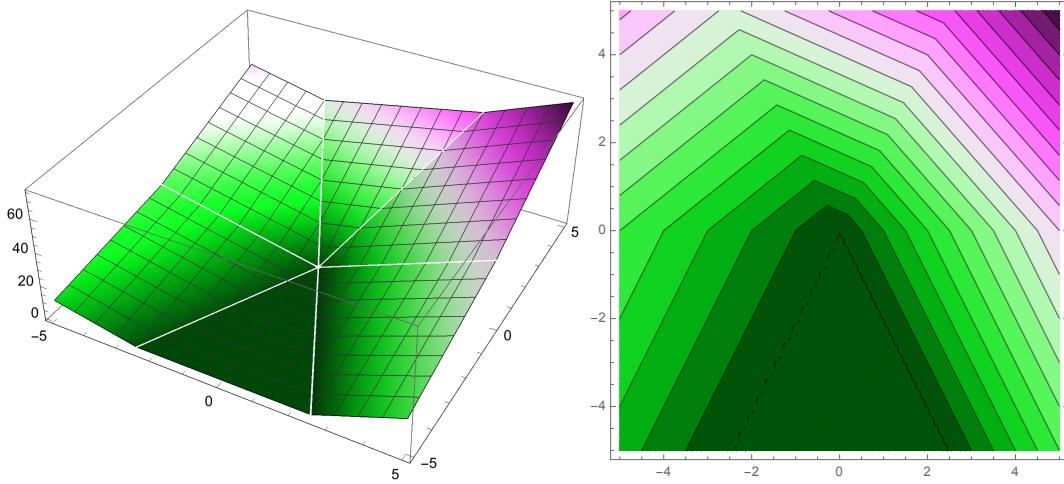


[Draw this by hand. [xwspace.pdf](#)]
[Observe that the x -space sample points are the normal vectors for the w -space lines. We can choose w to be anywhere in the shaded region.]

[For a sample point x in class C, w and x must be on the *same* side of the hyperplane that x transforms into. For a point x not in class C (marked by an X), w and x must be on *opposite* sides of the hyperplane that x transforms into. These rules determine the shaded region above, in which w must lie.]

[Again, what have we accomplished? We have switched from the problem of finding a hyperplane in x -space to the problem of finding a point in w -space. That's a much better fit to how we think about optimization algorithms.]

[Let's take a look at the risk function these three sample points create.]



`riskplot.pdf, riskiso.pdf` [Plot & isocontours of risk $R(w)$. Note how R 's creases match the dual chart above.]

[In this plot, we can choose w to be any point in the bottom pizza slice; all those points minimize R .]

[We have an optimization problem; we need an optimization algorithm to solve it.]

An optimization algorithm: gradient descent on R .

Given a starting point w , find gradient of R with respect to w ; this is the direction of steepest ascent. Take a step in the opposite direction. Recall [from your vector calculus class]

$$\nabla R(w) = \begin{bmatrix} \frac{\partial R}{\partial w_1} \\ \frac{\partial R}{\partial w_2} \\ \vdots \\ \frac{\partial R}{\partial w_d} \end{bmatrix} \quad \text{and} \quad \nabla(z \cdot w) = \begin{bmatrix} z_1 \\ z_2 \\ \vdots \\ z_d \end{bmatrix} = z$$

$$\nabla R(w) = \sum_{i \in V} \nabla -y_i X_i \cdot w = - \sum_{i \in V} y_i X_i$$

At any point w , we walk downhill in direction of steepest descent, $-\nabla R(w)$.

```

 $w \leftarrow$  arbitrary nonzero starting point (good choice is any  $y_i X_i$ )
while  $R(w) > 0$ 
   $V \leftarrow$  set of indices  $i$  for which  $y_i X_i \cdot w < 0$ 
   $w \leftarrow w + \epsilon \sum_{i \in V} y_i X_i$ 
return  $w$ 

```

$\epsilon > 0$ is the step size aka learning rate, chosen empirically. [Best choice depends on input problem!]

[Show plot of R again. Draw the typical steps of gradient descent.]

Problem: Slow! Each step takes $O(nd)$ time. [Can we improve this?]

Optimization algorithm 2: stochastic gradient descent

Idea: each step, pick **one** misclassified X_i ;
do gradient descent on loss fn $L(X_i \cdot w, y_i)$.

Called the perceptron algorithm. Each step takes $O(d)$ time.

[Not counting the time to search for a misclassified X_i .]

```
while some  $y_i X_i \cdot w < 0$ 
     $w \leftarrow w + \epsilon y_i X_i$ 
return  $w$ 
```

[Stochastic gradient descent is quite popular and we'll see it several times more this semester, especially for neural nets. However, stochastic gradient descent does not work for every problem that gradient descent works for. The perceptron risk function happens to have special properties that guarantee that stochastic gradient descent will always succeed.]

What if separating hyperplane doesn't pass through origin?

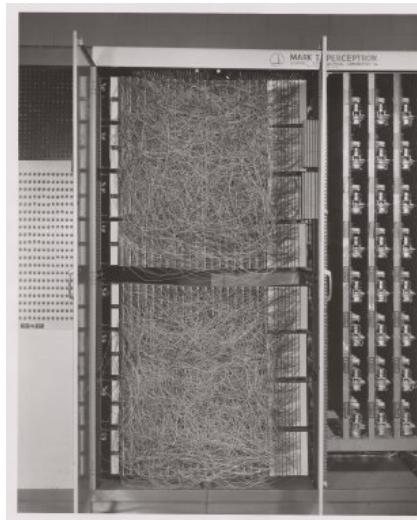
Add a fictitious dimension. Decision fn is

$$f(x) = w \cdot x + \alpha = [w_1 \ w_2 \ \alpha] \cdot \begin{bmatrix} x_1 \\ x_2 \\ 1 \end{bmatrix}$$

Now we have sample points in \mathbb{R}^{d+1} , all lying on hyperplane $x_{d+1} = 1$.

Run perceptron algorithm in $(d + 1)$ -dimensional space. [We are simulating a general hyperplane in d dimensions by using a hyperplane through the origin in $d + 1$ dimensions.]

[The perceptron algorithm was invented in 1957 by Frank Rosenblatt at the Cornell Aeronautical Laboratory. It was originally designed not to be a program, but to be implemented in hardware for image recognition on a 20×20 pixel image. Rosenblatt built a Mark I Perceptron Machine that ran the algorithm, complete with electric motors to do weight updates.]



Mark_I_perceptron.jpg (from Wikipedia, “Perceptron”) [The Mark I Perceptron Machine.
This is what it took to process a 20×20 image in 1957.]

[Then he held a press conference where he predicted that perceptrons would be “the embryo of an electronic computer that [the Navy] expects will be able to walk, talk, see, write, reproduce itself and be conscious of its existence.” We’re still waiting on that.]

[One interesting aspect of the perceptron algorithm is that it’s an “online algorithm,” which means that if new data points come in while the algorithm is already running, you can just throw them into the mix and keep looping.]

Perceptron Convergence Theorem: If data is linearly separable, perceptron algorithm will find a linear classifier that classifies all data correctly in at most $O(R^2/\gamma^2)$ iterations, where $R = \max |X_i|$ is “radius of data” and γ is the “maximum margin.”

[I’ll define “maximum margin” shortly.]

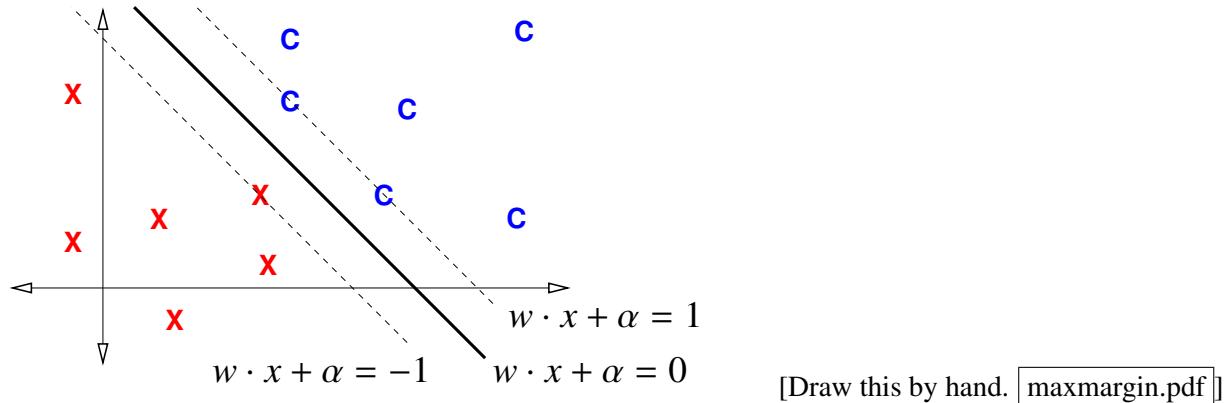
[We’re not going to prove this, because perceptrons are obsolete.]

[Although the step size/learning rate doesn’t appear in that big-O expression, it does have an effect on the running time, but the effect is hard to characterize. The algorithm gets slower if ϵ is too small because it has to take lots of steps to get down the hill. But it also gets slower if ϵ is too big for a different reason: it jumps right over the region with zero risk and oscillates back and forth for a long time.]

[Although stochastic gradient descent is faster for this problem than gradient descent, the perceptron algorithm is still slow. There’s no reliable way to choose a good step size ϵ . Fortunately, optimization algorithms have improved a lot since 1957. You can get rid of the step size by using any decent modern “line search” algorithm. Better yet, you can find a better decision boundary much more quickly by quadratic programming, which is what we’ll talk about next.]

MAXIMUM MARGIN CLASSIFIERS

The margin of a linear classifier is the distance from the decision boundary to the nearest sample point. What if we make the margin as wide as possible?



We enforce the constraints

$$y_i (w \cdot X_i + \alpha) \geq 1 \quad \text{for } i \in [1, n]$$

[Notice that the right-hand side is a 1, rather than a 0 as it was for the perceptron algorithm. It’s not obvious, but this a better way to formulate the problem, partly because it makes it impossible for the weight vector w to get set to zero.]

Recall: if $|w| = 1$, signed distance from hyperplane to X_i is $w \cdot X_i + \alpha$.

Otherwise, it's $\frac{w}{|w|} \cdot X_i + \frac{\alpha}{|w|}$. [We've normalized the expression to get a unit weight vector.]

Hence the margin is $\min_i \frac{1}{|w|} |w \cdot X_i + \alpha| \geq \frac{1}{|w|}$. [We get the inequality by substituting the constraints.]

There is a slab of width $\frac{2}{|w|}$ containing no sample points [with the hyperplane running along its middle].

To maximize the margin, minimize $|w|$. Optimization problem:

$$\begin{aligned} &\text{Find } w \text{ and } \alpha \text{ that minimize } |w|^2 \\ &\text{subject to } y_i(X_i \cdot w + \alpha) \geq 1 \quad \text{for all } i \in [1, n] \end{aligned}$$

Called a quadratic program in $d + 1$ dimensions and n constraints.

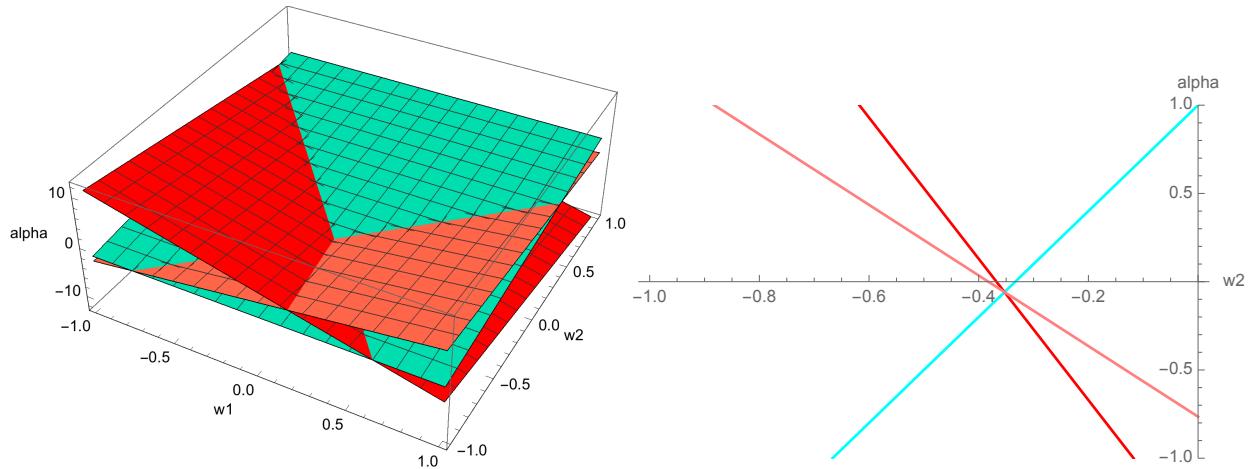
It has one unique solution! [If the points are linearly separable; otherwise, it has no solution.]

[A reason we use $|w|^2$ as an objective function, instead of $|w|$, is that the length function $|w|$ is not smooth at zero, whereas $|w|^2$ is smooth everywhere. This makes optimization easier.]

The solution gives us a maximum margin classifier, aka a hard margin support vector machine (SVM).

[Technically, this isn't really a support vector machine yet; it doesn't fully deserve that name until we add features and kernels, which we'll do in later lectures.]

[Let's see what these constraints look like in weight space.]



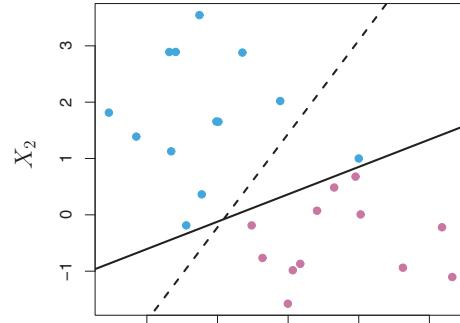
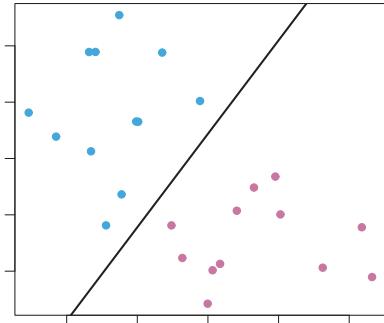
[weight3d.pdf](#), [weightcross.pdf](#) [This is an example of what the linear constraints look like in the 3D weight space (w_1, w_2, α) for an SVM with three training points. The SVM is looking for the point nearest the origin that lies above the blue plane (representing an in-class training point) but below the red and pink planes (representing out-of-class training points). In this example, that optimal point lies where the three planes intersect. At right we see a 2D cross-section $w_1 = 1/17$ of the 3D space, because the optimal solution lies in this cross-section. The constraints say that the solution must lie in the leftmost pizza slice, while being as close to the origin as possible, so the optimal solution is where the three lines meet.]

4 Soft-Margin Support Vector Machines; Features

SOFT-MARGIN SUPPORT VECTOR MACHINES (SVMs)

Solves 2 problems:

- Hard-margin SVMs fail if data not linearly separable.
- “ ” ” sensitive to outliers.



sensitive.pdf (ISL, Figure 9.5) [Example where one outlier moves the decision boundary a lot.]

Idea: Allow some points to violate the margin, with slack variables.

Modified constraint for point i :

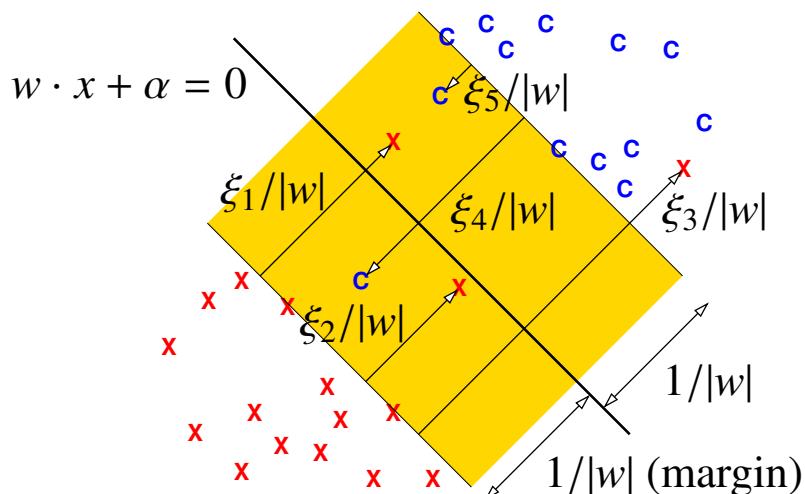
$$y_i(X_i \cdot w + \alpha) \geq 1 - \xi_i$$

[Observe that the only difference between these constraints and the hard margin constraints we saw last lecture is the extra slack term ξ_i .]

[We also impose new constraints, that the slack variables are never negative.]

$$\xi_i \geq 0$$

[This inequality ensures that all sample points that *don't* violate the margin are treated the same; they all have $\xi_i = 0$. Point i has nonzero ξ_i if and only if it violates the margin.]



slacker+.pdf [A margin where some points have slack.]

To prevent abuse of slack, we add a loss term to objective fn.

Optimization problem:

Find w , α , and ξ_i that minimize $ w ^2 + C \sum_{i=1}^n \xi_i$
subject to
$y_i(X_i \cdot w + \alpha) \geq 1 - \xi_i \quad \text{for all } i \in [1, n]$
$\xi_i \geq 0 \quad \text{for all } i \in [1, n]$

...a quadratic program in $d + n + 1$ dimensions and $2n$ constraints.

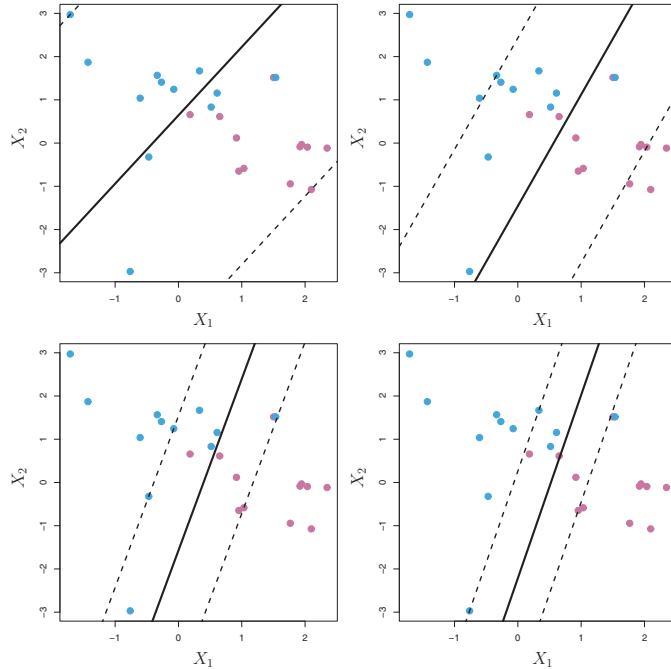
[It's a quadratic program because its objective function is quadratic and its constraints are linear inequalities.]

$C > 0$ is a scalar regularization hyperparameter that trades off:

	small C	big C
desire	maximize margin $1/ w $	keep most slack variables zero or small
danger	underfitting (misclassifies much training data)	overfitting (awesome training, awful test)
outliers	less sensitive	very sensitive
boundary	more “flat”	more sinuous

[The last row only applies to nonlinear decision boundaries, which we'll discuss next. Obviously, a linear decision boundary can't be “sinuous.”]

Use validation to choose C .



svmC.pdf (ISL, Figure 9.7) [Examples of how the slab varies with C . Smallest C at upper left; largest C at lower right.]

[One way to think about slack is to pretend that slack is money we can spend to buy permission for a sample point to violate the margin. The further a point penetrates the margin, the bigger the fine you have to pay. We want to make the margin as wide as possible, but we also want to spend as little money as possible. If the regularization parameter C is small, it means we're willing to spend lots of money on violations so we can get a wider margin. If C is big, it means we're cheap and we won't pay much for violations, even though we'll suffer a narrower margin. If C is infinite, we're back to a hard-margin SVM.]

FEATURES

Q: How to do nonlinear decision boundaries?

A: Make nonlinear features that lift points into a higher-dimensional space.

High- d linear classifier \rightarrow low- d nonlinear classifier.

[Features work with all classifiers—not only linear classifiers like perceptrons and SVMs, but also classifiers that are not linear.]

Example 1: The parabolic lifting map

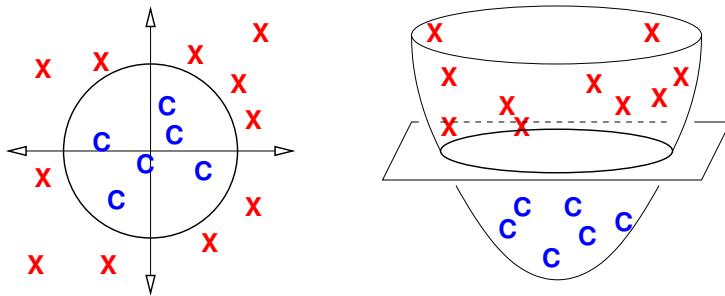
$$\Phi : \mathbb{R}^d \rightarrow \mathbb{R}^{d+1}$$

$$\Phi(x) = \begin{bmatrix} x \\ |x|^2 \end{bmatrix} \quad \leftarrow \text{lifts } x \text{ onto paraboloid } x_{d+1} = |x|^2$$

[We've added one new feature, $|x|^2$. Even though the new feature is just a function of other input features, it gives our linear classifier more power.]

Find a linear classifier in Φ -space.

It induces a sphere classifier in x -space.



[Draw this by hand. [circledec.pdf](#)]

Theorem: $\Phi(X_1), \dots, \Phi(X_n)$ are linearly separable iff X_1, \dots, X_n are separable by a hypersphere.

(Possibly an ∞ -radius hypersphere = hyperplane.)

Proof: Consider hypersphere in \mathbb{R}^d w/center c & radius ρ . Points inside:

$$\begin{aligned} |x - c|^2 &< \rho^2 \\ |x|^2 - 2c \cdot x + |c|^2 &< \rho^2 \\ \underbrace{[-2c^\top 1]}_{\text{normal vector in } \mathbb{R}^{d+1}} \underbrace{\begin{bmatrix} x \\ |x|^2 \end{bmatrix}}_{\Phi(x)} &< \rho^2 - |c|^2 \end{aligned}$$

Hence points inside sphere \rightarrow same side of hyperplane in Φ -space.

(Reverse implication works too.)

[Hyperspheres include hyperplanes as a special, degenerate case. A hyperplane is essentially a hypersphere with infinite radius. So hypersphere decision boundaries can do everything hyperplane decision boundaries can do, plus a lot more. With the parabolic lifting map, if you pick a hyperplane in Φ -space that is vertical, you get a hyperplane in x -space.]

Example 2: Axis-aligned ellipsoid/hyperboloid decision boundaries

[Draw examples of axis-aligned ellipse & hyperbola.]

In 3D, these have the formula

$$Ax_1^2 + Bx_2^2 + Cx_3^2 + Dx_1 + Ex_2 + Fx_3 = -G$$

[Here, the capital letters are scalars, not matrices.]

$$\Phi : \mathbb{R}^d \rightarrow \mathbb{R}^{2d}$$

$$\Phi(x) = [x_1^2 \quad \dots \quad x_d^2 \quad x_1 \quad \dots \quad x_d]^\top$$

$$\text{Hyperplane is } \underbrace{(A, B, C, D, E, F)}_w \cdot \Phi(x) = -G$$

[We've turned d input features into $2d$ features for our linear classifier. If the points are separable by an axis-aligned ellipsoid or hyperboloid, per the formula above, then the points lifted to Φ -space are separable by a hyperplane whose normal vector is (A, B, C, D, E, F) .]

Example 3: Ellipsoid/hyperboloid

[Draw example of non-axis-aligned ellipse.]

3D formula: [for a general ellipsoid or hyperboloid]

$$Ax_1^2 + Bx_2^2 + Cx_3^2 + Dx_1x_2 + Ex_2x_3 + Fx_3x_1 + Gx_1 + Hx_2 + Ix_3 = -J$$

$$\Phi(x) : \mathbb{R}^d \rightarrow \mathbb{R}^{(d^2+3d)/2}$$

[Now, our decision function can be any degree-2 polynomial.]

Isosurface defined by this equation is called a quadric. [In the special case of two dimensions, it's also known as a conic section. So our decision boundary can be an arbitrary conic section.]

[You'll notice that there is a quadratic blowup in the number of features, because every *pair* of input features creates a new feature in Φ -space. If the dimension is large, these feature vectors are getting huge, and that's going to impose a serious computational cost. But it might be worth it to find good classifiers for data that aren't linearly separable.]

Example 4: Decision fn is degree- p polynomial

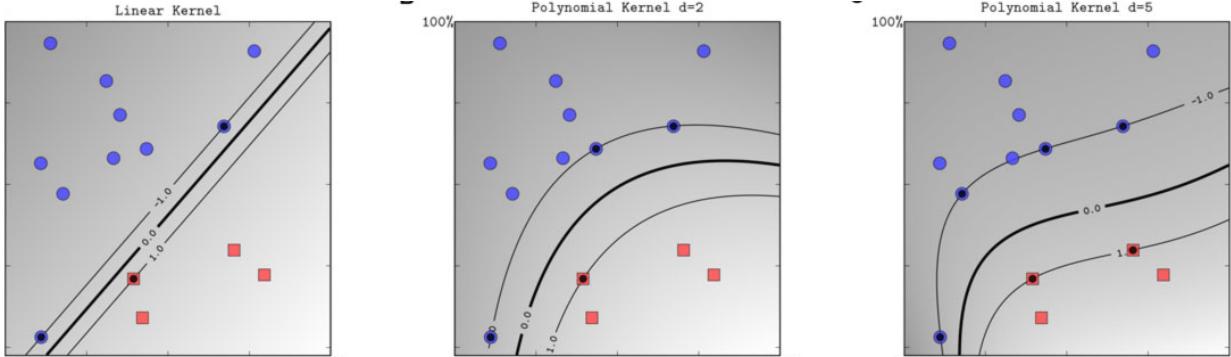
E.g. a cubic in \mathbb{R}^2 :

$$\Phi(x) = [x_1^3 \quad x_1^2 x_2 \quad x_1 x_2^2 \quad x_2^3 \quad x_1^2 \quad x_1 x_2 \quad x_2^2 \quad x_1 \quad x_2]^\top$$

$$\Phi(x) : \mathbb{R}^d \rightarrow \mathbb{R}^{O(d^p)}$$

[Now we're really blowing up the number of features! If you have, say, 100 features per sample point and you want to use degree-4 decision functions, then each lifted feature vector has a length of roughly 4 million, and your learning algorithm will take approximately forever to run.]

[However, later in the semester we will learn an extremely clever trick that allows us to work with these huge feature vectors very quickly, without ever computing them. It's called "kernelization" or "the kernel trick." So even though it appears now that working with degree-4 polynomials is computationally infeasible, it can actually be done quickly.]

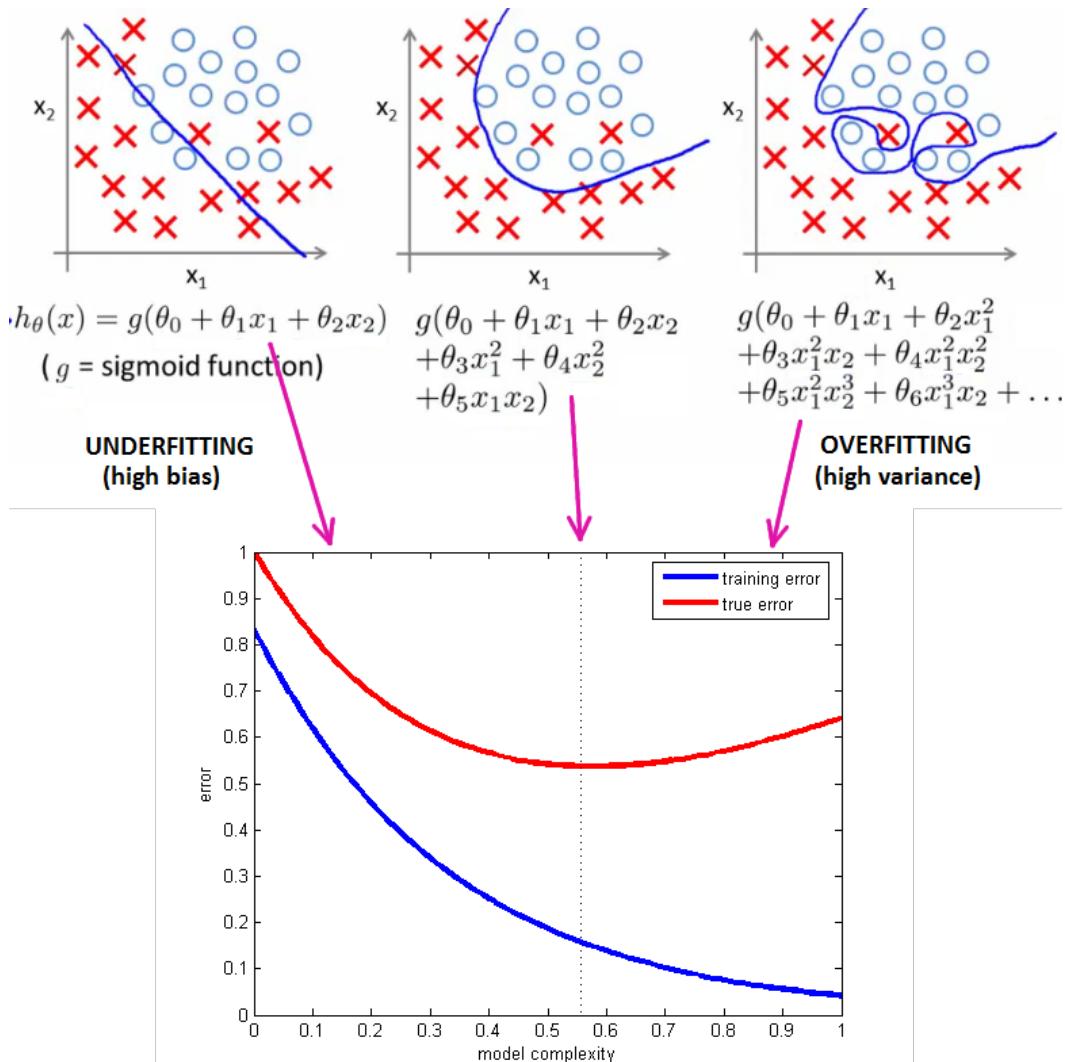


degree5.pdf [SVMs with degree 1/2/5 decision functions. Observe that the margin tends to get wider as the degree increases.]

[Increasing the degree like this accomplishes two things.

- First, the data might become linearly separable when you lift them to a high enough degree, even if the original data are not linearly separable.
- Second, raising the degree can increase the margin, so you might get a more robust decision boundary that generalizes better.

However, if you raise the degree too high, you will overfit the data and then generalization will get worse.]



[overfit.pdf](#) [Training vs. test error for degree 1/2/5 decision functions. (Artist's conception; these aren't actual calculations, just hand-drawn guesses. Please send me email if you know where to find figures like this with actual data.) In this example, a degree-2 decision gives the smallest test error.]

[Sometimes you should search for the ideal degree—not too small, not too big. It's a balancing act between underfitting and overfitting. The degree is an example of a *hyperparameter* that can be optimized by validation.]

[If you're using both polynomial features and a soft-margin SVM, now you have two hyperparameters: the degree and the regularization hyperparameter C . Generally, the optimal C will be different for every polynomial degree, so when you change the degree, you have to run validation again to find the best C for that degree.]

[So far I've talked only about polynomial features. But features can get much more complicated than polynomials, and they can be tailored to fit a specific problem. Let's consider a type of feature you might use if you wanted to implement, say, a handwriting recognition algorithm.]

Example 5: Edge detection

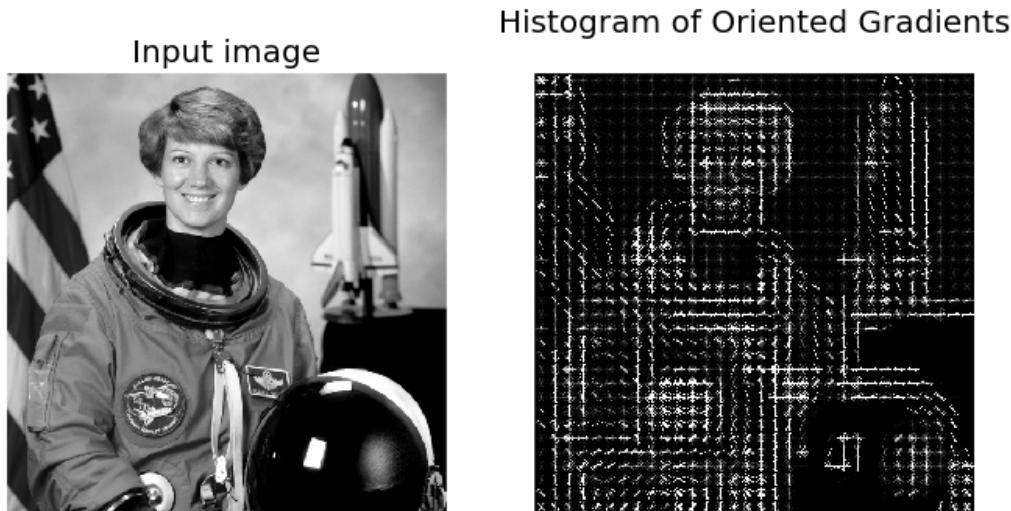
Edge detector: algorithm for approximating grayscale/color gradients in image, e.g.,

- tap filter
- Sobel filter
- oriented Gaussian derivative filter

[images are discrete, not continuous fields, so approximation of gradients is necessary.]

[See “Image Derivatives” on Wikipedia.]

Collect line orientations in local histograms (each having 12 orientation bins per region); use histograms as features (*instead* of raw pixels).



orientgrad.png [Image histograms.]

Paper: Maji & Malik, 2009.

[If you want to, optionally, use these features in future homeworks and try to win the Kaggle competition, this paper is a good online resource.]

[When they use a linear SVM on the raw pixels, Maji & Malik get an error rate of 15.38% on the test set. When they use a linear SVM on the histogram features, the error rate goes down to 2.64%.]

[Many applications can be improved by designing application-specific features. There’s no limit but your own creativity and ability to discern the structure hidden in your application.]

5 Machine Learning Abstractions and Numerical Optimization

ML ABSTRACTIONS [some meta comments on machine learning]

[When you write a large computer program, you break it down into subroutines and modules. Many of you know from experience that you need to have the discipline to impose strong abstraction barriers between different modules, or your program will become so complex you can no longer manage nor maintain it.]

[When you learn a new subject, it helps to have mental abstraction barriers, too, so you know when you can replace one approach with a different approach. I want to give you four levels of abstraction that can help you think about machine learning. It's important to make mental distinctions between these four things, and the code you write should have modules that reflect these distinctions as well.]

APPLICATION/DATA	
data labeled (classified) or not?	
yes: labels categorical (classification) or quantitative (regression)?	
no: similarity (clustering) or positioning (dimensionality reduction)?	
MODEL	[what kinds of hypotheses are permitted?]
e.g.:	
– decision fns: linear, polynomial, logistic, neural net, ...	
– nearest neighbors, decision trees	
– features	
– low vs. high capacity (affects overfitting, underfitting, inference)	
OPTIMIZATION PROBLEM	
– variables, objective fn, constraints	
e.g., unconstrained, convex program, least squares, PCA	
OPTIMIZATION ALGORITHM	
e.g., gradient descent, simplex, SVD	

[In this course, we focus primarily on the middle two levels. As a data scientist, you might be given an application, and your challenge is to turn it into an optimization problem that we know how to solve. We will talk a bit about optimization algorithms, but usually you'll use an optimization code that's faster and more robust than what you would write yourself.]

[The second level, the model, has a huge effect on the success of your learning algorithm. Sometimes you get a big improvement by tailoring the model or its features to fit the structure of your specific data. The model also has a big effect on whether you overfit or underfit. And if you want a model that you can interpret so you can do *inference*, the model has to have a simple structure. Lastly, you have to pick a model that leads to an optimization problem that can be solved. Some optimization problems are just too hard.]

[It's important to understand that when you change something in one level of this diagram, you probably have to change all the levels underneath it. If you switch your model from a linear classifier to a neural net, your optimization problem changes, and your optimization algorithm probably changes too.]

[Not all machine learning methods fit this four-level decomposition. Nevertheless, for everything you learn in this class, think about where it fits in this hierarchy. If you don't distinguish which math is part of the model and which math is part of the optimization algorithm, this course will be very confusing for you.]

OPTIMIZATION PROBLEMS

[I want to familiarize you with some types of optimization problems that can be solved reliably and efficiently, and the names of some of the optimization algorithms used to solve them. An important skill for you to develop is to be able to go from an application to a well-defined optimization problem. That skill depends on your ability to *recognize* well-studied types of optimization problems.]

Unconstrained

Goal: Find w that minimizes (or maximizes) a continuous objective fn $f(w)$.

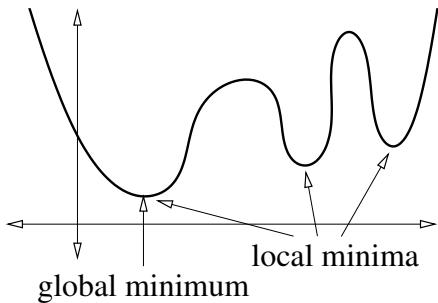
f is smooth if its gradient is continuous too.

A global minimum of f is a value w such that $f(w) \leq f(v)$ for every v .

A local minimum " " " " " " " "

for every v in a tiny ball centered at w .

[In other words, you cannot walk downhill from w .]

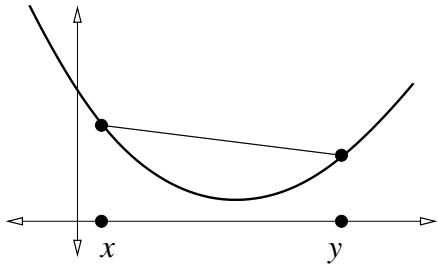


[Draw this by hand. [minima.pdf](#)]

Usually, finding a local minimum is easy;

finding the global minimum is hard. [or impossible]

Exception: A function is convex if for every $x, y \in \mathbb{R}^d$,
the line connecting $(x, f(x))$ to $(y, f(y))$ does not go below $f(\cdot)$.



[Draw this by hand. [convex.pdf](#)]

E.g. perceptron risk fn is convex and nonsmooth.

[When you sum together convex functions, you always get a convex function. The perceptron risk function is a sum of convex loss functions.]

A [continuous] convex function [on a closed, convex domain] has either

- no minimum (goes to $-\infty$), or
- just one local minimum, or
- a connected set of local minima that are all global minima with equal f .

[The perceptron risk function has the last of these three.]

[In the last two cases, if you walk downhill, you eventually converge to a global minimum.]

[However, there are many applications where you don't have a convex objective function, and your machine learning algorithm has to settle for finding a local minimum. For example, neural nets try to optimize an objective function that has *lots* of local minima; they almost never find a global minimum.]

Algs for smooth f :

- Gradient descent:
 - blind [with learning rate]
 - with line search:
 - secant method
 - Newton–Raphson (may need Hessian matrix of f)
 - stochastic (blind) [trains on one point per iteration, or a small batch]
- Nonlinear conjugate gradient [uses the secant or Newton–Raphson line search methods]
- Newton's method (needs Hessian matrix)

Algs for nonsmooth f :

- Gradient descent:
 - blind
 - with direct line search (e.g. golden section search)

These algs find a local minimum. [They don't reliably find a global minimum, because that's very hard.]

line search: finds a local minimum along the search direction by solving an optimization problem in 1D. [... instead of using a blind step size like the perceptron algorithm does. Solving a 1D problem is much easier than solving a higher-dimensional one.]

[Neural nets are unconstrained optimization problems with many, many local minima. They sometimes benefit from line search or second-order optimization algorithms, but when the input data set is very large, researchers often favor the dumb, blind, stochastic versions of gradient descent.]

[If you're optimizing over a d -dimensional space, the Hessian matrix is a $d \times d$ matrix and it's usually dense, so most methods that use the Hessian are computationally infeasible when d is large.]

Constrained Optimization (smooth equality constraints)

Goal: Find w that minimizes (maximizes) $f(w)$

subject to $g(w) = 0$

[\leftarrow observe that this is an isosurface]

where g is a smooth fn

[g may be a vector, encoding multiple constraints]

Alg: Use Lagrange multipliers.

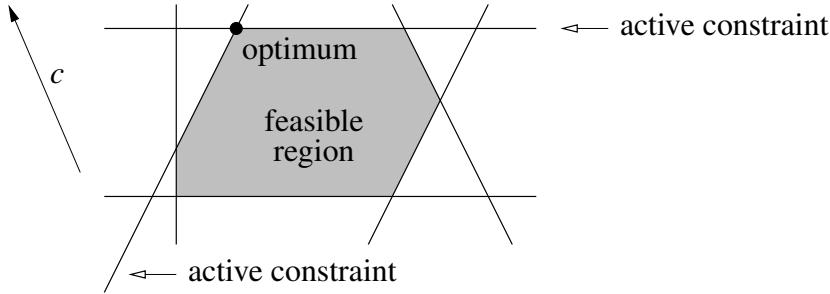
[to transform constrained to unconstrained optimization]

Linear Program

Linear objective fn + linear **inequality** constraints.

Goal: Find w that maximizes (or minimizes) $c \cdot w$
subject to $Aw \leq b$

where A is $n \times d$ matrix, $b \in \mathbb{R}^n$, expressing n linear constraints:
 $A_i w \leq b_i, \quad i \in [1, n]$



[Draw this by hand. [linprog.pdf](#)]

The set of points that satisfy all constraints is a convex polytope called the feasible region F [shaded].

The optimum is the point in F that is furthest in the direction c . [What does convex mean?]

A point set P is convex if for every $p, q \in P$, the line segment with endpoints p, q lies entirely in P .

[A polytope is just a polyhedron, generalized to higher dimensions.]

The optimum achieves equality for some constraints (but not most), called the active constraints of the optimum. [In the figure above, there are two active constraints. In an SVM, active constraints correspond to the sample points that touch or violate the slab, and they're also known as support vectors.]

[Sometimes, there is more than one optimal point. For example, in the figure above, if c pointed straight up, every point on the top horizontal edge would be optimal. The set of optimal points is always convex.]

Example: EVERY feasible point (w, α) gives a linear classifier:

Find w, α that maximizes 0
subject to $y_i(w \cdot X_i + \alpha) \geq 1 \quad \text{for all } i \in [1, n]$

IMPORTANT: The data are linearly separable iff the feasible region is not the empty set.

→ Also true for maximum margin classifier (quadratic program)

Algs for linear programming:

- Simplex (George Dantzig, 1947)

[Indisputably one of the most important and useful algorithms of the 20th century.]

[Walks along edges of polytope from vertex to vertex until it finds optimum.]

- Interior point methods

[Linear programming is very different from unconstrained optimization; it has a much more combinatorial flavor. If you knew which constraints would be the active constraints once you found the solution, it would be easy; the hard part is figuring out which constraints should be the active ones. There are exponentially many possibilities, so you can't afford to try them all. So linear programming algorithms tend to have a very discrete, computer science feeling to them, like graph algorithms, whereas unconstrained optimization algorithms tend to have a continuous, numerical mathematics feeling.]

[Linear programs crop up *everywhere* in engineering and science, but they're usually in disguise. An extremely useful talent you should develop is to recognize when a problem is a linear program.]

[A linear program solver can find a linear classifier, but it can't find the maximum margin classifier. We need something more powerful.]

Quadratic Program

Quadratic, convex objective fn + linear inequality constraints.

Goal: Find w that minimizes $f(w) = w^\top Qw + c^\top w$
subject to $Aw \leq b$

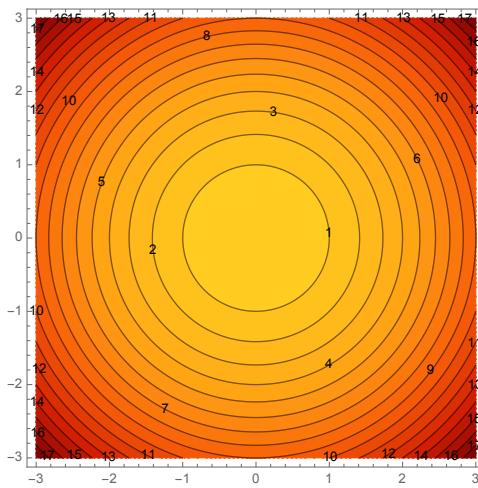
where Q is a symmetric, positive definite matrix.

A matrix is positive definite if $w^\top Qw > 0$ for all $w \neq 0$.

Only one local minimum! [Which is therefore the global minimum.]

[What if Q is not positive definite? If Q is indefinite, then f is not convex, the minimum is not always unique, and quadratic programming is NP-hard.]

Example: Find maximum margin classifier.



quadratic.pdf [Draw two polygons on these isocontours—one with one active constraint, and one with two—and show the constrained minimum for each polygon. “In an SVM, we are looking for the point in this polygon that’s closest to the origin.”]

Algs for quadratic programming:

- Simplex-like [commonly used for general-purpose quadratic programs, but not as good for SVMs as the following two algorithms that specifically exploit properties of SVMs]
- Sequential minimal optimization (SMO, used in LIBSVM)
- Coordinate descent (used in LIBLINEAR)

[One clever idea SMO uses is that it does a line search that uses the Hessian, but it's cheap to compute because SMO doesn't walk in the direction of steepest descent; instead it walks along just two coordinate axes at a time.]

Convex Program

Convex objective fn + convex inequality constraints.

[What I've given you here is, roughly, a sliding scale of optimization problems of increasing complexity, difficulty, and computation time. But even convex programs are relatively easy to solve. When you're trying to address the needs of real-world applications, it's not uncommon to devise an optimization problem with crazy inequalities and an objective function that's nowhere near convex. These are sometimes very, very hard to solve.]

Numerical optimization @ Berkeley: EECS 127/227AT/227BT/227C.

6 Decision Theory; Generative and Discriminative Models

DECISION THEORY aka Risk Minimization

[Today I'm going to talk about a style of classifier very different from SVMs. The classifiers we'll cover in the next few weeks are based on probability.]

[One aspect of probabilistic data is that sometimes a point in feature space doesn't have just one class. Suppose one borrower with income \$30,000 and debt \$15,000 defaults.

another " " " " " " " doesn't default.

So in your feature space, you have two feature vectors at the same point with different classes. Obviously, in that case, you can't draw a decision boundary that classifies all points with 100% accuracy.]

Multiple sample points with different classes could lie at same point:
we want a probabilistic classifier.

Suppose 10% of population has cancer, 90% doesn't.

Probability distributions for calorie intake, $P(X|Y)$: [caps here mean random variables, not matrices.]

calories	(X)	< 1,200	1,200–1,600	> 1,600
cancer	($Y = 1$)	20%	50%	30%
no cancer	($Y = -1$)	1%	10%	89%

[I made these numbers up. Please don't take them as medical advice.]

Recall: $P(X) = P(X|Y = 1)P(Y = 1) + P(X|Y = -1)P(Y = -1)$

$$P(1,200 \leq X \leq 1,600) = 0.5 \times 0.1 + 0.1 \times 0.9 = 0.14$$

You meet guy eating $x = 1,400$ calories/day. Guess whether he has cancer?

[If you're in a hurry, you might see that 50% of people with cancer eat 1,400 calories, but only 10% of people with no cancer do, and conclude that someone who eats 1,400 calories probably has cancer. But that would be wrong, because that reasoning fails to take the prior probabilities into account.]

Bayes' Theorem:

$$\begin{aligned} &\downarrow \text{posterior probability} \quad \downarrow \text{prior prob.} \quad \downarrow \text{for } 1,200 \leq X \leq 1,600 \\ P(Y = 1|X) &= \frac{P(X|Y = 1)P(Y = 1)}{P(X)} = \frac{0.05}{0.14} \\ P(Y = -1|X) &= \frac{P(X|Y = -1)P(Y = -1)}{P(X)} = \frac{0.09}{0.14} \end{aligned} \quad [\text{These two probs always sum to 1.}]$$

$$P(\text{cancer} | 1,200 \leq X \leq 1,600 \text{ cals}) = 5/14 \approx 36\%.$$

[So we probably shouldn't diagnose cancer.]

[BUT ... we're assuming that we want to maximize the chance of a correct prediction. But that's not always the right assumption. If you're developing a cheap screening test for cancer, you'd rather have more false positives and fewer false negatives. A false negative might mean somebody misses an early diagnosis and dies of a cancer that could have been treated if caught early. A false positive just means that you spend more money on more accurate tests.]

A loss function $L(z, y)$ specifies badness if classifier predicts z , true class is y .

$$\text{E.g., } L(z, y) = \begin{cases} 1 & \text{if } z = 1, y = -1 \quad \text{false positive is bad} \\ 5 & \text{if } z = -1, y = 1 \quad \text{false negative is BAAAAAAD} \\ 0 & \text{if } z = y \end{cases}$$

A 36% probability of loss 5 is worse than a 64% prob. of loss 1,
so we recommend further cancer screening.

Defs: loss fn above is asymmetrical.

The 0-1 loss function is 1 for incorrect predictions, [symmetrical]
0 for correct.

[Another example where you want a very asymmetrical loss function is for spam detection. Putting a good email in the spam folder is much worse than putting spam in your inbox.]

Let $r : \mathbb{R}^d \rightarrow \pm 1$ be a decision rule, aka classifier:
a fn that maps a feature vector x to 1 (“in class”) or -1 (“not in class”).

The risk for r is the expected loss over all values of x, y :

$$\begin{aligned} R(r) &= \mathbb{E}[L(r(X), Y)] \\ &= \sum_x \left(L(r(x), 1) P(Y = 1|X = x) + L(r(x), -1) P(Y = -1|X = x) \right) P(X = x) \\ &= P(Y = 1) \sum_x L(r(x), 1) P(X = x|Y = 1) + P(Y = -1) \sum_x L(r(x), -1) P(X = x|Y = -1) \end{aligned}$$

The Bayes decision rule aka Bayes classifier is the fn r^* that minimizes functional $R(r)$.

Assuming $L(z, y) = 0$ for $z = y$:

$$r^*(x) = \begin{cases} 1 & \text{if } L(-1, 1) P(Y = 1|X = x) > L(1, -1) P(Y = -1|X = x), \\ -1 & \text{otherwise} \end{cases}$$

When L is symmetric, [the big, key principle you should be remembering is]
pick the class with the biggest posterior probability.

[But if the loss function is asymmetric, then you must weight the posteriors with the losses.]

In cancer example, $r^*(x) = 1$ for $x \leq 1,600$; $r^*(x) = -1$ for $x > 1,600$.

The Bayes risk, aka optimal risk, is the risk of the Bayes classifier.

[In our cancer example, the last expression for risk R gives:]

$$R(r^*) = 0.1(5 \times 0.3) + 0.9(1 \times 0.01 + 1 \times 0.1) = 0.249$$

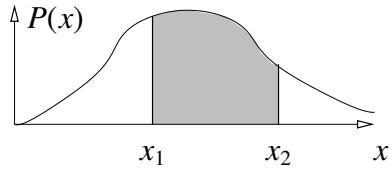
[It is interesting that, if we really know all these probabilities, we really can construct an ideal probabilistic classifier. But in real applications, we rarely know these probabilities; the best we can do is use statistical methods to estimate them.]

Deriving/using r^* is called risk minimization.

Continuous Distributions

Suppose X has a continuous probability density fn (PDF).

Review: [Go back to your CS 70 or stats notes if you don't remember this.]



[Draw this by hand. [integrate.pdf](#)]

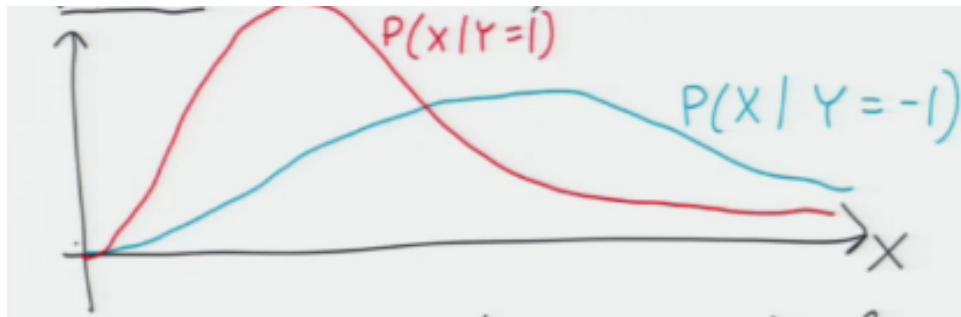
$$\text{prob. that random variable } X \in [x_1, x_2] = \int_{x_1}^{x_2} P(x) dx \quad [\text{shaded area}]$$

$$\text{area under whole curve} = 1 = \int_{-\infty}^{\infty} P(x) dx \quad \text{expected value of } f(x) : E[f(X)] = \int_{-\infty}^{\infty} f(x)P(x) dx$$

$$\text{mean } \mu = E[X] = \int_{-\infty}^{\infty} x P(x) dx$$

$$\text{variance } \sigma^2 = E[(X - \mu)^2] = E[X^2] - \mu^2$$

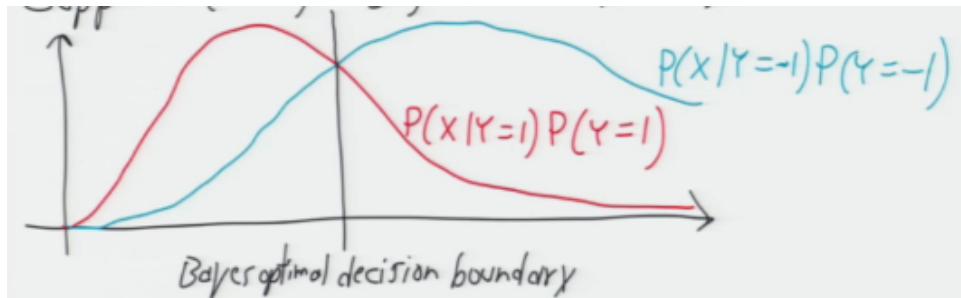
[Perhaps our cancer statistics look like this:]



[Draw this figure by hand (cancerconditional.png)] [The area under each curve is 1.]

[Let's go back to the 0-1 loss function for a moment. In other words, suppose you want a classifier that maximizes the chance of a correct prediction. The wrong answer would be to look where these two curves cross and make that be the decision boundary. As before, it's wrong because it doesn't take into account the prior probabilities.]

Suppose $P(Y = 1) = 1/3$, $P(Y = -1) = 2/3$, 0-1 loss:



[Draw this figure by hand (cancerposterior.png)]

[To maximize the chance you'll predict correctly whether somebody has cancer, the Bayes decision rule looks up x on this chart and picks the curve with the highest probability. In this example, that means you pick cancer when x is left of the optimal decision boundary, and no cancer when x is to the right.]

Define risk as before, replacing summations with integrals.

$$\begin{aligned} R(r) &= \mathbb{E}[L(r(X), Y)] \\ &= P(Y = 1) \int L(r(x), 1) P(X = x|Y = 1) dx + \\ &\quad P(Y = -1) \int L(r(x), -1) P(X = x|Y = -1) dx \end{aligned}$$

For Bayes decision rule, Bayes Risk is the area under minimum of functions above. [Shade it.]

Assuming $L(z, y) = 0$ for $z = y$:

$$R(r^*) = \int \min_{y=\pm 1} L(-y, y) P(X = x|Y = y) P(Y = y) dx$$

[If you want to use an asymmetrical loss function, just scale the curves vertically in the figure above.]

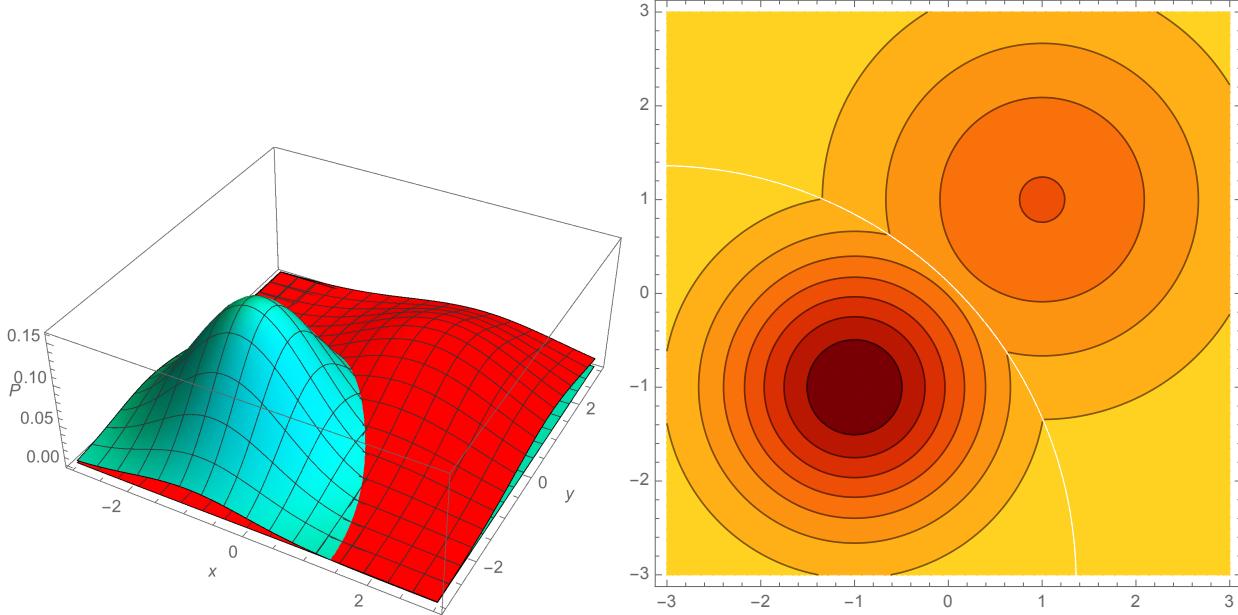
If L is 0-1 loss,

$$R(r) = P(r(x) \text{ is wrong})$$

and the Bayes optimal decision boundary is $\{x : \underbrace{P(Y = 1|X = x)}_{\text{decision fn}} = \underbrace{0.5}_{\text{iso value}}\}$

[then the risk has a particularly nice interpretation:]

[which makes sense, because R is the expected loss.]



[Notice that the accuracy of the probabilities is most important near the decision boundary. Far away from the decision boundary, a bit of error in the probabilities probably wouldn't change the classification.]

[You can also have multi-class classifiers, choosing among three or more classes. The Bayesian approach is a particularly convenient way to generate multi-class classifiers, because you can simply choose whichever class has the greatest posterior probability. Then the decision boundary lies wherever two or more classes are tied for the highest probability.]

3 WAYS TO BUILD CLASSIFIERS

- (1) Generative models (e.g. LDA) [We'll learn about LDA next lecture.]
 - Assume sample points come from probability distributions, different for each class.
 - Guess form of distributions
 - For each class C, fit distribution parameters to class C points, giving $P(X|Y = C)$
 - For each C, estimate $P(Y = C)$
 - Bayes' Theorem gives $P(Y|X)$
 - If 0-1 loss, pick class C that maximizes $P(Y = C|X = x)$ [posterior probability]
equivalently, maximizes $P(X = x|Y = C) P(Y = C)$
- (2) Discriminative models (e.g. logistic regression) [We'll learn about logistic regression in a few weeks.]
 - Model $P(Y|X)$ directly
- (3) Find decision boundary (e.g. SVM)
 - Model $r(x)$ directly (no posterior)

Advantage of (1 & 2): $P(Y|X)$ tells you probability your guess is wrong

[This is something SVMs don't do.]

Advantage of (1): you can diagnose outliers: $P(X)$ is very small

Disadvantages of (1): often hard to estimate distributions accurately;
real distributions rarely match standard ones.

[What I've written here doesn't actually define the phrases "generative model" or "discriminative model." The proper definitions accord with the way statisticians think about models. A generative model is a full probabilistic model of all variables, whereas a discriminative model provides a model only for the target variables that we want to predict.]

[It's important to remember that we rarely know precisely the value of any of these probabilities. There is usually error in all of these probabilities. In practice, generative models are most popular when you have phenomena that are well approximated by the normal distribution, and you have a lot of sample points, so you can approximate the shape of the distribution well.]

7 Gaussian Discriminant Analysis, including QDA and LDA

GAUSSIAN DISCRIMINANT ANALYSIS

Fundamental assumption: each class comes from normal distribution (Gaussian).

$$X \sim \mathcal{N}(\mu, \sigma^2) : P(x) = \frac{1}{(\sqrt{2\pi}\sigma)^d} \exp\left(-\frac{|x - \mu|^2}{2\sigma^2}\right) \quad [\mu \text{ & } x = \text{vectors}; \sigma = \text{scalar}; d = \text{dimension}]$$

For each class C , suppose we estimate mean μ_C , variance σ_C^2 , and prior $\pi_C = P(Y = C)$.

Given x , Bayes decision rule $r^*(x)$ returns class C that maximizes $P(X = x|Y = C)\pi_C$.

$\ln \omega$ is monotonically increasing for $\omega > 0$, so it is equivalent to maximize

$$Q_C(x) = \ln \left((\sqrt{2\pi})^d P(x) \pi_C \right) = -\frac{|x - \mu_C|^2}{2\sigma_C^2} - d \ln \sigma_C + \ln \pi_C$$

\uparrow quadratic in x . \uparrow normal distribution, estimates $P(X = x|Y = C)$

[In a 2-class problem, you can also incorporate an asymmetrical loss function the same way we incorporate the prior π_C . In a multi-class problem, it gets more complicated, because the penalty for guessing wrong might depend not just on the wrong guess, but also on the true class.]

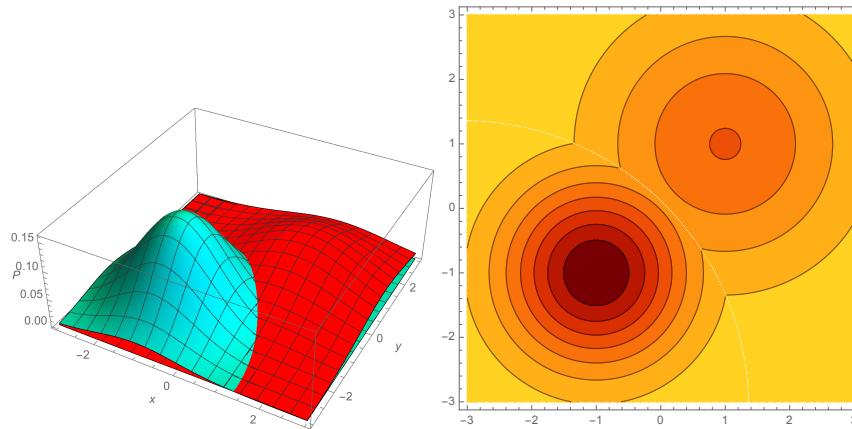
Quadratic Discriminant Analysis (QDA)

Suppose only 2 classes C, D. Then

$$r^*(x) = \begin{cases} C & \text{if } Q_C(x) - Q_D(x) > 0, \\ D & \text{otherwise.} \end{cases} \quad [\text{Pick the class with the biggest posterior probability}]$$

Decision fn is quadratic in x . Bayes decision boundary is $Q_C(x) - Q_D(x) = 0$.

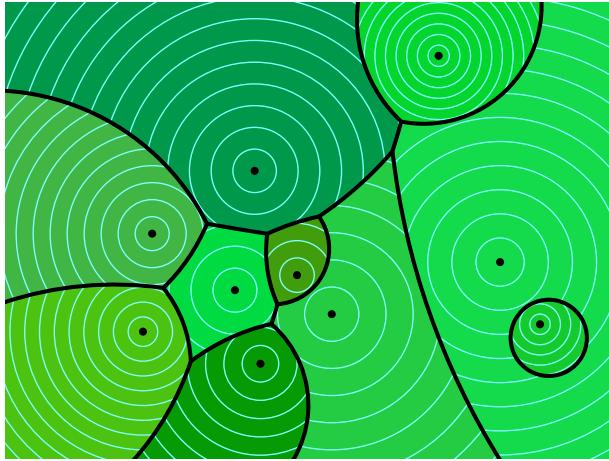
- In 1D, B.d.b. may have 1 or 2 points. [Solutions to a quadratic equation]
 - In d -D, B.d.b. is a quadric. [In 2D, that's a conic section]



`qda3d.pdf`, `qdacontour.pdf` [The same example I showed during the last lecture.]

[You're probably familiar with the Gaussian distribution where x and μ are scalars, but as I've written it, it applies equally well to a multi-dimensional feature space with isotropic Gaussians. Then x and μ are vectors, but the variance σ is still a scalar. Next lecture we'll look at anisotropic Gaussians where the variance is different along different directions.]

[QDA works very naturally with more than 2 classes.]



multiplicative.pdf [The feature space gets partitioned into regions. In two or more dimensions, you typically wind up with multiple decision boundaries that adjoin each other at joints. It looks like a sort of Voronoi diagram. In fact, it's a special kind of Voronoi diagram called a multiplicatively, additively weighted Voronoi diagram.]

[You might not be satisfied with just knowing how each point is classified. One of the great things about QDA is that you can also determine the probability that your classification is correct. Let's work that out.]

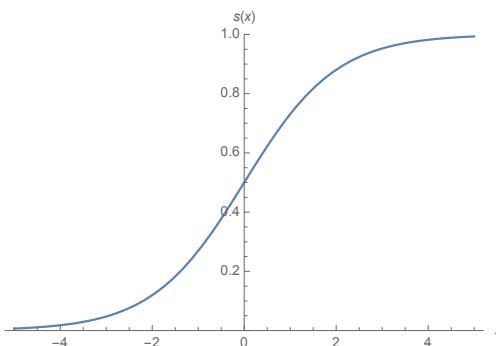
To recover posterior probabilities in 2-class case, use Bayes:

$$P(Y = C|X) = \frac{P(X|Y = C)\pi_C}{P(X|Y = C)\pi_C + P(X|Y = D)\pi_D}$$

recall $e^{Q_C(x)} = (\sqrt{2\pi})^d P(x)\pi_C$ [by definition of Q_C]

$$\begin{aligned} P(Y = C|X = x) &= \frac{e^{Q_C(x)}}{e^{Q_C(x)} + e^{Q_D(x)}} = \frac{1}{1 + e^{Q_D(x)-Q_C(x)}} \\ &= s(Q_C(x) - Q_D(x)), \quad \text{where} \end{aligned}$$

$$s(\gamma) = \frac{1}{1 + e^{-\gamma}} \iff \text{logistic fn aka sigmoid fn} \quad \text{[recall } Q_C - Q_D \text{ is the decision fn]}$$



logistic.pdf [The logistic function. Write beside it:] $s(0) = \frac{1}{2}$, $s(\infty) \rightarrow 1$, $s(-\infty) \rightarrow 0$, monotonically increasing.

[We interpret $s(0) = \frac{1}{2}$ as saying that on the decision boundary, there's a 50% chance of class C and a 50% chance of class D.]

Linear Discriminant Analysis (LDA)

[LDA is a variant of QDA with linear decision boundaries. It's less likely to overfit than QDA.]

Fundamental assumption: all the Gaussians have same variance σ .

[The equations simplify nicely in this case.]

$$Q_C(x) - Q_D(x) = \underbrace{\frac{(\mu_C - \mu_D) \cdot x}{\sigma^2}}_{w \cdot x} \underbrace{- \frac{|\mu_C|^2 - |\mu_D|^2}{2\sigma^2} + \ln \pi_C - \ln \pi_D}_{+\alpha}$$

[The quadratic terms in Q_C and Q_D cancelled each other out!]

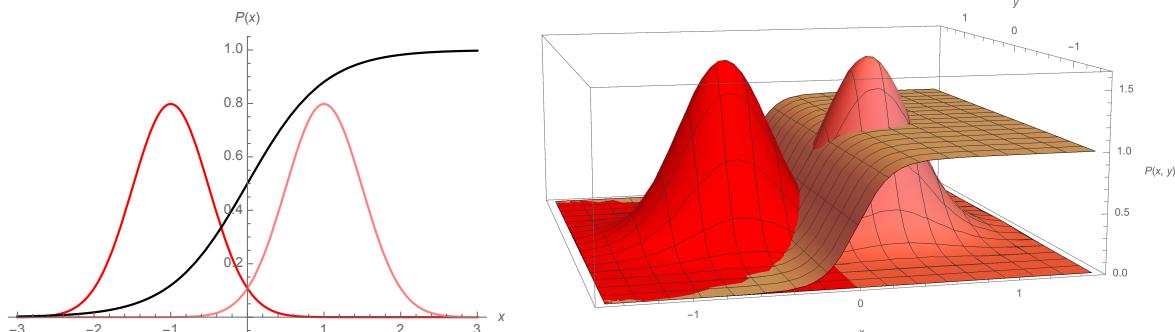
Now it's a linear classifier! Choose C that maximizes linear discriminant fn

$$\frac{\mu_C \cdot x}{\sigma^2} - \frac{|\mu_C|^2}{2\sigma^2} + \ln \pi_C \quad [\text{this works for any number of classes}]$$

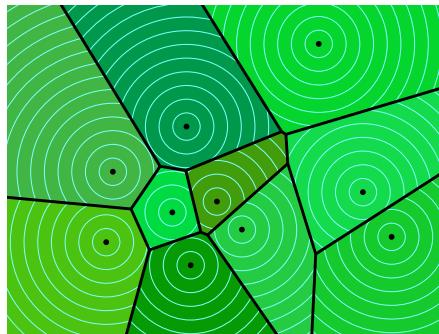
In 2-class case: decision boundary is $w \cdot x + \alpha = 0$

Bayes posterior is $P(Y = C|X = x) = s(w \cdot x + \alpha)$

[The effect of " $w \cdot x + \alpha$ " is to scale and translate the logistic fn in x -space. It's a linear transformation.]



lida1d.pdf [Two Gaussians (red) and the logistic function (black). The logistic function is the right Gaussian divided by the sum of the Gaussians. Observe that even when the Gaussians are 2D, the logistic function still looks 1D.]



voronoi.pdf [When you have many classes, their LDA decision boundaries form a classical Voronoi diagram if the priors π_C are equal. All the Gaussians have the same width.]

$$\text{If } \pi_C = \pi_D = \frac{1}{2} \Rightarrow (\mu_C - \mu_D) \cdot x - (\mu_C - \mu_D) \cdot \left(\frac{\mu_C + \mu_D}{2} \right) = 0$$

This is the centroid method!

MAXIMUM LIKELIHOOD ESTIMATION OF PARAMETERS (Ronald Fisher, circa 1912)

[To use Gaussian discriminant analysis, we must first fit Gaussians to the sample points and estimate the class prior probabilities. We'll do priors first—they're easier, because they involve a discrete distribution. Then we'll fit the Gaussians—they're less intuitive, because they're continuous distributions.]

Let's flip biased coins! Heads with probability p ; tails w/prob. $1 - p$.

10 flips, 8 heads, 2 tails. [Let me ask you a weird question.] What is the most likely value of p ?

Binomial distribution: $X \sim \mathcal{B}(n, p)$

$$P[X = x] = \binom{n}{x} p^x (1 - p)^{n-x} \quad [\text{this is the probability of getting exactly } x \text{ heads in } n \text{ coin flips}]$$

Our example: $n = 10$,

$$P[X = 8] = 45p^8(1 - p)^2 \stackrel{\text{def}}{=} \mathcal{L}(p)$$

Probability of 8 heads in 10 flips:

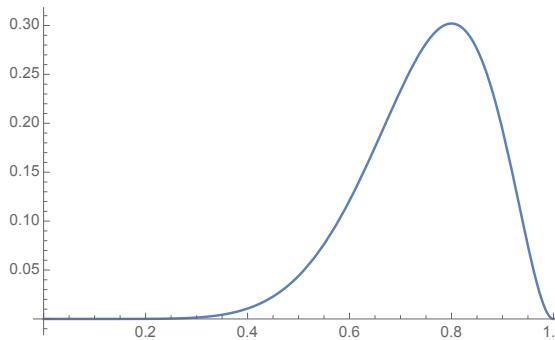
written as a fn $\mathcal{L}(p)$ of distribution parameter(s), this is the likelihood fn.

Maximum likelihood estimation (MLE): A method of estimating the parameters of a statistical model by picking the params that maximize [the likelihood function] \mathcal{L} .

...is one method of density estimation: estimating a PDF [probability density function] from data.

[Let's phrase it as an optimization problem.]

Find p that maximizes $\mathcal{L}(p)$.



binomlikelihood.pdf [Graph of $\mathcal{L}(p)$ for this example.]

Solve this example by setting derivative = 0:

$$\frac{d\mathcal{L}}{dp} = 360p^7(1 - p)^2 - 90p^8(1 - p) = 0$$

$$\Rightarrow 4(1 - p) - p = 0 \Rightarrow p = 0.8$$

[It shouldn't seem surprising that a coin that comes up heads 80% of the time is the coin most likely to produce 8 heads in 10 flips.]

[Note: $\frac{d^2\mathcal{L}}{dp^2} \doteq -18.9 < 0$ at $p = 0.8$, confirming it's a maximum.]

[Here's how this applies to prior probabilities. Suppose our data set is 10 sample points, and 8 of them are of class C and 2 are not. Then our estimated prior for class C will be $\pi_C = 0.8$.]

Likelihood of a Gaussian

Given sample points X_1, X_2, \dots, X_n , find best-fit Gaussian.

[Let's do this with a normal distribution instead of a binomial distribution. If you generate a random point from a normal distribution, what is the probability that it will be exactly at the mean of the Gaussian?]

[Zero. So it might seem like we have a problem here. With a continuous distribution, the probability of generating any particular point is zero. But we're just going to ignore that and do "likelihood" anyway.]

Likelihood of generating these points is

$$\mathcal{L}(\mu, \sigma; X_1, \dots, X_n) = P(X_1) P(X_2) \cdots P(X_n). \quad [\text{How do we maximize this?}]$$

The log likelihood $\ell(\cdot)$ is the \ln of the likelihood $\mathcal{L}(\cdot)$.

Maximizing likelihood \Leftrightarrow maximizing log likelihood.

$$\begin{aligned} \ell(\mu, \sigma; X_1, \dots, X_n) &= \ln P(X_1) + \ln P(X_2) + \dots + \ln P(X_n) \\ &= \sum_{i=1}^n \underbrace{\left(-\frac{|X_i - \mu|^2}{2\sigma^2} - d \ln \sqrt{2\pi} - d \ln \sigma \right)}_{\ln \text{ of normal distribution}} \end{aligned}$$

$$\text{Want to set } \nabla_\mu \ell = 0, \frac{\partial \ell}{\partial \sigma} = 0$$

$$\nabla_\mu \ell = \sum_{i=1}^n \frac{X_i - \mu}{\sigma^2} = 0 \quad \Rightarrow \quad \hat{\mu} = \frac{1}{n} \sum_{i=1}^n X_i \quad [\text{The hats } \hat{\cdot} \text{ mean "estimated"}]$$

$$\frac{\partial \ell}{\partial \sigma} = \sum_{i=1}^n \frac{|X_i - \mu|^2 - d\sigma^2}{\sigma^3} = 0 \quad \Rightarrow \quad \hat{\sigma}^2 = \frac{1}{dn} \sum_{i=1}^n |X_i - \mu|^2$$

We don't know μ exactly, so substitute $\hat{\mu}$ for μ to compute $\hat{\sigma}$.

I.e., we use mean & variance of points in class C to estimate mean & variance of Gaussian for class C.

For QDA: estimate conditional mean $\hat{\mu}_C$ & conditional variance $\hat{\sigma}_C^2$ of **each class C separately** [as above] & estimate the priors:

$$\hat{\pi}_C = \frac{n_C}{\sum_D n_D} \quad \Leftarrow \quad \text{total sample points in all classes} \quad [\hat{\pi}_C \text{ is the coin flip parameter}]$$

For LDA: same means & priors; one variance for all classes:

$$\hat{\sigma}^2 = \frac{1}{dn} \sum_C \sum_{\{i:y_i=C\}} |X_i - \hat{\mu}_C|^2 \quad \Leftarrow \quad \underline{\text{pooled within-class variance}}$$

[Notice that although LDA is computing one variance for all the data, each sample point contributes with respect to *its own class's mean*. This gives a very different result than if you simply use the global mean! It's usually smaller than the global variance. We say "within-class" because we use each point's distance from its class's mean, but "pooled" because we then pool all the classes together.]

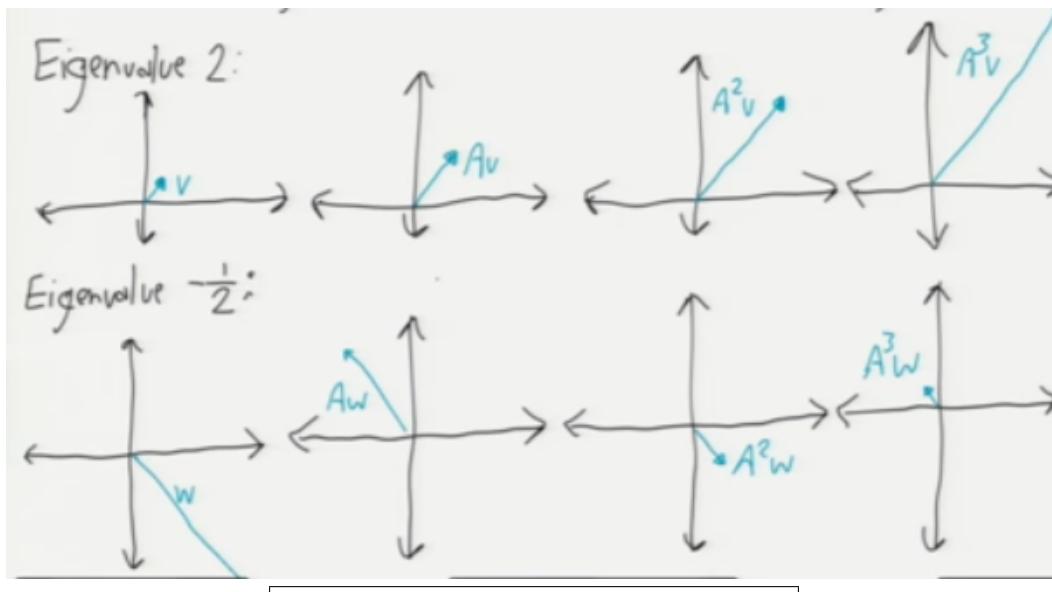
8 Eigenvectors and the Anisotropic Multivariate Normal Distribution

EIGENVECTORS

[I don't know if you were properly taught about eigenvectors here at Berkeley, but I sure don't like the way they're taught in most linear algebra books. So I'll start with a review. You all know the definition of an eigenvector:]

Given square matrix A , if $Av = \lambda v$ for some vector $v \neq 0$, scalar λ , then v is an eigenvector of A and λ is the eigenvalue of A associated w/ v .

[But what does that mean? It means that v is a magical vector that, after being multiplied by A , still points in the *same direction*, or in exactly the *opposite direction*.]



Draw this figure by hand (eigenvectors.png)

[For most matrices, most vectors don't have this property. So the ones that do are special, and we call them eigenvectors.]

[Clearly, when you scale an eigenvector, it's still an eigenvector. Only the direction matters, not the length. Let's look at a few consequences.]

Theorem: if v is eigenvector of A w/eigenvalue λ ,
then v is eigenvector of A^k w/eigenvalue λ^k [k is a +ve integer; we will use Theorem later]

Proof: $A^2v = A(\lambda v) = \lambda^2 v$, etc.

Theorem: moreover, if A is invertible,
then v is eigenvector of A^{-1} w/eigenvalue $1/\lambda$

Proof: $A^{-1}v = \frac{1}{\lambda}A^{-1}Av = \frac{1}{\lambda}v$

[look at the figures above, but go from right to left.]

[Stated simply: When you invert a matrix, the eigenvectors don't change, but the eigenvalues get inverted. When you square a matrix, the eigenvectors don't change, but the eigenvalues get squared.]

[Those theorems are pretty obvious. The next theorem is not obvious at all.]

Spectral Theorem: every real, symmetric $n \times n$ matrix has real eigenvalues and n eigenvectors that are mutually orthogonal, i.e., $v_i^\top v_j = 0$ for all $i \neq j$

[This takes about a page of math to prove.]

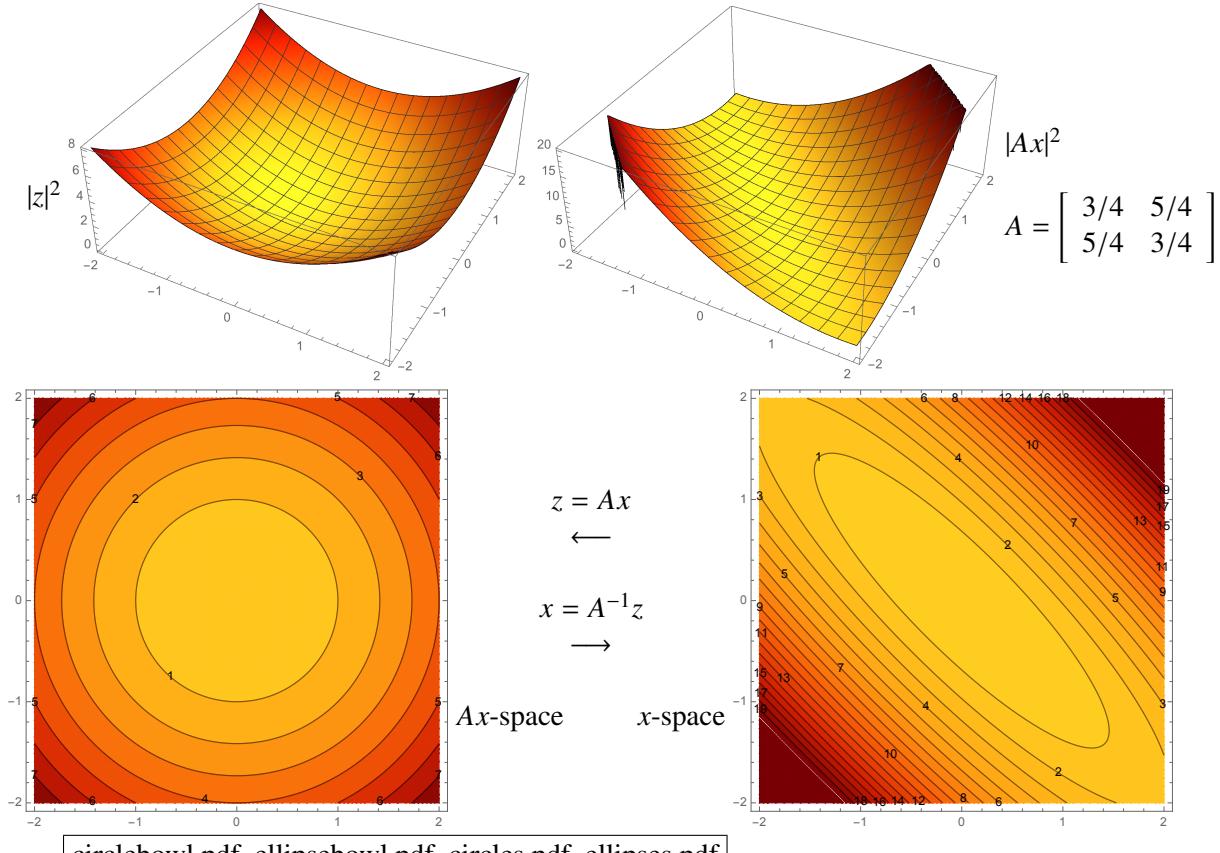
One minor detail is that a matrix can have more than n eigenvector directions. If two eigenvectors happen to have the same eigenvalue, then every linear combination of those eigenvectors is also an eigenvector. Then you have infinitely many eigenvector directions, but they all span the same plane. So you just arbitrarily pick two vectors in that plane that are orthogonal to each other. By contrast, the set of eigenvalues is always uniquely determined by a matrix, including the multiplicity of the eigenvalues.]

We can use them as a basis for \mathbb{R}^n .

Quadratic Forms

[My favorite way to visualize a symmetric matrix is to graph something called *the quadratic form*, which shows how applying the matrix affects the length of a vector. The following example uses the same two eigenvectors and eigenvalues as above.]

$$\begin{array}{ll} |z|^2 &= z^\top z \\ |Ax|^2 &= x^\top A^2 x \end{array} \quad \begin{array}{l} \Leftarrow \text{quadratic; isotropic; isosurfaces are spheres} \\ \Leftarrow \text{quadratic form of the matrix } A^2 \quad (\text{A symmetric}) \\ \Downarrow \text{anisotropic; isosurfaces are ellipsoids} \end{array}$$



[Both figures at left graph $|z|^2$, and both figures at right graph $|Ax|^2$.]

(Draw the stretch direction $(1, 1)$ with eigenvalue 2 and the shrink direction $(1, -1)$ with eigenvalue $-\frac{1}{2}$ on the ellipses at bottom right.)

[The matrix A maps the ellipses on the right to the circles on the left. They're stretching along the direction with eigenvalue 2, and shrinking along the direction with eigenvalue $-1/2$. You can also think of this process in reverse, if you remember that A^{-1} has the same eigenvectors but reciprocal eigenvalues. The matrix A^{-1} maps the circles to the ellipses, shrinking along the direction with eigenvalue 2, and stretching along the direction with eigenvalue $-1/2$. I can prove that.]

$|Ax|^2 = 1$ is an ellipsoid with axes v_1, v_2, \dots, v_n and radii $1/\lambda_1, 1/\lambda_2, \dots, 1/\lambda_n$

because if v_i has length $1/\lambda_i$, $|Av_i|^2 = |\lambda_i v_i|^2 = 1 \Rightarrow v_i$ lies on the ellipsoid

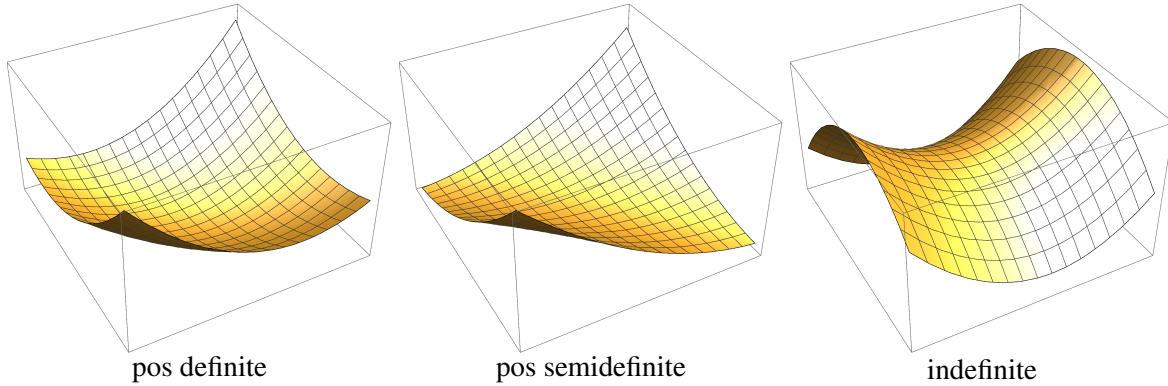
[The reason the ellipsoid radii are the reciprocals of the eigenvalues is that it is the matrix A^{-1} that maps the spheres to ellipsoids. So each axis of the spheres gets scaled by $1/\text{eigenvalue}$.]

bigger eigenvalue \Leftrightarrow steeper hill \Leftrightarrow shorter ellipsoid radius
 [↑ bigger curvature, to be precise]

Special case: A is diagonal \Leftrightarrow eigenvectors are coordinate axes
 \Leftrightarrow ellipsoids are axis-aligned

[Draw axis-aligned isocontours for a diagonal metric.]

A square matrix B is	<u>positive definite</u>	if $w^\top B w > 0$ for all $w \neq 0$.	\Leftrightarrow	all eigenvalues positive
	<u>positive semidefinite</u>	if $w^\top B w \geq 0$ for all w .	\Leftrightarrow	all eigenvalues nonnegative
	<u>indefinite</u>	if +ve eigenvalue & -ve eigenvalue		
	<u>invertible</u>	if no zero eigenvalue		



[posdef.pdf](#), [possemi.pdf](#), [indef.pdf](#)

[Examples of quadratic forms for positive definite, positive semidefinite, and indefinite matrices. Positive eigenvalues correspond to axes where the curvature goes up; negative eigenvalues correspond to axes where the curvature goes down. (Draw the eigenvector directions, and draw the flat trough in the positive semidefinite bowl.)]

What does this tell us about $x^\top A^2 x$?

[Recall that we've been visualizing the quadratic form of a matrix A^2 . The eigenvalues of A^2 are the squares of the eigenvalues of A , so A^2 cannot have a negative eigenvalue. Therefore, A^2 is positive semidefinite. A^2 might have an eigenvalue of zero, but if it does, the isosurfaces are cylinders instead of ellipsoids. We're going to use these ideas to define Gaussian distributions, and for that, we'll need a strictly positive definite covariance matrix.]

Building a Quadratic

[There are a lot of applications where you're given a matrix, and you want to extract the eigenvectors and eigenvalues. But when you're learning the math, I think it's more intuitive to go in the opposite direction. Suppose you have an ellipsoid in mind. Suppose you pick the ellipsoid axes and the radius along each axis, and you want to create the matrix that will embody the ellipsoid of your dreams.]

Choose n mutually orthogonal **unit** n -vectors v_1, \dots, v_n [so they specify an orthonormal coordinate system]

Let $V = [v_1 \ v_2 \ \dots \ v_n]$ $\Leftarrow n \times n$ matrix

Observe: $V^T V = I$

[off-diagonal 0's because the vectors are orthogonal]

[diagonal 1's because they're unit vectors]

$$\Rightarrow V^T = V^{-1} \Rightarrow VV^T = I$$

V is orthonormal matrix: acts like rotation (or reflection)

Choose some inverse radii λ_i :

$$\text{Let } \Lambda = \begin{bmatrix} \lambda_1 & 0 & \dots & 0 \\ 0 & \lambda_2 & & 0 \\ \vdots & & \ddots & \vdots \\ 0 & 0 & \dots & \lambda_n \end{bmatrix} \quad [\text{diagonal matrix of eigenvalues}]$$

Defn. of "eigenvector": $AV = VA$

[This is the same definition of eigenvector I gave you at the start of the lecture, but this is how we express it in matrix form, so we can cover all the eigenvectors in one statement.]

$$\Rightarrow AVV^T = V\Lambda V^T \quad [\text{which leads us to ...}]$$

Theorem: $A = V\Lambda V^T = \sum_{i=1}^n \lambda_i \underbrace{v_i v_i^T}_{\text{outer product}}$ has chosen eigenvectors/values

outer product: $n \times n$ matrix, rank 1

This is a matrix factorization called the eigendecomposition. [every real, symmetric matrix has one]

Λ is the diagonalized version of A .

V^T rotates the ellipsoid to be axis-aligned.

[So now, if you want a paraboloid with specified axes and radii, you can write it as $x^T A^2 x = 1$.]

Observe: $A^2 = V\Lambda V^T V\Lambda V^T = V\Lambda^2 V^T$

Given a symmetric PSD matrix M , we can find a symmetric square root $A = M^{1/2}$:

compute eigenvectors/values of M

take square roots of M 's eigenvalues

reassemble matrix A [with the same eigenvectors as M but changed eigenvalues]

[The first step of this algorithm—computing the eigenvectors and eigenvalues of a matrix—is much harder than the remaining two steps.]

ANISOTROPIC GAUSSIANS

[Let's revisit the multivariate Gaussian distribution, with different variances along different directions.]

$$X \sim \mathcal{N}(\mu, \Sigma)$$

[X and μ are d -vectors. X is random variable with mean μ .]

$$P(x) = \frac{1}{(\sqrt{2\pi})^d \sqrt{|\Sigma|}} \exp\left(-\frac{1}{2}(x - \mu)^\top \Sigma^{-1} (x - \mu)\right)$$

↑ determinant of Σ

Σ is the $d \times d$ SPD covariance matrix.

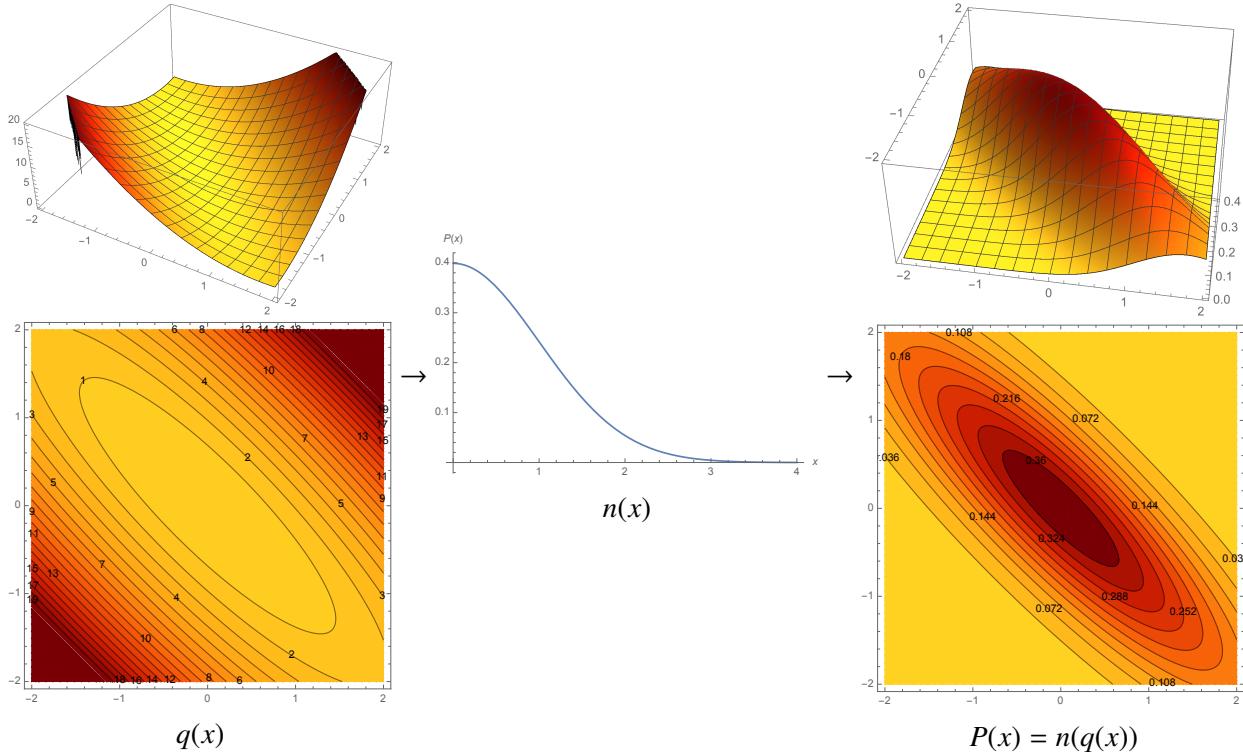
Σ^{-1} is the $d \times d$ SPD precision matrix.

Write $P(x) = n(q(x))$, where $q(x) = (x - \mu)^\top \Sigma^{-1} (x - \mu)$

↑ ↑
 $\mathbb{R} \rightarrow \mathbb{R}$, exponential $\mathbb{R}^d \rightarrow \mathbb{R}$, quadratic

[Now $q(x)$ is a function we understand—it's just a quadratic bowl centered at μ whose curvature is represented by the precision matrix Σ^{-1} . The other function $n(\cdot)$ is like a 1D Gaussian with a different normalization constant. This mapping $n(\cdot)$ does not change the isosurfaces.]

Principle: given monotonic $n : \mathbb{R} \rightarrow \mathbb{R}$, isosurfaces of $n(q(x))$ are same as $q(x)$ (different isovalues).



ellipsebowl.pdf, ellipses.pdf, gauss.pdf, gauss3d.pdf, gausscontour.pdf

[(Show this figure on a separate “whiteboard” for easy reuse next lecture.) A paraboloid (left) becomes a bivariate Gaussian (right) after you compose it with the univariate Gaussian (center).]

[One of the ideas I'm trying to get across is that if you understand the isosurfaces of a quadratic function, then you understand the isosurfaces of a Gaussian, because they're the same. The differences are in the isovalue—in particular, the Gaussian achieves its maximum at the mean, and decreases to zero as you move infinitely far away from the mean.]

$q(x)$ is the squared distance from $\Sigma^{-1/2}x$ to $\Sigma^{-1/2}\mu$. Consider the metric

$$d(x, \mu) = |\Sigma^{-1/2}x - \Sigma^{-1/2}\mu| = \sqrt{(x - \mu)^\top \Sigma^{-1}(x - \mu)} = \sqrt{q(x)}.$$

[So we think of the precision matrix as a “metric tensor” which defines a metric, a sort of warped distance from x to the mean μ .]

covariance: Let R, S be random variables—vectors or scalars

$$\text{Cov}(R, S) = E[(R - E[R])(S - E[S])^\top] = E[RS^\top] - \mu_R \mu_S^\top$$

$$\text{Var}(R) = \text{Cov}(R, R)$$

If R is a vector, covariance matrix for R is

$$\text{Var}(R) = \begin{bmatrix} \text{Var}(R_1) & \text{Cov}(R_1, R_2) & \dots & \text{Cov}(R_1, R_d) \\ \text{Cov}(R_2, R_1) & \text{Var}(R_2) & & \text{Cov}(R_2, R_d) \\ \vdots & & \ddots & \vdots \\ \text{Cov}(R_d, R_1) & \text{Cov}(R_d, R_2) & \dots & \text{Var}(R_d) \end{bmatrix} \quad [\text{symmetric; each } R_i \text{ is scalar}]$$

For a Gaussian $R \sim \mathcal{N}(\mu, \Sigma)$, one can show $\text{Var}(R) = \Sigma$.

[... by integrating the expectation in anisotropic spherical coordinates. It's a painful integral.]

[An important point is that statisticians didn't just arbitrarily decide to call Σ a covariance matrix. Rather, statisticians discovered that if you find the covariance of the normal distribution by integration, it turns out that the covariance is Σ . This is a happy discovery; it's rather elegant.]

R_i, R_j independent $\Rightarrow \text{Cov}(R_i, R_j) = 0$ [the reverse implication is not generally true, but ...]

$\text{Cov}(R_i, R_j) = 0$ AND multivariate normal dist. $\Rightarrow R_i, R_j$ independent

all features pairwise independent $\Rightarrow \text{Var}(R)$ is diagonal

$\text{Var}(R)$ is diagonal AND joint normal

\Leftrightarrow axis-aligned Gaussian; squared radii on diagonal of $\Sigma = \text{Var}(R)$

$$\Leftrightarrow \underbrace{P(x)}_{\text{multivariate univariate Gaussians}} = \underbrace{P(x_1) P(x_2) \cdots P(x_d)}$$

[So when the features are independent, you can write the multivariate Gaussian as a product of univariate Gaussians. When they aren't, you can do a change of coordinates to the eigenvector coordinate system, and write it as a product of univariate Gaussians in eigenvector coordinates.]

9 Anisotropic Gaussians, Maximum Likelihood Estimation, QDA, and LDA

ANISOTROPIC GAUSSIANS

[Recall from our last lecture the probability density function of the multivariate normal distribution in its full generality.]

$$\text{Normal PDF: } P(x) = \frac{1}{(\sqrt{2\pi})^d \sqrt{|\Sigma|}} \exp\left(-\frac{1}{2}(x - \mu)^\top \Sigma^{-1} (x - \mu)\right) \quad [x, \mu \text{ are } d\text{-vectors}]$$

\uparrow determinant of Σ

$$\text{Write } P(x) = n(q(x)), \text{ where } q(x) = (x - \mu)^\top \Sigma^{-1} (x - \mu)$$

$\uparrow \quad \uparrow$

$\mathbb{R} \rightarrow \mathbb{R}$, exponential $\mathbb{R}^d \rightarrow \mathbb{R}$, quadratic

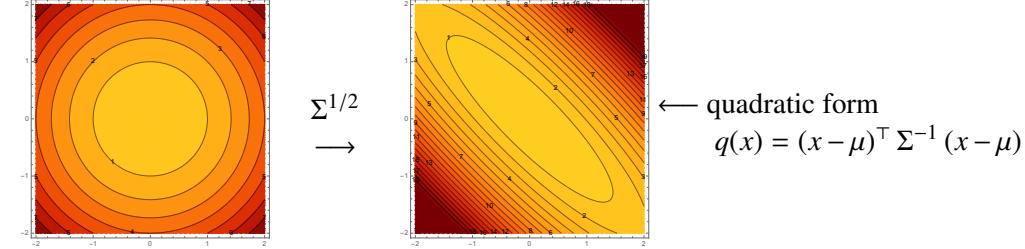
[The covariance matrix Σ and its symmetric square root and its inverse all play roles in our intuition about the multivariate normal distribution.]

$$\Sigma = V\Gamma V^\top \quad \text{covariance matrix}$$

\uparrow eigenvalues of Σ are variances along the eigenvectors, $\gamma_{11} = \sigma_1^2$

$$\Sigma^{1/2} = V\Gamma^{1/2}V^\top \quad \text{maps spheres to ellipsoids}$$

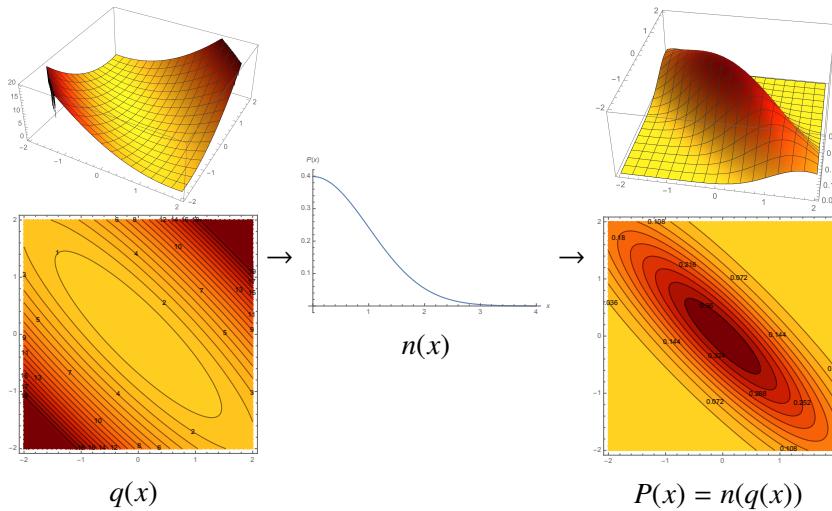
\uparrow eigenvalues of $\Sigma^{1/2}$ are Gaussian widths / ellipsoid radii / standard deviations, $\sqrt{\gamma_{ii}} = \sigma_i$



$$\Sigma^{-1} = V\Gamma^{-1}V^\top \quad \text{precision matrix (metric tensor)} \quad [\uparrow \text{quadratic form of } \Sigma^{-1} \text{ defines contours}]$$

$$\Sigma^{-1/2} = V\Gamma^{-1/2}V^\top \quad (\Sigma^{-1/2} \text{ was } A \text{ in last lecture})$$

[Recall from last lecture that the isocontours of the multivariate normal distribution are the same as the isocontours of the quadratic form of the precision matrix Σ^{-1} .]



Maximum Likelihood Estimation for Anisotropic Gaussians

Given sample points X_1, \dots, X_n and classes y_1, \dots, y_n , find best-fit Gaussians.

X_i is a column vector. [To fit our definition of the Gaussian distribution $P(x)$.]

[Once again, we want to fit the Gaussian that maximizes the likelihood of generating the sample points in a specified class. This time I won't derive the maximum-likelihood Gaussian; I'll just tell you the answer.]

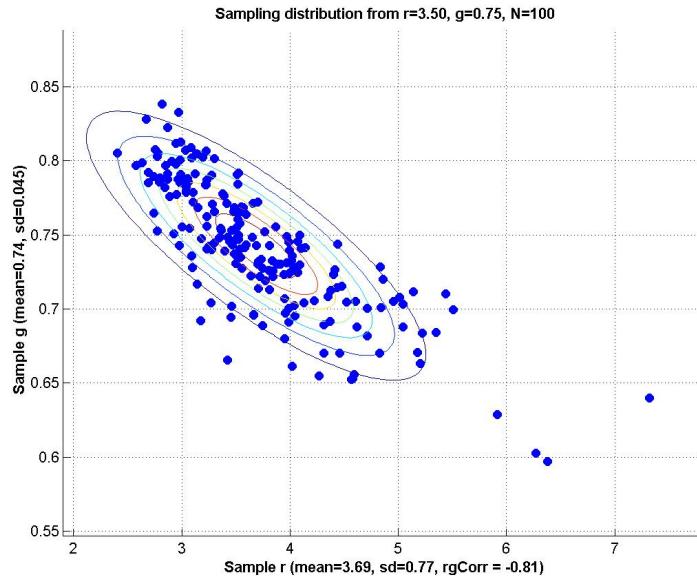
For QDA:

$$\hat{\Sigma}_C = \frac{1}{n_C} \sum_{i:y_i=C} \underbrace{(X_i - \hat{\mu}_C)(X_i - \hat{\mu}_C)^\top}_{\text{outer product matrix}} \quad \Leftarrow \text{conditional covariance for pts in class C}$$

[where n_C is the number of points in class C.]

Prior $\hat{\pi}_C$, mean $\hat{\mu}_C$: same as before

[$\hat{\pi}_C$ is number of points in class C \div total sample points; $\hat{\mu}_C$ is mean of sample points in class C.]



maxlike.jpg [Maximum likelihood estimation takes these points and outputs this Gaussian].

$\hat{\Sigma}_C$ is positive semidefinite, but not always definite!

[If there are some zero eigenvalues, the standard version of QDA just doesn't work. We can try to fix it by eliminating the zero-variance dimensions (eigenvectors). Homeworks 2 and 3 suggest two ways to do that.]

For LDA:

$$\hat{\Sigma} = \frac{1}{n} \sum_C \sum_{i:y_i=C} (X_i - \hat{\mu}_C)(X_i - \hat{\mu}_C)^\top \quad \Leftarrow \text{pooled within-class covariance matrix}$$

[Let's revisit QDA and LDA and see what has changed now that we know anisotropic Gaussians. The short answer is “not much has changed, but the graphs look cooler.” By the way, capital X once again represents a random variable.]

QDA

Choosing C that maximizes $P(X = x|Y = C)\pi_C$ is equivalent to maximizing the quadratic discriminant fn

$$Q_C(x) = \ln((\sqrt{2\pi})^d P_C(x)\pi_C) = -\frac{1}{2}(x - \mu_C)^\top \Sigma_C^{-1} (x - \mu_C) - \frac{1}{2} \ln |\Sigma_C| + \ln \pi_C$$

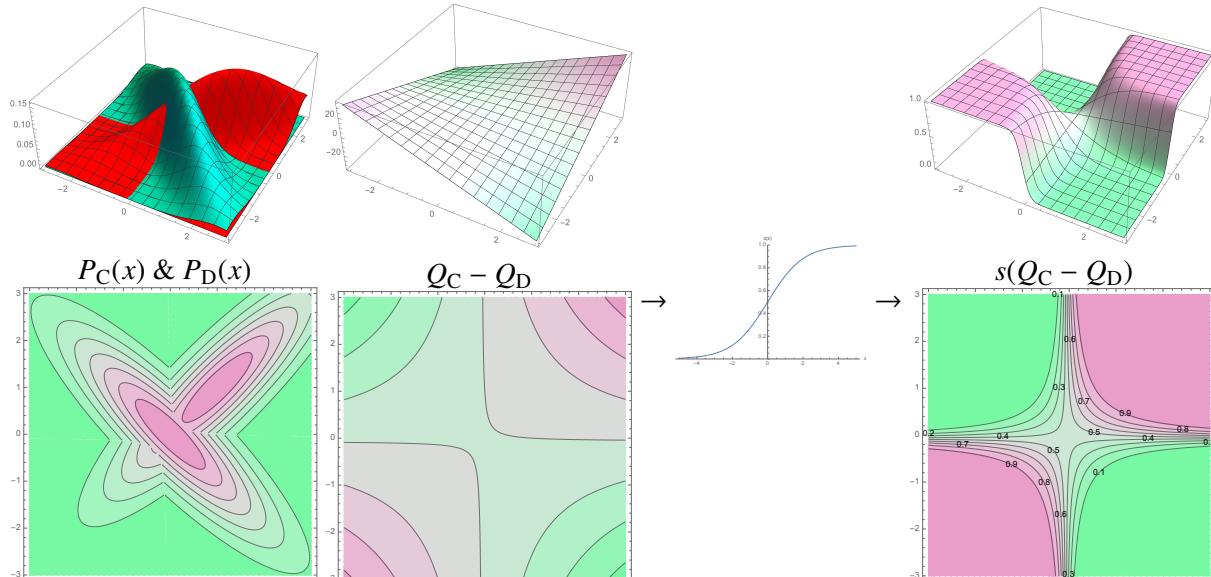
↑
Gaussian for C

[This works for any number of classes. In a multi-class problem, you just pick the class with the greatest quadratic discriminant for x .]

2 classes: Decision fn $Q_C(x) - Q_D(x)$ is quadratic, but may be indefinite

⇒ Bayes decision boundary is a quadric.

Posterior is $P(Y = C|X = x) = s(Q_C(x) - Q_D(x))$ where $s(\cdot)$ is logistic fn



[qdaaniso3d.pdf](#), [qdaanisocontour.pdf](#), [qdaanisodiff3d.pdf](#), [qdaanisodiffcontour.pdf](#),

[logistic.pdf](#), [qdaanisoposterior3d.pdf](#), [qdaanisoposteriorcontour.pdf](#)

[(Show this figure on a separate “whiteboard.”) An example where the decision boundary is a hyperbola—which is not possible with isotropic Gaussians. At left, two anisotropic Gaussians. Center left, the difference $Q_C - Q_D$. After applying the logistic function to this difference we obtain the posterior probabilities at right, which tells us the probability our prediction is correct. However, observe that we can see the decision boundary in both contour plots: it is $Q_C - Q_D = 0$ and $s(Q_C - Q_D) = 0.5$. We don’t need to apply the logistic function to find the decision boundary, but we do need to apply it if we want the posterior probabilities.]



[aniso.pdf](#) [When you have many classes, their QDA decision boundaries form an anisotropic Voronoi diagram. Interestingly, a cell of this diagram might not be connected.]

LDA

One $\hat{\Sigma}$ for all classes.

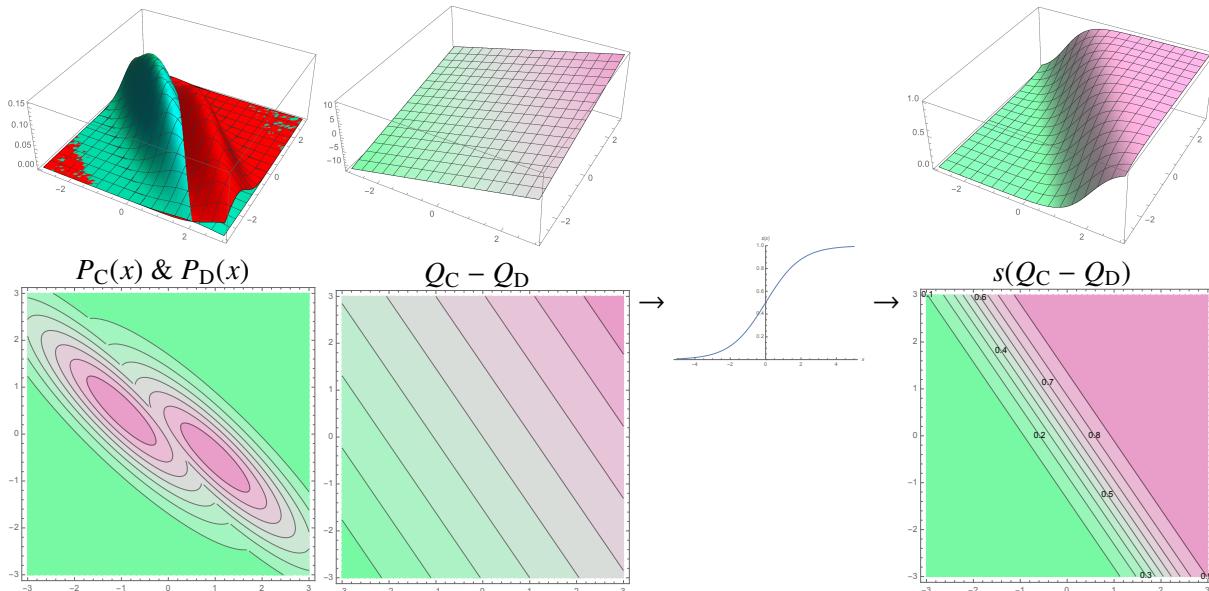
[Once again, the quadratic terms cancel each other out so the decision function is linear and the decision boundary is a hyperplane.]

$$Q_C(x) - Q_D(x) = \underbrace{(\mu_C - \mu_D)^\top \Sigma^{-1} x}_{w^\top x} - \underbrace{\frac{\mu_C^\top \Sigma^{-1} \mu_C - \mu_D^\top \Sigma^{-1} \mu_D}{2}}_{+\alpha} + \ln \pi_C - \ln \pi_D$$

Choose class C that maximizes the linear discriminant fn

$$\mu_C^\top \Sigma^{-1} x - \frac{1}{2} \mu_C^\top \Sigma^{-1} \mu_C + \ln \pi_C \quad [\text{works for any # of classes}]$$

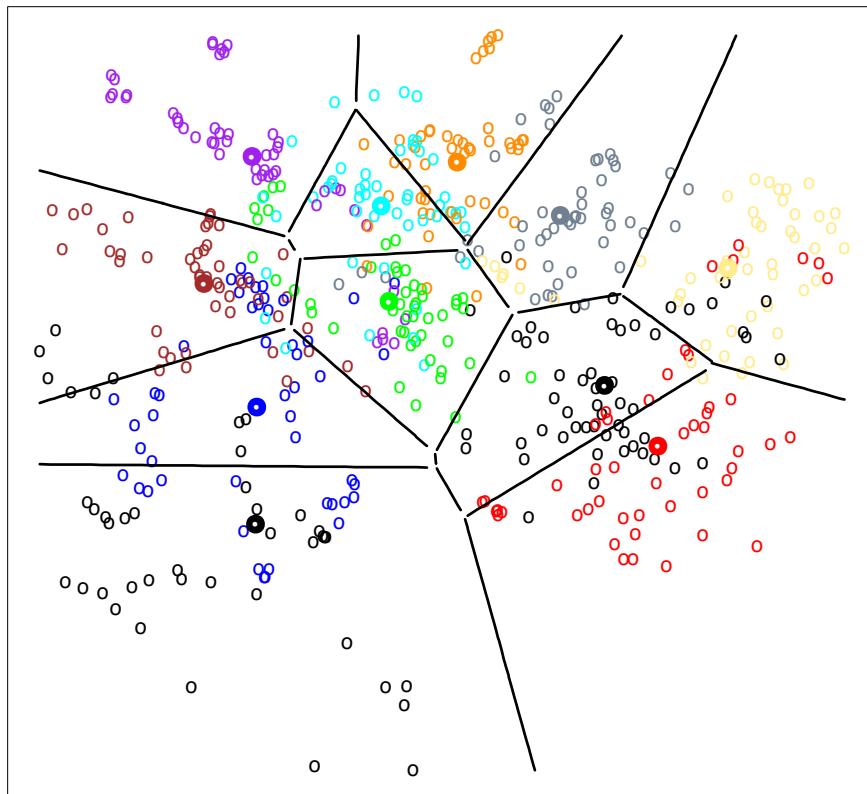
2 classes: Decision boundary is $w^\top x + \alpha = 0$
 Posterior is $P(Y = C|X = x) = s(w^\top x + \alpha)$



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logistic.pdf, lдаanisoposterior3d.pdf, lдаanisoposteriorcontour.pdf

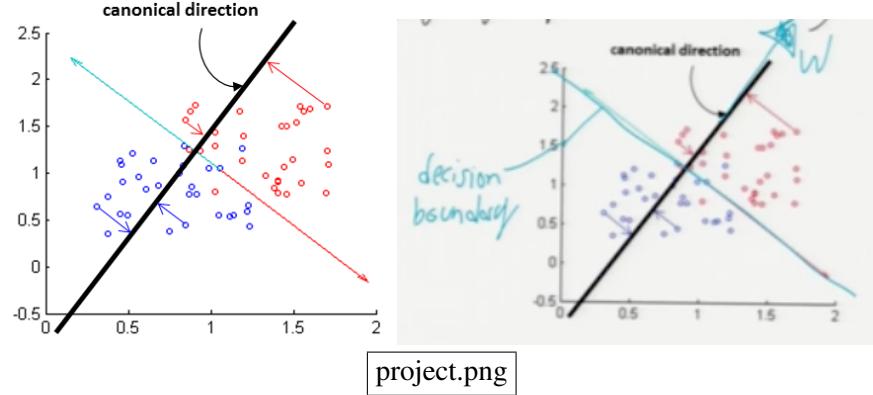
[Show this figure on a separate “whiteboard.”) In LDA, the decision boundary is always a hyperplane. Note that Mathematica messed up the top left plot a bit; there should be no red in the left corner, nor blue in the right corner.]



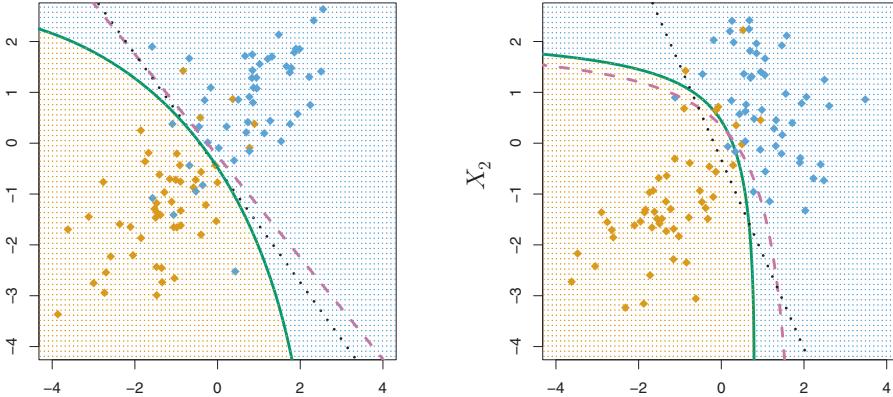
LDAdata.pdf (ESL, Figure 4.11) [An example of LDA with messy data. The points are not sampled from perfect Gaussians, but LDA still works reasonably well.]

Notes:

- LDA often interpreted as projecting points onto normal w ; cutting the line in half.



- For 2 classes,
 - LDA has $d + 1$ parameters (w, α);
 - QDA has $\frac{d(d+3)}{2} + 1$ params;
 - QDA more likely to overfit. [The danger is much bigger when the dimension d is large.]



ldaqda.pdf (ISL, Figure 4.9) [In these examples, the Bayes optimal decision boundary is purple (and dashed), the QDA decision boundary is green, the LDA decision boundary is black (and dotted). When the optimal boundary is linear, as at left, LDA gives a more stable fit whereas QDA may overfit. When the optimal boundary is curved, as at right, QDA often gives you a better fit.]

- With features, LDA can give nonlinear boundaries; QDA nonquadratic.
- We don't get **true** optimum Bayes classifier
 - estimate distributions from finite data
 - real-world data not perfectly Gaussian
- Changing priors or loss = adding constants to discriminant fns
[So it's very easy. In the 2-class case, it's equivalent to changing the isovalue . . .]
- Posterior gives decision boundaries for 10% probability, 50%, 90%, etc.
 - choosing isovalue = probability p is same as choosing asymmetrical loss p for false positive, $1 - p$ for false negative; OR as choosing $\pi_C = 1 - p, \pi_D = p$.

[LDA & QDA are the best method in practice for many applications. In the STATLOG project, either LDA or QDA were in the top three classifiers for 10 out of 22 datasets. But it's not because all those datasets are Gaussian. LDA & QDA work well when the data can only support simple decision boundaries such as linear or quadratic, because Gaussian models provide stable estimates. See ESL, Section 4.3]

Some Terms

Let X be $n \times d$ design matrix of sample pts

Each row i of X is a sample pt X_i^\top .

[Now I'm using capital X as a matrix instead of a random variable vector. I'm treating X_i as a column vector to match the standard convention for multivariate distributions like the Gaussian, but X_i^\top is a row of X .]

centering X : subtracting μ^\top from each row of X . $X \rightarrow \dot{X}$

$[\mu^\top$ is the mean of all the rows of X . Now the mean of all the rows of \dot{X} is zero.]

Let R be uniform distribution on sample pts. Sample covariance matrix is

$$\text{Var}(R) = \frac{1}{n} \dot{X}^\top \dot{X}.$$

[This is the simplest way to remember how to compute a covariance matrix for QDA. Imagine you have a design matrix X_C that contains only the sample points of class C; then you have $\hat{\Sigma}_C = \frac{1}{n_C} \dot{X}_C^\top \dot{X}_C$.]

[When we have points from an anisotropic Gaussian distribution, sometimes it's useful to perform a linear transformation that maps them to an axis-aligned distribution, or maybe even to an isotropic distribution. Recall that $\Sigma^{-1/2}$ maps ellipsoids to spheres.]

decorrelating \dot{X} : applying transform $Z = \dot{X}V$, where $\text{Var}(R) = V\Lambda V^\top$

[transforms the sample points to the eigenvector coordinate system]

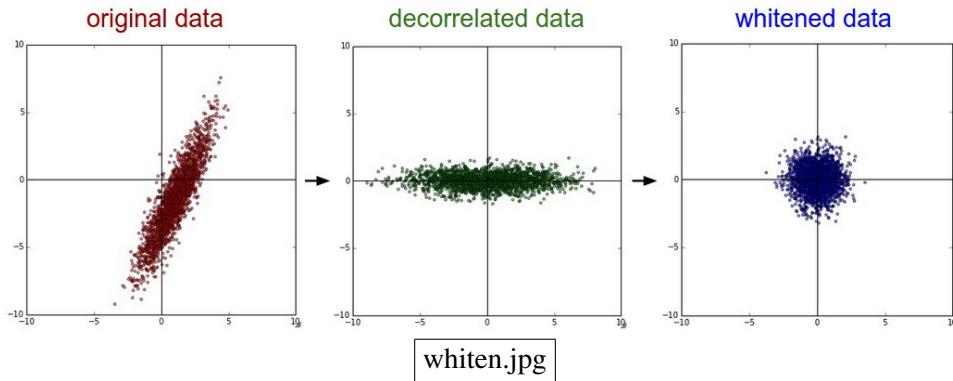
Then $\text{Var}(Z) = \Lambda$. [Z has diagonal covariance. If $X_i \sim \mathcal{N}(\mu, \Sigma)$, then approximately, $Z_i \sim \mathcal{N}(0, \Lambda)$.]

[Proof: $\text{Var}(Z) = (1/n)V^\top \dot{X}^\top \dot{X} V = V^\top \text{Var}(R) V = V^\top V \Lambda V^\top V = \Lambda$.]

spherling \dot{X} : applying transform $W = \dot{X} \text{Var}(R)^{-1/2}$

whitening X : centering + spherling, $X \rightarrow W$

Then W has covariance matrix I . [If $X_i \sim \mathcal{N}(\mu, \Sigma)$, then approximately, $W_i \sim \mathcal{N}(0, I)$.]



[Whitening input data is often used with other machine learning algorithms, like SVMs and neural networks. The idea is that some features may be much bigger than others—for instance, because they're measured in different units. SVMs penalize violations by large features more heavily than they penalize small features. Whitening the data before you run an SVM puts the features on an equal basis. Once nice thing about discriminant analysis is that whitening is built in.]

[Incidentally, what we've done here—computing a sample covariance matrix and its eigenvectors/values—is about 75% of an important unsupervised learning method called principal components analysis, or PCA, which we'll learn later in the semester.]

10 Regression, including Least-Squares Linear and Logistic Regression

REGRESSION aka Fitting Curves to Data

Classification: given point x , predict class (often binary)

Regression: given point x , predict a numerical value

[Classification gives a discrete prediction, whereas regression gives us a quantitative prediction, usually on a continuous scale.]

[We've already seen an example of regression in Gaussian discriminant analysis. QDA and LDA don't just give us a classifier; they also give us the probability that a particular class label is correct. So QDA and LDA implicitly do regression on probability values.]

- Choose form of regression fn $h(x; p)$ with parameters p ($h = \text{hypothesis}$)
 - like decision fn in classification [e.g. linear, quadratic, logistic in x]
- Choose a cost fn (objective fn) to optimize
 - usually based on a loss fn; e.g. risk fn = expected loss

Some regression fns:

- (1) linear: $h(x; w, \alpha) = w \cdot x + \alpha$
- (2) polynomial [equivalent to linear regression with added polynomial features]
- (3) logistic: $h(x; w, \alpha) = s(w \cdot x + \alpha)$ recall: logistic fn $s(\gamma) = \frac{1}{1+e^{-\gamma}}$

[The last choice is interesting. You'll recall that LDA produces a posterior probability function with this expression. So the logistic function seems to be a natural form for modeling certain probabilities. If we want to model class probabilities, sometimes we use LDA; but alternatively, we could skip fitting Gaussians to points, and instead just try to directly fit a logistic function to a set of probabilities.]

Some loss fns: let z be prediction $h(x)$; y be true value

- | | |
|---|--|
| (A) $L(z, y) = (z - y)^2$ | squared error |
| (B) $L(z, y) = z - y $ | absolute error |
| (C) $L(z, y) = -y \ln z - (1 - y) \ln(1 - z)$ | logistic loss, aka cross-entropy: $y \in [0, 1], z \in (0, 1)$ |

Some cost fns to minimize:

- | | |
|---|---|
| (a) $J(h) = \frac{1}{n} \sum_{i=1}^n L(h(X_i), y_i)$ | mean loss [you can leave out the " $\frac{1}{n}$ "] |
| (b) $J(h) = \max_{i=1}^n L(h(X_i), y_i)$ | maximum loss |
| (c) $J(h) = \sum_{i=1}^n \omega_i L(h(X_i), y_i)$ | weighted sum [some points are more important than others] |
| (d) $J(h) = \frac{1}{n} \sum_{i=1}^n L(h(X_i), y_i) + \lambda \ w\ ^2$ | ℓ_2 penalized/regularized |
| (e) $J(h) = \frac{1}{n} \sum_{i=1}^n L(h(X_i), y_i) + \lambda \ w\ _{\ell_1}$ | ℓ_1 penalized/regularized |

Some famous regression methods:

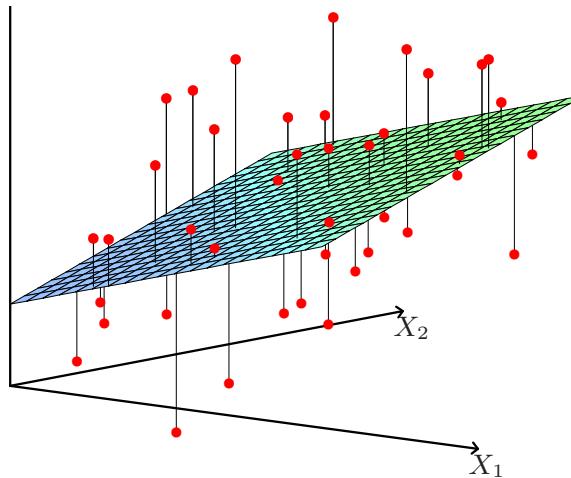
- | | | |
|---|--|-------------------------------------|
| Least-squares linear regr.: (1) + (A) + (a) | } | quadratic cost; minimize w/calculus |
| Weighted least-squ. linear: (1) + (A) + (c) | | |
| Ridge regression: (1) + (A) + (d) | convex cost; minimize w/gradient descent | |
| Logistic regr.: (3) + (C) + (a) | | |
| Lasso: (1) + (A) + (e) | quadratic program | |
| Least absolute deviations: (1) + (B) + (a) | | |
| Chebyshev criterion: (1) + (B) + (b) | linear program | |

[I have given you several choices of regression function form, several choices of loss function, and several choices of objective function. These are interchangeable parts where you can snap one part out and replace it with a different one. But the optimization algorithm and its speed depend crucially on which parts you pick. Let's consider some examples.]

LEAST-SQUARES LINEAR REGRESSION (Gauss, 1801)

Linear regression fn (1) + squared loss fn (A) + cost fn (a).

$$\text{Find } w, \alpha \text{ that minimizes } \sum_{i=1}^n (X_i \cdot w + \alpha - y_i)^2$$



`linregress.pdf` (ISL, Figure 3.4) [An example of linear regression.]

Convention: X is $n \times d$ design matrix of sample pts
 y is n -vector of scalar labels

$$\left[\begin{array}{cccccc} X_{11} & X_{12} & \dots & X_{1j} & \dots & X_{1d} \\ X_{21} & X_{22} & & X_{2j} & & X_{2d} \\ \vdots & & & & & \\ X_{i1} & X_{i2} & & X_{ij} & & X_{id} \\ \vdots & & & & & \\ X_{n1} & X_{n2} & & X_{nj} & & X_{nd} \end{array} \right] \leftarrow \text{point } X_i^\top \quad \left[\begin{array}{c} y_1 \\ y_2 \\ \vdots \\ y_n \end{array} \right] \quad \begin{matrix} \uparrow \\ \text{feature column } X_{*j} \\ \uparrow \\ y \end{matrix}$$

Usually $n > d$. [But not always.]

Recall fictitious dimension trick [from Lecture 3]: rewrite $h(x) = x \cdot w + \alpha$ as

$$[x_1 \ x_2 \ 1] \cdot \begin{bmatrix} w_1 \\ w_2 \\ \alpha \end{bmatrix}.$$

Now X is an $n \times (d+1)$ matrix; w is a $(d+1)$ -vector. [We've added a column of all-1's to the end of X .] [We rewrite the optimization problem above:]

$$\text{Find } w \text{ that minimizes } |Xw - y|^2 = \text{RSS}(w), \text{ for residual sum of squares}$$

Optimize by calculus:

$$\begin{aligned}\text{minimize RSS}(w) &= w^\top X^\top X w - 2y^\top X w + y^\top y \\ \nabla \text{RSS} &= 2X^\top X w - 2X^\top y = 0 \\ \Rightarrow \underbrace{X^\top X}_{(d+1) \times (d+1)} \underbrace{w}_{(d+1)-\text{vectors}} &= X^\top y \quad \Leftarrow \text{the normal equations [} w \text{ unknown; } X \& y \text{ known]}\end{aligned}$$

If $X^\top X$ is singular, problem is underconstrained

[because the sample points all lie on a common hyperplane. Notice that $X^\top X$ is always positive semidefinite.]

We use a linear solver to find $w = \underbrace{(X^\top X)^{-1} X^\top y}_{X^+, \text{ the pseudoinverse of } X, (d+1) \times n}$ [never actually invert the matrix!]

[We never compute X^+ directly, but we are interested in the fact that w is a linear transformation of y .] [X is usually not square, so X can't have an inverse. However, every X has a pseudoinverse X^+ , and if $X^\top X$ is invertible, then X^+ is a "left inverse."]

Observe: $X^+ X = (X^\top X)^{-1} X^\top X = I \Leftarrow (d+1) \times (d+1)$ [which explains the name "left inverse"]

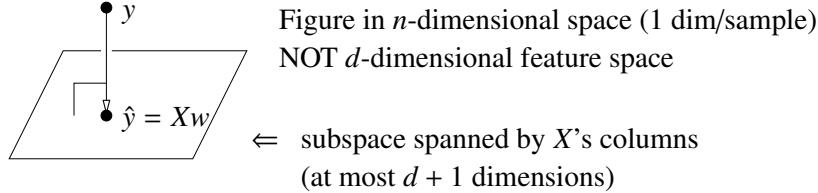
Observe: the predicted values of y are $\hat{y}_i = w \cdot X_i \Rightarrow \hat{y} = Xw = XX^+y = Hy$

where $\underbrace{H}_{n \times n} = XX^+$ is called the hat matrix because it puts the hat on y

[Ideally, H would be the identity matrix and we'd have a perfect fit, but if $n > d + 1$, then H is singular.]

Interpretation as a projection:

- $\hat{y} = Xw \in \mathbb{R}^n$ is a linear combination of columns of X (one column per feature)
- For fixed X , varying w , Xw is subspace of \mathbb{R}^n spanned by columns



- Minimizing $|\hat{y} - y|^2$ finds point \hat{y} nearest y on subspace
 - ⇒ projects y orthogonally onto subspace
 - [the vertical line is the direction of projection and the error vector]
- Error is smallest when line is perpendicular to subspace: $X^\top(Xw - y) = 0$
 - ⇒ the normal equations!
- Hat matrix H does the projecting. [H is sometimes called the projection matrix.]

Advantages:

- Easy to compute; just solve a linear system.
- Unique, stable solution. [... except when the problem is underconstrained.]

Disadvantages:

- Very sensitive to outliers, because errors are squared!
- Fails if $X^\top X$ is singular.

[Apparently, least-squares linear regression was first posed and solved in 1801 by the great mathematician Carl Friedrich Gauss, who used least-squares regression to predict the trajectory of the planetoid Ceres. A paper he wrote on the topic is regarded as the birth of modern linear algebra.]

LOGISTIC REGRESSION (David Cox, 1958)

Logistic regression fn (3) + logistic loss fn (C) + cost fn (a).
Fits “probabilities” in range (0, 1).

Usually used for classification. The input y_i 's *can* be probabilities, but in most applications they're all 0 or 1.

QDA, LDA: generative models

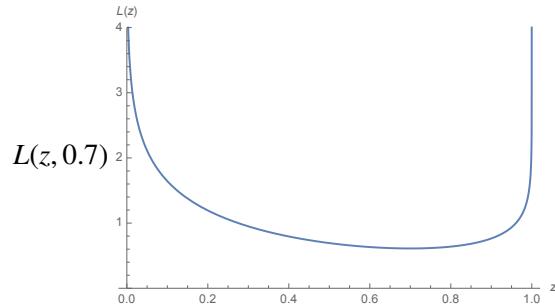
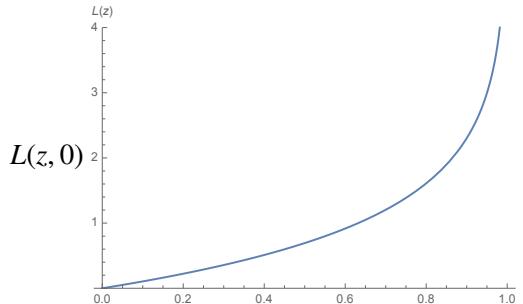
logistic regression: discriminative model

[We've learned from LDA that in classification, the posterior probabilities are often modeled well by a logistic function. So why not just try to fit a logistic function directly to the data, skipping the Gaussians?]

With X and w including the fictitious dimension; α is w 's last component ...

Find w that minimizes

$$J = - \sum_{i=1}^n \left(y_i \ln s(X_i \cdot w) + (1 - y_i) \ln (1 - s(X_i \cdot w)) \right)$$

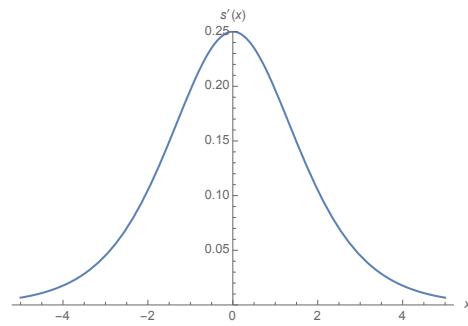
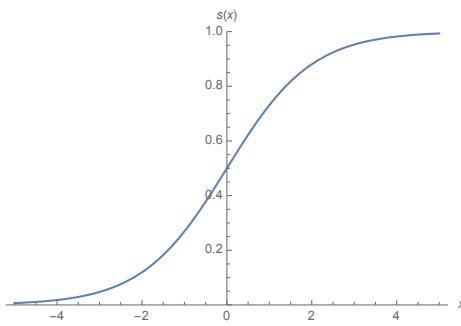


logloss0.pdf, loglosspt7.pdf [Plots of the loss $L(z, y)$ for $y = 0$ (left) and $y = 0.7$ (right). As you might guess, the left function is minimized at $z = 0$, and the right function is minimized at $z = 0.7$. These loss functions are always convex.]

$J(w)$ is convex! Solve by gradient descent.

[To do gradient descent, we'll need to compute some derivatives.]

$$\begin{aligned} s'(\gamma) &= \frac{d}{d\gamma} \frac{1}{1 + e^{-\gamma}} = \frac{e^{-\gamma}}{(1 + e^{-\gamma})^2} \\ &= s(\gamma)(1 - s(\gamma)) \end{aligned}$$



logistic.pdf, dlogistic.pdf [Plots of $s(\gamma)$ (left) and $s'(\gamma)$ (right).]

Let $s_i = s(X_i \cdot w)$

$$\begin{aligned}
 \nabla_w J &= -\sum \left(\frac{y_i}{s_i} \nabla s_i - \frac{1-y_i}{1-s_i} \nabla s_i \right) \\
 &= -\sum \left(\frac{y_i}{s_i} - \frac{1-y_i}{1-s_i} \right) s_i (1-s_i) X_i \\
 &= -\sum (y_i - s_i) X_i \\
 &= -X^\top (y - s(Xw)) \quad \text{where } s(Xw) = \begin{bmatrix} s_1 \\ s_2 \\ \vdots \\ s_n \end{bmatrix} \quad [\text{applies } s \text{ component-wise to } Xw]
 \end{aligned}$$

Gradient descent rule: $w \leftarrow w + \epsilon X^\top (y - s(Xw))$

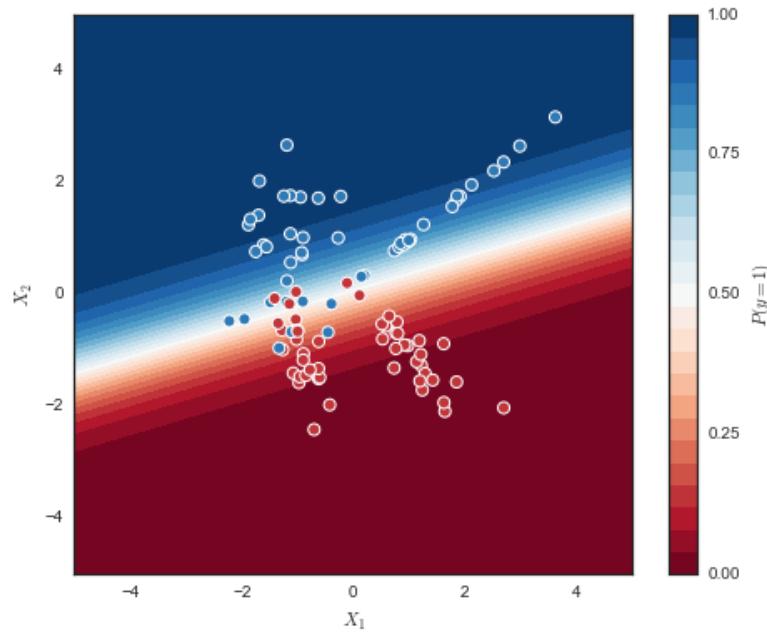
Stochastic gradient descent: $w \leftarrow w + \epsilon (y_i - s(X_i \cdot w)) X_i$

Works best if we shuffle points in random order, process one by one.

For very large n , sometimes converges before we visit all points!

[This looks a lot like the perceptron learning rule. The only difference is that the “ $-s_i$ ” part is new.]

Starting from $w = 0$ works well in practice.



problogistic.png, by “mwascom” of Stack Overflow

<http://stackoverflow.com/questions/28256058/plotting-decision-boundary-of-logistic-regression>

[An example of logistic regression.]

11 More Regression; Newton's Method; ROC Curves

LEAST-SQUARES POLYNOMIAL REGRESSION

Replace each X_i with feature vector $\Phi(X_i)$ with all terms of degree $0 \dots p$

$$\text{e.g., } \Phi(X_i) = [X_{i1}^2 \ X_{i1}X_{i2} \ X_{i2}^2 \ X_{i1} \ X_{i2} \ 1]^\top$$

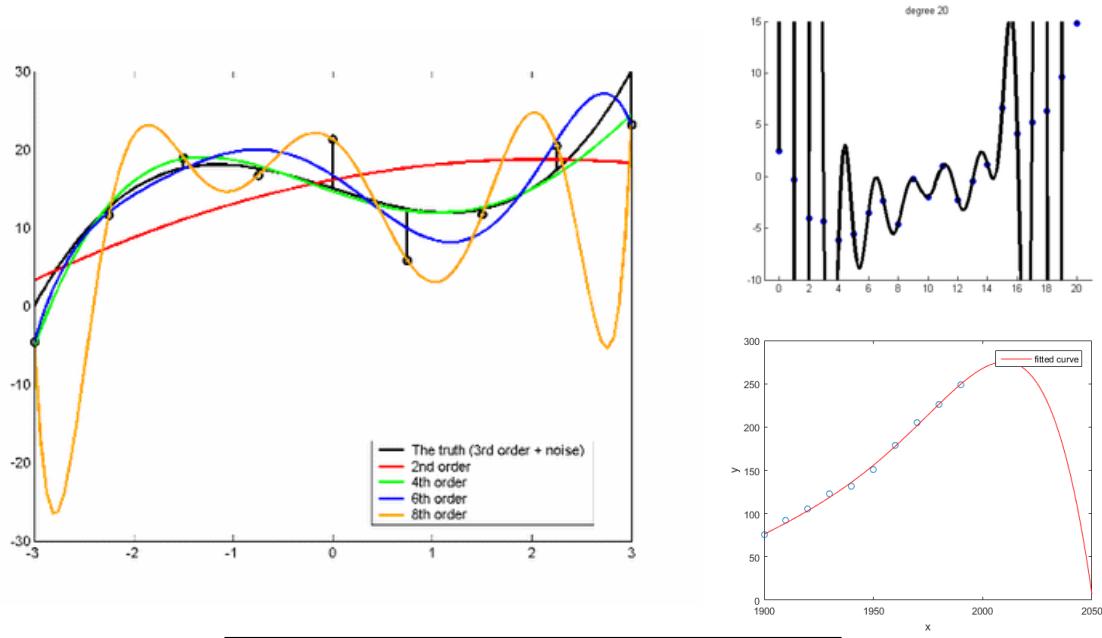
[Notice that we've added the fictitious dimension "1" here, so we don't need to add it again to do linear or logistic regression. This basis covers all polynomials quadratic in X_{i1} and X_{i2} .]

Can also use non-polynomial features (e.g., edge detectors).

Otherwise just like linear or logistic regression.

Log. reg. + quadratic features = same logistic posteriors as QDA.

Very easy to overfit!



overunder.png, degree20.png, UScensusquartic.png

[Here are some examples of polynomial overfitting, to show the importance of choosing the polynomial degree very carefully. At left, we have sampled points from a degree-3 curve (black) with added noise. We show best-fit polynomials of degrees 2, 4, 6, and 8 found by regression of the black points. The degree-4 curve (green) fits the true curve (black) well, whereas the degree-2 curve (red) underfits and the degree-6 and 8 curves (blue, yellow) overfit the noise and oscillate. The oscillations in the yellow degree-8 curve are a characteristic problem of polynomial interpolation.]

[At upper right, a degree-20 curve shows just how insane high-degree polynomial oscillations can get. It takes a great deal of densely spaced data to tame the oscillations in a high degree curve, and there isn't nearly enough data here.]

[At lower right, somebody has regressed a degree-4 curve to U.S. census population numbers. The curve doesn't oscillate, but can you nevertheless see a flaw? This shows the difficulty of *extrapolation* outside the range of the data. As a general rule, extrapolation is much harder than interpolation. The k -nearest neighbor classifier is one of the few that does extrapolation decently without occasionally returning crazy values.]

WEIGHTED LEAST-SQUARES REGRESSION

Linear regression fn (1) + squared loss fn (A) + cost fn (c).

[The idea of weighted least-squares is that some sample points might be more trusted than others, or there might be certain points you want to fit particularly well. So you assign those more trusted points a higher weight. If you suspect some points of being outliers, you can assign them a lower weight.]

Assign each sample pt a weight ω_i ; collect them in $n \times n$ diagonal matrix Ω .

Greater $\omega_i \rightarrow$ work harder to minimize $|\hat{y}_i - y_i|^2$ recall: $\hat{y} = Xw$ [\hat{y}_i is predicted label for X_i]

$$\boxed{\text{Find } w \text{ that minimizes } (Xw - y)^\top \Omega (Xw - y)} = \sum_{i=1}^n \omega_i (X_i \cdot w - y_i)^2$$

[As with ordinary least-squares regression, we find the minimum by setting the gradient to zero, which leads us to the normal equations.]

Solve for w in normal equations: $X^\top \Omega X w = X^\top \Omega y$

[Once again, you can interpret this method as a projection of y .]

Note: $\Omega^{1/2}\hat{y}$ is the pt nearest $\Omega^{1/2}y$ on subspace spanned by columns of $\Omega^{1/2}X$.

[If you stretch the n -dimensional space by applying the linear transformation $\Omega^{1/2}$, \hat{y} is an orthogonal projection of y onto the stretched subspace.]

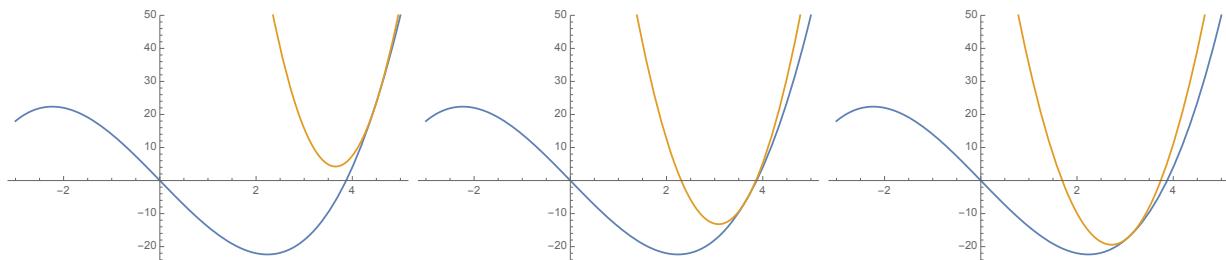
NEWTON'S METHOD

Iterative optimization method for smooth fn $J(w)$.

Often much faster than gradient descent. [We'll use Newton's method for logistic regression.]

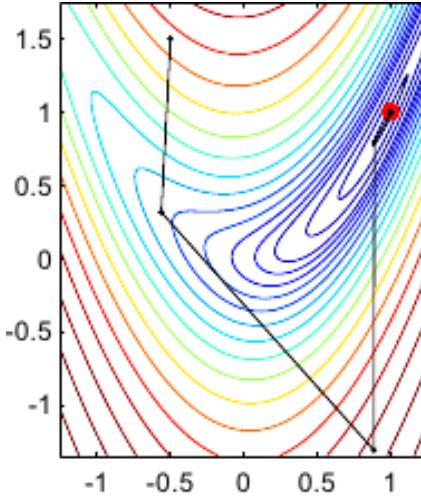
Idea: You're at point v . Approximate $J(w)$ near v by quadratic fn.

Jump to its unique critical pt. Repeat until bored.



[newton1.pdf](#), [newton2.pdf](#), [newton3.pdf](#)

[Three iterations of Newton's method in one-dimensional space. We seek the minimum of the blue curve, J . Each brown curve is a local quadratic approximation to J . Each iteration, we jump to the bottom of the brown parabola.]



newton2D.png [Steps taken by Newton's method in two-dimensional space.]

Taylor series about v :

$$\nabla J(w) = \nabla J(v) + (\nabla^2 J(v))(w - v) + O(|w - v|^2)$$

where $\nabla^2 J(v)$ is the Hessian matrix of J at v .

Find critical pt w by setting $\nabla J(w) = 0$:

$$w = v - (\nabla^2 J(v))^{-1} \nabla J(v)$$

[This is an iterative update rule you can repeat until it converges to a solution. As usual, we don't really want to compute a matrix inverse. Instead, we solve a linear system of equations, typically by Cholesky factorization or the conjugate gradient method.]

Newton's method:

```

pick starting point w
repeat until convergence
    e ← solution to linear system  $(\nabla^2 J(w)) e = -\nabla J(w)$ 
    w ← w + e
  
```

Warning: Doesn't know difference between minima, maxima, saddle pts.

Starting pt must be "close enough" to desired solution.

[If the objective function J is actually quadratic, Newton's method needs only one step to find the correct answer. The closer J is to quadratic, the faster Newton's method tends to converge.]

[Newton's method is superior to blind gradient descent for some optimization problems for several reasons. First, it tries to find the right step length to reach the minimum, rather than just walking an arbitrary distance downhill. Second, rather than follow the direction of steepest descent, it tries to choose a better descent direction.]

[Nevertheless, it has some major disadvantages. The biggest one is that computing the Hessian can be quite expensive, and it has to be recomputed every iteration. It can work well for low-dimensional weight spaces, but you would never use it for a neural network, because there are too many weights. Newton's method also doesn't work for most nonsmooth functions. It particularly fails for the perceptron risk function, whose Hessian is zero, except where the Hessian is not even defined.]

LOGISTIC REGRESSION (continued)

[Let's use Newton's method to solve logistic regression faster.]

$$\text{Recall: } s'(\gamma) = s(\gamma)(1 - s(\gamma)), \quad s_i = s(X_i \cdot w), \quad s = \begin{bmatrix} s_1 \\ s_2 \\ \vdots \\ s_n \end{bmatrix},$$

$$\nabla_w J = - \sum_{i=1}^n (y_i - s_i) X_i = -X^\top(y - s)$$

[Now let's derive the Hessian too, so we can use Newton's method.]

$$\nabla_w^2 J(w) = \sum_{i=1}^n s_i(1 - s_i) X_i X_i^\top = X^\top \Omega X \quad \text{where } \Omega = \begin{bmatrix} s_1(1 - s_1) & 0 & \dots & 0 \\ 0 & s_2(1 - s_2) & & 0 \\ \vdots & & \ddots & \vdots \\ 0 & 0 & \dots & s_n(1 - s_n) \end{bmatrix}$$

Ω is +ve definite $\forall w \Rightarrow X^\top \Omega X$ is +ve semidefinite $\forall w \Rightarrow J(w)$ is convex.

[The logistic regression cost function is convex, so Newton's method finds a globally optimal point if it converges at all.]

Newton's method:

```
w ← 0
repeat until convergence
  e ← solution to normal equations  $(X^\top \Omega X) e = X^\top(y - s)$            Recall:  $\Omega, s$  are fns of  $w$ 
  w ← w + e
```

[Notice that this looks a lot like weighted least squares, but the weight matrix Ω and the right-hand-side vector $y - s$ change every iteration. So we call it ...]

An example of iteratively reweighted least squares.

Ω prioritizes points with s_i near 0.5; tunes out pts near 0/1.

[In other words, sample points near the decision boundary have the biggest effect on the iterations. Meanwhile, the iterations move the decision boundary; in turn, that movement may change which points have the most influence. In the end, only the points near the decision boundary make a big contribution to the logistic fit.]

Idea: If n very large, save time by using a random subsample of the pts per iteration. Increase sample size as you go.

[The principle is that the first iteration isn't going to take you all the way to the optimal point, so why waste time looking at *all* the sample points? Whereas the last iteration should be the most accurate one.]

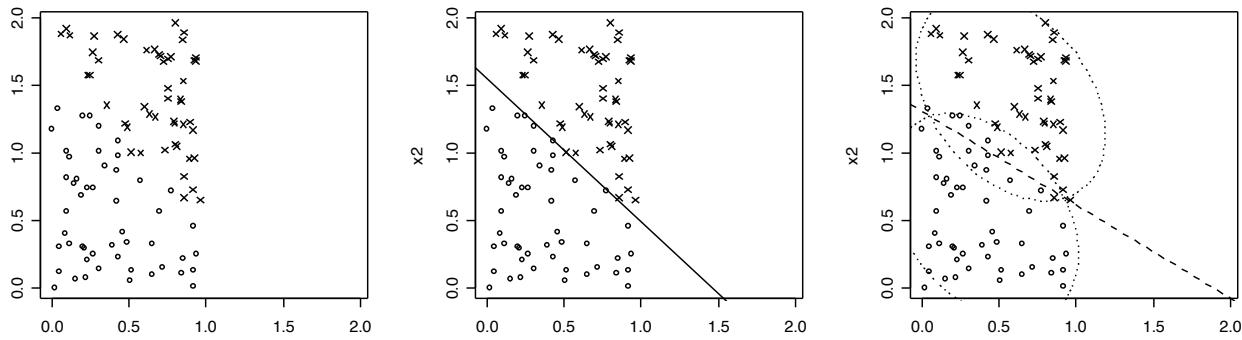
LDA vs. Logistic Regression

Advantages of LDA:

- For well-separated classes, LDA stable; log. reg. surprisingly unstable
- > 2 classes easy & elegant; log. reg. needs modifying (softmax regression)
- LDA slightly more accurate when classes nearly normal, especially if n is small

Advantages of log. reg.:

- More emphasis on decision boundary [though not as much as a hard-margin SVM]
 [When we analyzed Newton's method, we learned that points near the decision boundary have greater influence on the boundary. By contrast, LDA gives all the sample points equal weight when fitting Gaussians to them.]

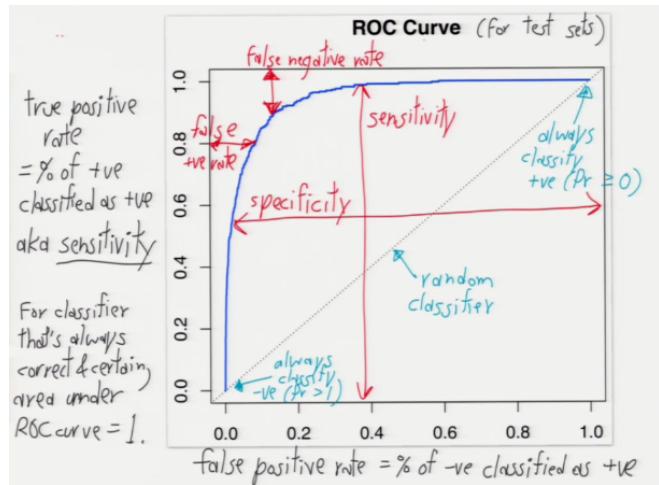
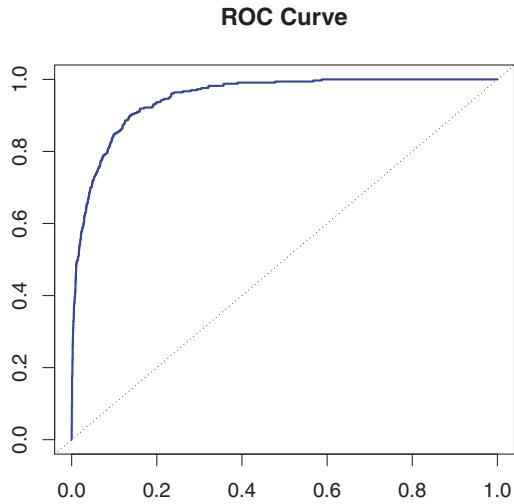


logregvsLDAuni.pdf [Logistic regression vs. LDA for a linearly separable data set with a very narrow margin. Logistic regression (center) succeeds in separating these classes, because points near the decision boundary have more influence. (Some luck is involved too; unlike a hard-margin SVM, logistic regression doesn't always separate separable points.)
 In this example, LDA (right) misclassifies some of the training points.]

- Hence less sensitive to most outliers
 [including outliers far out on the wrong side of the decision boundary ... compare with SVMs]
- Easy & elegant treatment of “partial” class membership; LDA pts are all-or-nothing
 [because logistic regression accommodates labels between 0 and 1]
- More robust on some non-Gaussian distributions (e.g., dists. w/large skew)

[When you use logistic regression with quadratic features, you get a quadric decision boundary, just as you do with QDA. Based on what I've said here, do you think logistic regression with quadratic features gives you exactly the same classifier as QDA?]

ROC CURVES (for test sets)



ROC.pdf

[This is a ROC curve. That stands for receiver operating characteristics, which is an awful name but we're stuck with it for historical reasons.]

A ROC curve is a way to evaluate your classifier after it is trained.

It is made by running a classifier on the test set or validation set.

It shows the rate of false positives vs. true positives for a range of settings.

We assume there is a knob we can turn to trade off these two types of error. For our purposes, that knob is the posterior probability threshold for Gaussian discriminant analysis or logistic regression.

However, neither axis of this plot is that knob.]

x-axis: “false positive rate = % of –ve classified as +ve”

y-axis: “true positive rate = % of +ve classified as +ve aka sensitivity”

“false negative rate”: vertical distance from curve to top [1 – sensitivity]

“specificity”: horizontal distance from curve to right [1 – false positive rate; “true negative rate”]

[You generate this curve by trying *every* probability threshold; for each threshold, measure the false positive & true positive rates and plot a point.]

upper right corner: “always classify +ve ($\text{Pr} \geq 0$)”

lower left corner: “always classify –ve ($\text{Pr} > 1$)”

diagonal: “random classifiers”

[A rough measure of a classifier’s effectiveness is the area under the curve. For a classifier that is always correct, the area under the curve is one. For the random classifier, the area under the curve is 1/2, so you’d better do better than that.]

[IMPORTANT: In practice, the trade-off between false negatives and false positives is usually negotiated by choosing a point on this plot, based on real test data, and NOT by taking the choice of threshold that’s best in theory.]

12 Statistical Justifications; the Bias-Variance Decomposition

STATISTICAL JUSTIFICATIONS FOR REGRESSION

[So far, I've talked about regression as a way to fit curves to points. Recall that early in the semester I divided machine learning into 4 levels: the application, the model, the optimization problem, and the optimization algorithm. My last two lectures about regression were at the bottom two levels: optimization. The cost functions that we optimize seem somewhat arbitrary. Today, let's take a step up to the second level, the model. I will describe some models, how they lead to those optimization problems, and how they contribute to underfitting or overfitting.]

Typical model of reality:

- sample points come from unknown prob. distribution: $X_i \sim D$
- y -values are sum of unknown, non-random fn + random noise:
 $\forall X_i, \quad y_i = f(X_i) + \epsilon_i, \quad \epsilon_i \sim D', \quad D' \text{ has mean zero}$

[We are positing that reality is described by a function f . We don't know f , but f is not a random variable; it represents a consistent relationship between X and y that we can estimate. We add to f a random variable ϵ , which represents measurement errors and all the other sources of statistical error when we measure real-world phenomena. Notice that the noise is independent of X . That's a pretty big assumption, and often it does not apply in practice, but that's all we'll have time to deal with this semester. Also notice that this model leaves out systematic errors, like when your measuring device adds one to every measurement, because we usually can't diagnose systematic errors from data alone.]

Goal of regression: find h that estimates f .

Ideal approach: choose $h(x) = \underbrace{\mathbb{E}_Y[Y|X=x]}_{\text{expectation}} = f(x) + \mathbb{E}[\epsilon] = f(x)$

[If this expectation exists at all, it partly justifies our model of reality. We can retroactively define f to be this expectation.]

[Draw figure showing example f , distribution for a fixed x .]

Least-Squares Regression from Maximum Likelihood

Suppose $\epsilon_i \sim \mathcal{N}(0, \sigma^2)$; then $y_i \sim \mathcal{N}(f(X_i), \sigma^2)$

Recall that log likelihood for normal dist. is

$$\ln P(y_i) = -\frac{(y_i - \mu)^2}{2\sigma^2} - \text{constant}, \quad \Leftrightarrow \mu = f(X_i)$$

$$\ell(f; X, y) = \ln(P(y_1)P(y_2)\dots P(y_n)) = \ln P(y_1) + \dots + \ln P(y_n) = -\frac{1}{2\sigma^2} \sum (y_i - f(X_i))^2 - \text{constant}$$

[We can use maximum likelihood to decide what to do, but now the "parameter" we want to estimate is f .]

Takeaway: Max likelihood on "parameter" $f \Rightarrow$ estimate f by least-squares

[So if the noise is normally distributed, maximum likelihood justifies using the least-squares cost function.] [However, I've told you in previous lectures that least-squares is very sensitive to outliers. If the error is truly normally distributed, that's not a big deal, especially when you have a lot of sample points. But in the real world, the distribution of outliers often isn't normal. Outliers might come from wrongly measured measurements, data entry errors, anomalous events, or just not having a normal distribution. When you have a heavy-tailed distribution, for example, least-squares isn't a good choice.]

Empirical Risk

The risk for hypothesis h is expected loss $R(h) = E[L]$ over all $x \in \mathbb{R}^d, y \in \mathbb{R}$.

Discriminative model: we don't know X 's dist. D . How can we minimize risk?

[If we have a generative model, we can estimate the joint probability distribution for X and Y and derive the expected loss. That's what we did for Gaussian discriminant analysis. But today I'm assuming we don't have a generative model, so we don't know those probabilities. Instead, we're going to approximate the distribution in a very crude way: we pretend that the sample points *are* the distribution.]

Empirical distribution: the **discrete** uniform distribution over the sample pts

Empirical risk: expected loss under empirical distribution

$$\hat{R}(h) = \frac{1}{n} \sum_{i=1}^n L(h(X_i), y_i)$$

[The hat on the R indicates it's only a cheap approximation of the true, unknown statistical risk we really want to minimize. Often, this is the best we can do. For many but not all distributions, it converges to the true risk in the limit as $n \rightarrow \infty$. Choosing h that minimizes \hat{R} is called empirical risk minimization.]

Takeaway: this is why we [usually] minimize the sum of loss fns.

Logistic Loss from Maximum Likelihood

What cost fn should we use for probabilities?

Actual probability pt X_i is in the class is y_i ; predicted prob. is $h(X_i)$.

Imagine β duplicate copies of X_i : $y_i\beta$ are in the class, $(1 - y_i)\beta$ are not.

[The size of β isn't very important, but imagine that $y_i\beta$ and $(1 - y_i)\beta$ are both integers for all i .]

[If we use maximum likelihood estimation to choose the weights most likely to generate these duplicated samples, we get the following likelihood.]

$$\text{Likelihood is } \mathcal{L}(h; X, y) = \prod_{i=1}^n h(X_i)^{y_i\beta} (1 - h(X_i))^{(1-y_i)\beta}$$

$$\begin{aligned} \text{Log likelihood is } \ell(h) &= \ln \mathcal{L}(h) \\ &= \beta \sum_i \left(y_i \ln h(X_i) + (1 - y_i) \ln(1 - h(X_i)) \right) \\ &= -\beta \sum \text{logistic loss fn } L(h(X_i), y_i). \end{aligned}$$

Max log likelihood \Rightarrow minimize $\sum L(h(X_i), y_i)$.

[So the principle of maximum likelihood explains where the weird logistic loss function comes from.]

THE BIAS-VARIANCE DECOMPOSITION

There are 2 sources of error in a hypothesis h :

bias: error due to inability of hypothesis h to fit f perfectly

e.g. fitting quadratic f with a linear h

variance: error due to fitting random noise in data

e.g. we fit linear f with a linear h , yet $h \neq f$.

Model: $X_i \sim D, \epsilon_i \sim D', y_i = f(X_i) + \epsilon_i$ [remember that D' has mean zero]

fit hypothesis h to X, y

Now h is a random variable; i.e. its weights are random

Consider arbitrary pt $z \in \mathbb{R}^d$ (not necessarily a sample pt!) & $\gamma = f(z) + \epsilon$.

[So z is *arbitrary*, whereas γ is *random*.]

Note: $E[\gamma] = f(z)$; $\text{Var}(\gamma) = \text{Var}(\epsilon)$ [the mean comes from f , and the variance comes from ϵ]

Risk fn when loss = squared error:

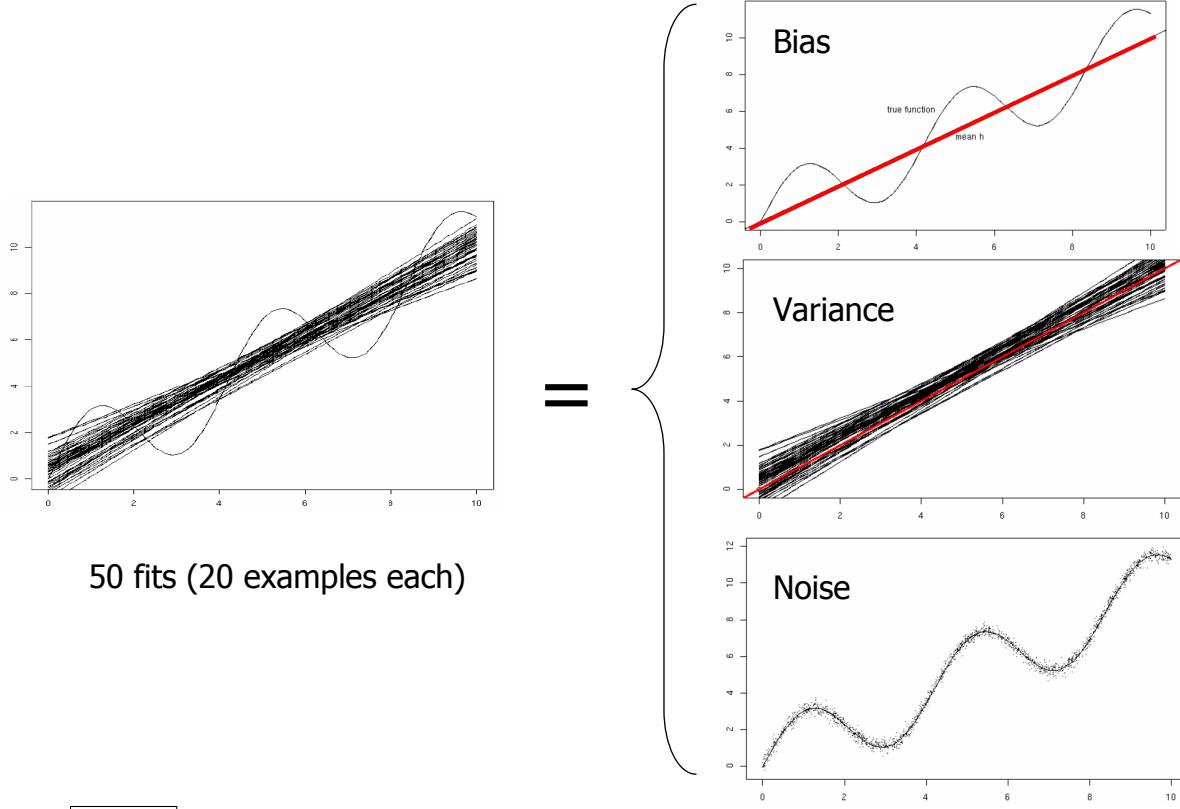
$$R(h) = E[L(h(z), \gamma)]$$

↑ take expectation over possible training sets X, y & values of γ

[Stop and take a close look at this expectation. Remember that the hypothesis h is a random variable. We are taking a mean over the probability distribution of hypotheses. That seems pretty weird if you've never seen it before. But remember, the training data X and y come from probability distributions. We use the training data to choose weights, so the weights that define h also come from some probability distribution. It might be hard to work out what that distribution is, but it exists. This “ $E[\cdot]$ ” is integrating the loss over all possible values of the weights.]

$$\begin{aligned} &= E[(h(z) - \gamma)^2] \\ &= E[h(z)^2] + E[\gamma^2] - 2E[\gamma h(z)] \quad [\text{Observe that } \gamma \text{ and } h(z) \text{ are independent}] \\ &= \text{Var}(h(z)) + E[h(z)]^2 + \text{Var}(\gamma) + E[\gamma]^2 - 2E[\gamma]E[h(z)] \\ &= (E[h(z)] - E[\gamma])^2 + \text{Var}(h(z)) + \text{Var}(\gamma) \\ &= \underbrace{(E[h(z)] - f(z))^2}_{\text{bias}^2 \text{ of method}} + \underbrace{\text{Var}(h(z))}_{\text{variance of method}} + \underbrace{\text{Var}(\epsilon)}_{\text{irreducible error}} \end{aligned}$$

[This is called the *bias-variance decomposition* of the risk function. Let's look at an intuitive interpretation of these three parts.]



bvn.pdf [In this example, we're trying to fit a sine wave with lines, which obviously aren't going to be accurate. At left, we have generated 50 different hypotheses (lines). At upper right, the red line is the *expected hypothesis*—an average over infinitely many hypotheses. The black curve illustrates test points on the true function f . We see that most test points have a large bias (squared difference between the black and red curves), because lines don't fit sine waves well. However, some of the test points happen to have a small bias—where the sine wave crosses the red line. At center right, the variance is the expected squared difference between a random black line and the red line. At lower right, the irreducible error is the expected squared difference between a random test point and the sine wave.]

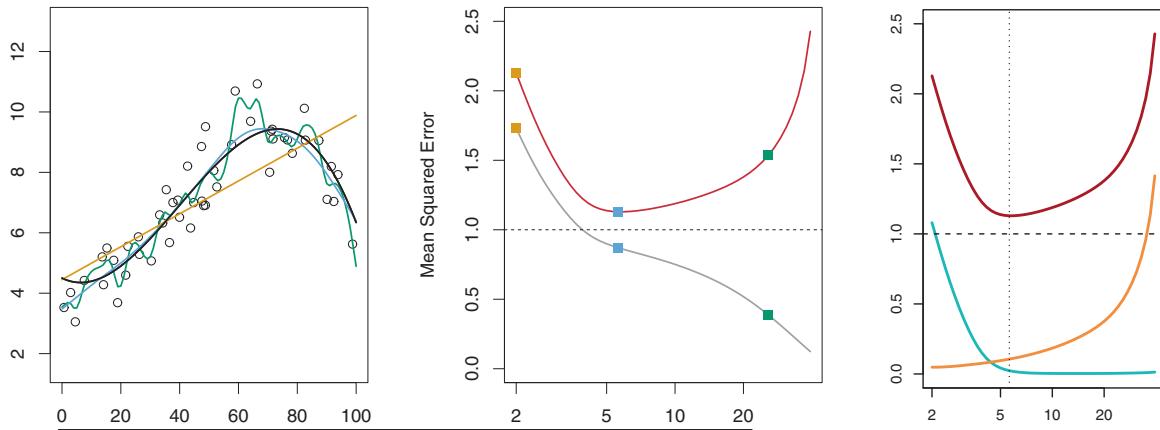
This is pointwise version [of the bias-variance decomposition.]

Mean version: let $z \sim D$ be random variable; take mean over D of bias², variance.

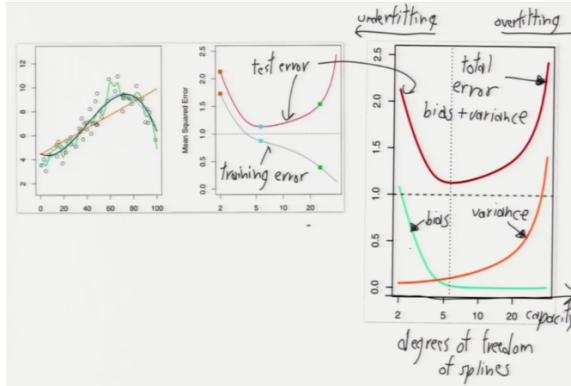
[So you can decompose one test point's error into these three components, or you can decompose the error of the hypothesis over its entire range into three components, which tells you roughly how big they'll be on a large test set.]

[Now I will write down a list of consequences of what we've just learned.]

- Underfitting = too much bias
- Most overfitting caused by too much variance
- Training error reflects bias but not variance; test error reflects both [which is why low training error can fool you when you've overfitted]
- For many distributions, variance $\rightarrow 0$ as $n \rightarrow \infty$
- If h can fit f exactly, for many distributions bias $\rightarrow 0$ as $n \rightarrow \infty$
- If h cannot fit f well, bias is large at “most” points
- Adding a good feature reduces bias; adding a bad feature rarely increases it
- Adding a feature usually increases variance [don't add a feature unless it reduces bias more]
- Can't reduce irreducible error [hence its name]
- Noise in test set affects only $\text{Var}(\epsilon)$;
noise in training set affects only bias & $\text{Var}(h)$
- For real-world data, f is rarely knowable [& noise model might be wrong]
[so we can't actually put numbers to the bias-variance decomposition on real-world data]
- But we can test learning algs by choosing f & making synthetic data



splinefit.pdf, biasvarspline.pdf (ISL, Figures 2.9 and 2.12) [At left, a data set is fit with splines having various degrees of freedom. The synthetic data is taken from the black curve with added noise. At center, we plot training error (gray) and test error (red) as a function of the number of degrees of freedom. At right, we plot the squared test error as a sum of squared bias (blue) and variance (orange). As the number of degrees of freedom increases, the training and test errors both decrease up to degree 6 because the bias decreases, but for higher degrees the test error increases because the variance increases.]



Example: Least-Squares Linear Reg.

For simplicity, no fictitious dimension.

[This implies that our linear regression function has to be zero at the origin.]

Model: $f(z) = v^T z$ (reality is linear)

[So we could fit f perfectly with a linear h if not for the noise in the training set.]

Let e be noise n -vector, $e_i \sim \mathcal{N}(0, \sigma^2)$

Training labels: $y = Xv + e$

[X & y are the inputs to linear regression. We don't know v or e .]

Lin. reg. computes weights

$$w = X^+y = X^+(Xv + e) = v + \underbrace{X^+e}_{\text{noise in weights}} \quad [\text{We want } w = v, \text{ but the noise in } y \text{ becomes noise in } w.]$$

BIAS is $E[h(z)] - f(z) = E[w^T z] - v^T z = E[z^T X^+ e] = z^T X^+ E[e] = 0$

Warning: This does not mean $h(z) - f(z)$ is always 0!

Sometimes +ve, sometimes -ve, mean over training sets is 0.

[Those deviations from the mean are captured in the variance.]

[When the bias is zero, a perfect fit is possible. But when a perfect fit is possible, not all learning methods give you a bias of zero; here it's a benefit of the squared error loss function. With a different loss function, we might have a nonzero bias even fitting a linear h to a linear f .]

VARIANCE is $\text{Var}(h(z)) = \text{Var}(w^T z) = \text{Var}(z^T v + z^T X^+ e) = \text{Var}(z^T X^+ e)$

[This is the dot product of a vector $z^T X^+$ with an isotropic, normally distributed vector e . The dot product reduces it to a one-dimensional Gaussian along the direction $z^T X^+$, so this variance is just the variance of the 1D Gaussian times the squared length of the vector $z^T X^+$.]

$$\begin{aligned} &= \sigma^2 |z^T X^+|^2 = \sigma^2 z^T (X^T X)^{-1} X^T X (X^T X)^{-1} z \\ &= \sigma^2 z^T (X^T X)^{-1} z \end{aligned}$$

If we choose coordinate system so $E[X] = 0$,

then $X^T X \rightarrow n \text{Cov}(D)$ as $n \rightarrow \infty$, so one can show that for $z \sim D$,

$$\text{Var}(h(z)) \approx \sigma^2 \frac{d}{n}$$

[where d is the dimension—the number of features per sample point.]

Takeaways: Bias can be zero when hypothesis function can fit the real one!

[This is a nice property of the squared error loss function.]

Variance portion of RSS (overfitting) decreases as $1/n$ (sample points),
increases as d (features)
or $O(d^p)$ if you use degree- p polynomials.

[I've used linear regression because it's a relatively simple example. But this bias-variance trade-off applies to many learning algorithms, including classification as well as regression. But for most learning algorithms, the math gets a lot more complicated than this, if you can do it at all. Sometimes there are multiple competing bias-variance models and no consensus on which is the right one.]

13 Shrinkage: Ridge Regression, Subset Selection, and Lasso

RIDGE REGRESSION aka Tikhonov Regularization

(1) + (A) + ℓ_2 penalized mean loss (d).

$$\boxed{\text{Find } w \text{ that minimizes } |Xw - y|^2 + \lambda |w'|^2} = J(w)$$

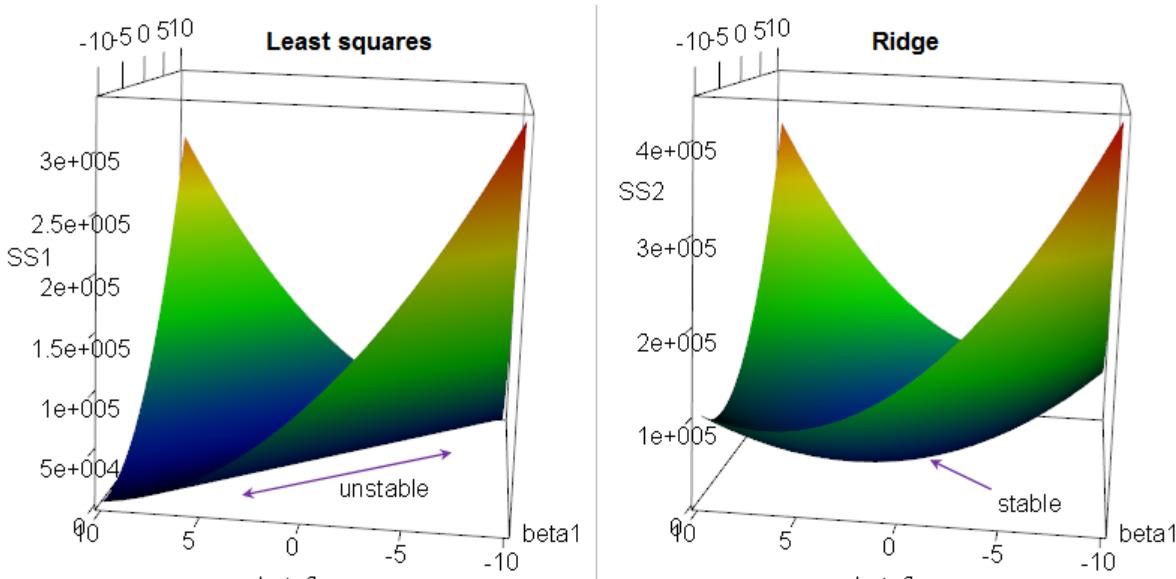
where w' is w with component α replaced by 0.

X has fictitious dimension but we DON'T penalize α .

Adds a regularization term, aka a penalty term, for shrinkage: to encourage small $|w'|$. Why?

- Guarantees positive definite normal eq'n's; always unique solution.

[Standard least-squares linear regression yields singular normal equations when the sample points lie on a common hyperplane in feature space.] E.g., when $d > n$.



`ridgequad.png` [The cost function $J(w)$ with and without regularization.]

[At left, we see a quadratic form for a positive semidefinite cost function associated with least-squares regression. This cost function has many minima, and the regression problem is said to be ill-posed. By adding a small penalty term, we obtain a positive definite quadratic form (right), which has one unique minimum. The term “regularization” implies that we are turning an ill-posed problem into a well-posed problem.]

[That was the original motivation, but the next has become more important ...]

- Reduces overfitting by reducing variance. Why?

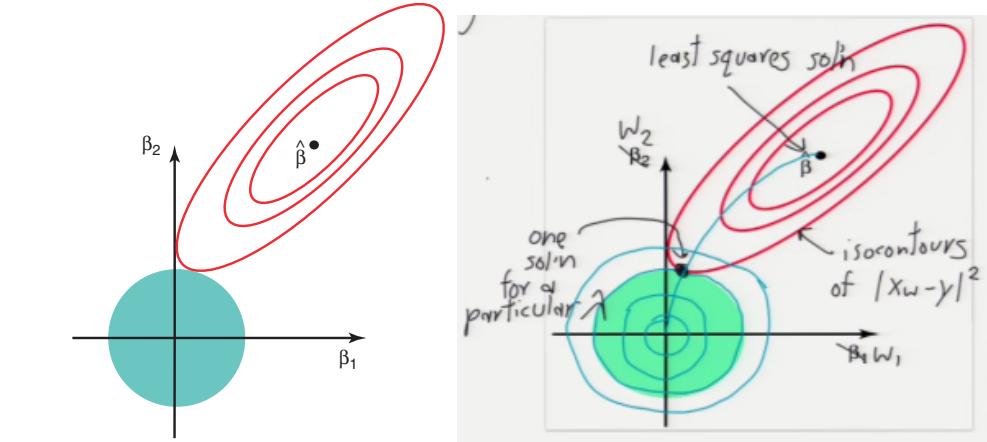
Imagine: $500x_1 - 500x_2$ is best fit for well-separated points with $y_i \in [0, 1]$.

Small change in $x \Rightarrow$ big change in y !

[Given that all the y values in the data are small, it's a sure sign of overfitting if tiny changes in x cause huge changes in y .]

So we penalize large weights.

[This use of regularization is closely related to the first one. When you have large variance and a lot of overfitting, it implies that your problem is *close to* being ill-posed, even though technically it might be well-posed.]



ridgeterms.pdf (ISL, Figure 6.7) [In this plot of weight space, $\hat{\beta}$ (read as “ \hat{w} ”) is the least-squares solution. The red ellipses are the isocontours of $|Xw - y|^2$. The isocontours of $|w|$ are circles centered at the origin (blue). The solution lies where a red isocontour just touches a blue isocontour tangentially. As λ increases, the solution will occur at a more outer red isocontour and a more inner blue isocontour. This helps to reduce overfitting.]

Setting $\nabla J = 0$ gives normal eq’ns

$$(X^T X + \lambda I') w = X^T y$$

where I' is identity matrix w/bottom right set to zero. [Don’t penalize the bias term α .]

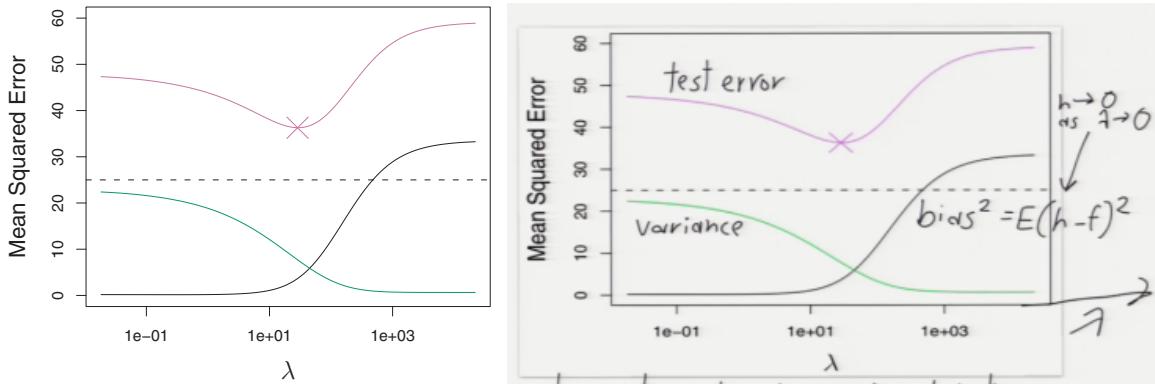
Algorithm: Solve for w . Return $h(z) = w^T z$.

Increasing $\lambda \Rightarrow$ more regularization; smaller $|w'|$

Recall [from the previous lecture] our data model $y = Xv + e$, where e is noise.

Variance of ridge reg. is $\text{Var}(z^T(X^T X + \lambda I')^{-1} X^T e)$.

As $\lambda \rightarrow \infty$, variance $\rightarrow 0$, but bias increases.



ridgebiasvar.pdf (ISL, Figure 6.5) [Plot of bias² & variance as λ increases.]

[So, as usual for the bias-variance trade-off, the test error as a function of λ is a U-shaped curve. We find the bottom by validation.]

λ is a hyperparameter; tune by (cross-)validation.

Ideally, features should be “normalized” to have same variance.

Alternative: use asymmetric penalty by replacing I' w/other diagonal matrix. [For example, if you use polynomial features, you could use different penalties for monomials of different degrees.]

Bayesian Justification for Ridge Reg.

Assign a prior probability on w' : a Gaussian $P(w') \sim \mathcal{N}(0, \sigma^2)$. Apply MLE to the posterior prob.
[This prior probability says that we think weights close to zero are more likely to be correct.]

$$\begin{aligned} \text{Bayes' Theorem: posterior } P(w|X, y) &= \frac{P(y|X, w) \times \text{prior } P(w')}{P(y|X)} = \frac{\mathcal{L}(w) P(w')}{P(y|X)} \\ \text{Maximize log posterior} &= \ln \mathcal{L}(w) + \ln P(w') - \text{const} \\ &= -\text{const} |Xw - y|^2 - \text{const} |w'|^2 - \text{const} \\ &\Rightarrow \text{Minimize } |Xw - y|^2 + \lambda |w'|^2 \end{aligned}$$

This method (using likelihood, but maximizing posterior) is called maximum *a posteriori* (MAP).
[A prior probability on the weights is another way to understand regularizing ill-posed problems.]

SUBSET SELECTION

[Some of you may have noticed as early as Homework 1 that you can sometimes get better performance on a spam classifier simply by dropping some useless features.]

All features increase variance, but not all features reduce bias.

Idea: Identify poorly predictive features, ignore them (weight zero).

Less overfitting, smaller test errors.

2nd motivation: Inference. Simpler models convey interpretable wisdom.

Useful in all classification & regression methods.

Sometimes it's hard: Different features can partly encode same information.

Combinatorially hard to choose best feature subset.

Alg: Best subset selection. Try all $2^d - 1$ nonempty subsets of features. [Train one classifier per subset.]
Choose the best classifier by (cross-)validation. Slow.

[Obviously, best subset selection isn't feasible if we have a lot of features. But it gives us an "ideal" algorithm to compare practical algorithms with. If d is large, there is no algorithm that's both guaranteed to find the best subset and that runs in acceptable time. But heuristics often work well.]

Heuristic 1: Forward stepwise selection.

Start with null model (0 features); repeatedly add best feature until validation errors start increasing (due to overfitting) instead of decreasing. At each outer iteration, inner loop tries every feature and chooses the best by validation. Requires training $O(d^2)$ models instead of $O(2^d)$.

Not perfect: e.g., won't find the best 2-feature model if neither of those
features yields the best 1-feature model. [That's why it's a heuristic.]

Heuristic 2: Backward stepwise selection.

Start with all d features; repeatedly remove feature whose removal gives best reduction in validation error.

Also trains $O(d^2)$ models.

Additional heuristic: Only try to remove features with small weights.

Q: small relative to what?

Recall: variance of least-squ. regr. is proportional to $\sigma^2(X^\top X)^{-1}$

z-score of weight w_i is $z_i = \frac{w_i}{\sigma \sqrt{v_i}}$ where v_i is i th diagonal entry of $(X^\top X)^{-1}$.

Small z-score hints "true" w_i could be zero.

[Forward stepwise is a better choice when you suspect only a few features will be good predictors. Backward stepwise is better when you suspect you'll keep most of the features. If you're lucky, you'll stop early.]

LASSO (Robert Tibshirani, 1996)

Regression w/regularization: (1) + (A) + ℓ_1 penalized mean loss (e).

“Least absolute shrinkage and selection operator”

[This is a regularized regression method similar to ridge regression, but it has the advantage that it often naturally sets some of the weights to zero.]

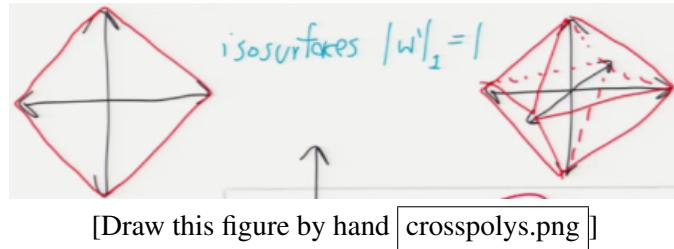
Find w that minimizes $|Xw - y|^2 + \lambda \|w'\|_1$

where $\|w'\|_1 = \sum_{i=1}^d |w_i|$ (Don't penalize α .)

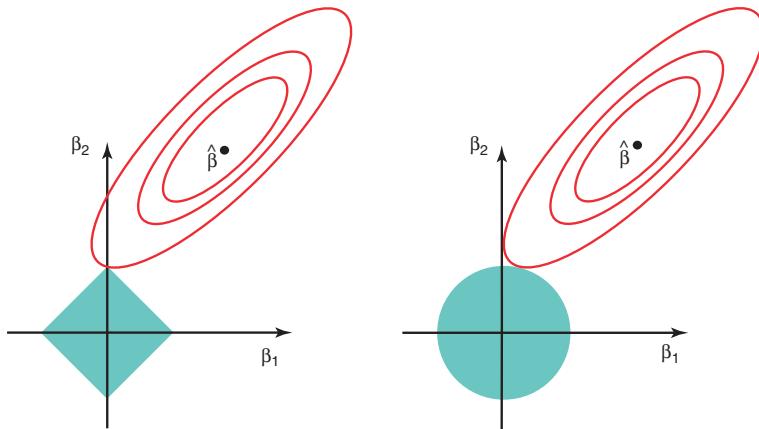
Recall ridge regr.: isosurfaces of $|w'|^2$ are hyperspheres.

The isosurfaces of $\|w'\|_1$ are cross-polytopes.

The unit cross-polytope is the convex hull of all the positive & negative unit coordinate vectors.



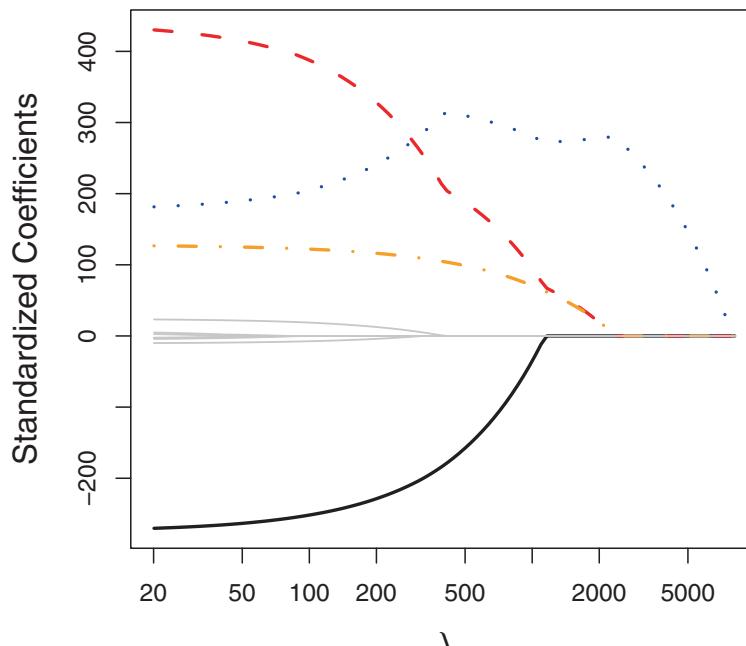
[You get larger and smaller cross-polytope isosurfaces by scaling these.]



[lassoridge.pdf](#) [Isocontours of the terms of the objective function for the Lasso appear at left. The ridge regression terms appear at right for comparison.]

[The red ellipses are the isocontours of $|Xw - y|^2$, and the least-squares solution lies at their center. The isocontours of $\|w'\|_1$ are diamonds centered at the origin (blue). The solution lies where a red isocontour just touches a blue diamond. What's interesting here is that in this example, the red isocontour touches just the tip of the diamond. So the weight w_1 gets set to zero. That's what we want to happen to weights that don't have enough influence. This doesn't always happen—for instance, the red isosurface could touch a side of the diamond instead of a tip of the diamond.]

[When you go to higher dimensions, you might have several weights set to zero. For example, in 3D, if the red isosurface touches a sharp vertex of the cross-polytope, two of the three weights get set to zero. If it touches a sharp edge of the cross-polytope, one weight gets set to zero. If it touches a flat side of the cross-polytope, no weight is zero.]



lassoweights.pdf (ISL, Figure 6.6) [Weights as a function of λ .]

[This shows the weights for a typical linear regression problem with about 10 variables. You can see that as λ increases, more and more of the weights become zero. Only four of the weights are really useful for prediction; they're in color. Statisticians used to choose λ by looking at a chart like this and trying to eyeball a spot where there aren't too many predictors and the weights aren't changing too fast. But nowadays they prefer cross-validation.]

Sometimes sets some weights to zero, especially for large λ .

Algs: subgradient descent, least-angle regression (LARS), forward stagewise

[Lasso can be reformulated as a quadratic program, but it's a quadratic program with 2^d constraints, because a d -dimensional cross-polytope has 2^d facets. In practice, special-purpose optimization methods have been developed for Lasso. I'm not going to teach you one, but if you need one, look up the last two of these algorithms. LARS is built into the R Programming Language for statistics.]

[As with ridge regression, you should probably normalize the features first before applying Lasso.]

14 The Kernel Trick

KERNELS

Recall: with d input features, degree- p polynomials blow up to $O(d^p)$ features.

[When d is large, this gets computationally intractable really fast.]

As I said in Lecture 4, if you have 100 features per feature vector and you want to use degree-4 decision functions, then each lifted feature vector has a length of roughly 4 million.]

Today, magically, we use those features without computing them!

Observation: In many learning algos,

- the weights can be written as a linear combo of sample points, &
- we can use inner products of $\Phi(x)$'s only \Rightarrow don't need to compute $\Phi(x)$!

$$\text{Suppose } w = X^\top a = \sum_{i=1}^n a_i X_i \quad \text{for some } a \in \mathbb{R}^n.$$

Substitute this identity into alg. and optimize n dual weights a (aka dual parameters) instead of $d + 1$ (or d^p) primal weights w .

Kernel Ridge Regression

Center X and y so their means are zero: $X_i \leftarrow X_i - \mu_X$, $y_i \leftarrow y_i - \mu_y$

This lets us replace I' with I in normal equations:

$$(X^\top X + \lambda I)w = X^\top y$$

[To dualize ridge regression, we need the weights to be a linear combination of the sample points. Unfortunately, that only happens if we penalize the bias term $w_{d+1} = \alpha$, as these normal equations do. Fortunately, when we center X and y , the “expected” value of the intercept is zero. The actual value won’t usually be exactly zero, but it will often be close enough that we won’t do much harm by penalizing the intercept.]

$$w = \frac{1}{\lambda}(X^\top y - X^\top Xw) = X^\top a \quad \text{where } a = \frac{1}{\lambda}(y - Xw)$$

This shows that w is a linear combo of sample points! To compute a :

$$\lambda a = y - Xw = y - XX^\top a \Rightarrow a = (XX^\top + \lambda I)^{-1}y$$

a is a dual solution; solves the dual form of ridge regression:

Find a that minimizes $|XX^\top a - y|^2 + \lambda|X^\top a|^2$

[We obtain this dual form by substituting $w = X^\top a$ into the original ridge regression cost function.]

Regression fn is

$$h(z) = w^\top z = a^\top Xz = \sum_{i=1}^n a_i (X_i^\top z) \quad \Leftarrow \text{weighted sum of inner products}$$

Let $k(x, z) = x^\top z$ be kernel fn.

[Later, we'll replace x and z with $\Phi(x)$ and $\Phi(z)$, and that's where the magic will happen.]

Let $K = XX^\top$ be $n \times n$ kernel matrix. Note $K_{ij} = k(X_i, X_j)$.

K is singular if $n > d + 1$. [And sometimes even if it's not.]

In that case, no solution if $\lambda = 0$. [Then we must choose a positive λ . But that's okay.]

Dual ridge reg. alg:

$$\begin{array}{lll} \forall i, j, & K_{ij} \leftarrow k(X_i, X_j) & \Leftarrow O(n^2d) \text{ time} \\ \text{Solve } (K + \lambda I)a = y \text{ for } a & & \Leftarrow O(n^3) \text{ time} \\ \text{for each test pt } z & & \\ h(z) \leftarrow \sum_{i=1}^n a_i k(X_i, z) & & \Leftarrow O(nd) \text{ time} \end{array}$$

Does not use X_i directly! Only k . [This will become important soon.]

Dual: solve $n \times n$ linear system, $O(n^3 + n^2d)$ time

Primal: " $d \times d$ ", $O(d^3 + d^2n)$ time

We prefer dual when $d > n$. [Moreover, if we add polynomial terms as new features, the d in the primal running time increases, but we will see that the d in the kernelized dual running time does not increase.]

The Kernel Trick (aka Kernelization)

[Here's the magic part. We will see that we can compute a polynomial kernel that involves many monomial terms without actually computing those terms.]

The polynomial kernel of degree p is $k(x, z) = (x^\top z + 1)^p$

Theorem: $(x^\top z + 1)^p = \Phi(x)^\top \Phi(z)$ where $\Phi(x)$ contains every monomial in x of degree $0 \dots p$.

Example for $d = 2, p = 2$:

$$\begin{aligned} (x^\top z + 1)^2 &= x_1^2 z_1^2 + x_2^2 z_2^2 + 2x_1 z_1 x_2 z_2 + 2x_1 z_1 + 2x_2 z_2 + 1 \\ &= [x_1^2 \quad x_2^2 \quad \sqrt{2}x_1 x_2 \quad \sqrt{2}x_1 \quad \sqrt{2}x_2 \quad 1] [z_1^2 \quad z_2^2 \quad \sqrt{2}z_1 z_2 \quad \sqrt{2}z_1 \quad \sqrt{2}z_2 \quad 1]^\top \\ &= \Phi(x)^\top \Phi(z) \end{aligned} \quad [\text{This is how we're defining } \Phi.]$$

[Notice the factors of $\sqrt{2}$. If you try a higher polynomial degree p , you'll see a wider variety of these constants. We have no control of the constants that appear in $\Phi(x)$, but they don't matter very much, because the primal weights w will scale themselves to compensate. Even though we won't be directly computing the primal weights ... they still implicitly exist.]

Key win: compute $\Phi(x)^\top \Phi(z)$ in $O(d)$ time instead of $O(d^p)$, even though $\Phi(x)$ has length $O(d^p)$.

Kernel ridge regression replaces X_i with $\Phi(X_i)$:

$$\text{Let } k(x, z) = \Phi(x)^\top \Phi(z) \quad \text{But don't compute } \Phi(x) \text{ or } \Phi(z); \text{ compute } k(x, z) = (x^\top z + 1)^p !$$

[I think what we've done here is pretty mind-blowing: we can now do polynomial regression with an exponentially long, high-order polynomial in less time than it would take even to compute the final polynomial. The running time is sublinear, actually much smaller than linear, in the length of the Φ vectors.]

Kernel Perceptrons

Featurized perceptron alg:

```
w ← y1 Φ(X1) [starting point is arbitrary, but can't be zero]
while some yi Φ(Xi) · w < 0
    w ← w + ε yi Φ(Xi)
for each test pt z
    h(z) ← w · Φ(z)
```

Let $\Phi(X)$ be $n \times D$ matrix with rows $\Phi(X_i)^\top$, $D = \text{length of } \Phi(\cdot)$, $K = \Phi(X) \Phi(X)^\top$

Dualize with $w = \Phi(X)^\top a$. Then the code “ $a_i \leftarrow a_i + \epsilon y_i$ ” has same effect as “ $w \leftarrow w + \epsilon y_i \Phi(X_i)$ ”

[So the dual weight a_i records what multiple of sample point i the perceptron algorithm has added to w .]

$$\Phi(X_i) \cdot w = (\Phi(X)w)_i = (\Phi(X) \Phi(X)^\top a)_i = (Ka)_i$$

Dual perceptron alg:

```
a ← [y1 0 ... 0]^\top [starting point is arbitrary, but can't be zero]
∀i, j, Kij ← k(Xi, Xj) ⇐ O(n2d) time (kernel trick)
while some yi (Ka)i < 0
    ai ← ai + ε yi ⇐ O(1) time; update Ka in O(n) time
for each test pt z
    h(z) ← ∑j=1n aj k(Xj, z) ⇐ O(nd) time [kernel trick]
```

[A big deal is that the running times depend on the original d , not on the length D of $\Phi(\cdot)$!]

OR we can compute $w = \Phi(X)^\top a$ once in $O(nD)$ time & evaluate test pts in $O(D)$ time/pt

[... which is a win if the numbers of training points and test points both exceed D/d .]

Kernel Logistic Regression

[The stochastic gradient descent step for logistic regression is just a small modification of the step for perceptrons. But recall that we're no longer looking for misclassified sample points. Instead, we apply the gradient descent rule to sample points in a stochastic, random order—or, alternatively, to all the points at once.]

Stochastic gradient descent step:

$$a_i \leftarrow a_i + \epsilon (y_i - s((Ka)_i)) \quad [\text{where } s \text{ is the logistic function}]$$

[Just like with perceptrons, every time you update one weight a_i , you can update Ka in $O(n)$ time so you don't have to compute it from scratch on the next iteration. If you prefer batch gradient descent ...]

Batch gradient descent step:

$$a \leftarrow a + \epsilon (y - s(Ka)) \quad \Leftarrow \text{applying } s \text{ component-wise to vector } Ka$$

For each test pt z:

$$h(z) \leftarrow s \left(\sum_{j=1}^n a_j k(X_j, z) \right)$$

[If you're using logistic regression as a classifier and you don't care about the posterior probabilities, you can skip the logistic function and just compute the summation, like in the perceptron algorithm.]

The Gaussian Kernel

[Mind-blowing as the polynomial kernel is, I think our next trick is even more mind-blowing. Since we can now do fast computations in spaces with exponentially large dimensions, why don't we go all the way and generate feature vectors in infinite-dimensional space?]

Gaussian kernel, aka radial basis fn kernel: there exists a $\Phi(x)$ such that

$$k(x, z) = \exp\left(-\frac{|x - z|^2}{2\sigma^2}\right) \quad [\text{This kernel takes } O(d) \text{ time to compute.}]$$

[In case you're curious, here's the feature vector that gives you this kernel, for the case where you have only one input feature per sample point.]

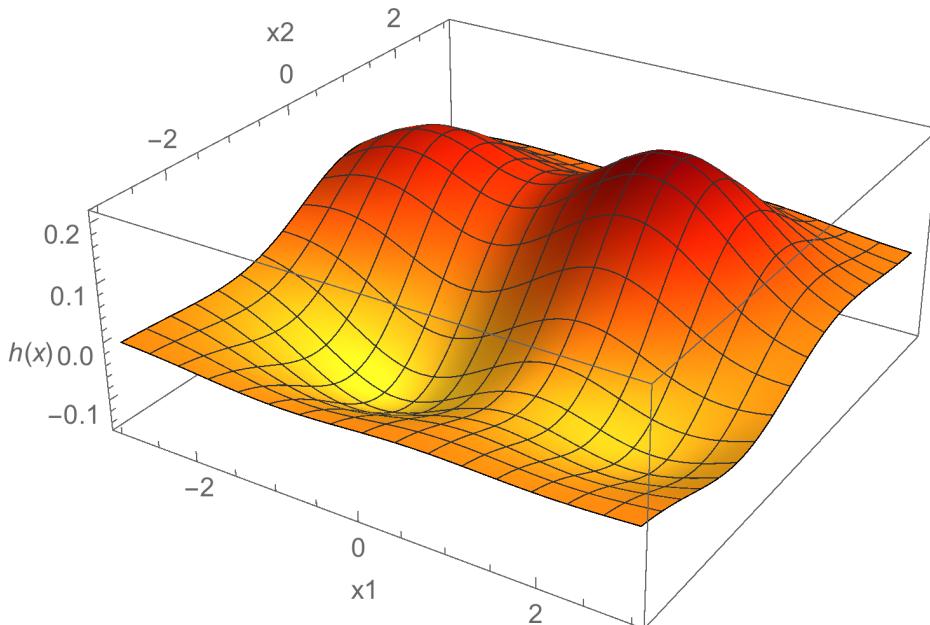
e.g. for $d = 1$,

$$\Phi(x) = \exp\left(-\frac{x^2}{2\sigma^2}\right) \left[1, \frac{x}{\sigma\sqrt{1!}}, \frac{x^2}{\sigma^2\sqrt{2!}}, \frac{x^3}{\sigma^3\sqrt{3!}}, \dots\right]^\top$$

[This is an infinite vector, and $\Phi(x) \cdot \Phi(z)$ is a series that converges to $k(x, z)$. Nobody actually uses this value of $\Phi(x)$ directly, or even cares about it; they just use the kernel function $k(\cdot, \cdot)$.]

[At this point, it's best *not* to think of points in a high-dimensional space. It's no longer a useful intuition. Instead, think of the kernel k as a measure of how similar or close together two points are to each other.]

Key observation: hypothesis $h(z) = \sum_{j=1}^n a_j k(X_j, z)$ is a linear combo of Gaussians centered at sample pts.
 [The dual weights are the coefficients of the linear combination.]
 [The Gaussians are a basis for the hypothesis.]



gausskernel.pdf [A hypothesis h that is a linear combination of Gaussians centered at four sample points, two with positive weights and two with negative weights. If you use ridge regression with a Gaussian kernel, your “linear” regression will look something like this.]

Very popular in practice! Why?

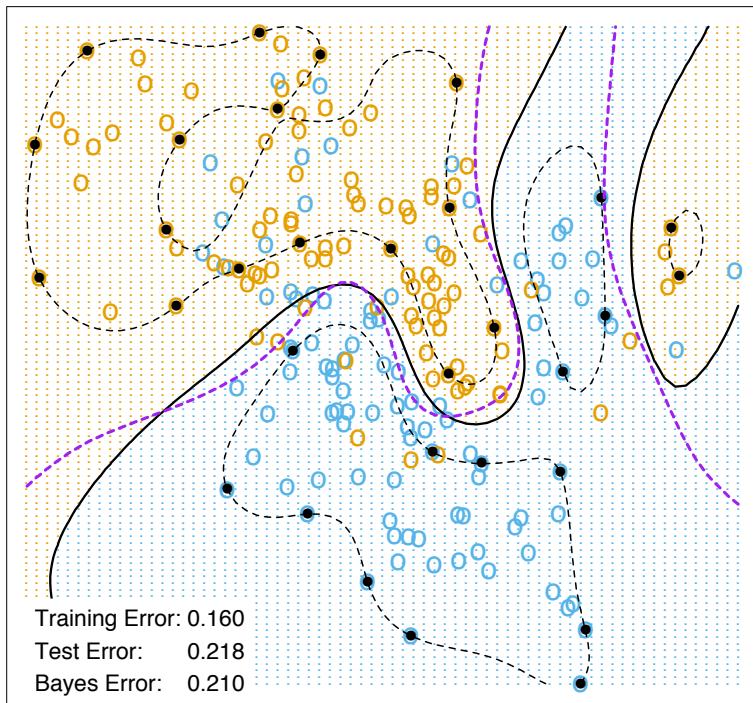
- Gives very smooth h [In fact, h is infinitely differentiable; it's C^∞ -continuous.]
- Behaves somewhat like k -nearest neighbors, but smoother
- Oscillates less than polynomials (depending on σ)
- $k(x, z)$ interpreted as a similarity measure. Maximum when $z = x$; goes to 0 as distance increases.
- Sample points “vote” for value at z , but closer points get weightier vote.

[The “standard” kernel $x \cdot z$ assigns more weight to vectors that point in roughly the same direction as z . By contrast, the Gaussian kernel assigns more weight to sample points near z .]

σ trades off bias vs. variance:

larger $\sigma \rightarrow$ wider Gaussians, smoother $h \rightarrow$ more bias, less variance

Choose σ by (cross-)validation.



gausskernelsvm.pdf (ESL, Figure 12.3) [The decision boundary (solid black) of a soft-margin SVM with a Gaussian kernel. Observe that in this example, it comes reasonably close to the Bayes optimal decision boundary (dashed purple). The dashed black curves are the boundaries of the margin. The small black disks are the support vectors that lie on the margin boundary.]

[By the way, there are many other kernels that, like the Gaussian kernel, are defined directly as kernel functions without worrying about Φ . But not every function can be a kernel function. A function is qualified only if it always generates a positive semidefinite kernel matrix, for every sample. There is an elaborate theory about how to construct valid kernel functions. However, you probably won't need it. The polynomial and Gaussian kernels are the two most popular by far.]

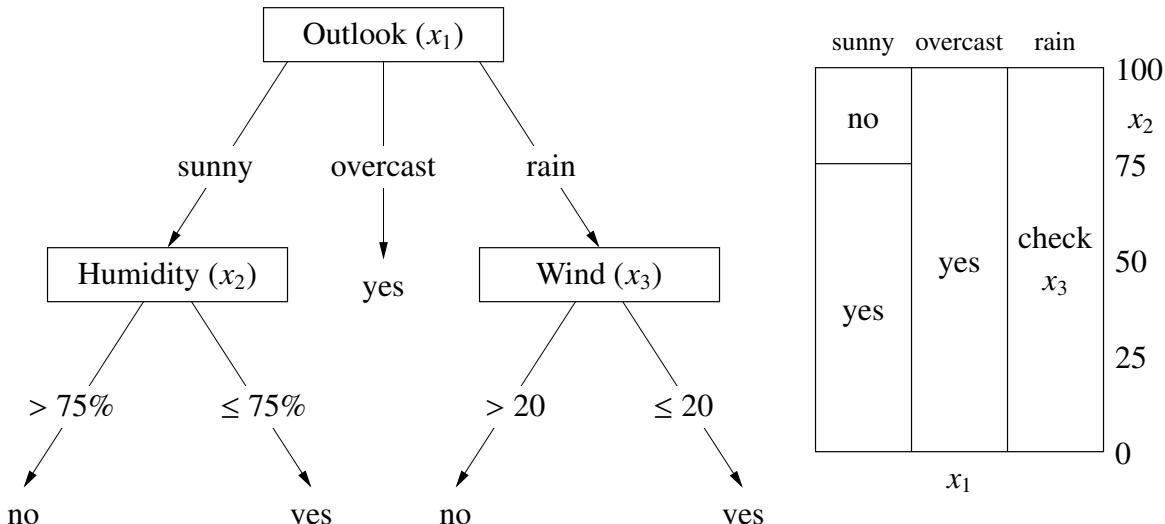
15 Decision Trees

DECISION TREES

Nonlinear method for classification and regression.

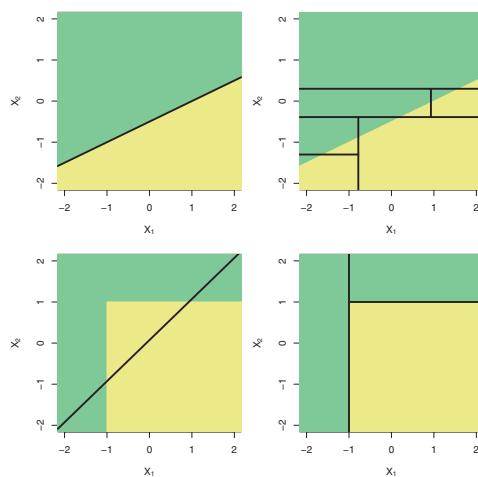
Uses tree with 2 node types:

- internal nodes test feature values (usually just one) & branch accordingly
- leaf nodes specify class $h(x)$



[Draw this by hand. [decree.pdf](#)] Deciding whether to go out for a picnic.]

- Cuts x -space into rectangular cells
- Works well with both categorical and quantitative features
- Interpretable result (inference)
- Decision boundary can be arbitrarily complicated



[treelinearcompare.pdf](#) (ISL, Figure 8.7) [Comparison of linear classifiers (left) vs. decision trees (right) on 2 examples.]

Consider classification first. Greedy, top-down learning heuristic:

[This algorithm is more or less obvious, and has been rediscovered many times. It's naturally recursive. I'll show how it works for classification first; later I'll talk about how it works for regression.]

Let $S \subseteq \{1, 2, \dots, n\}$ be set of sample point indices.

Top-level call: $S = \{1, 2, \dots, n\}$.

$\text{GrowTree}(S)$

```

if ( $y_i = C$  for all  $i \in S$  and some class C) then {
    return new leaf( $C$ )           [We say the leaves are pure]
} else {
    choose best splitting feature  $j$  and splitting value  $\beta$  (*)  

     $S_l = \{i : X_{ij} < \beta\}$           [Or you could use  $\leq$  and  $>$ ]  

     $S_r = \{i : X_{ij} \geq \beta\}$   

    return new node( $j, \beta$ ,  $\text{GrowTree}(S_l)$ ,  $\text{GrowTree}(S_r)$ )
}

```

(*) How to choose best split?

- Try all splits. [All features, and all splits within a feature.]
- For a set S , let $J(S)$ be the cost of S .
- Choose the split that minimizes $J(S_l) + J(S_r)$; or,
the split that minimizes weighted average $\frac{|S_l|J(S_l) + |S_r|J(S_r)}{|S_l| + |S_r|}$.

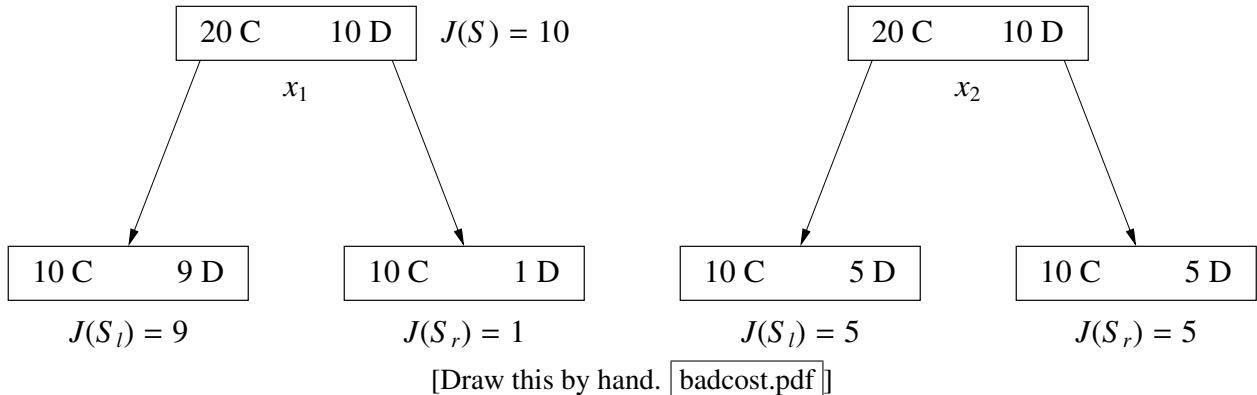
[Here, I'm using the vertical brackets $|\cdot|$ to denote set cardinality.]

How to choose cost $J(S)$?

[I'm going to start by suggesting a mediocre cost function, so you can see why it's mediocre.]

Idea 1 (bad): Label S with the class C that labels the most points in S .

$$J(S) \leftarrow \# \text{ of points in } S \text{ not in class } C.$$



Problem: Left split much better, but $J(S_l) + J(S_r) = 10$ for both. Weighted avg prefers right split!

[There are many different splits that all have the same total cost. We want a cost function that better distinguishes between them.]

Idea 2 (good): Measure the entropy.

[An idea from information theory.]

Let Y be a random class variable, and suppose $P(Y = C) = p_C$.

The surprise of Y being class C is $-\log_2 p_C$.

[Always nonnegative.]

- event w/prob. 1 gives us zero surprise.

- event w/prob. 0 gives us infinite surprise!

[In information theory, the surprise is equal to the expected number of bits of information we need to transmit which events happened to a recipient who knows the probabilities of the events. Often this means using fractional bits, which may sound crazy, but it makes sense when you're compiling lots of events into a single message; e.g., a sequence of biased coin flips.]

The entropy of an index set S is the average surprise

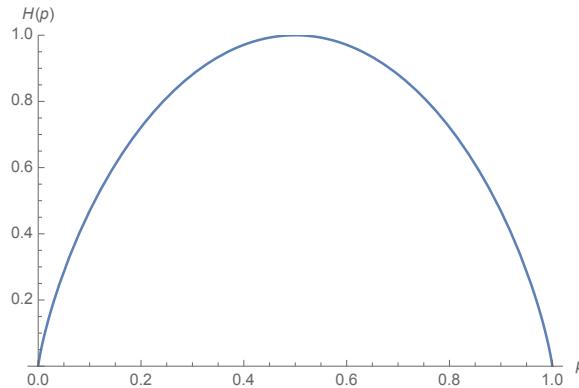
$$H(S) = - \sum_C p_C \log_2 p_C, \quad \text{where } p_C = \frac{|\{i \in S : y_i = C\}|}{|S|}. \quad \begin{array}{l} \text{[The proportion of points in } S \\ \text{that are in class } C.\text{]} \end{array}$$

If all points in S belong to same class? $H(S) = -1 \log_2 1 = 0$.

Half class C, half class D? $H(S) = -0.5 \log_2 0.5 - 0.5 \log_2 0.5 = 1$.

n points, all different classes? $H(S) = -\log_2 \frac{1}{n} = \log_2 n$.

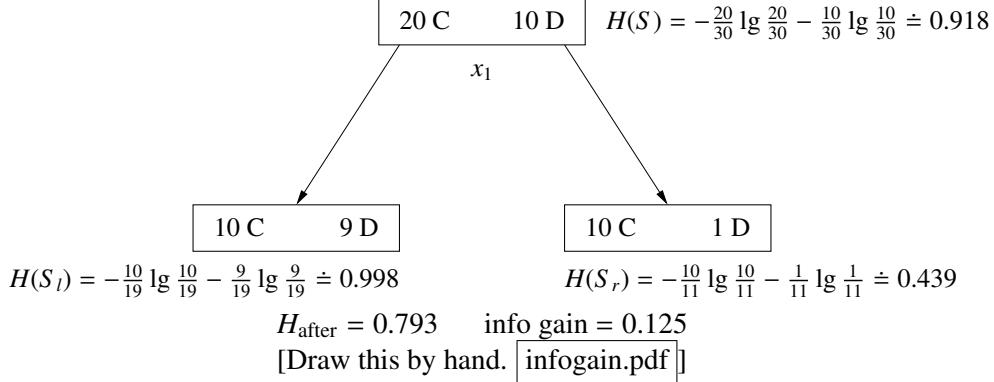
[The entropy is the expected number of bits of information we need to transmit to identify the class of a sample point in S chosen uniformly at random. It makes sense that it takes 1 bit to specify C or D when each class is equally likely. And it makes sense that it takes $\log_2 n$ bits to specify one of n classes when each class is equally likely.]



entropy.pdf [Plot of the entropy $H(p_C)$ when there are only two classes. The probability of the second class is $p_D = 1 - p_C$, so we can plot the entropy with just one dependent variable. If you have > 2 classes, you would need a multidimensional chart to graph the entropy, but the entropy is still strictly concave.]

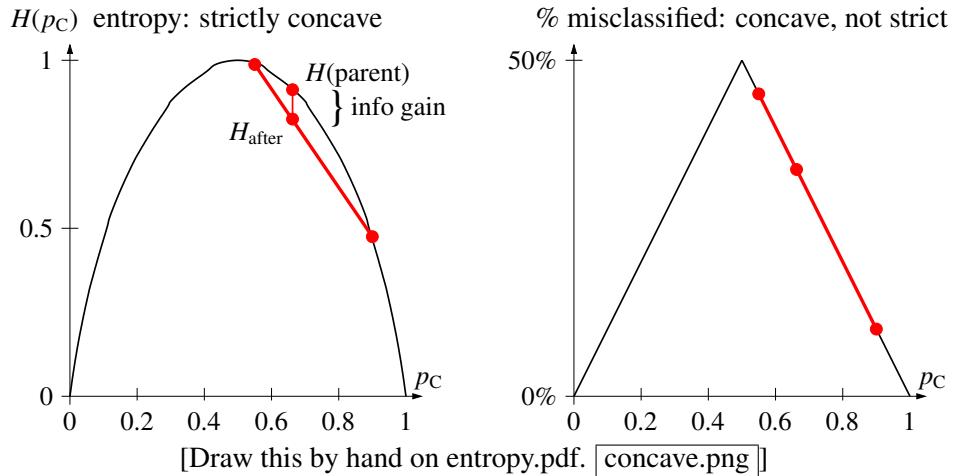
Weighted avg entropy after split is $H_{\text{after}} = \frac{|S_l| H(S_l) + |S_r| H(S_r)}{|S_l| + |S_r|}$.

Choose split that maximizes information gain $H(S) - H_{\text{after}}$. [Which is just the same as minimizing H_{after} .]



Info gain always positive except when one child is empty or for all C, $P(y_i = C|i \in S_l) = P(y_i = C|i \in S_r)$. [Which is the case for the second split we considered.]

[Recall the graph of the entropy.]



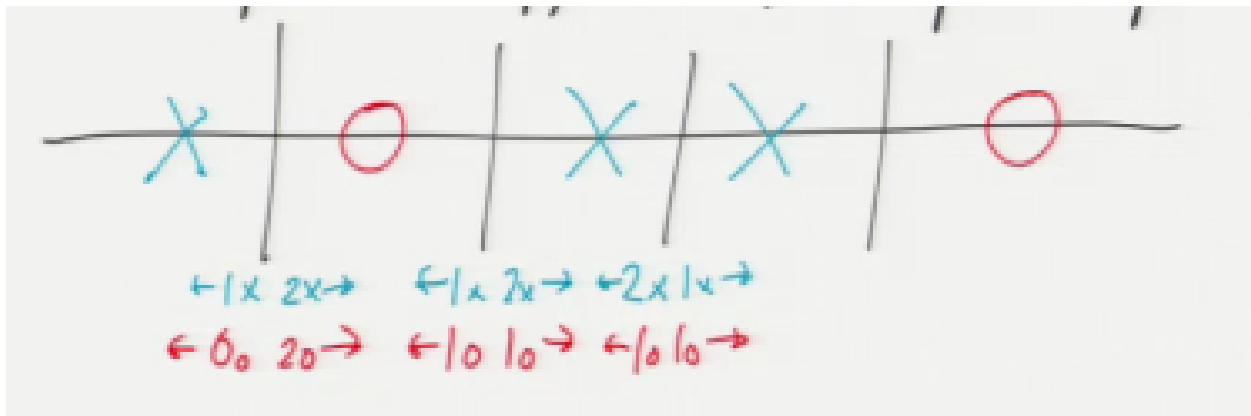
[Suppose we pick two points on the entropy curve, then draw a line segment connecting them. Because the entropy curve is strictly concave, the interior of the line segment is strictly below the curve. Any point on that segment represents a weighted average of the two entropies for suitable weights. If you unite the two sets into one parent set, the parent set's value p_C is the weighted average of the children's p_C 's. Therefore, the point directly above that point on the curve represents the parent's entropy. The information gain is the vertical distance between them. So the information gain is positive unless the two child sets both have exactly the same p_C and lie at the same point on the curve.]

[On the other hand, for the graph on the right, plotting the % misclassified, if we draw a line segment connecting two points on the curve, the segment might lie entirely on the curve. In that case, uniting the two child sets into one, or splitting the parent set into two, changes neither the total misclassified sample points nor the weighted average of the % misclassified. The bigger problem, though, is that many different splits will get the same weighted average cost; this test doesn't distinguish the quality of different splits well.]

[By the way, the entropy is not the only function that works well. Many concave functions work fine, including the simple polynomial $p(1-p)$.]

More on choosing a split:

- For binary feature x_i : children are $x_i = 0$ & $x_i = 1$.
- If x_i has 3+ discrete values: split depends on application.
[Sometimes it makes sense to use multiway splits; sometimes binary splits.]
- If x_i is quantitative: sort x_i values in S ; try splitting between each pair of *unequal* consecutive values.
[We can radix sort the points in linear time, and if n is huge we should.]
Clever bit: As you scan sorted list from left to right, you can update entropy in $O(1)$ time per point!⁵
[This is important for obtaining a fast tree-building time.]
[Draw a row of X's and O's; show how we update the # of X's and # of O's in each of S_L and S_R as we scan from left to right.]



scan.png

Algs & running times:

- Test point: Walk down tree until leaf. Return its label.
Worst-case time is $O(\text{tree depth})$.
For binary features, that's $\leq d$. [Quantitative features may go deeper.]
Usually (not always) $\leq O(\log n)$.
- Training: For binary features, try $O(d)$ splits at each node.
For quantitative features, try $O(n'd)$ splits; n' = points in node
Either way $\Rightarrow O(n'd)$ time at this node
[Quantitative features are asymptotically just as fast as binary features because of our clever way of computing the entropy for each split.]
Each point participates in $O(\text{depth})$ nodes, costs $O(d)$ time in each node.
Running time $\leq O(nd \text{ depth})$.
[As nd is the size of the design matrix X , and the depth is often logarithmic, this is a surprisingly reasonable running time.]

⁵Let C be the number of class C sample points to the left of a potential split and c be the number to the right of the split. Let D be the number of class not-C points to the left of the split and d be the number to the right of the split. Update C , c , D , and d at each split as you move from left to right. At each potential split, calculate the entropy of the left set as $-\frac{C}{C+D} \log_2 \frac{C}{C+D} - \frac{D}{C+D} \log_2 \frac{D}{C+D}$ and the entropy of the right set as $-\frac{c}{c+d} \log_2 \frac{c}{c+d} - \frac{d}{c+d} \log_2 \frac{d}{c+d}$. Note: $\log 0$ is undefined, but this formula works if we use the convention $0 \log 0 = 0$.

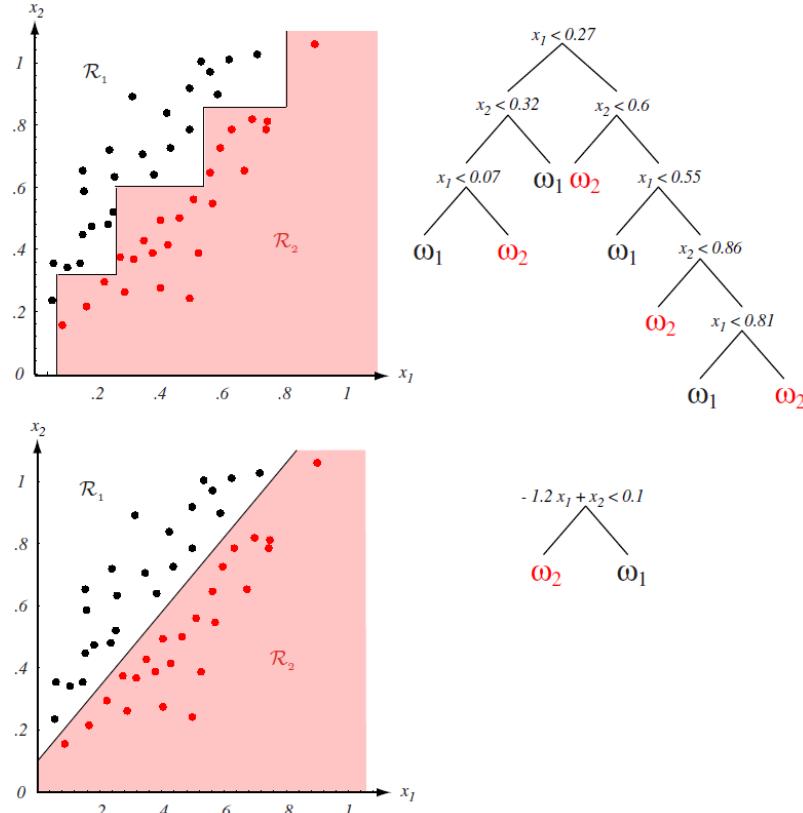
It follows that the weighted average of the two entropies is $-\frac{1}{n'} \left(C \log_2 \frac{C}{C+D} + D \log_2 \frac{D}{C+D} + c \log_2 \frac{c}{c+d} + d \log_2 \frac{d}{c+d} \right)$, where $n' = C + D + c + d$ is the total number of sample points stored in this treenode. Choose the split that minimizes this weighted average.

16 More Decision Trees, Ensemble Learning, and Random Forests

DECISION TREES (continued)

Multivariate Splits

Find non-axis-aligned splits with other classification algs or by generating them randomly.



multivariate.pdf [An example where an ordinary decision tree needs many splits to approximate a diagonal linear decision boundary, but a single multivariate split takes care of it.]

[Here you can use classification algorithms such as SVMs, logistic regression, and Gaussian discriminant analysis. Decision trees permit these algorithms to find more complicated decision boundaries by making them hierarchical.]

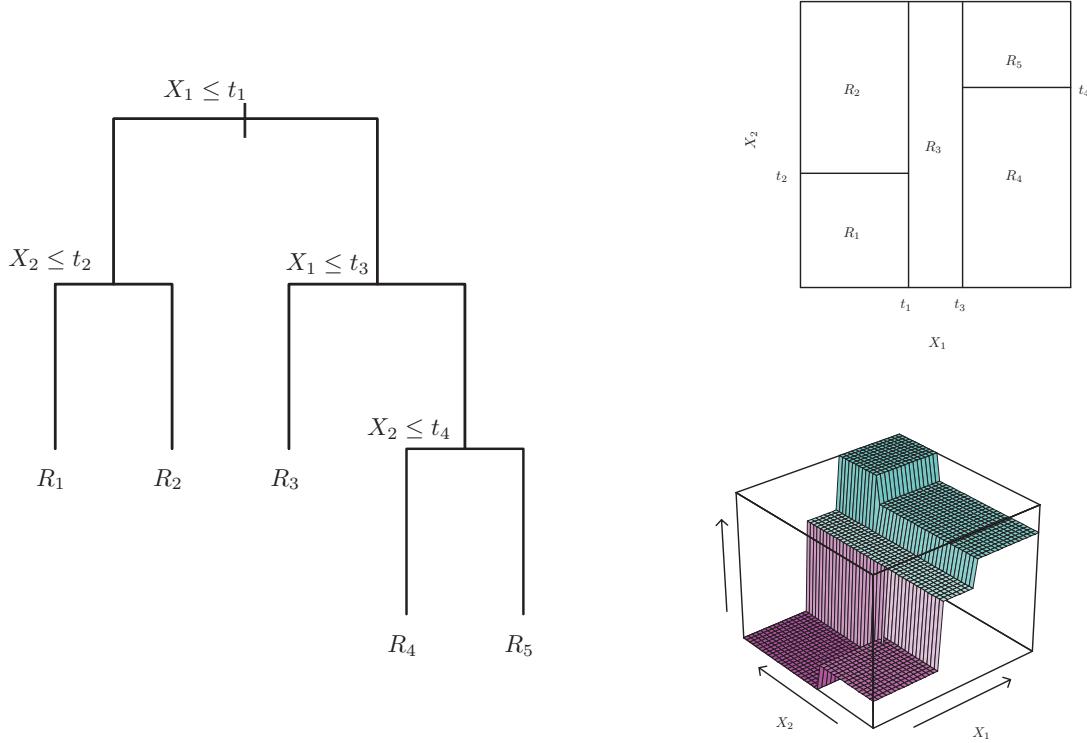
May gain better classifier at cost of worse interpretability or speed.

[One of the reasons decision trees are typically so fast is that they check only one feature at each treenode. But if there are hundreds of features, and you have to check all of them at every level of the tree to classify a point, it slows down classification a lot. So it sometimes pays to consider methods like forward stepwise selection when you're learning so that when you classify, you only need to check a few features.]

Can limit # of features per split: forward stepwise selection, Lasso.

Decision Tree Regression

Creates a piecewise constant regression fn.



regressstree.pdf, regressstreefn.pdf (ISL, Figure 8.3) [Decision tree regression.]

Cost $J(S) = \frac{1}{|S|} \sum_{i \in S} (y_i - \bar{y})^2$, where \bar{y} is the mean y_i for sample points in subset S .

[So if all the points in a node have the same y -value, then the cost is zero.]

[We choose the split that minimizes the weighted average of the costs of the children after the split.]

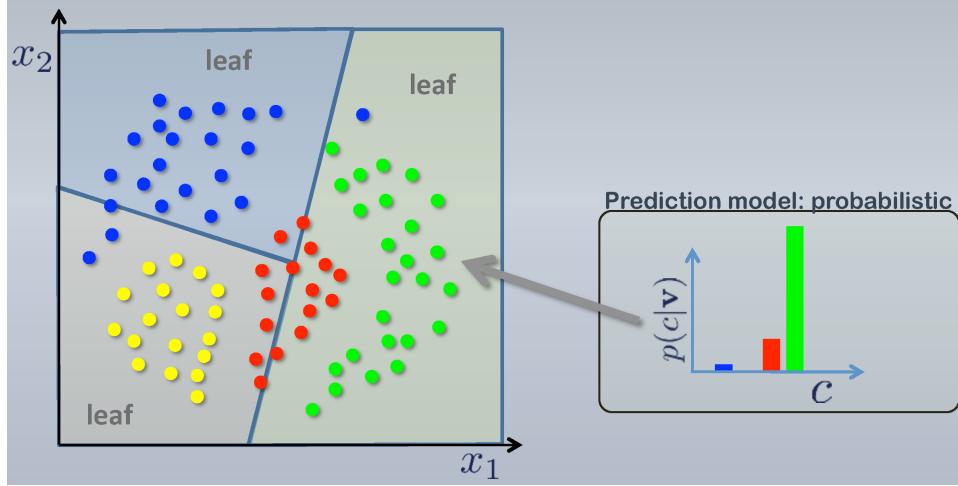
Stopping Early

[The basic version of the decision tree algorithm keeps subdividing treenodes until every leaf is pure. We don't have to do that; sometimes we prefer to stop subdividing treenodes earlier.]

Why?

- Limit tree depth (for speed)
- Limit tree size (big data sets)
- Complete tree may overfit
- Given noise or overlapping distributions, purity of leaves is counterproductive; better to estimate posterior probs

[When you have overlapping class distributions, refining the tree down to one sample point per leaf is absolutely guaranteed to overfit. It's better to stop early, then classify each leaf node by taking a vote of its sample points. Alternatively, you can use the points to estimate a posterior probability for each leaf. If there are many points in each leaf, the posterior probabilities might be reasonably accurate.]



leaf.pdf [In the decision tree at left, each leaf has multiple classes. Instead of returning the majority class, each leaf could return a posterior probability histogram, as illustrated at right.]

How? Select stopping condition(s):

- Next split doesn't reduce entropy/error enough (dangerous; pruning is better)
- Most of node's points (e.g. > 95%) have same class [to deal with outliers]
- Node contains few sample points (e.g. < 10)
- Cell's edges are all tiny
- Depth too great
- Use (cross-)validation to compare

[The last is the slowest but most effective way to know when to stop: use validation to decide whether splitting the node is a win or a loss. But if your goal is to avoid overfitting, it's generally even more effective to grow the tree a little too large and then use validation to prune it back. We'll talk about that next.]

Leaves with multiple points return

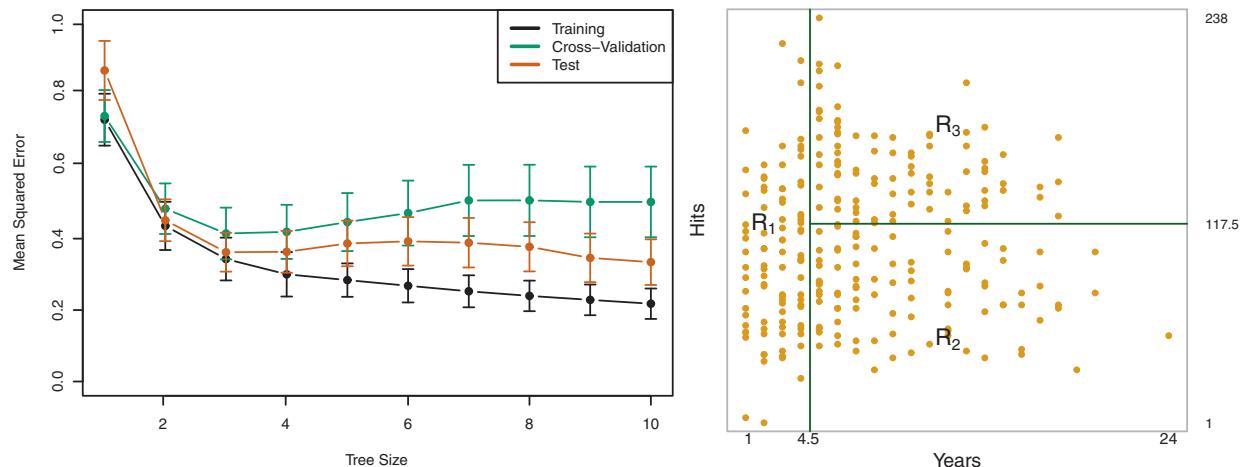
- a majority vote or class posterior probs (classification) or
- an average (regression).

Pruning

Grow tree too large; greedily remove each split whose removal improves cross-validation performance. More reliable than stopping early.

[We have to do cross-validation once for each split that we're considering removing. But you can do that pretty cheaply. What you *don't* do is reclassify every sample point from scratch. Instead, you keep track of which points in the validation set end up at which leaf. When you are deciding whether to remove a split, you just look at the validation points in the two leaves you're thinking of removing, and see how they will be reclassified and how that will change the error rate. You can do this very quickly.]

[The reason why pruning often works better than stopping early is because often a split that doesn't seem to make much progress is followed by a split that makes a lot of progress. If you stop early, you'll never find out. Pruning is a simple idea, but it's highly recommended when you have the time.]



`prunehitters.pdf`, `prunedhitters.pdf` (ISL, Figures 8.5 & 8.2) [At left, a plot of decision tree size vs. errors for baseball hitter data. At right, the best decision tree has three leaves. Players' salaries: R1 = \$165,174, R2 = \$402,834, R3 = \$845,346.]

[In this example, a 10-node decision tree was constructed to predict the salaries of baseball players, based on their years in the league and average hits per season. Then the tree was pruned by cross-validation. The best decision tree on the validation data turned out to have just three leaves.]

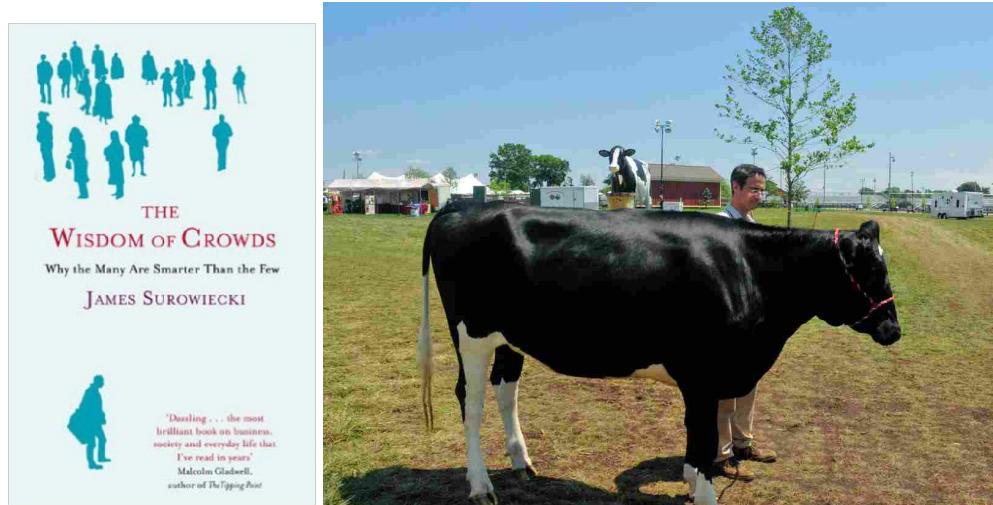
ENSEMBLE LEARNING

Decision trees are fast, simple, interpretable, easy to explain, invariant under scaling/translation, robust to irrelevant features.

But not the best at prediction. [Compared to previous methods we've seen.]
High variance.

[For example, suppose we take a training data set, split it into two halves, and train two decision trees, one on each half of the data. It's not uncommon for the two trees to turn out very different. In particular, if the two trees pick different features for the very first split at the top of the tree, then it's quite common for the trees to be completely different. So decision trees tend to have high variance.]

[So let's think about how to fix this. As an analogy, imagine that you are generating random numbers from some distribution. If you generate just one number, it might have high variance. But if you generate n numbers and take their average, then the variance of that average is n times smaller. So you might ask yourself, can we reduce the variance of decision trees by taking an average answer of a bunch of decision trees? Yes we can.]



wisdom.jpg, penelope.jpg [James Surowiecki's book "The Wisdom of Crowds" and Penelope the cow. Surowiecki tells us this story ...]

[A 1906 county fair in Plymouth, England had a contest to guess the weight of an ox. A scientist named Francis Galton was there, and he did an experiment. He calculated the median of everyone's guesses. The median guess was 1,207 pounds, and the true weight was 1,198 pounds, so the error was less than 1%. Even the cattle experts present didn't estimate it that accurately.]

[NPR repeated the experiment in 2015 with a cow named Penelope whose photo they published online. They got 17,000 guesses, and the average guess was 1,287 pounds. Penelope's actual weight was 1,355 pounds, so the crowd got it to within 5 percent.]

[The main idea is that sometimes the average opinion of a bunch of idiots is better than the opinion of one expert. And so it is with learning algorithms. We call a learning algorithm a weak learner if it does better than guessing randomly. And we combine a bunch of weak learners to get a strong one.]

[Incidentally, James Surowiecki, the author of the book, guessed 725 pounds for Penelope. So he was off by 87%. He's like a bad decision tree who wrote a book about how to benefit from bad decision trees.]

We can take average of output of

- different learning algs
- same learning alg on many training sets [if we have tons of data]
- bagging: same learning alg on many random subsamples of one training set
- random forests: randomized decision trees on random subsamples

[These last two are the most common ways to use averaging, because usually we don't have enough training data to use fresh data for every learner.]

[Averaging is not specific to decision trees; it can work with many different learning algorithms. But it works particularly well with decision trees.]

Regression algs: take median or mean output

Classification algs: take majority vote OR average posterior probs

[Apology to readers: I show some videos in this lecture, which cannot be included in this report.]

[Show averageaxis.mov] [Here's a simple classifier that takes an average of "stumps," trees of depth 1. Observe how good the posterior probabilities look.]

[Show averageaxistree.mov] [Here's a 4-class classifier with depth-2 trees.]

[The Netflix Prize was an open competition for the best collaborative filtering algorithm to predict user ratings for films, based on previous ratings. It ran for three years and ended in 2009. The winners used an extreme ensemble method that took an average of many different learning algorithms. In fact, a couple of top teams combined into one team so they could combine their methods. They said, “Let’s average our models and split the money,” and that’s what happened.]

Use learners with low bias (e.g. deep decision trees).

High variance & some overfitting are okay. Averaging reduces the variance!

[Each learner may overfit, but each overfits in its own unique way.]

Averaging sometimes reduces the bias & increases flexibility;

e.g. creating nonlinear decision boundary from linear classifiers.

Hyperparameter settings usually different than 1 learner.

[Because averaging learners reduces their variance. But averaging rarely reduces bias as much as it reduces variance, so you want to get the bias nice and small before you average.]

of trees is another hyperparameter.

Bagging = Bootstrap AGGregatING (Leo Breiman, 1994)

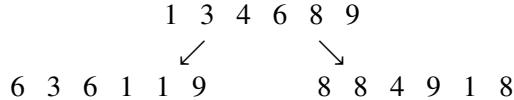
[Leo Breiman was a statistics professor right here at Berkeley. He did his best work after he retired in 1993. The bagging algorithm was published the following year, and then he went on to co-invent random forests as well. Unfortunately, he died in 2005.]



breiman.gif [Leo Breiman]

[Bagging is a randomized method for creating many different learners from the same data set. It works well with many different learning algorithms. One exception seems to be k -nearest neighbors; bagging mildly degrades it.]

Given n -point training sample, generate random subsample of size n' by sampling *with replacement*. Some points chosen multiple times; some not chosen.



If $n' = n$, $\sim 63.2\%$ are chosen. [On average; this fraction varies randomly.]

Build learner. Points chosen j times have greater weight:

[If a point is chosen j times, we want to treat it the same way we would treat j different points all bunched up infinitesimally close together.]

- Decision trees: j -time point has $j \times$ weight in entropy.
- SVMs: j -time point incurs $j \times$ penalty to violate margin.
- Regression: j -time point incurs $j \times$ loss.

Repeat until T learners. Metalearner takes test point, feeds it into all T learners, returns average/majority output.

Random Forests

Random sampling isn't random enough!

[With bagging, often the decision trees look very similar. Why is that?] One really strong predictor → same feature split at top of every tree.

[For example, if you're building decision trees to identify spam, the first split might always be "viagra." Random sampling might not change that. If the trees are very similar, then taking their average doesn't reduce the variance much.]

Idea: At each treenode, take random sample of m features (out of d).

Choose best split from m features.

[We're not allowed to split on the other $d - m$ features!]

Different random sample for each treenode.

$m \approx \sqrt{d}$ works well for classification; $m \approx d/3$ for regression.

[So if you have 100-dimensional feature space, you randomly choose 10 features and pick the one of those 10 that gives the best split. But m is a hyperparameter, and you might get better results by tuning it for your particular application. These values of m are good starting guesses.]

Smaller $m \rightarrow$ more randomness, less tree correlation, more bias

[One reason this works is if there's a really strong predictor, only a fraction of the trees can choose that predictor as the first split. That fraction is m/d . So the split tends to "decorrelate" the trees. And that means that when you take the average of the trees, you'll have less variance.]

[You have to be careful, though, because you don't want to dumb down the trees too much in your quest for decorrelation. Averaging works best when you have very strong learners that are also diverse. But it's hard to create a lot of learners that are very different yet all very smart. The Netflix Prize winners did it, but it was a huge amount of work.]

Sometimes test error reduction up to 100s or even 1,000s of decision trees!

Disadvantage: loses interpretability/inference.

[But the compensation is it's more accurate than a single decision tree.]

Variation: generate m random multivariate splits (oblique lines, quadrics); choose best split.

[You have to be a bit clever about how you generate random decision boundaries; I'm not going to discuss that today. I'll just show lots of examples.]

[Show treesidesdeep.mov] [Lots of good-enough conic random decision trees.]

[Show averageline.mov]

[Show averageconic.mov]

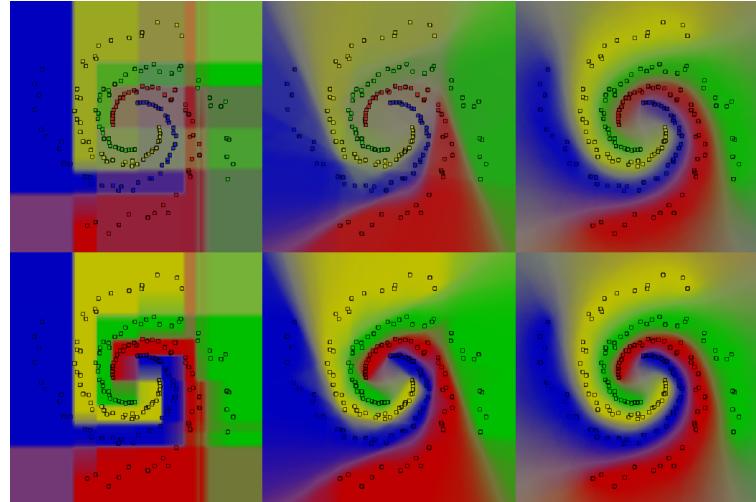
[Show square.mov] [Depth 2; look how good the posterior probabilities look.]

[Show squaresmall.mov] [Depth 2; see the uncertainty away from the center.]

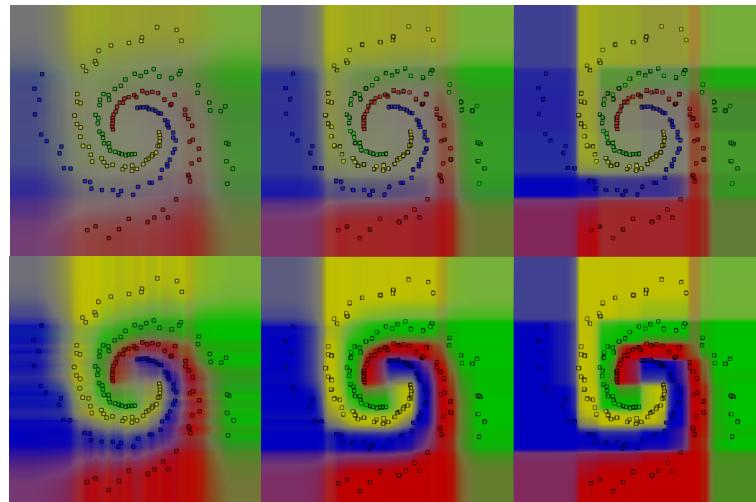
[Show spiral2.mov] [Doesn't look like a decision tree at all, does it?]

[Show overlapdepth14.mov] [Overlapping classes. This example overfits!]

[Show overlapdepth5.mov] [Better fit.]



500.pdf [Random forest classifiers for 4-class spiral data. Each forest takes the average of 400 trees. The top row uses trees of depth 4. The bottom row uses trees of depth 12. From left to right, we have axis-aligned splits, splits with lines with arbitrary rotations, and splits with conic sections. Each split is chosen to be the best of 500 random choices.]



randomness.pdf [Random forest classifiers for the same data. Each forest takes the average of 400 trees. In these examples, *all* the splits are axis-aligned. The top row uses trees of depth 4. The bottom row uses trees of depth 12. From left to right, we choose each split from 1, 5, or 50 random choices. The more choices, the less bias and the better the classifier.]

17 Neural Networks

NEURAL NETWORKS

Can do both classification & regression.

[They tie together several ideas from the course: perceptrons, logistic regression, ensembles of learners, and stochastic gradient descent. They also tie in the idea of lifting sample points to a higher-dimensional feature space, but with a new twist: neural nets can learn features themselves.]

[I want to begin by reminding you of the story I told you at the beginning of the semester, about Frank Rosenblatt's invention of perceptrons in 1957. Remember that he held a press conference where he predicted that perceptrons would be "the embryo of an electronic computer that [the Navy] expects will be able to walk, talk, see, write, reproduce itself and be conscious of its existence."]

[Perceptron research continued until something monumental happened in 1969. Marvin Minsky, one of the founding fathers of AI, and Seymour Papert published a book called "Perceptrons." Sounds good, right? Well, part of the book was devoted to things perceptrons can't do. And one of those things is XOR.]

		x_1	
		XOR	
		0	1
x_2	0	0	1
	1	1	0

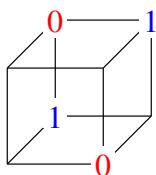
[Think of the four outputs here as sample points in two-dimensional space. Two of them are in class 1, and two of them are in class 0. We want to find a linear classifier that separates the 1's from the 0's. Can we do it? No.]

[So Minsky and Papert were basically saying, "Frank. You're telling us this machine is going to be conscious of its own existence but it can't do XOR?"] []

[The book had a devastating effect on the field. After its publication, almost no research was done on neural net-like ideas for a decade, a time we now call the "AI Winter." Shortly after the book was published, Frank Rosenblatt died.] []

[One thing I don't understand is why the book was so fatal when there are several almost obvious ways to get around the XOR problem. Here's the easiest.]

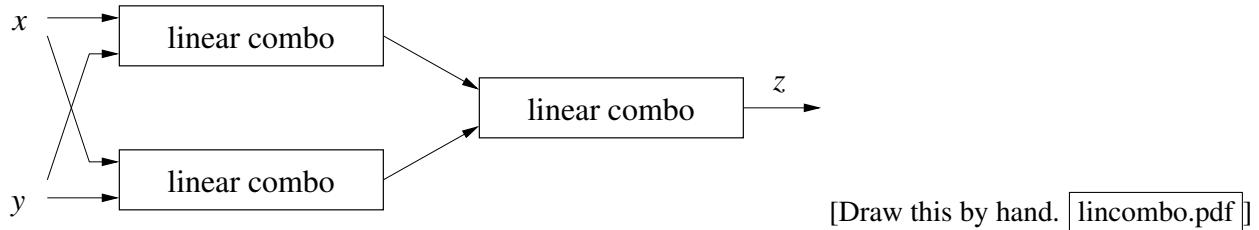
If you add one new quadratic feature, $x_1 x_2$, XOR is linearly separable in 3D.



[Draw this by hand. [xorcube.pdf](#)]

[Now we can find a plane that cuts through the cube obliquely and separates the 0's from the 1's.]

[However, there's an even more powerful way to do XOR. The idea is to design linear classifiers whose output is the input to other linear classifiers. That way, you should be able to emulate arbitrarily logical circuits. Suppose I put together some linear decision functions like this.]



[Draw this by hand. [lincombo.pdf](#)]

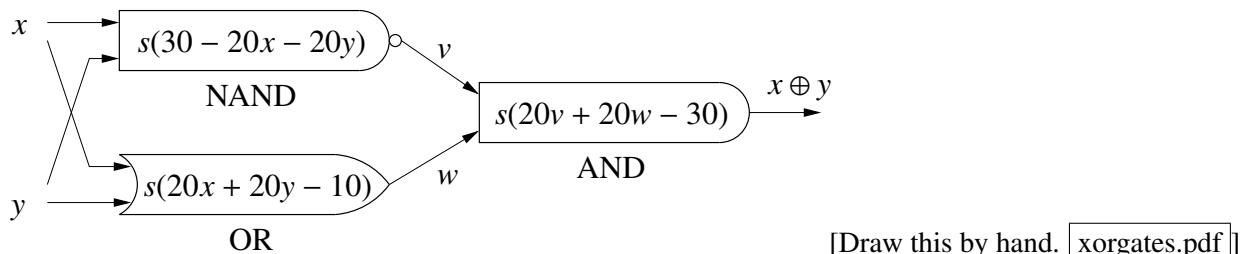
[If I interpret the output as 1 if z is positive or 0 if z is negative, can I do XOR with this?]

A linear combo of a linear combo is a linear combo . . . only works for linearly separable points.

[We need one more idea to make neural nets. We need to add some sort of nonlinearity between the linear combinations. Let's call these boxes that compute linear combinations "neurons." If a neuron runs the linear combination it computes through some nonlinear function before sending it on to other neurons, then the neurons can act somewhat like logic gates. The nonlinearity could be as simple as clamping the output so it can't go below zero. And that's what people usually use in practice these days.]

[The traditional choice has been to use the logistic function. The logistic function can't go below zero or above one, which is nice because it can't ever get huge and oversaturate the other neurons it's sending information to. The logistic function is also smooth, which means it has well-defined gradients and Hessians we can use in gradient descent.]

[With logistic functions between the linear combinations, here's a two-level perceptron that computes the XOR function.]



[Draw this by hand. [xorgates.pdf](#)]

[Note that the logistic function at the output is optional; we could just take the sign of the output instead.]

Network with 1 Hidden Layer

Input layer: $x_1, \dots, x_d ; x_{d+1} = 1$

Hidden units: $h_1, \dots, h_m ; h_{m+1} = 1$

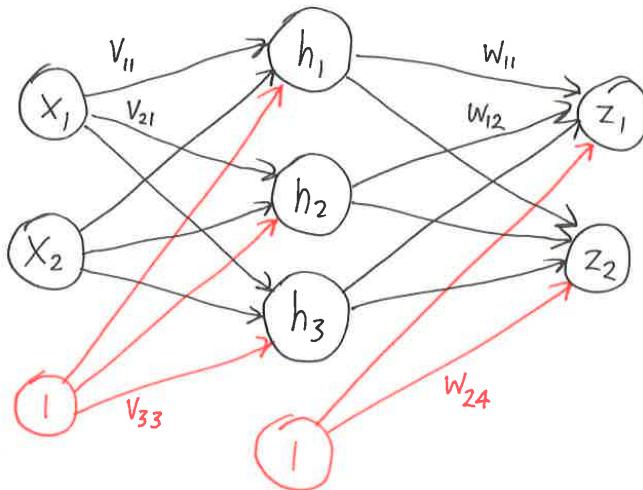
Output layer: z_1, \dots, z_k

Layer 1 weights: $m \times (d + 1)$ matrix V V_i^\top is row i

Layer 2 weights: $k \times (m + 1)$ matrix W W_i^\top is row i

Recall [logistic function] $s(\gamma) = \frac{1}{1+e^{-\gamma}}$. Other nonlinear fns can be used.

For vector v , $s(v) = \begin{bmatrix} s(v_1) \\ s(v_2) \\ \vdots \end{bmatrix}$. [We apply s to a vector component-wise.]



[Draw this by hand. [1hiddenlayer.pdf](#)]

$$h = s_1(Vx)$$

$$\dots \text{that is, } h_i = s\left(\sum_{j=1}^{d+1} V_{ij}x_j\right)$$

$$z = s(Wh) = s(Ws_1(Vx))$$

↑
add a 1 to end of s vector

[We might have more than one output so that we can build multiple classifiers that share hidden units. One of the interesting advantages of neural nets is that if you train multiple classifiers simultaneously, sometimes some of them come out better because they can take advantage of particularly useful hidden units that first emerged to support one of the other classifiers.]

[We can add more hidden layers, and for image recognition tasks it's common to have 12 to 40 hidden layers. There are many variations you can experiment with—for instance, you can have connections that go forward more than one layer.]

Training

Usually stochastic or batch gradient descent.

Pick loss fn $L(z, y)$

↑↑

predictions true values (could be vectors)

Cost fn is $J(h) = \frac{1}{n} \sum_{i=1}^n L(h(X_i), Y_i)$

[I'm using a capital Y here because now Y is a matrix with one row for each sample point and one column for each output of the neural net. Sometimes there is just one output, but many neural net applications have more.]

Usually there are many local minima!

[The cost function for a neural net is, generally, not even close to convex. For that reason, it's possible to wind up in a bad minimum. We'll talk later about some clever ways to coax neural nets into better minima.]

[Now let me ask you this. Suppose we start by setting all the weights to zero, and then we do gradient descent on the weights. What will go wrong?]

[This neural network has a symmetry: there's really no difference between one hidden unit and any other hidden unit. The gradient descent algorithm has no way to break the symmetry between hidden units. You can get stuck in a situation where all the weights out of an input unit have the same value, and all the weights into an output unit have the same value, and they have no way to become different from each other. To avoid this problem, and in the hopes of finding a better local minimum, we start with random weights.]

Let w be a vector containing all the weights in V & W . Batch gradient descent:

```
w ← vector of random weights
repeat
    w ← w - ε ∇J(w)
```

[We've just rewritten all the weights as a vector for notational convenience. When you actually write the code, you should probably operate directly on the weight matrices V and W .]

[It's important to make sure the random weights aren't too big, because if a unit's output gets too close to zero or one, it can get "stuck," meaning that a modest change in the input values causes barely any change in the output value. Stuck units tend to stay stuck because in that operating range, the gradient $s'(\cdot)$ of the logistic function is close to zero.]

[Instead of batch gradient descent, we can use stochastic gradient descent, which means we use the gradient of one sample point's loss function at each step. Typically, we shuffle the points in a random order, or just pick one randomly at each step.]

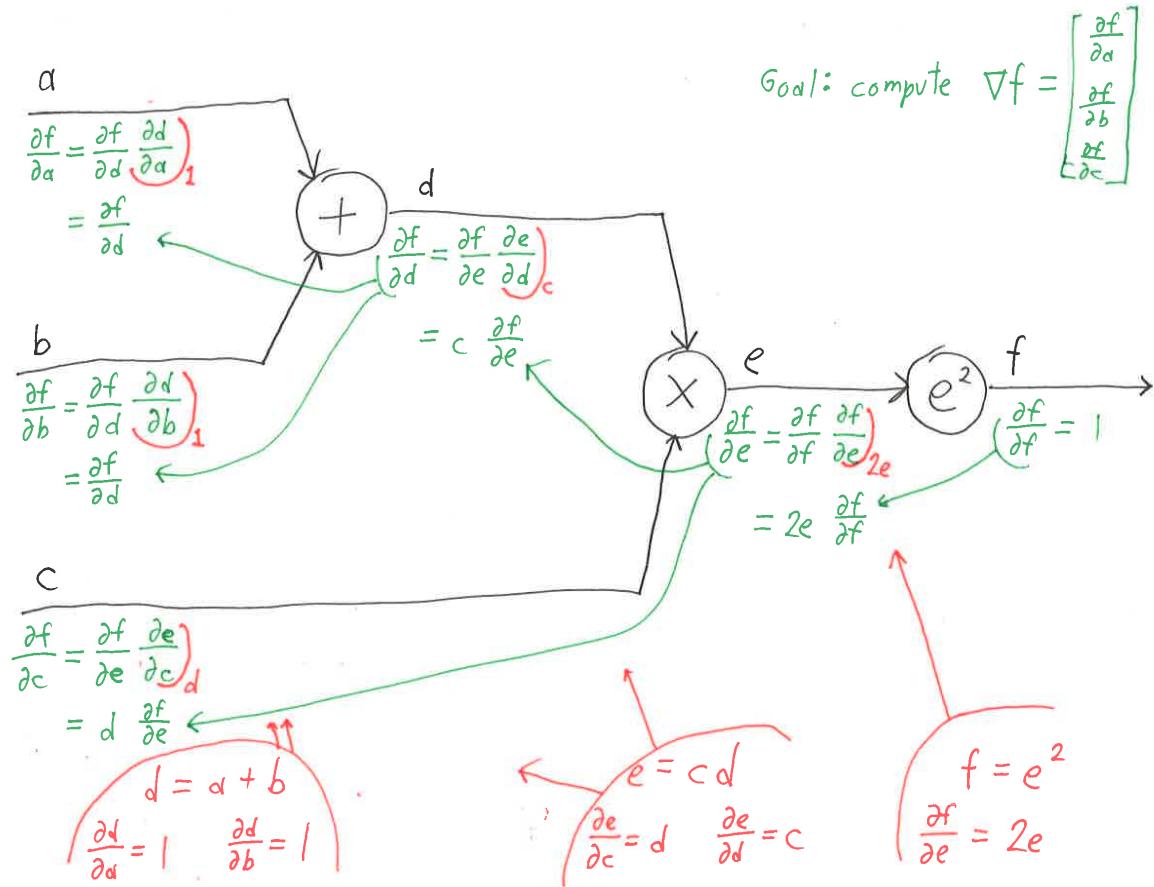
[The hard part of this algorithm is computing the gradient. If you simply derive one derivative for each weight, you'll find that for a network with multiple layers of hidden units, it takes time linear in the number of edges in the neural network to compute a derivative for one weight. Multiply that by the number of weights. We're going to spend some time learning to improve the running time to linear in the number of edges.]

Naive gradient computation: $O(\text{edges}^2)$ time

Backpropagation: $O(\text{edges})$ time

Computing Gradients for Arithmetic Expressions

[Let's see what it takes to compute the gradient of an arithmetic expression. It turns into repeated applications of the chain rule from calculus.]



Each value z gives partial derivative of the form

$$\frac{\partial f}{\partial z} = \left(\frac{\partial f}{\partial n} \frac{\partial n}{\partial z} \right)$$

computed during forward pass

computed during backward pass

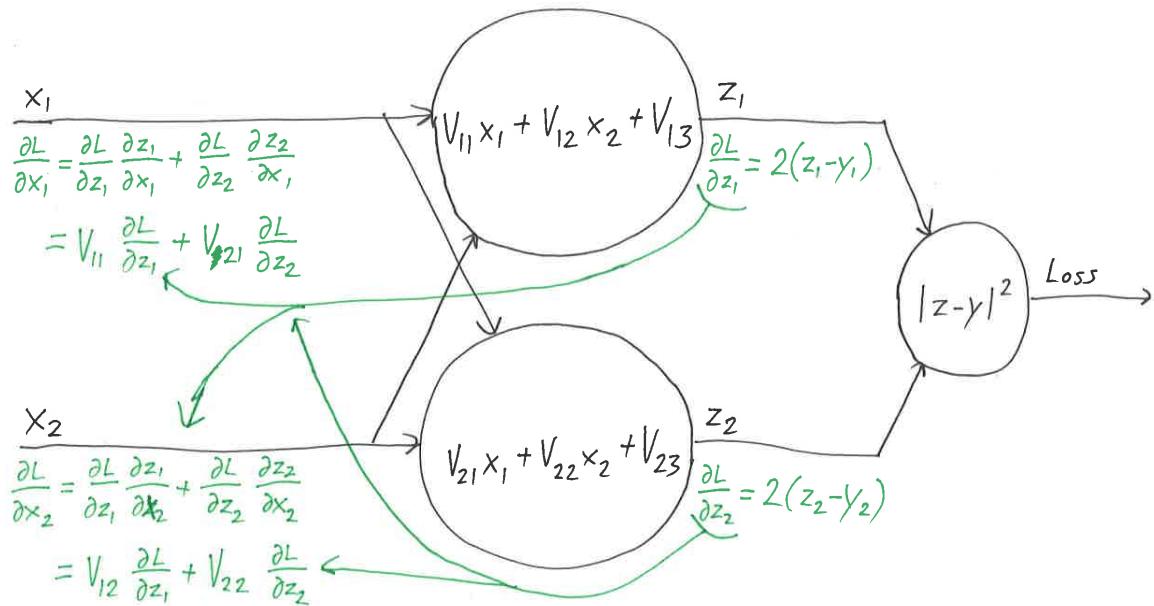
after forward pass

"backpropagation"

where z is input to n .

[Draw this by hand. [gradients.pdf](#) Draw the black diagram first. Then the goal (upper right). Then the green and red expressions, from left to right, leaving out the green arrows. Then the green arrows, starting at the right side of the page and moving left. Lastly, write the text at the bottom. (Use the same procedure for the next two figures.)]

[What if a unit's output goes to more than one unit? Then we need to understand a more complicated version of the chain rule. Let's try it with an expression that's similar to what you'll encounter in a neural net.]



[Draw this by hand. [gradientspartial.pdf](#)]

[Here we're using a standard rule of multivariate calculus:]

$$\frac{\partial}{\partial \tau} L(z_1(\tau), z_2(\tau)) = \frac{\partial L}{\partial z_1} \frac{\partial z_1}{\partial \tau} + \frac{\partial L}{\partial z_2} \frac{\partial z_2}{\partial \tau} = \nabla_z L \cdot \frac{\partial}{\partial \tau} z$$

[Observe that we're doing dynamic programming here. We're computing the solutions of subproblems, then using each solution to compute the solutions of several bigger problems.]

The Backpropagation Alg.

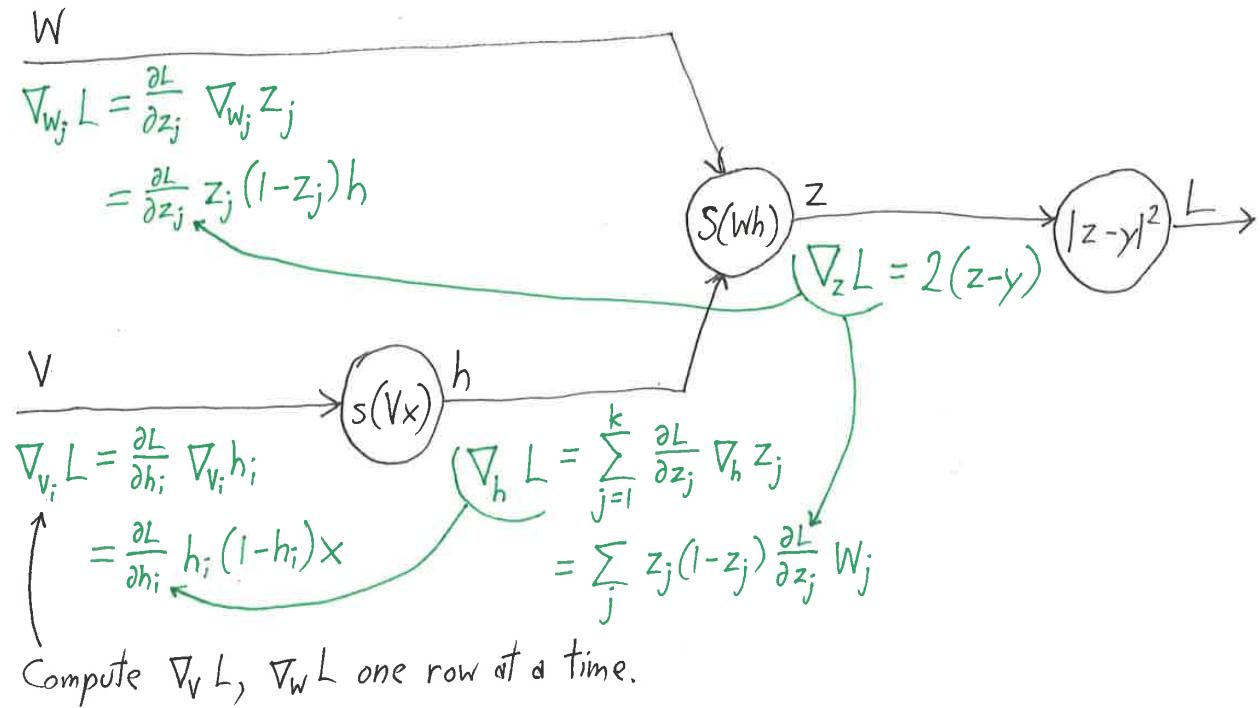
[Backpropagation is a dynamic programming algorithm for computing the gradients we need to do neural net stochastic gradient descent in time linear in the number of weights.]

V_i^\top is row i of weight matrix V [and likewise for rows of W]

Recall $s'(\gamma) = s(\gamma)(1 - s(\gamma))$

$$\begin{aligned} h_i &= s(V_i \cdot x), \text{ so} & \nabla_{V_i} h_i &= s'(V_i \cdot x) x = h_i(1 - h_i)x \\ z_j &= s(W_j \cdot h), \text{ so} & \nabla_{W_j} z_j &= s'(W_j \cdot h) h = z_j(1 - z_j)h & \nabla_h z_j &= z_j(1 - z_j)W_j \end{aligned}$$

[Here is the arithmetic expression for the same neural network I drew for you three illustrations ago. It looks very different when you depict it like this, but don't be fooled; it's exactly the same network I started with.]



18 Neurons; Variations on Neural Networks

NEURONS

[The field of artificial intelligence started with some wrong premises. The early AI researchers attacked problems like chess and theorem proving, because they thought those exemplified the essence of intelligence. They didn't pay much attention at first to problems like vision and speech understanding. Any four-year-old can do those things, and so researchers underestimated their difficulty. Today, we know better. Computers can effortlessly beat four-year-olds at chess, but they still can't play with toys well. We've come to realize that rule-based symbol manipulation is not the primary defining mark of intelligence. Even rats do computations that we're hard pressed to match with our computers. We've also come to realize that these are different classes of problems that require very different styles of computation. Brains and computers have very different strengths and weaknesses, which reflect their different computing styles.]

[Neural networks are partly inspired by the workings of actual brains. Let's take a look at a few things we know about biological neurons, and contrast them with both neural nets and traditional computation.]

- CPUs: largely sequential, nanosecond gates, fragile if gate fails
superior for arithmetic, logical rules, perfect key-based memory
- Brains: very parallel, millisecond neurons, fault-tolerant

[Neurons are continually dying. You've probably lost a few since this lecture started. But you probably didn't notice. And that's interesting, because it points out that our memories are stored in our brains in a diffuse representation. There is no one neuron whose death will make you forget that $2 + 2 = 4$. Artificial neural nets often share that resilience. Brains and neural nets seem to superpose memories on top of each other, all stored together in the same weights, sort of like a hologram.]

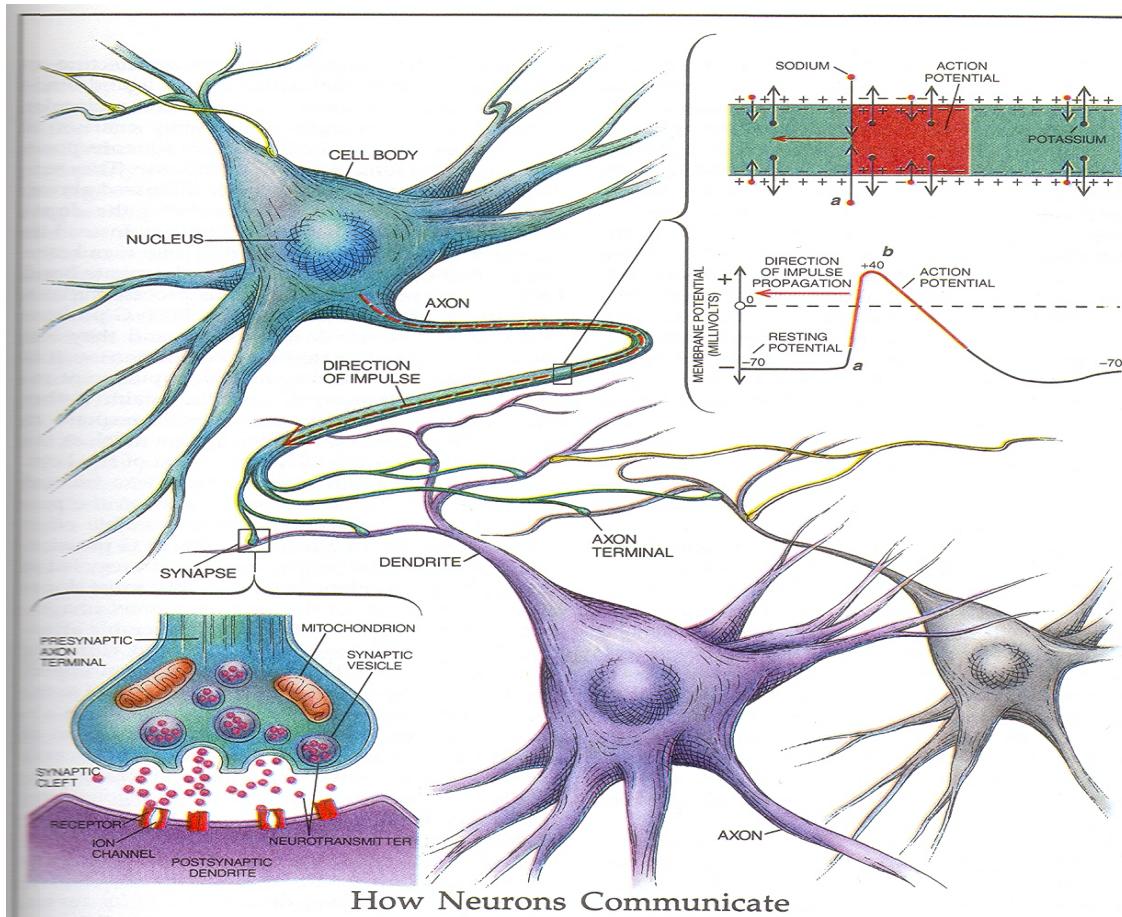
[In the 1920's, the psychologist Karl Lashley conducted experiments to identify where in the brain memories are stored. He trained rats to run a maze, and then made lesions in different parts of the cerebral cortex, trying to erase the memory trace. Lashley failed; his rats could still find their way through the maze, no matter where he put lesions. He concluded that memories are not stored in any one area of the brain, but are distributed throughout it. Neural networks, properly trained, can duplicate this property.]

superior for vision, speech, associative memory

[By "associative memory," I mean noticing connections between things. One thing our brains are very good at is retrieving a pattern if we specify only a portion of the pattern.]

[It's impressive that even though a neuron needs a few milliseconds to transmit information to the next neurons downstream, we can perform very complex tasks like interpreting a visual scene in a tenth of a second. This is possible because neurons run in parallel, but also because of their computation style.]

[Neural nets try to emulate the parallel, associative thinking style of brains, and they are among the best techniques we have for many fuzzy problems, including some problems in vision and speech. Not coincidentally, neural nets are also inferior at many traditional computer tasks such as multiplying 10-digit numbers or compiling source code.]



[neurons.pdf](#)

- Neuron: A cell in brain/nervous system for thinking/communication
- Action potential or spike: An electrochemical impulse fired by a neuron to communicate w/other neurons
- Axon: The limb(s) along which the action potential propagates; “output”
[Most axons branch out eventually, sometimes profusely near their ends.]
[It turns out that giant squids have a very large axon they use for fast water jet propulsion. The mathematics of action potentials was first characterized in these giant squid axons, and that work won a Nobel Prize in Physiology in 1963.]
- Dendrite: Smaller limbs by which neuron receives info; “input”
- Synapse: Connection from one neuron’s axon to another’s dendrite
[Some synapses connect axons to muscles or glands.]
- Neurotransmitter: Chemical released by axon terminal to stimulate dendrite

[When an action potential reaches an axon terminal, it causes tiny containers of neurotransmitter, called vesicles, to empty their contents into the space where the axon terminal meets another neuron’s dendrite. That space is called the synaptic cleft. The neurotransmitters bind to receptors on the dendrite and influence the next neuron’s body voltage. This sounds incredibly slow, but it all happens in 1 to 5 milliseconds.]

You have about 10^{11} neurons, each with about 10^4 synapses.

[]

Analogies: [between artificial neural networks and brains]

- Output of unit \leftrightarrow firing rate of neuron

[An action potential is “all or nothing”—all action potentials have the same shape and size. The output of a neuron is not signified by voltage like the output of a transistor. The output of a neuron is the frequency at which it fires. Some neurons can fire at nearly 1,000 times a second, which you might think of as a strong “1” output. Conversely, some types of neurons can go for minutes without firing. But some types of neurons never stop firing, and for those you might interpret a firing rate of 10 times per second as “0”.]

- Weight of connection \leftrightarrow synapse strength

- Positive weight \leftrightarrow excitatory neurotransmitter (e.g. glutamine)

- Negative weight \leftrightarrow inhibitory neurotransmitter (e.g. GABA, glycine) [Gamma aminobutyric acid.]

[A typical neuron is either excitatory at all its axon terminals, or inhibitory at all its terminals. It can't switch from one to the other. Artificial neural nets have an advantage here.]

- Linear combo of inputs \leftrightarrow summation

[A neuron fires when the sum of its inputs, integrated over time, reaches a high enough voltage. However, the neuron body voltage also decays slowly with time, so if the action potentials are coming in slowly enough, the neuron might not fire at all.]

- Logistic/sigmoid fn \leftrightarrow firing rate saturation

[A neuron can't fire more than 1,000 times a second, nor less than zero times a second. This limits its ability to overpower downstream neurons without help. We accomplish the same thing with the sigmoid function.]

- Weight change/learning \leftrightarrow synaptic plasticity

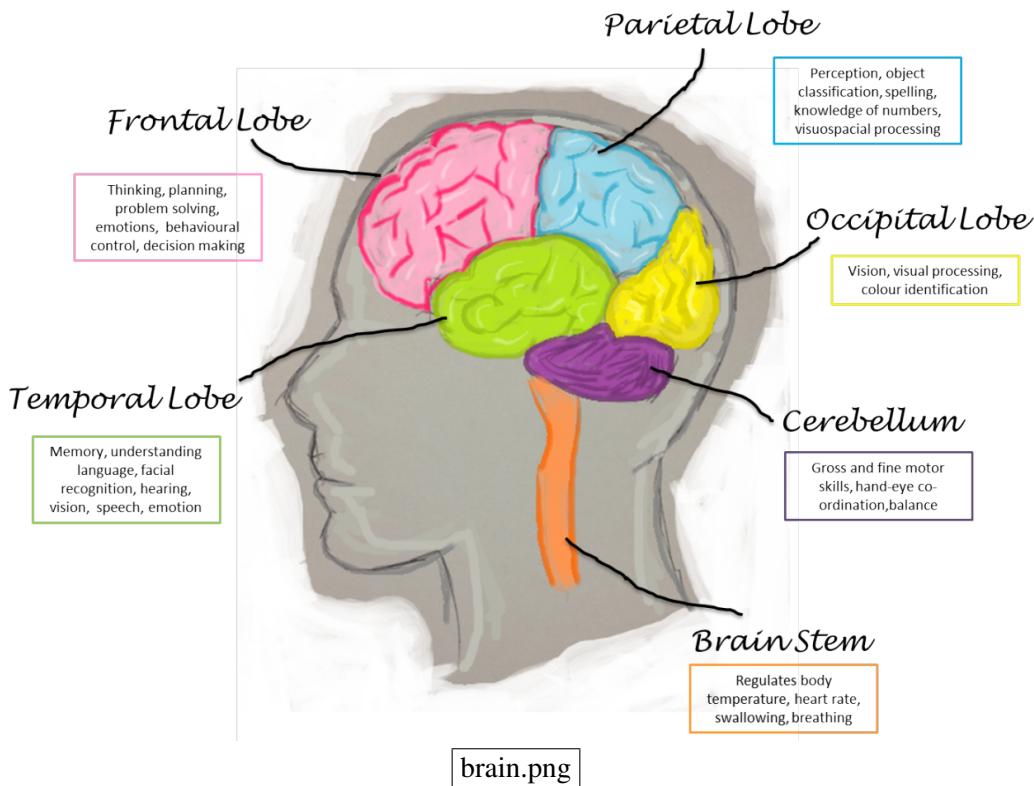
Hebb's rule (1949): “Cells that fire together, wire together.”

[This doesn't mean that the cells have to fire at exactly the same time. But if one cell's firing tends to make another cell fire more often, their excitatory synaptic connection tends to grow stronger. There's a reverse rule for inhibitory connections. And there are ways for neurons that aren't even connected to grow connections.]

[There are simple computer learning algorithms based on Hebb's rule. It can work, but it's generally not nearly as fast or effective as backpropagation.]

[Backpropagation is one part of artificial neural networks for which there is no analogy in the brain. Brains probably do not do backpropagation . . . though some people do speculate otherwise.]

[The brain is very modular.]



brain.png

[(The following items are all spoken, not written ...)]

- The part of our brain we think of as most characteristically human is the cerebral cortex, the seat of self-awareness, language, and abstract thinking.

But the brain has a lot of other parts that take the load off the cortex.

- Our brain stem regulates functions like heartbeat, breathing, and sleep.
- Our cerebellum governs fine coordination of motor skills. When we talk about “muscle memory,” much of that is in the cerebellum, and it saves us from having to consciously think about how to walk or talk or brush our teeth, so the cortex can focus on where to walk and what to say. []
- Our limbic system is the center of emotion and motivation, and as such, it makes a lot of the big decisions. I sometimes think that 90% of the job of our cerebral cortex is to rationalize decisions that have already been made by the limbic system. []
- Our visual cortex (in the occipital lobe) performs a lot of processing on the input from your eyes to change it into a more useful form. Neuroscientists and computer scientists are particularly interested in the visual cortex for several reasons. Vision is an important problem for computers. The visual cortex is one of the easier parts of the brain to study in lab animals. The visual cortex is largely a feedforward network with few neurons going backward, so it’s easier for us to train computers to behave like the visual cortex.]

[Although the brain has lots of specialized modules, one thing that's interesting about the frontal lobe is that it seems to be made of general-purpose neural tissue that looks more or less the same everywhere, at least before it's trained. If you experience damage to part of the frontal lobe early enough in life, while your brain is still growing, the functions will just relocate to a different part of the frontal lobe, and you'll probably never notice the difference.]

[As computer scientists, our primary motivation for studying neurology is to try to get clues about how we can get computers to do tasks that humans are good at. But neurologists and psychologists have also been part of the study of neural nets from the very beginning. Their motivations are scientific: they're curious how humans think, and how we can do the things we do.]

NEURAL NET VARIATIONS

[I want to show you a few basic variations on the standard neural network I showed you last class, and how some of these variations change backpropagation.]

Regression: usually linear output unit(s)—omit sigmoid fn.

[If you make that change, the gradient changes too, and you have to change the derivation of backprop. The derivation gets simpler, so I'll leave it as an exercise.]

Classification: to choose from $k \geq 3$ classes, use softmax fn. [Here we deploy k separate output units.]

Let $t = Wh$ be k -vector of linear combos in final layer.

$$\text{Softmax output is } z_j(t) = \frac{e^{t_j}}{\sum_{i=1}^k e^{t_i}}. \quad \frac{\partial z_j}{\partial t_j} = z_j \cdot (1 - z_j), \quad \frac{\partial z_j}{\partial t_i} = -z_j z_i, j \neq i, \quad \nabla_h z_j = z_j \cdot (W_j - W^\top z).$$

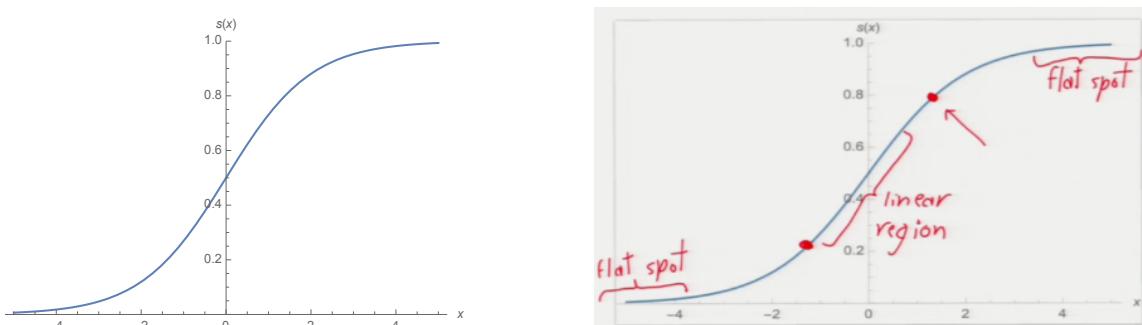
Each $z_j \in (0, 1)$; their sum is 1.

[Interpret z_j as the probability of the input belonging to class j . For example, in the digit recognition problem, we might have 10 softmax output units, one for each digit class.]

[If you have only 2 classes, just use one sigmoid output; it's equivalent to 2-way softmax.]

Unit Saturation

Problem: When unit output s is close to 0 or 1 for most training points, $s' = s(1 - s) \approx 0$, so gradient descent changes s very slowly. Unit is “stuck.” Slow training & bad local minima.



logistic.pdf [Draw flat spots, “linear” region, & maximum curvature points (at $s(\lambda) \doteq 0.21$ and $s(\lambda) \doteq 0.79$) of the sigmoid function. Ideally, we would stay away from the flat spots.]

[Wikipedia calls this the “vanishing gradient problem.”]

[The more layers your network has, the more problematic this problem becomes. Most of the early attempts to train deep, many-layered neural nets failed.]

Mitigation: [None of these are complete cures.]

- (1) Initial weight of edge into unit with fan-in η :
random with mean zero, std. dev. $1/\sqrt{\eta}$.

[The bigger the fan-in of a unit, the easier it is to saturate it. So we choose smaller random initial weights for gates with bigger fan-in.]

- (2) Set target values to 0.15 & 0.85 instead of 0 & 1.

[Recall that the sigmoid function can never be 0 or 1; it can only come close. Relaxing the target values helps prevent the output units from getting saturated. The numbers 0.15 and 0.85 are reasonable because the sigmoid function achieves its greatest curvature when its output is near 0.21 or 0.79. But experiment to find the best values.]

[Option (2) helps to avoid stuck output units, but not stuck hidden units. So ...]

- (3) Modify backprop to add small constant (typically ~ 0.1) to s' .

[This hacks the gradient so a unit can't get stuck. We're not doing *steepest* descent any more, because we're not using the real gradient. But often we're finding a better descent direction that will get us to a minimum faster. This hack originates with Scott Fahlman's *Quickprop* algorithm.]

- (4) Cross-entropy loss fn (aka logistic loss) instead of squared error.

$$L(z, y) = - \sum_i \left(y_i \ln z_i + (1 - y_i) \ln(1 - z_i) \right) \quad [\text{for sigmoid outputs only}]$$

$\begin{array}{c} \uparrow \text{true values} \\ \uparrow \text{prediction} \end{array} \quad \left. \right\} \text{vectors}$

For backprop, we derive $\frac{\partial L}{\partial z_j} = \frac{z_j - y_j}{z_j(1 - z_j)}$. This simplifies some other [previously derived] gradients:

$$\begin{aligned} \nabla_{W_j} L &= \frac{\partial L}{\partial z_j} z_j (1 - z_j) h = (z_j - y_j) h; \Rightarrow \nabla_W L = (z - y) h^\top \quad [\text{an outer product matrix}] \\ \nabla_h L &= \sum_{j=1}^k \frac{\partial L}{\partial z_j} z_j (1 - z_j) W_j = \sum (z_j - y_j) W_j = W^\top (z - y) \end{aligned}$$

[Notice that the denominator of $\partial L / \partial z_j$ cancels out the derivative of the sigmoid, $z_j(1 - z_j)$! This saves the unit from getting stuck when the derivative is small. It is related to the fact that the logistic loss goes to infinity as the predicted value z approaches zero or one. The vanishing gradient of the sigmoid is compensated for by the huge gradient of the logistic loss.]

For k -class softmax output, cross-entropy is $L(z, y) = - \sum_i^k y_i \ln z_i$.

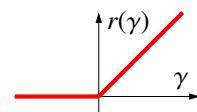
[Cross-entropy losses are only for sigmoid and softmax outputs. By contrast, for regression we typically use linear outputs, which don't saturate, so the squared error loss is better for them.]

[Like option (2), cross-entropy loss helps to avoid stuck output units, but not stuck hidden units.]

(5) Replace sigmoids with ReLUs: rectified linear units.

ramp fn aka hinge fn: $r(\gamma) = \max\{0, \gamma\}$

$$r'(\gamma) = \begin{cases} 1 & \gamma \geq 0 \\ 0 & \gamma < 0 \end{cases}$$



[The derivative is not defined at zero, but we just pretend it is.]

Popular for many-layer networks with large training sets.

[One nice thing about ramp functions is that they and their gradients are very fast to compute. Computers compute exponentials slowly. Even though ReLUs are linear in each half of their range, they're still nonlinear enough to easily compute functions like XOR.]

[Obviously, the gradient is sometimes zero, so you might wonder if ReLUs can get stuck too. Fortunately, it's rare for a ReLU's gradient to be zero for *all* the training data; it's usually zero for just some sample points. But yes, ReLUs sometimes get stuck too; just not as often as sigmoids.]

[The output of a ReLU can be arbitrarily large, creating the danger that it might overwhelm units downstream. This is called the “exploding gradient problem,” and it is not a big problem in shallow networks, but it becomes a big problem in deep or recurrent networks.]

[Note that option (5) makes options (2)–(4) irrelevant.]

[Now I will show you how to derive the backprop equations for a softmax output, the cross-entropy loss function, and ℓ_2 regularization—which helps to reduce overfitting, just like in ridge regression.]

$$\nabla_{W_i} L = \sum_{j=1}^k \frac{\partial L}{\partial z_j} \nabla_{W_i} z_j + 2\lambda w'_i$$

$$= (-y_i + z_i \sum_{j=1}^k y_j) h + 2\lambda w'_i$$

$$z_j = \frac{e^{w_j \cdot h}}{\sum_{i=1}^k e^{w_i \cdot h}}$$

$$\frac{\partial L}{\partial z_j} = -\frac{y_i}{z_j} + \lambda |w'_i|^2$$

$$\nabla_h L = \sum_{j=1}^k \frac{\partial L}{\partial z_j} \nabla_h z_j$$

$$= \sum_j y_j (w^T z - w_j)$$

l_2 regularization (optional)

[w' is a vector containing all the non-bias weights in matrices V & W .]

[Draw this by hand. [gradsoftmax4.pdf](#)]

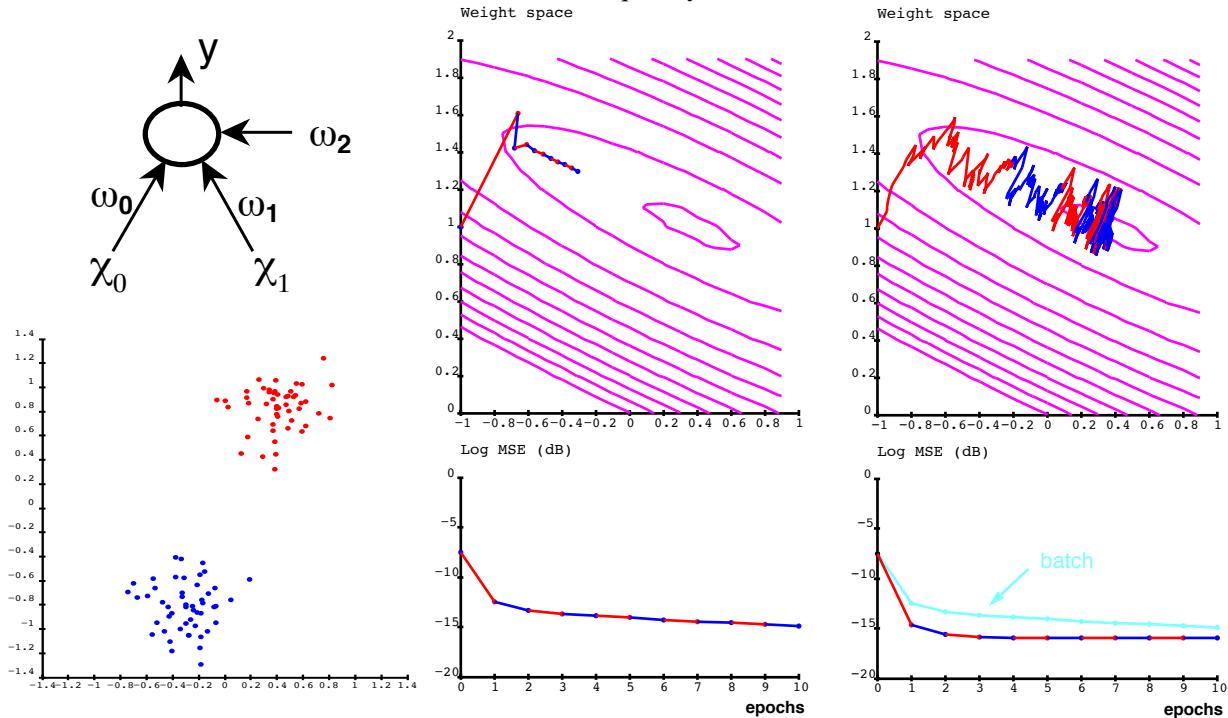
19 Better Neural Network Training; Convolutional Neural Networks

[I'm going to talk about a bunch of heuristics that make gradient descent faster, or make it find better local minima, or prevent it from overfitting. I suggest you implement vanilla stochastic backprop first, and experiment with the other heuristics only after you get that working.]

Heuristics for Faster Training

[A big disadvantage of neural nets is that they take a long, long time to train compared to other classification methods we've studied. Here are some ways to speed them up. Unfortunately, you usually have to experiment with techniques and hyperparameters to find which ones will help with your particular application.]

- (1)–(5) above. [Fix the unit saturation problem.]
- Stochastic gradient descent: faster than batch on large, redundant data sets.
[Whereas batch gradient descent walks downhill on one cost function, stochastic descent takes a very short step downhill on one point's loss function and then another short step on another point's loss function. The cost function is the sum of the loss functions over all the sample points, so one batch step behaves similarly to n stochastic steps and takes the same amount of time. But if you have many different examples of the digit "9", they contain much redundant information, and stochastic gradient descent learns the redundant information more quickly.]



batchvsstoch.pdf (LeCun et al., “Efficient BackProp”)
[Left: a simple neural net with only three weights, and its 2D training data. Center: batch gradient descent makes only a little progress each epoch. (Epochs alternate between red and blue.) Right: stochastic descent decreases the error much faster than batch descent.]

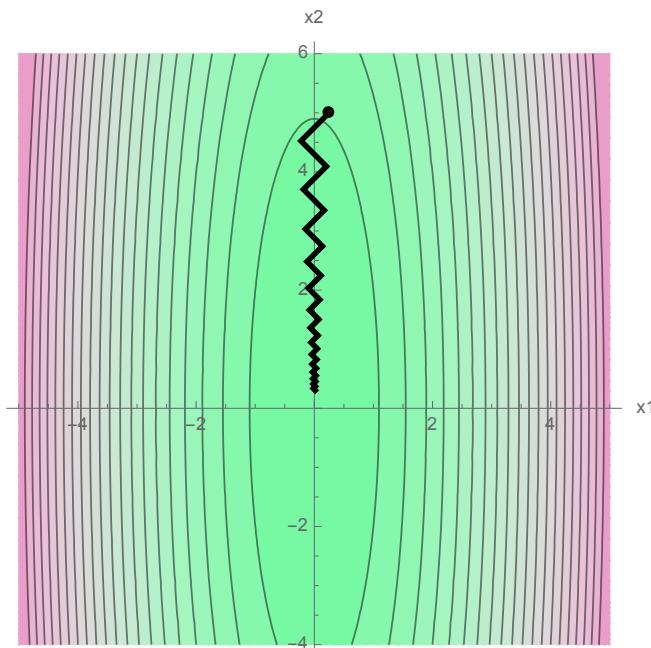
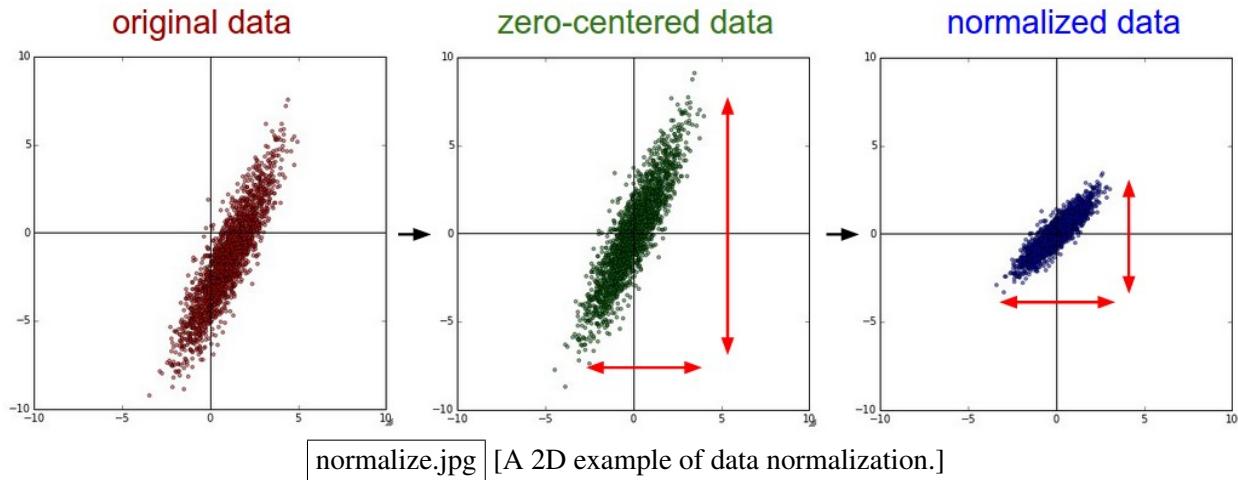
One epoch presents every training example once. Training usually takes many epochs, but if sample is huge [and carries lots of redundant information] it can take less than one.

– Normalizing the data.

– Center each feature so mean is zero.

– Then scale each feature so variance ≈ 1 .

[The first step seems to make it easier for hidden units to get into a good operating region of the sigmoid or ReLU. The second step makes the objective function better conditioned, so gradient descent converges faster.]



illcondition.pdf [Skewed data often leads to an objective function with an ill-conditioned (highly eccentric) Hessian. Gradient descent in these functions can be painfully slow, as this figure shows. Normalizing helps by reducing the eccentricity. Whitening reduces the eccentricity even more, but it's much more expensive. Another thing that helps with elongated troughs like this is momentum, which we'll discuss shortly. It eventually gets us moving fast along the long axis.]

- “Centering” the hidden units helps too.

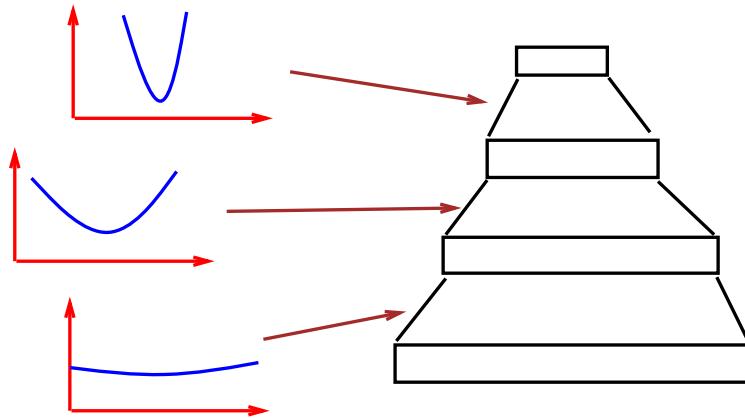
Replace sigmoids with $\tanh \gamma = \frac{e^\gamma - e^{-\gamma}}{e^\gamma + e^{-\gamma}} = 2 s(2\gamma) - 1$.

[This function ranges from -1 to 1 instead of from 0 to 1 .]

[If you use \tanh units, don’t forget that you also need to change backprop to replace s' with the derivative of \tanh . Also, good output target values change to roughly -0.7 and 0.7 .]

- Different learning rate for each layer of weights.

Earlier layers have smaller gradients, need larger learning rate.



curvaturelayers.pdf [In this illustration, the inputs are at the bottom, and the outputs at the top. The derivatives tend to be smaller at the earlier layers.]

- Emphasizing schemes.

[Neural networks learn the most redundant examples quickly, and the most rare examples slowly. So we try to emphasize the uncommon examples.]

- Present examples from rare classes more often, or w/bigger ϵ .
- Same for misclassified examples.

[Be forewarned that emphasizing schemes can backfire if you have really bad outliers.]

- Second-order optimization.

[Unfortunately, Newton’s method is completely impractical, because the Hessian is too large and expensive to compute. There have been a lot of attempts to incorporate curvature information into neural net learning in cheaper ways, but none of them are popular yet.]

- Nonlinear conjugate gradient: works well for small nets + small data + regression.
Batch descent only! → Too slow with redundant data.
- Stochastic Levenberg Marquardt: approximates a diagonal Hessian.
[The authors claim convergence is typically three times faster than well-tuned stochastic gradient descent. The algorithm is complicated.]

- Acceleration schemes: RMSprop, Adam, AMSGrad.

[These are quite popular. Look them up online if you’re curious.]

Heuristics for Avoiding Bad Local Minima

- (1)–(5) above. [Fix the unit saturation problem.]
- Stochastic gradient descent. “Random” motion gets you out of shallow local minima.
[Even if you’re at a local minimum of the cost function, each sample point’s loss function will be trying to push you in a different direction. Stochastic gradient descent looks like a random walk or Brownian motion, which can shake you out of a shallow minimum.]
- Momentum. Gradient descent changes “velocity” Δw slowly.
Carries us through shallow local minima to deeper ones.

```

 $\Delta w \leftarrow -\epsilon \nabla J(w)$ 
repeat
   $w \leftarrow w + \Delta w$ 
   $\Delta w \leftarrow -\epsilon \nabla J(w) + \beta \Delta w$ 

```

Good for both batch & stochastic. Choose hyperparameter $\beta < 1$.

[The hyperparameter β specifies how much momentum persists from iteration to iteration.]

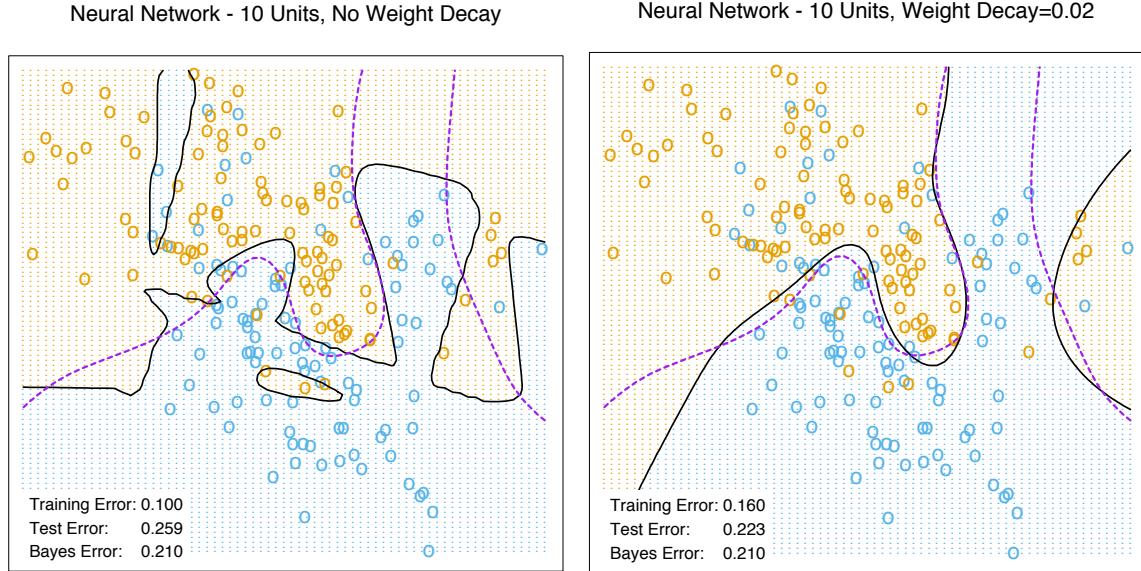
[I’ve seen conflicting advice on β . Some researchers set it to 0.9; some set it close to zero. Geoff Hinton suggests starting at 0.5 and slowly increasing it to 0.9 or higher as the gradients get small.]

[If β is large, you should usually choose ϵ smaller to compensate, and you might also want to change the first line so the initial velocity is larger.]

[A problem with momentum is that once it gets close to a good minimum, it oscillates around the minimum. But it’s more likely to get close to a good minimum in the first place.]

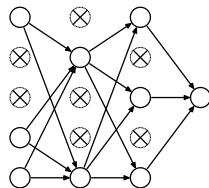
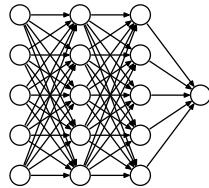
Heuristics to Avoid Overfitting

- Ensemble of neural nets. Random initial weights + bagging.
[We saw how well ensemble learning works for decision trees. It works well for neural nets too. The combination of random initial weights and bagging helps ensure that each neural net comes out differently. Obviously, ensembles of neural nets are slow to train.]
- ℓ_2 regularization, aka weight decay.
Add $\lambda |w'|^2$ to the cost/loss fn, where w' is vector of all weights except biases.
[w' includes all the weights in matrices V and W except the last column of each, rewritten as a vector.]
[We do this for the same reason we do it in ridge regression: penalizing large weights reduces overfitting by reducing the variance of the method.]
Effect: $-\epsilon \frac{\partial L}{\partial w_i}$ has extra term $-2\epsilon\lambda w_i$
Weight decays by factor $1 - 2\epsilon\lambda$ if not reinforced by training.]



weightdecayoff.pdf, weightdecayon.pdf (ESL, Figure 11.4) Write “softmax + cross-entropy loss”. [Examples of 2D classification without (left) and with (right) weight decay. Observe that the second example better approximates the Bayes optimal boundary (dashed purple curve).]

- Dropout emulates an ensemble in one network.



dropout.pdf

[During training, we temporarily disable a random subset of the units, along with all the edges in and out of those units. It seems to work well to disable each hidden unit with probability 0.5, and to disable input units with a smaller probability. We do stochastic gradient descent and we frequently change which random subset of units is disabled. The authors claim that their method gives even better generalization than ℓ_2 regularization. It gives some of the advantages of an ensemble, but it's faster to train.]

- Fewer hidden units.

[The number of hidden units is a hyperparameter you can use to adjust the bias-variance tradeoff. If there's too few, you can't learn well, but if there's too many, you may overfit. ℓ_2 regularization and dropout make it safer to have too many hidden units, so it's less critical to find just the right number.]

CONVOLUTIONAL NEURAL NETWORKS (ConvNets; CNNs)

[Convolutional neural nets have caused a big resurgence of interest in neural nets in the last decade. Often you'll hear the buzzword deep learning, which refers to neural nets with many layers. Many of the most successful deep networks have been convolutional. Just two weeks ago, the ACM announced that the 2018 Alan M. Turing Award is awarded to Geoff Hinton, Yann LeCun, and Yoshua Bengio for their work on deep neural networks.]

Vision: inputs are large images. 200×200 image = 40,000 pixels.

If we connect them all to 40,000 hidden units \rightarrow 1.6 billion connections.

Neural nets are often overparametrized: too many weights, too little data.

[As a rule of thumb, if you have a huge number of weights, you need a huge amount of data, or you risk overfitting. Another problem with having billions of weights is that the network becomes very slow to train or even to use.]

[Researchers have addressed these problems by taking inspiration from the neurology of the visual system. Remember that early in the semester, I told you that you can get better performance on the handwriting recognition task by using edge detectors. Edge detectors have two interesting properties. First, each edge detector looks at just one small part of the image. Second, the edge detection computation is the same no matter which part of the image you apply it to. So let's apply these two properties to neural net design.]

ConvNet ideas:

- (1) Local connectivity: A hidden unit (in early layer) connects only to small patch of units in previous layer.

[This improves the overparametrization problem, and speeds up both the forward pass and the training considerably.]

- (2) Shared weights: Groups of hidden units share same set of input weights, called a mask aka filter aka kernel. [No relation to the kernels of Lecture 14.] We learn several masks.

[Each mask operates on every patch of image.]

Masks \times patches = hidden units in first hidden layer.

If one patch learns to detect edges, *every* patch has an edge detector.

[Because the mask that detects edges is applied to every patch.]

ConvNets exploit repeated structure in images, audio.

Convolution: the same linear transformation applied to different parts of the input by shifting.

[Shared weights improve the overparametrization problem even more, because shared weights means fewer weights. It's a kind of regularization.]

[But shared weights have another big advantage. Suppose that gradient descent starts to develop an edge detector. That edge detector is being trained on *every* part of every image, not just on one spot. And that's good, because edges appear at different locations in different images. The location no longer matters; the edge detector can learn from edges in every part of the image.]

[In a neural net, you can think of hidden units as features that we learn, as opposed to features that you code up yourself. Convolutional neural nets take them to the next level by learning features from multiple patches simultaneously and then applying those features everywhere, not just in the patches where they were originally learned.]

[By the way, although local connectivity is inspired by our visual systems, shared weights obviously don't happen in biology.]

[Show slides on computing in the visual cortex and ConvNets, available from the CS 189 web page at <https://people.eecs.berkeley.edu/~jrs/189/lec/cnn.pdf>. Sorry, readers, there are too many images to include here. The narration is below.]

[Neurologists can stick needles into individual neurons in animal brains. After a few hours the neuron dies, but until then they can record its action potentials. In this way, biologists quickly learned how some of the neurons in the retina, called retinal ganglion cells, respond to light. They have interesting receptive fields, illustrated in the slides, which show that each ganglion cell receives excitatory stimulation from receptors in a small patch of the retina but inhibitory stimulation from other receptors around it.]

[The signals from these cells propagate to the V1 visual cortex in the occipital lobe at the back of your skull. The V1 cells proved much harder to understand. David Hubel and Torsten Wiesel of the Johns Hopkins University put probes into the V1 visual cortex of cats, but they had a very hard time getting any neurons to fire there. However, a lucky accident unlocked the secret and ultimately won them the 1981 Nobel Prize in Physiology.]

[Show video HubelWiesel.mp4, taken from YouTube: <https://www.youtube.com/watch?v=IOHayh06LJ4>]

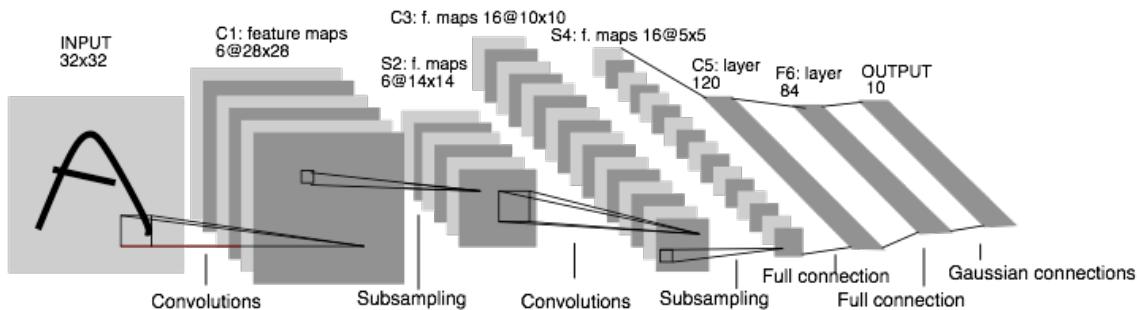
[The glass slide happened to be at the particular orientation the neuron was sensitive to. The neuron doesn't respond to other orientations; just that one. So they were pretty lucky to catch that.]

[The simple cells act as line detectors and/or edge detectors by taking a linear combination of inputs from retinal ganglion cells.]

[The complex cells act as location-independent line detectors by taking inputs from many simple cells, which are location dependent.]

[Later researchers showed that local connectivity runs through the V1 cortex by projecting certain images onto the retina and using radioactive tracers in the cortex to mark which neurons had been firing. Those images show that the neural mapping from the retina to V1 is "retinotopic," i.e., locality preserving. This is a big part of the inspiration for convolutional neural networks!]

[Unfortunately, as we go deeper into the visual system, layers V2 and V3 and so on, we know less and less about what processing the visual cortex does.]



LeNet5.png Architecture of LeNet5.

[ConvNets were first popularized by the success of Yann LeCun's "LeNet 5" handwritten digit recognition software. LeNet 5 has six hidden layers! Layers 1 and 3 are convolutional layers in which groups of units share weights. Layers 2 and 4 are pooling layers that make the image smaller. These are just hardcoded max-functions with no weights and nothing to train. Layers 5 and 6 are just regular layers of hidden units with no shared weights. A great deal of experimentation went into figuring out the number of layers and their sizes. At its peak, LeNet 5 was responsible for reading the zip codes on 10% of US Mail. Another Yann LeCun system was deployed in ATMs and check reading machines and was reading 10 to 20% of all the checks in the US by the late 90's. LeCun is one of the Turing Award winners I told you about earlier.]

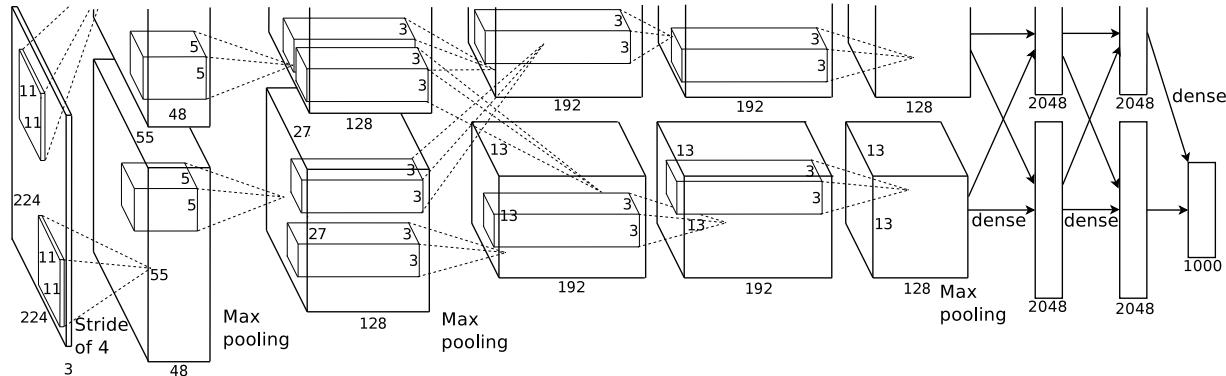
[Show Yann LeCun's video LeNet5.mov, illustrating LeNet 5.]

[When ConvNets were first applied to image analysis, researchers found that some of the learned masks are edge detectors or line detectors, similar to the ones that Hubel and Wiesel discovered! This created a lot of excitement in both the computer learning community and the neuroscience community. The fact that a neural net can naturally learn the same features as the mammalian visual cortex is impressive.]

[I told you two lectures ago that neural nets research was popular in the 60's, but the 1969 book *Perceptrons* killed interest in them throughout the 70's. They came back in the 80's, but interest was partly killed off a second time in the 00's by ... guess what? By support vector machines. SVMs work well for a lot of tasks, they're much faster to train, and they more or less have only one hyperparameter, whereas neural nets take a lot of work to tune.]

[Neural nets are now in their third wave of popularity. The single biggest factor in bringing them back is probably big data. Thanks to the internet, we now have absolutely huge collections of images to train neural nets with, and researchers have discovered that neural nets often give better performance than competing algorithms when you have huge amounts of data to train them with. In particular, convolutional neural nets are now learning better features than hand-tuned features. That's a recent change.]

[One event that brought attention back to neural nets was the ImageNet Image Classification Challenge in 2012. The winner of that competition was a neural net, and it won by a huge margin, about 10%. It's called AlexNet, and it's surprisingly similarly to LeNet 5, in terms of how its layers are structured. However, there are some new innovations that led to their prize-winning performance, in addition to the fact that the training set had 1.4 million images: they used ReLUs, dropout, and GPUs for training.]



[alexnet.pdf](#) Architecture of AlexNet.

[If you want to learn more about deep neural networks, there's a whole new undergraduate class at Berkeley just for you: CS 182.]

20 Unsupervised Learning and Principal Components Analysis

UNSUPERVISED LEARNING

We have sample points, but no labels!

No classes, no y -values, nothing to predict.

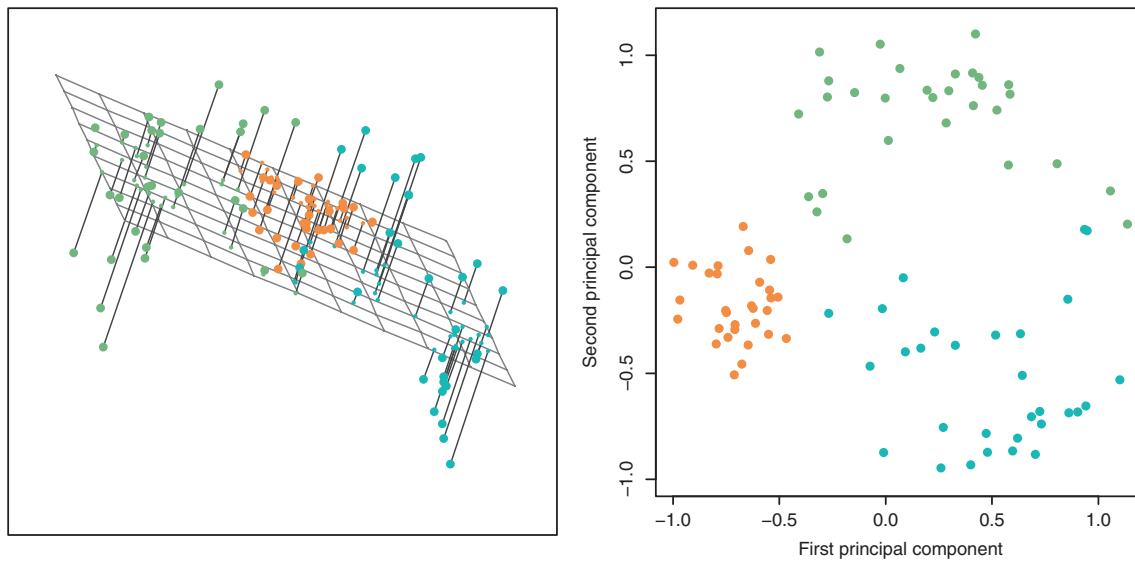
Goal: Discover structure in the data.

Examples:

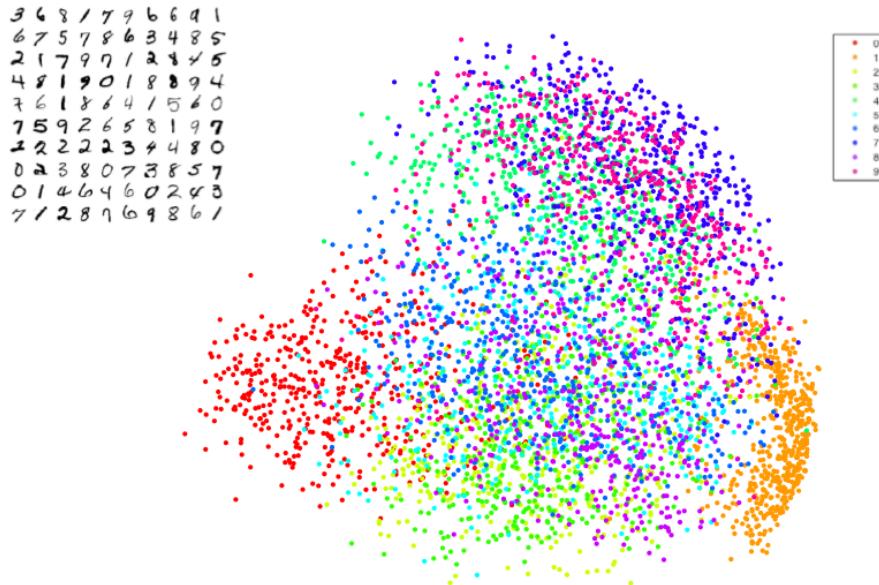
- Clustering: partition data into groups of similar/nearby points.
- Dimensionality reduction: data often lies near a low-dimensional subspace (or manifold) in feature space; matrices have low-rank approximations.
[Whereas clustering is about grouping similar sample points, dimensionality reduction is more about identifying a continuous variation from sample point to sample point.]
- Density estimation: fit a continuous distribution to discrete data.
[When we use maximum likelihood estimation to fit Gaussians to sample points, that's density estimation, but we can also fit functions more complicated than Gaussians, with more local variation.]

PRINCIPAL COMPONENTS ANALYSIS (PCA) (Karl Pearson, 1901)

Goal: Given sample points in \mathbb{R}^d , find k directions that capture most of the variation. (Dimensionality reduction.)



3dpca.pdf [Example of 3D points projected to 2D by PCA.]



[pcadigits.pdf] [The (high-dimensional) MNIST digits projected to 2D. Two dimensions aren't enough to fully separate the digits, but observe that the digits 0 (red) and 1 (orange) are well on their way to being separated.]

Why?

- Find a small basis for representing variations in complex things, e.g. faces, genes.
 - Reducing # of dimensions makes some computations cheaper, e.g. regression.
 - Remove irrelevant dimensions to reduce overfitting in learning alg.
- Like subset selection, but the “features” aren't axis-aligned;
they're linear combos of input features.

[Sometimes PCA is used as a preprocess before regression or classification for the last two reasons.]

Let X be $n \times d$ design matrix. [No fictitious dimension.]

From now on, assume X is centered: mean X_i is zero.

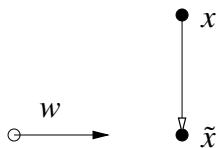
[As usual, we can center the data by computing the mean x -value, then subtracting the mean from each sample point.]

[Let's start by seeing what happens if we pick just one principal direction.]

Let w be a unit vector.

The orthogonal projection of point x onto vector w is $\tilde{x} = (x \cdot w) w$

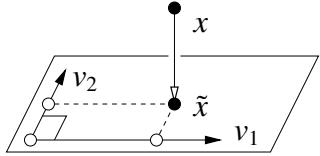
If w not unit, $\tilde{x} = \frac{x \cdot w}{|w|^2} w$



[The idea is that we're going to pick the best direction w , then project all the data down onto w so we can analyze it in a one-dimensional space. Of course, we lose a lot of information when we project down from d dimensions to just one. So, suppose we pick several directions. Those directions span a subspace, and we want to project points orthogonally onto the subspace. This is easy if the directions are orthogonal to each other.]

Given orthonormal directions v_1, \dots, v_k , $\tilde{x} = \sum_{i=1}^k (x \cdot v_i) v_i$

[The word “orthonormal” implies they’re mutually orthogonal and length 1.]



[Usually we don’t actually want the projected point in \mathbb{R}^d ; usually we want the principle coordinates $x \cdot v_i$ in principal components space.]

$X^\top X$ is square, symmetric, positive semidefinite, $d \times d$ matrix. [As it’s symmetric, its eigenvalues are real.]

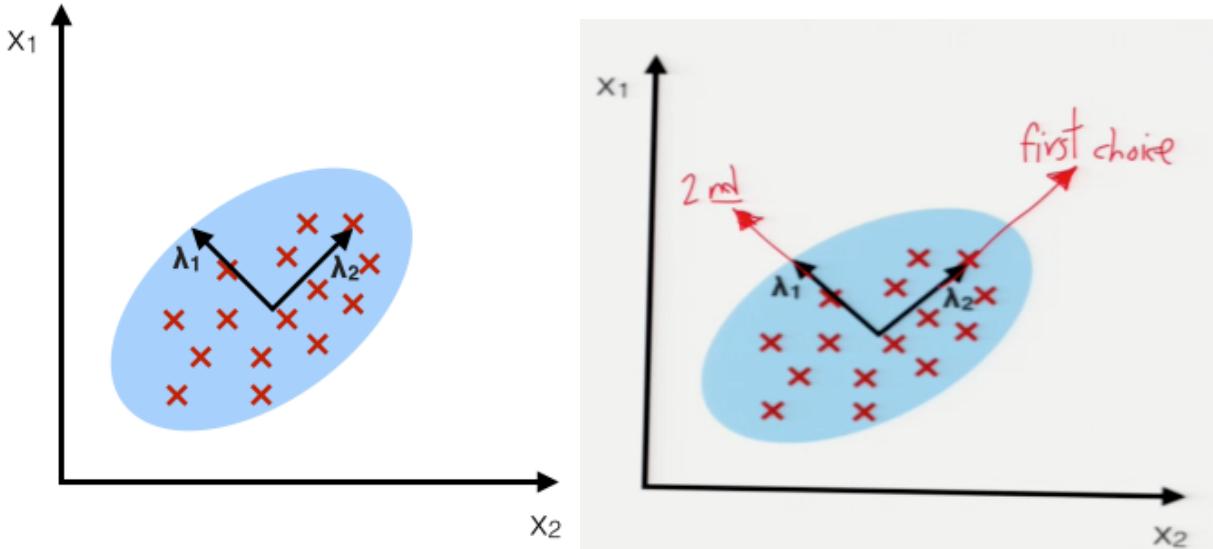
Let $0 \leq \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_d$ be its eigenvalues. [sorted]

Let v_1, v_2, \dots, v_d be corresponding orthogonal **unit** eigenvectors.

[It turns out that the principal directions will be these eigenvectors, and the most important ones will be the ones with the greatest eigenvalues. I will show you this in three different ways.]

PCA derivation 1: Fit a Gaussian to data with maximum likelihood estimation.

Choose k Gaussian axes of greatest variance.



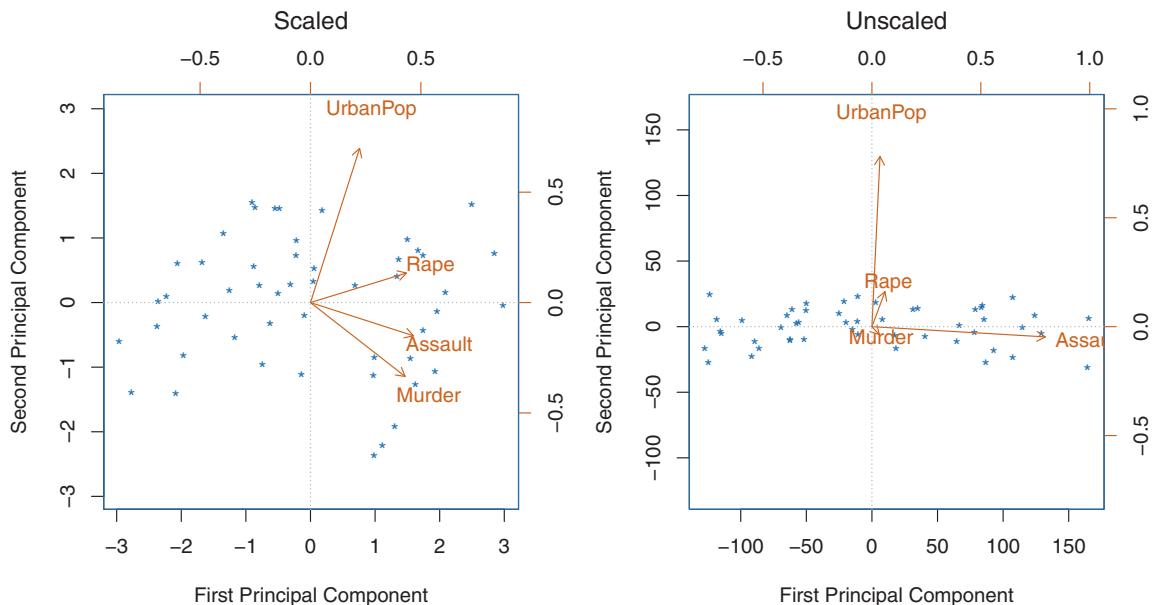
gaussfitpca.png [A Gaussian fitted to sample points.]

Recall that MLE estimates a covariance matrix $\hat{\Sigma} = \frac{1}{n} X^\top X$. [Presuming X is centered.]

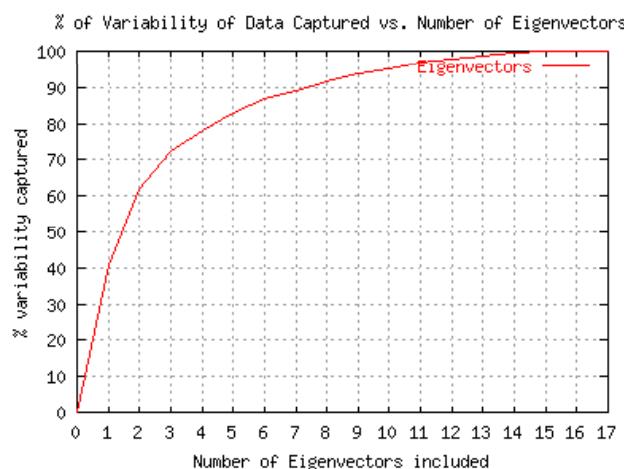
PCA Alg:

- Center X .
- Optional: Normalize X . Units of measurement different?
 - Yes: Normalize.
[Bad for principal components to depend on arbitrary choice of scaling.]
 - No: Usually don’t.
[If several features have the same unit of measurement, but some of them have smaller variance than others, that difference is usually meaningful.]
- Compute unit eigenvectors/values of $X^\top X$.

- Optional: choose k based on the eigenvalue sizes.
- For the best k -dimensional subspace, pick eigenvectors v_{d-k+1}, \dots, v_d .
- Compute the coordinates $x \cdot v_i$ of training/test data in principal components space.
[When we do this projection, we have two choices: we can un-center the input training data before projecting it, OR we can translate the test data by the same vector we used to translate the training data when we centered it.]



normalize.pdf [Projection of 4D data onto a 2D subspace. Normalized data at left; unnormalized data at right. The arrows show the four original axes projected on the two principal components. When the data are not normalized, rare occurrences like murder have little influence on the principal directions. Which is better? It depends on whether you think that low-frequency events like murder and rape should have a disproportionate influence.]

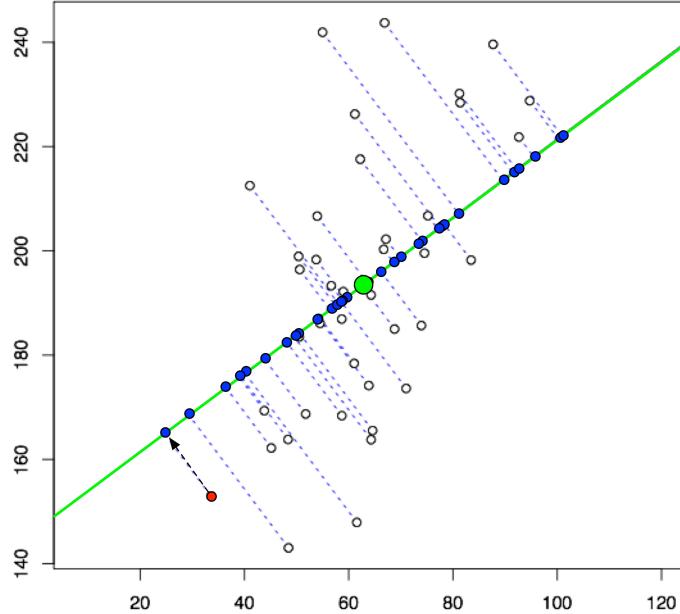


$$\% \text{ of variability} = \frac{\sum_{i=1}^k \lambda_i}{\sum_{i=1}^n \lambda_i}$$

variance.pdf [Plot of # of eigenvectors vs. percentage of sample variance captured for a 17D data set. In this example, just 3 eigenvectors capture 70% of the variance.]

[If you are using PCA as a preprocess for a supervised learning algorithm, there's a more effective way to choose k : (cross-)validation.]

PCA derivation 2: Find direction w that maximizes sample variance of projected data
 [In other words, when we project the data down, we don't want it all to bunch up; we want to keep it as spread out as possible.]



project.jpg [Points projected on a line. We wish to choose the orientation of the green line to maximize the sample variance of the blue points.]

$$\text{Maximize } \text{Var}(\{\tilde{X}_1, \tilde{X}_2, \dots, \tilde{X}_n\}) = \frac{1}{n} \sum_{i=1}^n \left(X_i \cdot \frac{w}{|w|} \right)^2 = \frac{1}{n} \frac{|Xw|^2}{|w|^2} = \frac{1}{n} \underbrace{\frac{w^\top X^\top X w}{w^\top w}}_{\text{Rayleigh quotient of } X^\top X \text{ and } w}$$

[This fraction is a well-known construction called the Rayleigh quotient. When you see it, you should smell eigenvectors nearby. How do we maximize this?]

If w is an eigenvector v_i , Ray. quo. = λ_i

→ of all eigenvectors, v_d achieves maximum variance λ_d/n .

One can show v_d beats every other vector too.

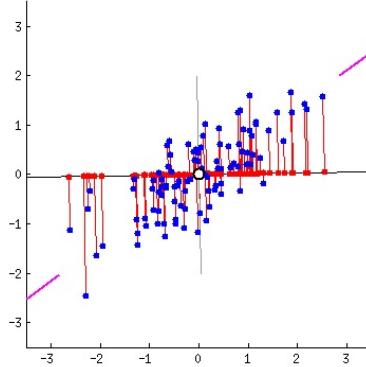
[Because every vector w is a linear combination of eigenvectors, and so its Rayleigh quotient will be a convex combination of eigenvalues. It's easy to prove this, but I don't have the time. For the proof, look up "Rayleigh quotient" in Wikipedia.]

[So the top eigenvector gives us the best direction. But we typically want k directions. After we've picked one direction, then we have to pick a direction that's orthogonal to the best direction. But subject to that constraint, we again pick the direction that maximizes the sample variance.]

What if we constrain w to be orthogonal to v_d ? Then v_{d-1} is optimal.

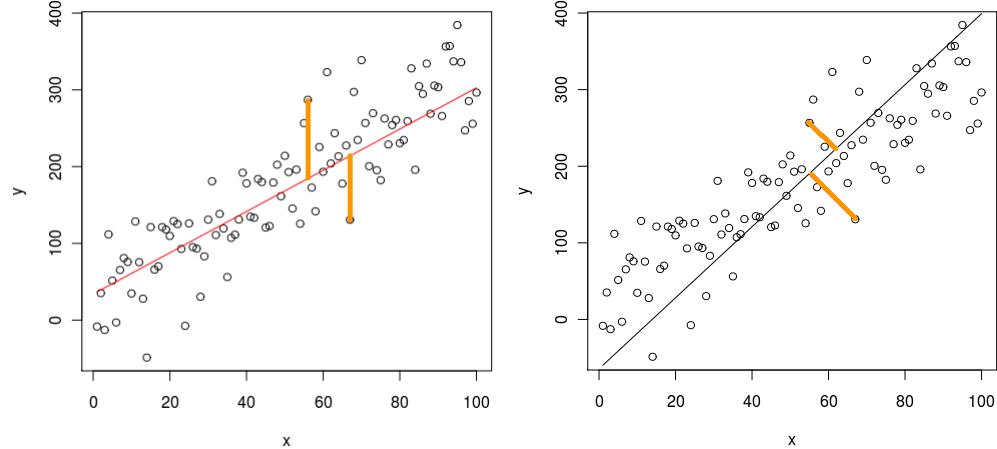
[And if we need a third direction orthogonal to v_d and v_{d-1} , the optimal choice is v_{d-2} . And so on.]

PCA derivation 3: Find direction w that minimizes “projection error”



PCAnimation.gif [This is an animated GIF; unfortunately, the animation can't be included in the PDF lecture notes. Find the direction of the black line for which the sum of squares of the lengths of the red lines is smallest.]

[You can think of this as a sort of least-squares linear regression, with one subtle but important change. Instead of measuring the error in a fixed vertical direction, we're measuring the error in a direction orthogonal to the principal component direction we choose.]

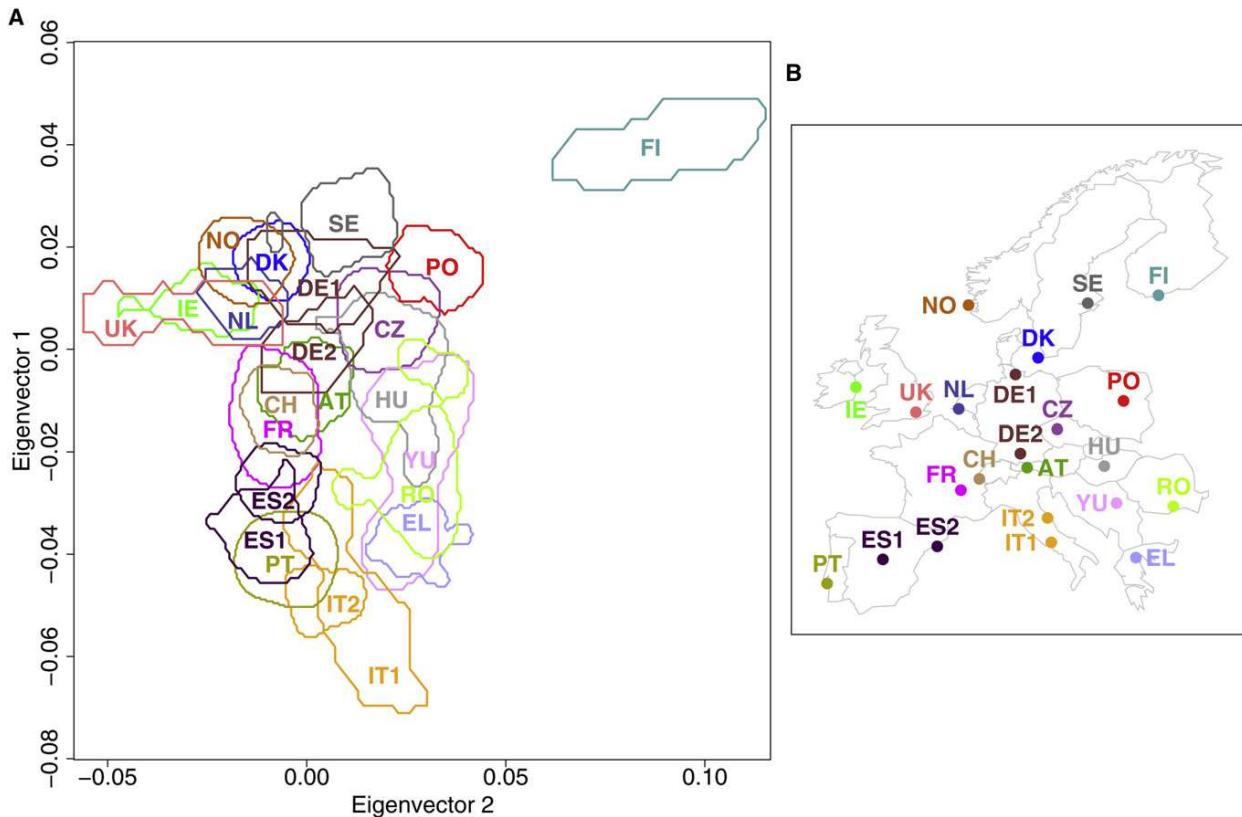


projlsq.png, projpca.png [Least-squares linear regression vs. PCA. In linear regression, the projection direction is always vertical; whereas in PCA, the projection direction is orthogonal to the projection hyperplane. In both methods, however, we minimize the sum of the squares of the projection distances.]

$$\begin{aligned} \text{Minimize } \sum_{i=1}^n |X_i - \tilde{X}_i|^2 &= \sum_{i=1}^n \left| X_i - \frac{X_i \cdot w}{|w|^2} w \right|^2 = \sum_{i=1}^n \left(|X_i|^2 - \left(X_i \cdot \frac{w}{|w|} \right)^2 \right) \\ &= \text{constant} - n \text{ (variance from derivation 2).} \end{aligned}$$

Minimizing projection error = maximizing variance.

[From this point, we carry on with the same reasoning as derivation 2.]



europegenetics.pdf (Lao et al., Current Biology, 2008.) [Illustration of the first two principal components of the single nucleotide polymorphism (SNP) matrix for the genes of various Europeans. The input matrix has 2,541 people from these locations in Europe (right), and 309,790 SNPs. Each SNP is binary, so think of it as 309,790 dimensions of zero or one. The output (left) shows spots on the first two principal components where there was a high density of projected people from a particular national type. What's amazing about this is how closely the projected genotypes resemble the geography of Europe.]

Eigenfaces

X contains n images of faces, d pixels each.

[If we have a 200×200 image of a face, we represent it as a vector of length 40,000, the same way we represent the MNIST digit data.]

Face recognition: Given a query face, compare it to all training faces; find nearest neighbor in \mathbb{R}^d .

[This works best if you have several training photos of each person you want to recognize, with different lighting and different facial expressions.]

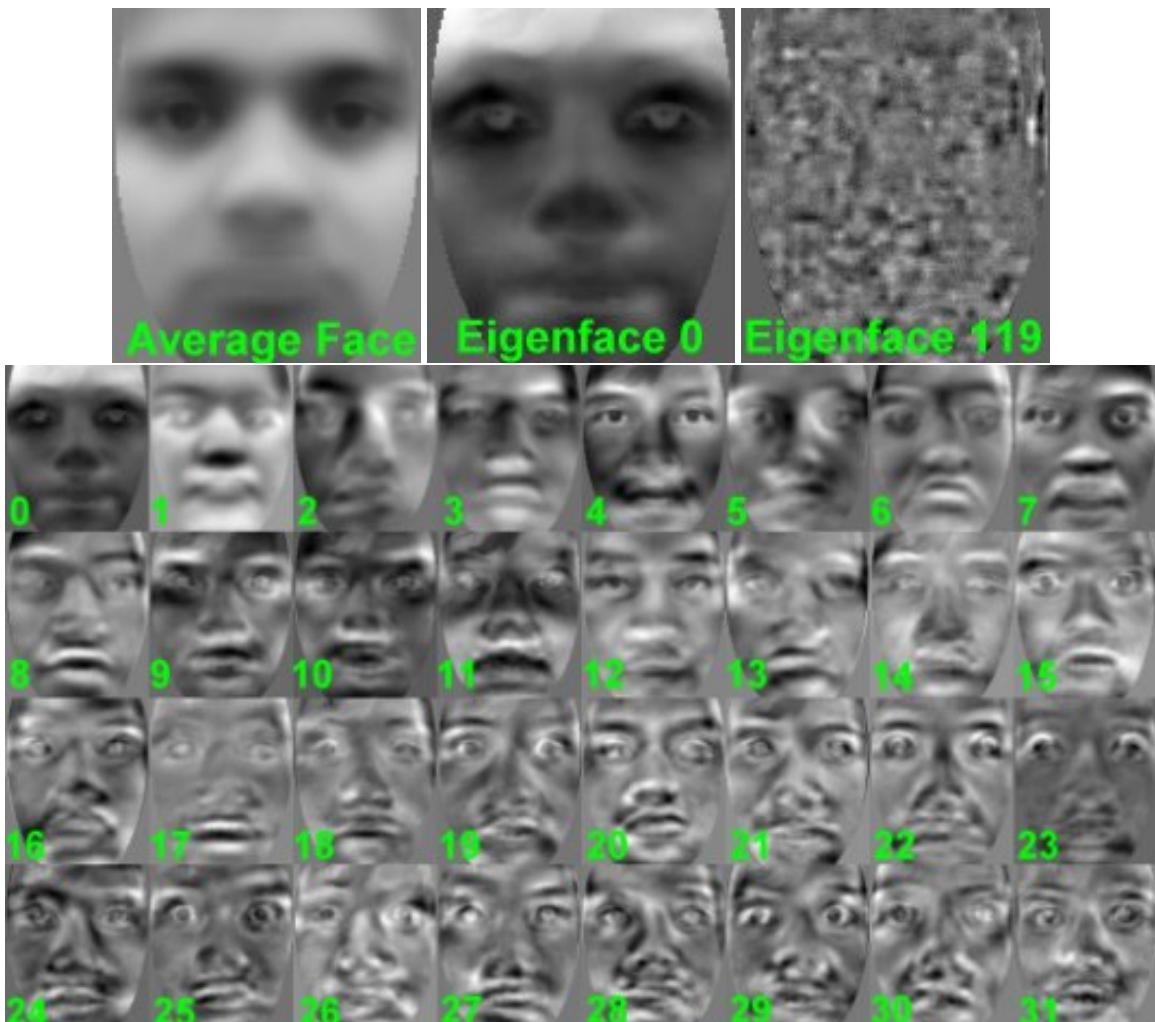
Problem: Each query takes $\Theta(nd)$ time.

Solution: Run PCA on faces. Reduce to much smaller dimension d' .

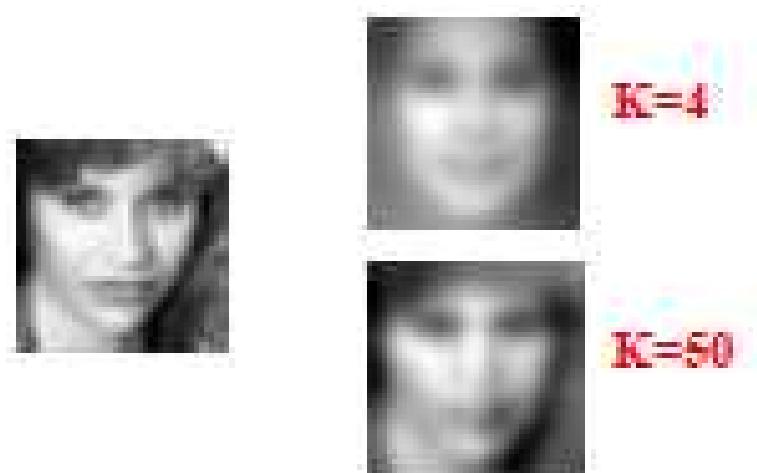
Now nearest neighbor takes $O(nd')$ time.

[Possibly even less. We'll talk about speeding up nearest-neighbor search at the end of the semester. If the dimension is small enough, you can sometimes do better than linear time.]

[If you have 500 stored faces with 40,000 pixels each, and you reduce them to 40 principal components, then each query face requires you to read 20,000 stored principle coordinates instead of 20 million pixels.]



[*facerecaverage.jpg*, *facereceigen0.jpg*, *facereceigen119.jpg*, *facereceigen.jpg*] [Images of the eigenfaces. The “average face” is the mean used to center the data.]



[*eigenfaceproject.pdf*] [Images of a face (left) projected onto the first 4 and 50 eigenvectors, with the average face added back. These last image is blurry but good enough for face recognition.]

For best results, equalize the intensity distributions first.



facerecequalize.jpg [Image equalization.]

[Eigenfaces are not perfect. They encode both face shape *and* lighting. Ideally, we would have some way to factor out lighting and analyze face shape only, but that's harder. Some people say that the first 3 eigenfaces are usually all about lighting, and you sometimes get better facial recognition by dropping the first 3 eigenfaces.]

[Show Blanz–Vetter face morphing video (morphmod.mpg).]

[Blanz and Vetter use PCA in a more sophisticated way for 3D face modeling. They take 3D scans of people's faces and find correspondences between peoples' faces and an idealized model. For instance, they identify the tip of your nose, the corners of your mouth, and other facial features, which is something the original eigenface work did not do. Instead of feeding an array of pixels into PCA, they feed the 3D locations of various points on your face into PCA. This works more reliably.]

21 The Singular Value Decomposition; Clustering

The Singular Value Decomposition (SVD) [and its Application to PCA]

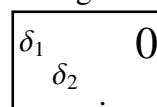
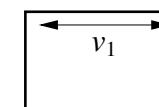
Problems: Computing $X^\top X$ takes $\Theta(nd^2)$ time.

$X^\top X$ is poorly conditioned \rightarrow numerically inaccurate eigenvectors.

[The SVD improves both these problems.]

[Earlier this semester, we learned about the eigendecomposition of a square, symmetric matrix. Unfortunately, nonsymmetric matrices don't eigendecompose nearly as nicely, and non-square matrices don't have eigenvectors at all. Happily, there is a similar decomposition that works for all matrices, even if they're not symmetric and not square.]

Fact: If $n \geq d$, we can find a singular value decomposition $X = UDV^\top$

X	=	U	D diagonal	V^\top	$= \sum_{i=1}^d \underbrace{\delta_i u_i v_i^\top}_{\text{rank 1}} \text{outer product matrix}$
	=	 u_1 u_d	 δ_1 δ_2 \ddots δ_d	 v_1 v_d	
$n \times d$	=	$n \times d$	$d \times d$	$d \times d$	

$U^\top U = I$

orthonormal u_i 's are left singular vectors of X

$V^\top V = I$

orthonormal v_i 's are right singular vectors of X

Diagonal entries $\delta_1, \dots, \delta_d$ of D are nonnegative singular values of X .

[Some of the singular values might be zero. The number of nonzero singular values is equal to the rank of the centered design matrix X . If all the sample points lie on a line, there is only one nonzero singular value. If the points span a subspace of dimension r , there are r nonzero singular values.]

[If $n < d$, an SVD still exists, but now U is square and V is not.]

Fact: v_i is an eigenvector of $X^\top X$ w/eigenvalue δ_i^2 .

$$\text{Proof: } X^T X = V D U^T U D V^T = V D^2 V^T$$

which is an eigendecomposition of $X^\top X$.

[The columns of V are the eigenvectors of $X^T X$, which is what we need for PCA. If $n < d$, V will omit some of the eigenvectors that have eigenvalue zero, but those are useless for PCA. The SVD also tells us the eigenvalues, which are the squares of the singular values. By the way, that's related to why the SVD is more numerically stable: the ratios between singular values are smaller than the ratios between eigenvalues.]

Fact: We can find the k greatest singular values & corresponding vectors in $O(ndk)$ time.
 [So we can save time by computing some of the singular vectors without computing all of them.]
 [There are approximate, randomized algorithms that are even faster, producing an approximate SVD in $O(nd \log k)$ time. These are starting to become popular in algorithms for very big data.]
 [<https://code.google.com/archive/p/redsvd/>]

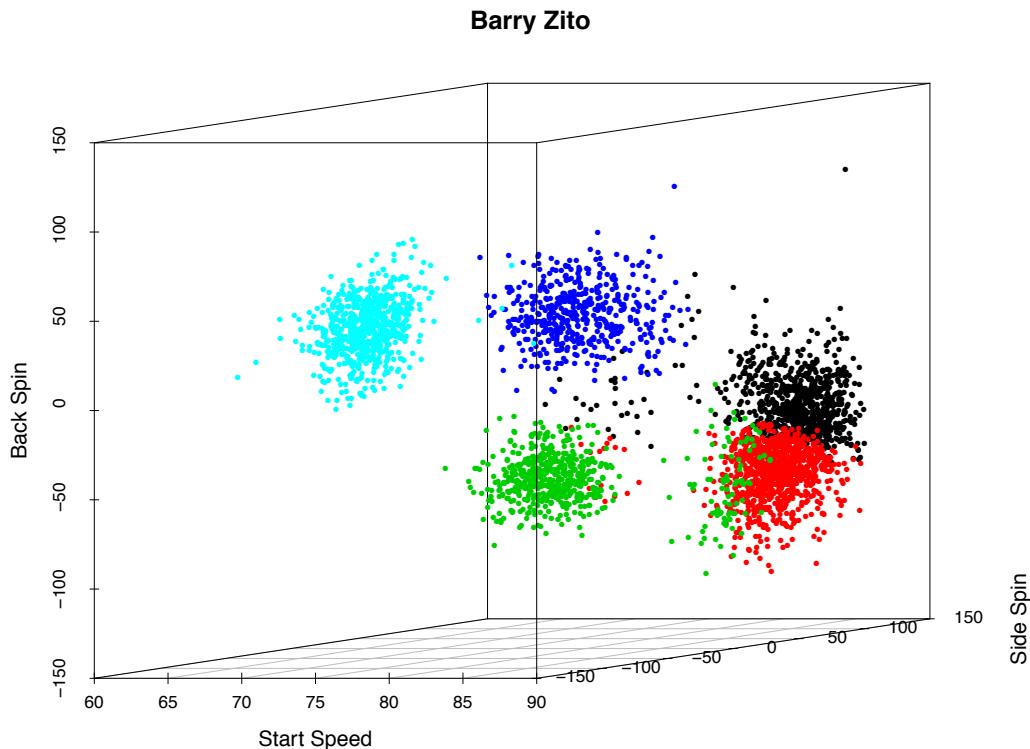
Important: Row i of UD gives the coordinates of sample point X_i in principal components space (i.e. $X_i \cdot v_j$ for each j).
 [So we don't need to explicitly project the input points onto principal components space; the SVD has already done it for us.]

CLUSTERING

Partition data into clusters so points in a cluster are more similar than across clusters.

Why?

- Discovery: Find songs similar to songs you like; determine market segments
- Hierarchy: Find good taxonomy of species from genes
- Quantization: Compress a data set by reducing choices
- Graph partitioning: Image segmentation; find groups in social networks



4-Seam Fastball	2-Seam Fastball	Changeup	Slider	Curveball
Black	Red	Green	Blue	Light Blue

zito.pdf (from a talk by Michael Pane) [k -means clusters that classify Barry Zito's baseball pitches. Here we discover that there really are distinct classes of baseball pitches.]

k -Means Clustering aka Lloyd's Algorithm (Stuart Lloyd, 1957)

Goal: Partition n points into k disjoint clusters.

Assign each input point X_i a cluster label $y_i \in [1, k]$.

Cluster i 's mean is $\mu_i = \frac{1}{n_i} \sum_{y_j=i} X_j$, given n_i points in cluster i .

Find y that minimizes $\sum_{i=1}^k \sum_{y_j=i} X_j - \mu_i ^2$	[Sum of the squared distances from points to their cluster means.]
---	--

NP-hard. Solvable in $O(nk^n)$ time. [Try every partition.]

k -means heuristic: Alternate between

(1) y_j 's are fixed; update μ_i 's

(2) μ_i 's are fixed; update y_j 's

Halt when step (2) changes no assignments.

[So, we have an assignment of points to clusters. We compute the cluster means. Then we reconsider the assignment. A point might change clusters if some other's cluster's mean is closer than its own cluster's mean. Then repeat.]

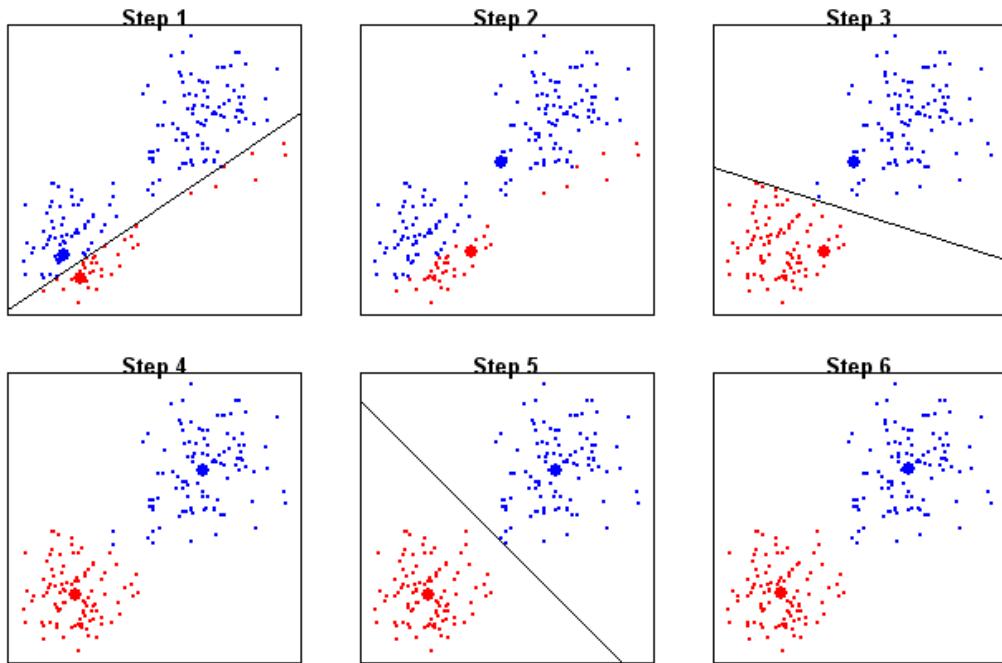
Step (1): One can show (calculus) the optimal μ_i is the mean of the points in cluster i .

[This is easy calculus, so I leave it as a short exercise.]

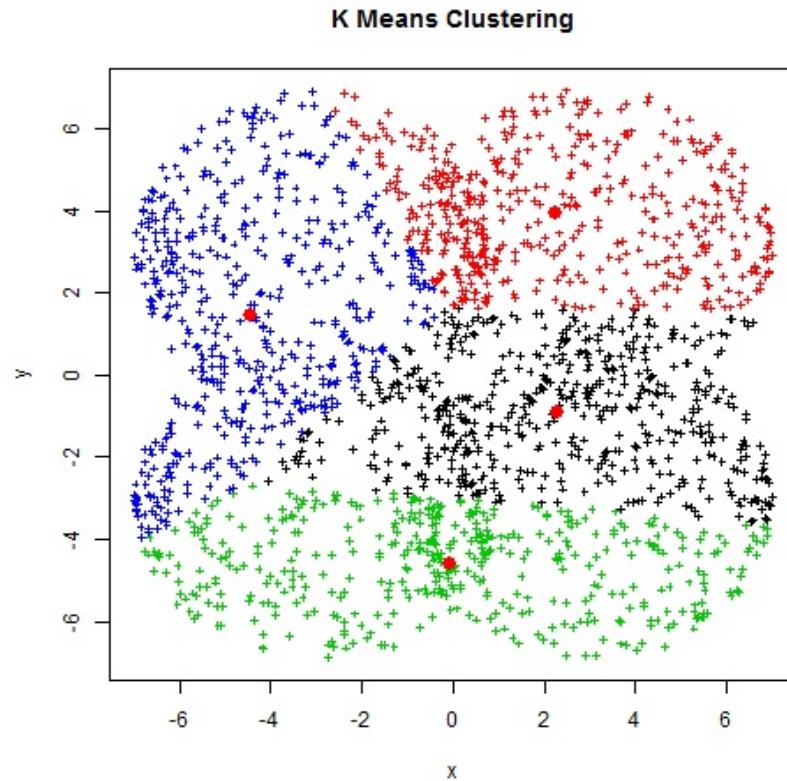
Step (2): The optimal y assigns each point X_j to the closest center μ_i .

[This should be even more obvious than step (1).]

[If there's a tie, and one of the choices is for X_j to stay in the same cluster as the previous iteration, always take that choice.]



2means.png [An example of 2-means. Odd-numbered steps reassigned the data points. Even-numbered steps compute new means.]



4meansanimation.gif [This is an animated GIF of 4-means with many points. Unfortunately, the animation can't be included in the PDF lecture notes.]

Both steps decrease objective fn *unless* they change nothing.

[Therefore, the algorithm never returns to a previous assignment.]

Hence alg. must terminate. [As there are only finitely many assignments.]

[This argument doesn't say anything optimistic about the running time, because we might see $O(k^n)$ different assignments before we halt. In theory, one can actually construct point sets in the plane that take an exponential number of iterations, but those don't come up in practice.]

Usually very fast in practice. Finds a local minimum, often not global.

[... which is not surprising, as this problem is NP-hard.]



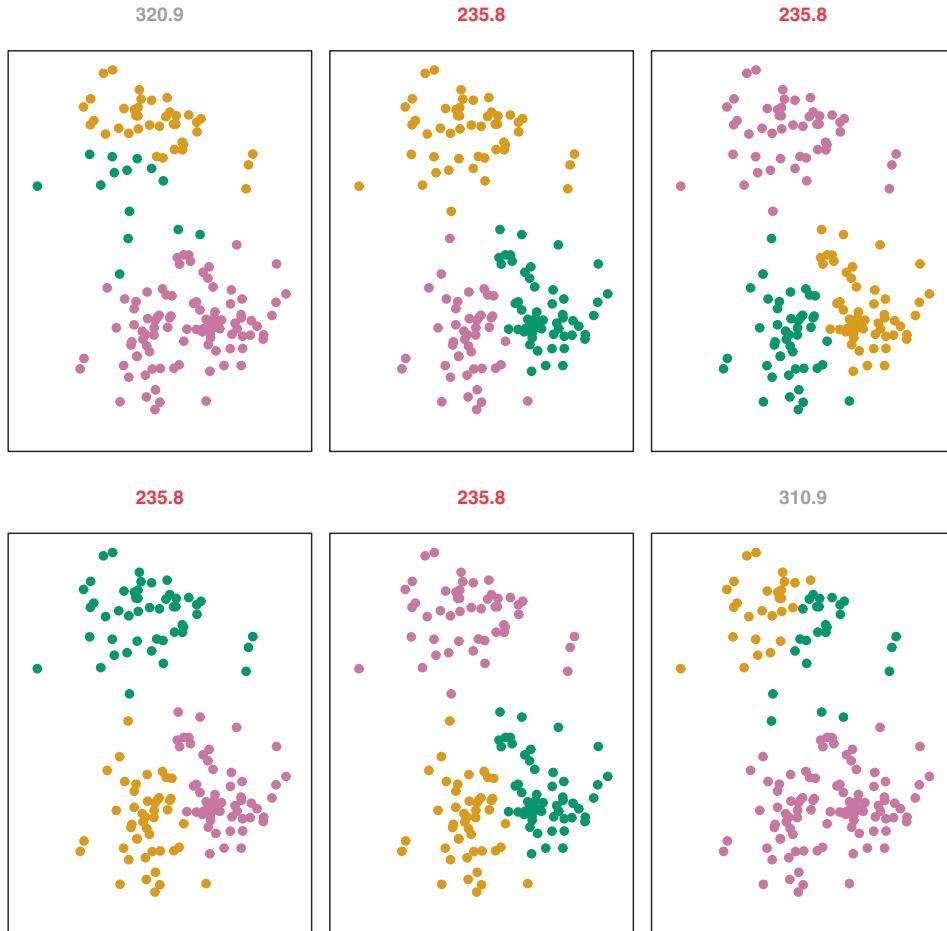
4meansbad.png [An example where 4-means clustering fails.]

Getting started:

- Forgy method: choose k random sample points to be initial μ_i 's; go to (2).
- Random partition: randomly assign each sample point to a cluster; go to (1).
- k -means++: like Forgy, but biased distribution. [Each center is chosen with a preference for points far from previous centers.]

[k -means++ is a little more work, but it works well in practice and theory. Forgy seems to be better than random partition, but Wikipedia mentions some variants of k -means for which random partition is better.]

For best results, run k -means multiple times with random starts.



`kmeans6times.pdf` (ISL, Figure 10.7) [Clusters found by running 3-means 6 times on the same sample points. The algorithm finds three different local minima.]

[Why did we choose that particular objective function to minimize? Partly because it is equivalent to minimizing the following function.]

Equivalent objective fn: the within-cluster variation

$$\text{Find } y \text{ that minimizes } \sum_{i=1}^k \frac{1}{n_i} \sum_{y_j=i} \sum_{y_m=i} |X_j - X_m|^2$$

[At the minimizer, this objective function is equal to twice the previous one. It's a worthwhile exercise to show that—it's harder than it looks. The nice thing about this expression is that it doesn't include the means; it's a function purely of the input points and the clusters we assign them to.]

Normalize the data? [before applying k -means]

Same advice as for PCA. Sometimes yes, sometimes no.

[If some features are much larger than others, they will tend to dominate the Euclidean distance. So if you have features in different units of measurement, you probably should normalize them. If you have features in the same unit of measurement, you usually shouldn't, but it depends on context.]

***k*-Medoids Clustering**

Generalizes k -means beyond Euclidean distance. [Means aren't optimal for other distance metrics.]

Specify a distance fn $d(x, y)$ between points x, y , aka dissimilarity.

Can be arbitrary; ideally satisfies triangle inequality $\overline{d(x, y)} \leq d(x, z) + d(z, y)$.

[Sometimes people use the ℓ_1 norm or the ℓ_∞ norm. Sometimes people specify a matrix of pairwise distances between the input points.]

[Suppose you have a database that tells you how many of each product each customer bought. You'd like to cluster together customers who buy similar products for market analysis. But if you cluster customers by Euclidean distance, you'll get a big cluster of all the customers who have only ever bought one thing. So Euclidean distance is not a good measure of dissimilarity. Instead, it makes more sense to treat each customer as a vector and measure the *angle* between two customers. If there's a large angle between customers, they're dissimilar.]

Replace mean with medoid, the sample point that minimizes total distance to other points in same cluster. [So the medoid of a cluster is always one of the input points.]

[One difficulty with k -means is that you have to choose the number k of clusters before you start, and there isn't any reliable way to guess how many clusters will best fit the data. The next method, hierarchical clustering, has the advantage in that respect. By the way, there is a whole Wikipedia article on "Determining the number of clusters in a data set."]

Hierarchical Clustering

Creates a tree; every subtree is a cluster.

[So some clusters contain smaller clusters.]

Bottom-up, aka agglomerative clustering:

start with each point a cluster; repeatedly fuse pairs.

[Draw figure of points in the plane; pair clusters together until all points are in one top-level cluster.]

Top-down, aka divisive clustering:

start with all pts in one cluster; repeatedly split it.

[Draw figure of points in the plane; divide points into subsets hierarchically until each point is in its own subset.]

[When the input is a point set, agglomerative clustering is used much more in practice than divisive clustering. But when the input is a graph, it's the other way around: divisive clustering is more common.]

We need a distance fn for clusters A, B :

complete linkage: $d(A, B) = \max\{d(w, x) : w \in A, x \in B\}$

single linkage: $d(A, B) = \min\{d(w, x) : w \in A, x \in B\}$

average linkage: $d(A, B) = \frac{1}{|A||B|} \sum_{w \in A} \sum_{x \in B} d(w, x)$

centroid linkage: $d(A, B) = d(\mu_A, \mu_B)$ where μ_S is mean of S

[The first three of these linkages work for any distance function, even if the input is just a matrix of distances between all pairs of points. The centroid linkage only really makes sense if we're using the Euclidean distance. But there's a variation of the centroid linkage that uses the medoids instead of the means, and medoids are defined for any distance function. Moreover, medoids are more robust to outliers than means.]

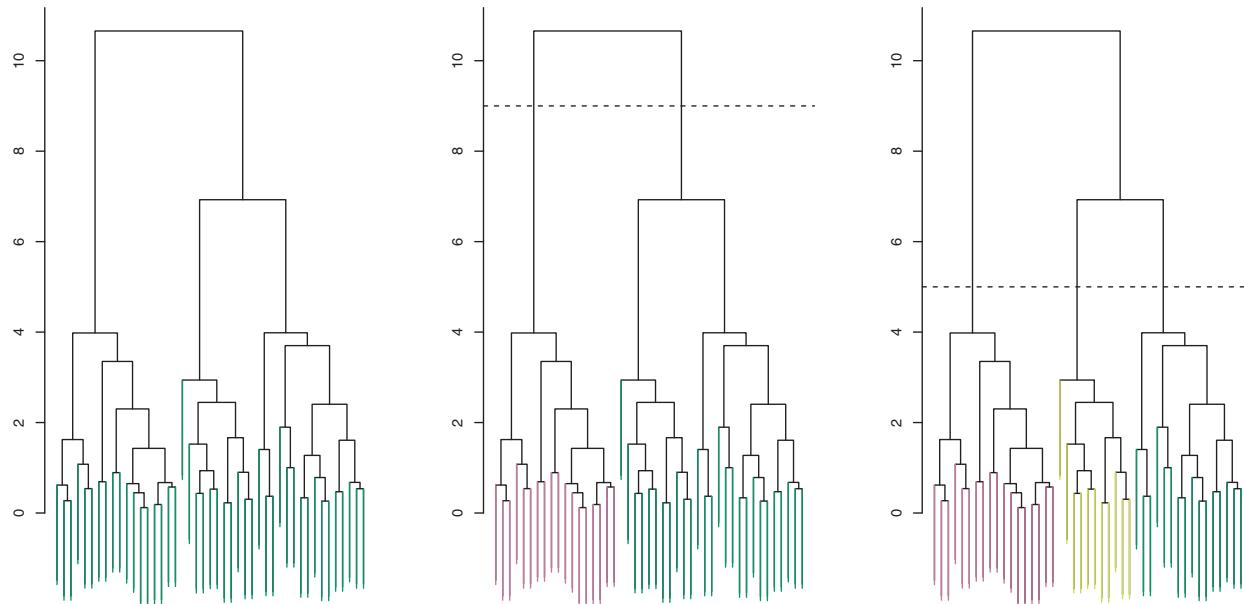
Greedy agglomerative alg.:

Repeatedly fuse the two clusters that minimize $d(A, B)$

Naively takes $O(n^3)$ time.

[But for complete and single linkage, there are more sophisticated algorithms called CLINK and SLINK, which run in $O(n^2)$ time. A package called ELKI has publicly available implementations.]

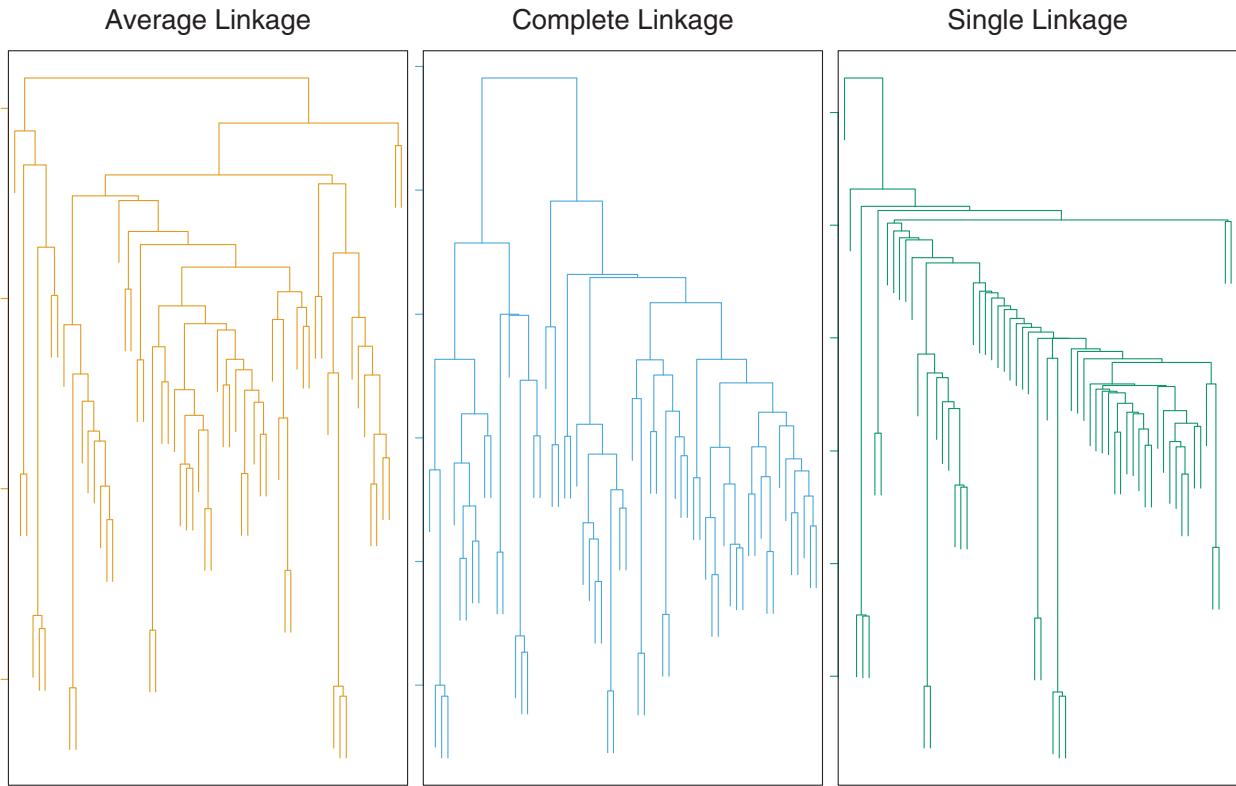
Dendrogram: Illustration of the cluster hierarchy (tree) in which the vertical axis encodes all the linkage distances.



dendrogram.pdf (ISL, Figure 10.9) [Example of a dendrogram cut into 1, 2, or 3 clusters.]

Cut dendrogram into clusters by horizontal line according to your choice of # of clusters OR intercluster distance.

[It's important to be aware that the horizontal axis of a dendrogram has no meaning. You could swap some node's left children and right children and it would still be the same dendrogram. It doesn't mean anything that two leaves happen to be next to each other.]



[linksages.pdf (ISL, Figure 10.12)] [Comparison of average, complete (max), and single (min) linkages. Observe that the complete linkage gives the best-balanced dendrogram, whereas the single linkage gives a very unbalanced dendrogram that is sensitive to outliers (especially near the top of the dendrogram).]

[Probably the worst of these is the single linkage, because it's very sensitive to outliers. Notice that if you cut this example into three clusters, two of them have only one node. It also tends to give you a very unbalanced tree.]

[The complete linkage tends to be the best balanced, because when a cluster gets large, the furthest point in the cluster is always far away. So large clusters are more resistant to growth than small ones. If balanced clusters are your goal, this is your best choice.]

[In most cases you probably want the average or complete linkage.]

Warning: centroid linkage can cause inversions where a parent cluster is fused at a lower height than its children.

[So statisticians don't like it, but nevertheless, centroid linkage is popular in genomics.]

[As a final note, all the clustering algorithms we've studied so far are unstable, in the sense that deleting a few input points can sometimes give you very different results. But these unstable heuristics are still the most commonly used clustering algorithms. And it's not clear to me whether a truly stable clustering algorithm is even possible.]

22 Spectral Graph Clustering

SPECTRAL GRAPH CLUSTERING

Input: Weighted, undirected graph $G = (V, E)$. No self-edges.

w_{ij} = weight of edge $(i, j) = (j, i)$; zero if $(i, j) \notin E$.

[Think of the edge weights as a similarity measure. A big weight means that the two vertices want to be in the same cluster. So the circumstances are the opposite of the last lecture on clustering. Then, we had a distance or dissimilarity function, so small numbers meant that points wanted to stay together. Today, big numbers mean that vertices want to stay together.]

Goal: Cut G into 2 (or more) pieces G_i of similar sizes,
but don't cut too much edge weight.

[That's a vague goal. There are many ways to make this precise.

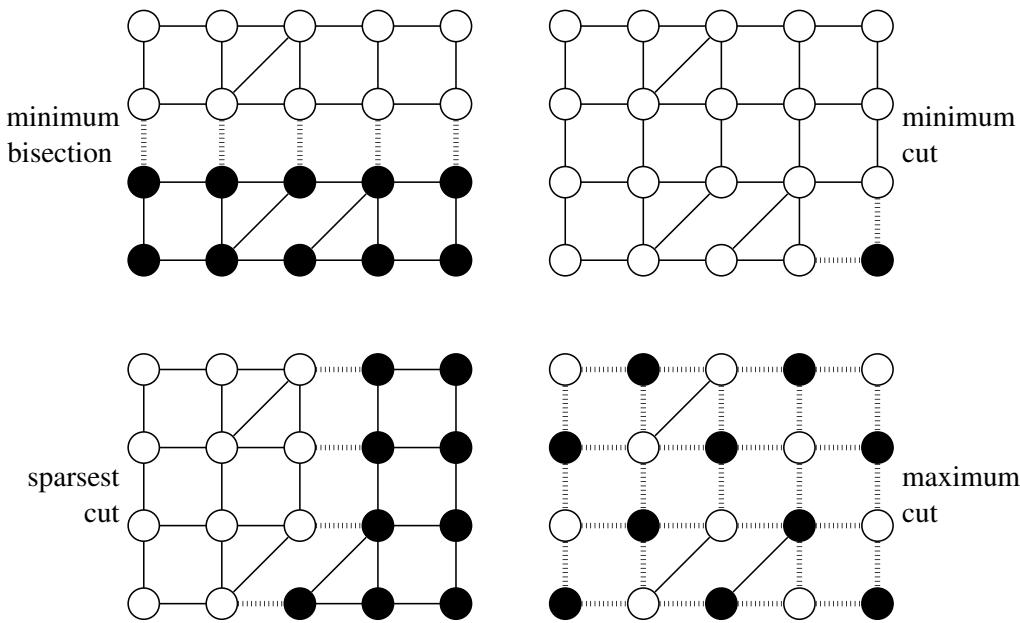
Here's a typical goal, which we'll solve approximately.]

e.g., Minimize the sparsity $\frac{\text{Cut}(G_1, G_2)}{\text{Mass}(G_1)\text{Mass}(G_2)}$, aka cut ratio

where $\text{Cut}(G_1, G_2)$ = total weight of cut edges

$\text{Mass}(G_1)$ = # of vertices in G_1 OR assign masses to vertices

[The denominator “ $\text{Mass}(G_1)\text{Mass}(G_2)$ ” penalizes imbalanced cuts.]



[graph.pdf] [Four cuts. All edges have weight 1.

Upper left: the minimum bisection; a bisection is perfectly balanced.

Upper right: the minimum cut. Usually very unbalanced; not what we want.

Lower left: the sparsest cut, which is good for many applications.

Lower right: the maximum cut; in this case also the maximum bisection.]

Sparsest cut, min bisection, max cut all NP-hard.

[Today we will look for an approximate solution to the sparsest cut problem.]

[We will turn this combinatorial graph cutting problem into algebra.]

Let $n = |V|$. Let $y \in \mathbb{R}^n$ be an indicator vector:

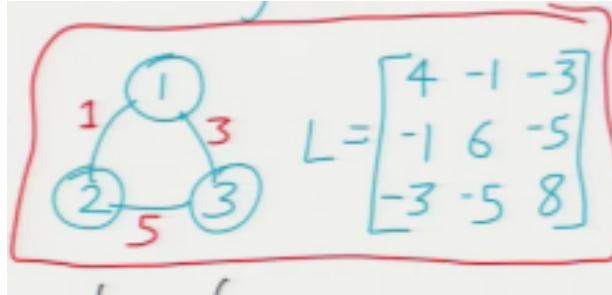
$$y_i = \begin{cases} 1 & \text{vertex } i \in G_1, \\ -1 & \text{vertex } i \in G_2. \end{cases}$$

Then $w_{ij} \frac{(y_i - y_j)^2}{4} = \begin{cases} w_{ij} & (i, j) \text{ is cut,} \\ 0 & (i, j) \text{ is not cut.} \end{cases}$

$$\begin{aligned} \text{Cut}(G_1, G_2) &= \sum_{(i,j) \in E} w_{ij} \frac{(y_i - y_j)^2}{4} && [\text{This is quadratic, so let's try to write it with a matrix.}] \\ &= \frac{1}{4} \sum_{(i,j) \in E} (w_{ij} y_i^2 - 2w_{ij} y_i y_j + w_{ij} y_j^2) \\ &= \frac{1}{4} \left(\underbrace{\sum_{(i,j) \in E} -2w_{ij} y_i y_j}_{\text{off-diagonal terms}} + \underbrace{\sum_{i=1}^n y_i^2 \sum_{k \neq i} w_{ik}}_{\text{diagonal terms}} \right) \\ &= \frac{y^T L y}{4}, \end{aligned}$$

where $L_{ij} = \begin{cases} -w_{ij}, & i \neq j, \\ \sum_{k \neq i} w_{ik}, & i = j. \end{cases}$

L is symmetric, $n \times n$ Laplacian matrix for G .



[Draw this by hand [graphexample.png](#)]

L is effectively a matrix representation of G . For the purpose of partitioning a graph, there is no need to distinguish edges of weight zero from edges that are not in the graph.]

[We see that minimizing the weight of the cut is equivalent to minimizing the Laplacian quadratic form $y^T L y$. This lets us turn graph partitioning into a problem in matrix algebra.]

[Usually we assume there are no negative weights, in which case $\text{Cut}(G_1, G_2)$ can never be negative, so it follows that L is positive semidefinite.]

If $y = \mathbf{1} = [1 \ 1 \ \dots \ 1]^T$, then $\text{Cut}(G_1, G_2) = 0$, so

$\mathbf{1}$ is an eigenvector of L with eigenvalue 0.

[If G is a connected graph and all the edge weights are positive, then this is the only zero eigenvalue. But if G is not connected, L has one zero eigenvalue for each connected component of G . It's easy to prove, but time prevents me.]

Bisection: exactly $n/2$ vertices in G_1 , $n/2$ in G_2 . Write $\mathbf{1}^\top \mathbf{y} = 0$.

[So we have reduced graph bisection to this constrained optimization problem.]

Minimum bisection:

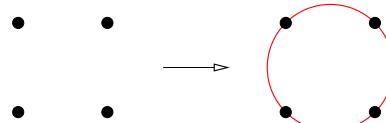
Find \mathbf{y} that minimizes $\mathbf{y}^\top L \mathbf{y}$ subject to $\forall i, y_i = 1$ or $y_i = -1$ and $\mathbf{1}^\top \mathbf{y} = 0$	\leftarrow binary constraint \leftarrow balance constraint
--	---

Also NP-hard. We relax the binary constraint. \rightarrow fractional vertices!

[A very common approach in combinatorial optimization algorithms is to relax some of the constraints so a discrete problem becomes a continuous problem. Intuitively, this means that you can put $1/3$ of vertex 7 in graph G_1 and the other $2/3$ of vertex 7 in graph G_2 . You can even put $-1/2$ of vertex 7 in graph G_1 and $3/2$ of vertex 7 in graph G_2 . This sounds crazy, but the continuous problem is much easier to solve than the combinatorial problem. After we solve it, we will round the vertex values to $+1/-1$, and we'll hope that our solution is still close to optimal.]

[But we can't just drop the binary constraint. We still need *some* constraint to rule out the solution $\mathbf{y} = 0$.]

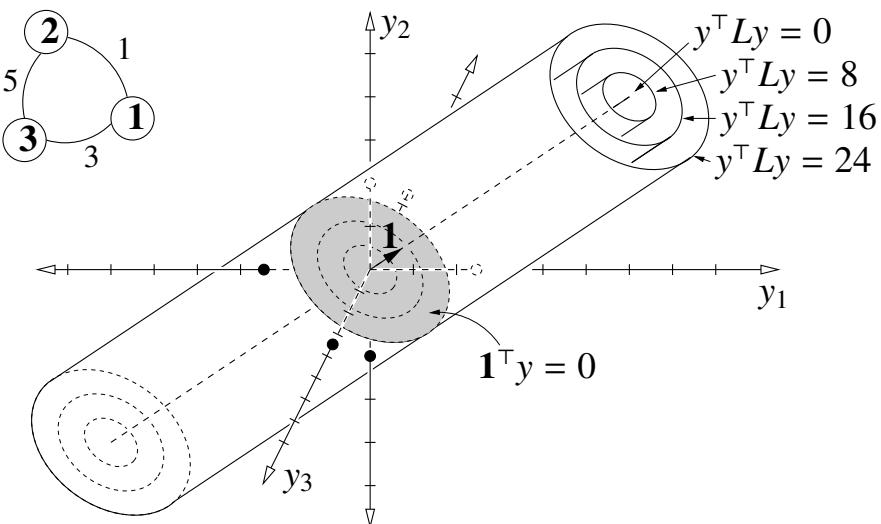
New constraint: \mathbf{y} must lie on hypersphere of radius \sqrt{n} .



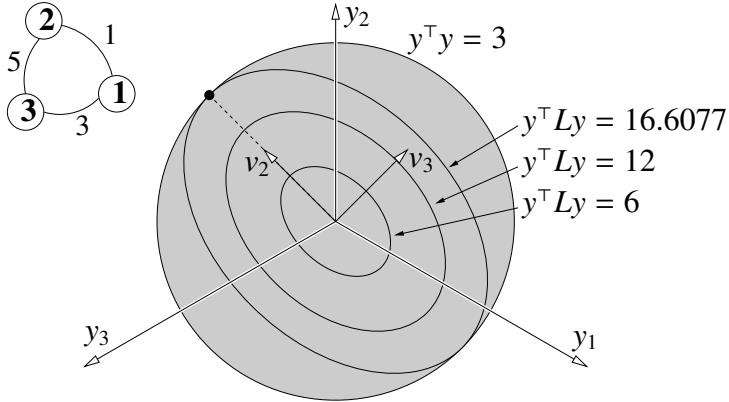
[Draw this by hand. [circle.pdf](#)] [Instead of constraining \mathbf{y} to lie at a vertex of the hypercube, we constrain \mathbf{y} to lie on the hypersphere through those vertices.]

Relaxed problem:

Minimize $\mathbf{y}^\top L \mathbf{y}$ subject to $\mathbf{y}^\top \mathbf{y} = n$ and $\mathbf{1}^\top \mathbf{y} = 0$	$\left\} = \text{Minimize } \frac{\mathbf{y}^\top L \mathbf{y}}{\mathbf{y}^\top \mathbf{y}} = \text{Rayleigh quotient of } L \text{ & } \mathbf{y}$
--	---



[cylinder.pdf](#) [The isosurfaces of $\mathbf{y}^\top L \mathbf{y}$ are elliptical cylinders. The gray cross-section is the hyperplane $\mathbf{1}^\top \mathbf{y} = 0$. We seek the point that minimizes $\mathbf{y}^\top L \mathbf{y}$, subject to the constraints that it lies on the gray cross-section and that it lies on a sphere centered at the origin.]



[endview.pdf] [The same isosurfaces restricted to the hyperplane $\mathbf{1}^\top \mathbf{y} = 0$. The solution is constrained to lie on the outer circle.]

[You should remember this Rayleigh quotient from the lecture on PCA. As I said then, when you see a Rayleigh quotient, you should smell eigenvectors nearby. The \mathbf{y} that minimizes this Rayleigh quotient is the eigenvector with the smallest eigenvalue. We already know what that eigenvector is: it's **1**. But that violates our balance constraint. As you should recall from PCA, when you've used the most extreme eigenvector and you need an orthogonal one, the next-best optimizer of the Rayleigh quotient is the next eigenvector.]

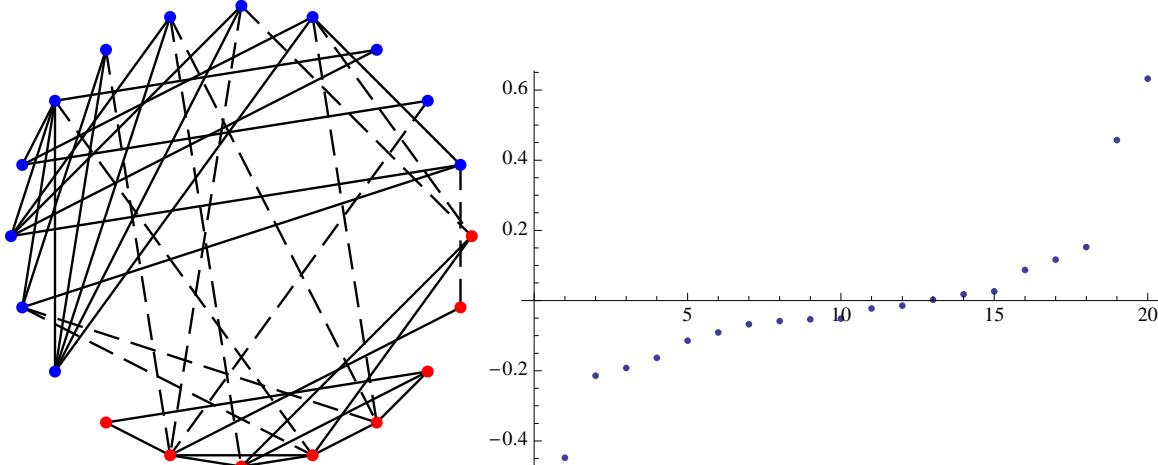
Let λ_2 = second-smallest eigenvalue of L .

Eigenvector v_2 is the Fiedler vector.

[It would be wonderful if every component of the Fiedler vector was 1 or -1 , but that happens more or less never. So we round v_2 . The simplest way is to round all positive entries to 1 and all negative entries to -1 . But in both theory and practice, it's better to choose the threshold as follows.]

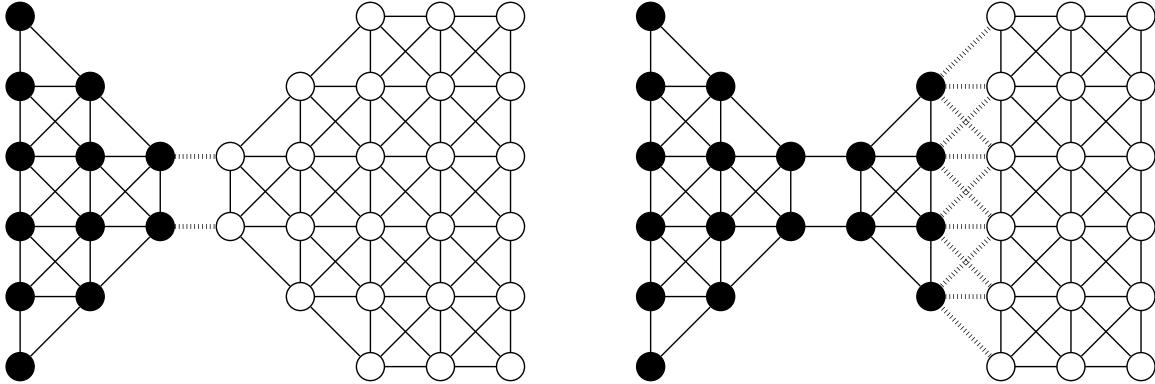
Spectral partitioning alg:

- Compute Fiedler vector v_2 of L
- Round v_2 with a sweep cut:
 - = Sort components of v_2 .
 - = Try the $n - 1$ cuts between successive components. Choose min-sparsity cut.
[If we're clever about updating the sparsity, we can try all these cuts in time linear in the number of edges in G .]



[specgraph.pdf, specvector.pdf] [Left: example of a graph partitioned by the sweep cut.
Right: what the un-rounded Fiedler vector looks like.]

[One consequence of relaxing the binary constraint is that the balance constraint no longer forces an exact bisection. But that's okay; we're cool with a slightly off-balance constraint if it means we cut fewer edges.]



[lopsided.pdf](#) [A graph for which an off-balance cut (left) is sparser than a balanced one (right).]

Vertex Masses

[Sometimes you want the notion of balance to accord more prominence to some vertices than others. We can assign masses to vertices.]

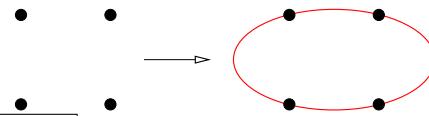
Let M be diagonal matrix with vertex masses on diagonal.

New balance constraint: $\mathbf{1}^\top M \mathbf{y} = 0$.

[This new balance constraint says that G_1 and G_2 should each have the same total mass. It turns out that this new balance constraint is easier to satisfy if we also revise the sphere constraint a little bit.]

New ellipsoid constraint: $\mathbf{y}^\top M \mathbf{y} = \text{Mass}(G) = \sum M_{ii}$.

[Instead of a sphere, now we constrain y to lie on an axis-aligned ellipsoid.]



[Draw this by hand. [ellipse.pdf](#)] [The constraint ellipsoid passes through the points of the hypercube.]

Now we want Fiedler vector of generalized eigensystem $Lv = \lambda Mv$.

[Most algorithms for computing eigenvectors and eigenvalues of symmetric matrices can easily be adapted to compute eigenvectors and eigenvalues of symmetric generalized eigensystems too.]

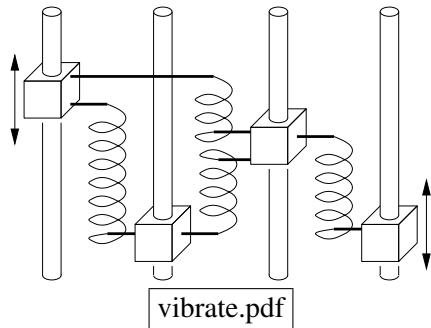
[For the grad students, here's the most important theorem in spectral graph partitioning.]

Fact: Sweep cut finds a cut w/sparsity $\leq \sqrt{2\lambda_2 \max_i \frac{L_{ii}}{M_{ii}}}$: Cheeger's inequality.

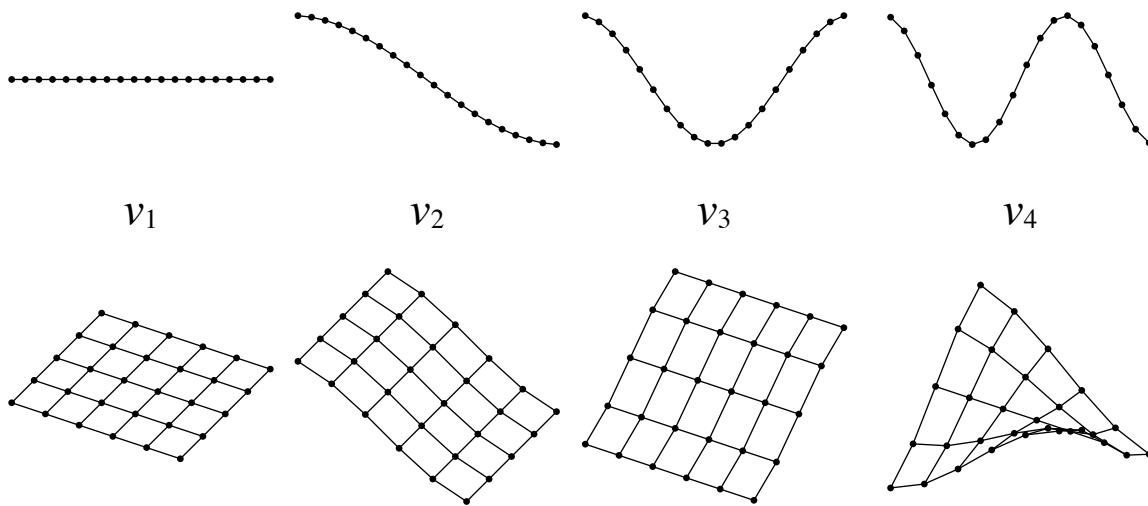
The optimal cut has sparsity $\geq \lambda_2/2$.

[So the spectral partitioning algorithm is an approximation algorithm, albeit not one with a constant factor of approximation. Cheeger's inequality is a very famous result in spectral graph theory, because it's one of the most important cases where you can relax a combinatorial optimization problem to a continuous optimization problem, round the solution, and still have a provably decent solution to the original combinatorial problem.]

Vibration Analogy



[For intuition about spectral partitioning, think of the eigenvectors as vibrational modes in a physical system of springs and masses. Each vertex models a point mass that is constrained to move freely along a vertical rod. Each edge models a vertical spring with rest length zero and stiffness proportional to its weight, pulling two point masses together. The masses are free to oscillate sinusoidally on their rods. The eigenvectors of the generalized eigensystem $Lv = \lambda Mv$ are the vibrational modes of this physical system, and their eigenvalues are proportional to their frequencies.]



grids.pdf [Vibrational modes in a path graph and a grid graph.]

[These illustrations show the first four eigenvectors for two simple graphs. On the left, we see that the first eigenvector is the eigenvector of all 1's, which represents a vertical translation of all the masses in unison. That's not really a vibration, which is why the eigenvalue is zero. The second eigenvector is the Fiedler vector, which represents the vibrational mode with the lowest frequency. Each component indicates the amplitude with which the corresponding point mass oscillates. At any point in time as the masses vibrate, roughly half the mass is moving up while half is moving down. So it makes sense to cut between the positive components and the negative components. The third eigenvector also gives us a nice bisection of the grid graph, entirely different from the Fiedler vector. Some more sophisticated graph clustering algorithms use multiple eigenvectors.]

[I want to emphasize that spectral partitioning takes a global view of a graph. It looks at the whole gestalt of the graph and finds a good cut. By comparison, the clustering algorithms we saw last lecture were much more local in nature, so they're easier to fool.]

Greedy Divisive Clustering

Partition G into 2 subgraphs; recursively cluster them.

[The sparsity is a good criterion for graph clustering. Use G 's sparsest cut to divide it into two subgraphs, then recursively cut them. You can stop when you have the right number of clusters, or you could keep going until each subgraph is a single vertex and create a dendrogram.]

Can form a dendrogram, but it may have inversions.

[There's no reason to expect that the sparsity of a subgraph is smaller than the sparsity of the parent graph, so the dendrogram can have inversions. But it's still useful for getting an arbitrary number of clusters on demand.]

The Normalized Cut

Set vertex i 's mass $M_{ii} = L_{ii}$. [Sum of edge weights adjoining vertex i .]

[That is how we define a “normalized cut,” which turns out to be a good choice for many different applications.]

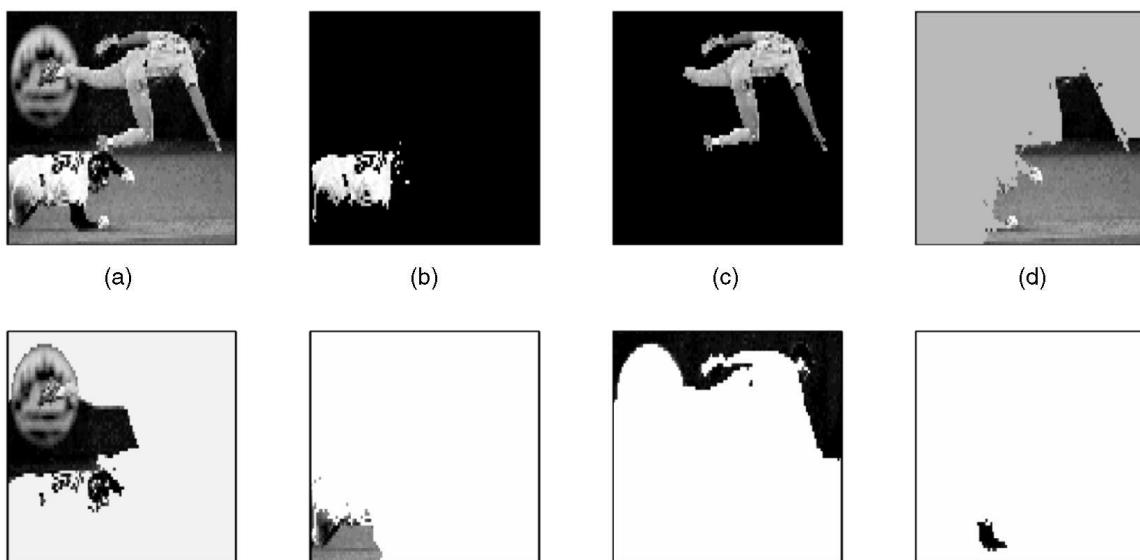
Popular for image segmentation.

[Image segmentation is the problem of looking at a photograph and separating it into different objects. To do that, we define a graph on the pixels.]

For pixels with location w_i , brightness b_i , use graph weights

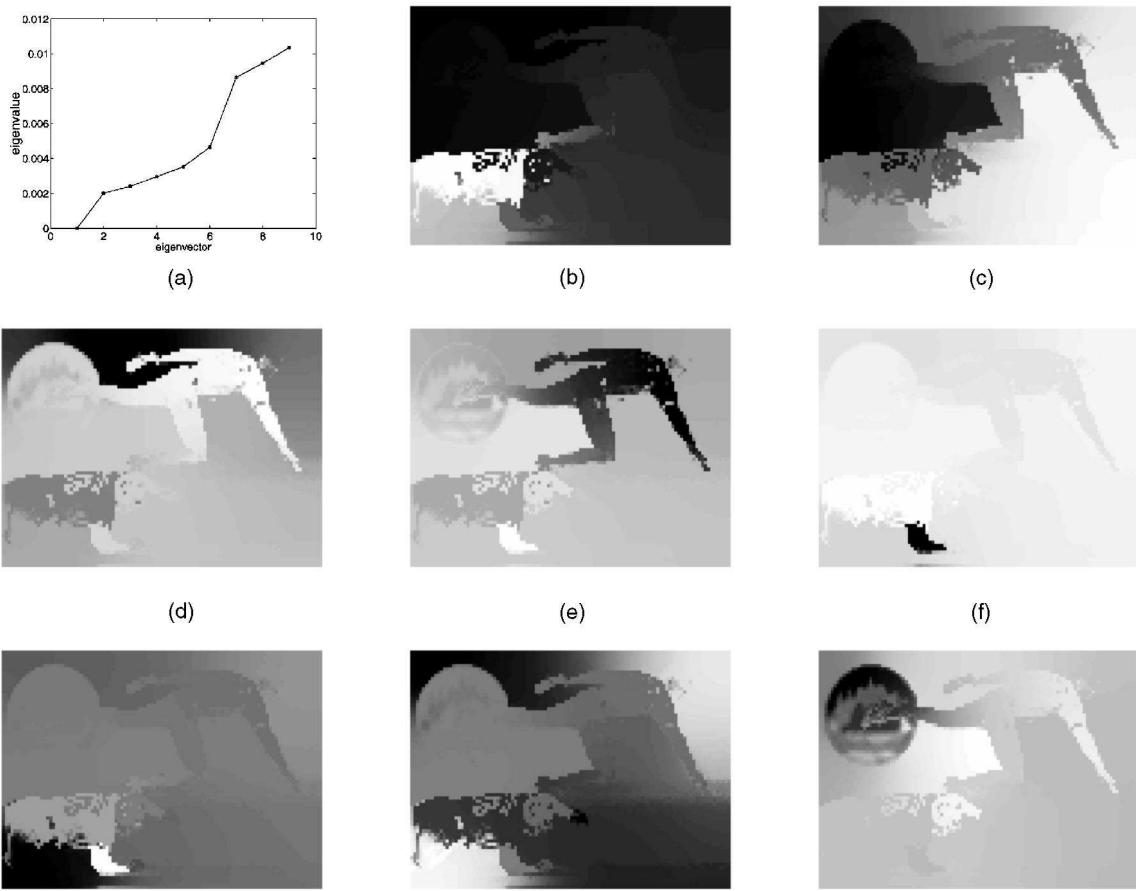
$$w_{ij} = \exp\left(-\frac{|w_i - w_j|^2}{\alpha} - \frac{|b_i - b_j|^2}{\beta}\right) \quad \text{or zero if } |w_i - w_j| \text{ large.}$$

[We choose a distance threshold, typically less than 4 to 10 pixels apart. Pixels that are far from each other aren't connected. α and β are empirically chosen constants. It often makes sense to choose β proportional to the variance of the brightness values.]



baseballsegment.pdf (Shi and Malik, “Normalized Cut and Image Segmentation”)

[A segmentation of a photo of a scene from a baseball game (upper left). The other figures show segments of the image extracted by recursive spectral partitioning.]



baseballvectors.pdf (Shi and Malik) [Eigenvectors 2–9 from the baseball image.]

Invented by [our own] Prof. Jitendra Malik and his student Jianbo Shi.

23 Multiple Eigenvectors; Random Projection; Latent Factor Analysis

Clustering w/Multiple Eigenvectors

[When we use the Fiedler vector for spectral graph clustering, it tells us how to divide a graph into two graphs. If we want more than two clusters, we can use divisive clustering: we repeatedly cut the subgraphs into smaller subgraphs by computing their Fiedler vectors. However, there are several other methods to subdivide a graph into k clusters in one shot that use multiple eigenvectors rather than just the Fiedler vector v_2 . These methods are usually faster and sometimes give better results. They use k eigenvectors in a natural way to cluster a graph into k subgraphs.]

For k clusters, compute first k eigenvectors $v_1 = \mathbf{1}, v_2, \dots, v_k$ of generalized eigensystem $Lv = \lambda Mv$.

$$V = \begin{array}{c|c} \begin{matrix} & \uparrow \\ 1 & \\ 1 & \\ 1 & \\ v_2 & \\ 1 & \\ 1 & \\ 1 & \\ 1 & \\ 1 & \\ 1 & \\ 1 & \\ 1 & \\ \downarrow & \end{matrix} & = \end{array} \begin{array}{c} V_1 \\ \vdots \\ V_n \end{array}$$

[V 's columns are the eigenvectors with the k smallest eigenvalues.]
 [Yes, we do include the all-1's vector v_1 as one of the columns of V .]

[Draw this by hand. [eigenvectors.pdf](#)]

Row V_i is spectral vector [my name] for vertex i . [The rows are vectors in a k -dimensional space I'll call the "spectral space." When we were using just one eigenvector, it made sense to cluster vertices together if their components were close together. When we use more than one eigenvector, it turns out that it makes sense to cluster vertices together if their spectral vectors point in similar directions.]

Normalize each row V_i to unit length.

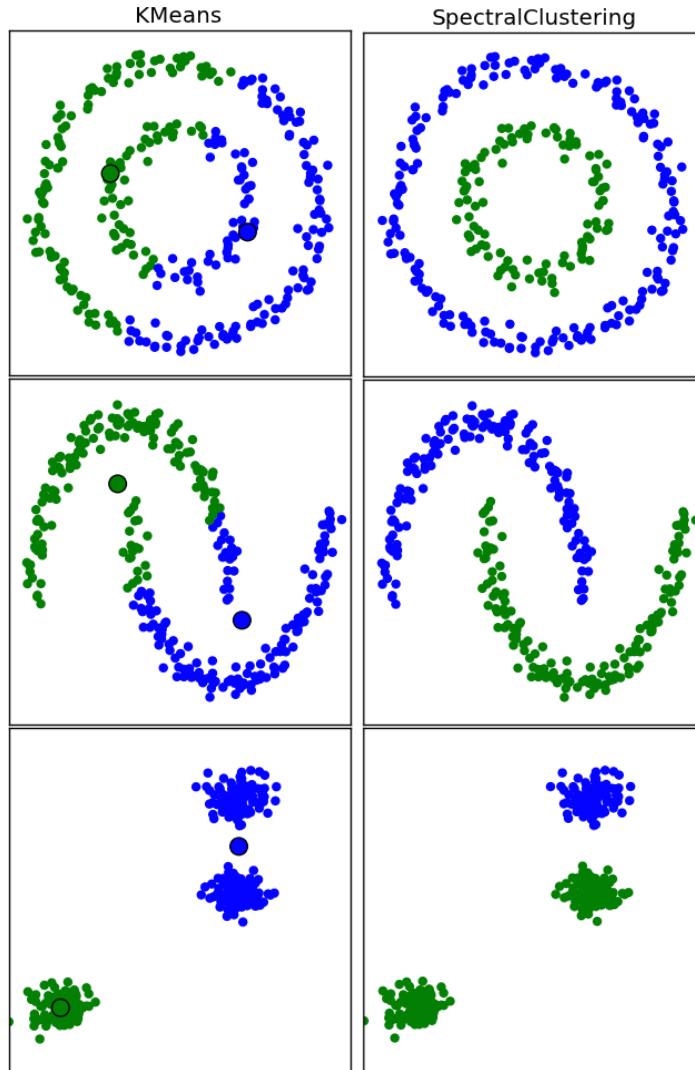
[Now you can think of the spectral vectors as points on a unit sphere centered at the origin.]



[Draw this by hand [vectorclusters.png](#)] [A 2D example showing two clusters on a circle. If the graph has k components, the points in each cluster will have identical spectral vectors that are exactly orthogonal to all the other components' spectral vectors (left). If we modify the graph by connecting these components with small-weight edges, we get vectors more like those at right—not exactly orthogonal, but still tending toward distinct clusters.]

k -means cluster these vectors.

[Because all the spectral vectors lie on the sphere, k -means clustering will cluster together vectors that are separated by small angles.]



`compkmeans.png, compspectral.png` [Comparison of point sets clustered by k -means—just k -means by itself, that is—vs. a spectral method. To create a graph for the spectral method, we use an exponentially decaying function to assign weights to pairs of points, like we used for image segmentation but without the brightnesses.]

Invented by [our own] Prof. Michael Jordan, Andrew Ng [when he was still a student at Berkeley], Yair Weiss.

[This wasn't the first algorithm to use multiple eigenvectors for spectral clustering, but it has become one of the most popular.]

RANDOM PROJECTION

A cheap alternative to PCA as preprocess for clustering, classification, regression.

Approximately preserves distances between points!

[We project onto a random subspace instead of the “best” subspace, but take a fraction of the time of PCA. It works best when you project a very high-dimensional space to a medium-dimensional space. Because we roughly preserve the distances, algorithms like k -means clustering and nearest neighbor classifiers will give similar results to what they would give in high dimensions, but they run much faster.]

Pick a small ϵ , a small δ , and a random subspace $S \subset \mathbb{R}^d$ of dimension k , where $k = \frac{2 \ln(1/\delta)}{\epsilon^2/2 - \epsilon^3/3}$.

For any pt q , let \hat{q} be orthogonal projection of q onto S , multiplied by $\sqrt{\frac{d}{k}}$.

[The multiplication by $\sqrt{d/k}$ helps preserve the distances between points after you project.]

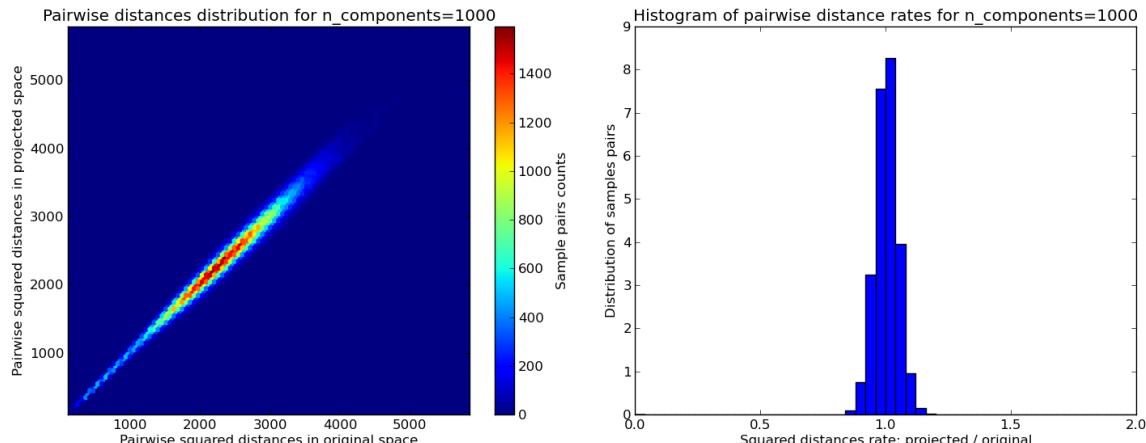
Johnson–Lindenstrauss Lemma (modified):

For any two pts $q, w \in \mathbb{R}^d$, $(1 - \epsilon)|qw|^2 \leq |\hat{q}\hat{w}|^2 \leq (1 + \epsilon)|qw|^2$ with probability $\geq 1 - 2\delta$.

Typical values: $\epsilon \in [0.02, 0.5]$, $\delta \in [1/n^3, 0.05]$.

[So the distance between two points after projecting might change by 2 to 50%. In practice, you can experiment with k to find the best speed-accuracy tradeoff. If you want most inter-point distances to be accurate, you should set δ smaller than $1/n^2$, so you need a subspace of dimension $\Theta(\log n)$. Reducing δ doesn’t cost much, but reducing ϵ costs more. You can bring 1,000,000 sample points down to a 10,000-dimensional space with a 10% error in the distances.]

[What is remarkable about this result is that the dimension d of the input points doesn’t matter.]



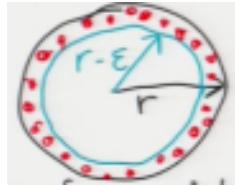
[100000to1000.pdf](#) [Comparison of inter-point distances before and after projecting points in 100,000-dimensional space down to 1,000 dimensions.]

[Why does this work? A random projection of a vector is like taking a random vector and selecting k components. The mean of the squares of those k components approximates the mean for the whole population.]

[How do you get a uniformly distributed random projection direction? You can choose each component from a univariate Gaussian distribution, then normalize the vector to unit length. How do you get a random subspace? You can choose k random vectors, then use Gram–Schmidt orthogonalization to make them mutually orthonormal. Interestingly, Indyk and Motwani show that if you skip the expensive normalization and Gram–Schmidt steps, random projection still works almost as well, because random vectors in a high-dimensional space are nearly equal in length and nearly orthogonal to each other with high probability.]

THE GEOMETRY OF HIGH-DIMENSIONAL SPACES

Consider shell between spheres of radii r & $r - \epsilon$.



[Draw this by hand [concentric.png](#)] [Concentric balls. In high dimensions, almost every point chosen uniformly at random in the outer ball lies outside the inner ball.]

Volume of outer ball $\propto r^d$

Volume of inner ball $\propto (r - \epsilon)^d$

Ratio of inner ball volume to outer =

$$\frac{(r - \epsilon)^d}{r^d} = \left(1 - \frac{\epsilon}{r}\right)^d \approx \exp\left(-\frac{\epsilon d}{r}\right) \quad \text{which is small for large } d.$$

E.g. if $\frac{\epsilon}{r} = 0.1$ & $d = 100$, inner ball has $90\%^{100} = 0.0027\%$ of volume.

Random points from uniform distribution in ball: nearly all are in outer shell.

" " " Gaussian " : nearly all are in some shell.

If the dimension is very high, the majority of the random points generated from an isotropic Gaussian distribution are approximately at the same distance from the center. So they lie in a thin shell. Why? Consider a d -dimensional normal distribution with mean zero. By Pythagoras' Theorem, the squared distance from a random point P to the mean is]

$$\|P\|^2 = P_1^2 + P_2^2 + \dots + P_d^2.$$

[Each component P_i is sampled independently from a univariate normal distribution with mean zero. When you add d independent random numbers, you scale the mean by d but you scale the standard deviation by only \sqrt{d} .]

$$E[\|P\|^2] = d E[P_1^2].$$

$$\text{std}(\|P\|^2) = \sqrt{d} \text{ std}(P_1^2).$$

[So when d is large, the distance from P to the mean is concentrated in a narrow shell whose radius is proportional to \sqrt{d} with a standard deviation proportional to $\sqrt[4]{d}$.]

[The principle here is that when you take the mean of a very large sample, you get a very accurate estimate of the population mean. When you sample one point from a high-dimensional normal distribution, it's like sampling d different scalars from one-dimensional normal distributions.]

Lessons:

- In high dimensions, sometimes nearest neighbor and farthest neighbor don't differ much.
- k -means clustering and nearest neighbor classifiers [and many other classifiers] are less effective for large d .

[Your head TA, Marc Khoury, has a nice short essay entitled “Counterintuitive Properties of High Dimensional Space”, which you can read at]

<https://marckhoury.github.io/counterintuitive-properties-of-high-dimensional-space/>

LATENT FACTOR ANALYSIS [aka Latent Semantic Indexing]

[You can think of this as dimensionality reduction for matrices.]

Suppose X is a term-document matrix: [aka bag-of-words model]

row i represents document i ; column j represents term j . [Term = word.]

[Term-document matrices are usually sparse, meaning most entries are zero.]

X_{ij} = occurrences of term j in doc i

better: $\log(1 + \text{occurrences})$ [So frequent words don't dominate.]

[Better still is to weight the entries so rare words give big entries and common words like "the" give small entries. To do that, you need to know how frequently each word occurs in general. I'll omit the details, but this is the common practice.]

Recall SVD $X = UDV^\top = \sum_{i=1}^d \delta_i u_i v_i^\top$. Suppose $\delta_i \leq \delta_j$ for $i \geq j$.

Unlike PCA, we usually don't center X .

For large δ_i , u_i and v_i represent a cluster of documents & terms.

- Large components in u_i mark books using similar/related terms, i.e., a genre.
- " " " v_i mark frequent terms in that genre.
- E.g., u_1 might have large components for the romance novels,
- v_1 " " " " for terms "passion," "ravish," "bodice" ...

[... and δ_1 would give us an idea how much bigger the romance novel market is than the markets for every other genre of books.]

[v_1 and u_1 tell us that there is a large subset of books that tend to use the same large subset of words. We can read off the words by looking at the larger components of v_1 , and we can read off the books by looking at the larger components of u_1 .]

[The property of being a romance novel is an example of a latent factor. So is the property of being the sort of word used in romance novels. There's nothing in X that tells you explicitly that romance novels exist, but the similar vocabulary is a hidden connection between them that gives them a large singular value. The vector u_1 reveals which books have that genre, and v_1 reveals which words are emphasized in that genre.]

Like clustering, but clusters overlap: if u_1 picks out romances &
 u_2 picks out histories, they both pick out historical romances.

[So you can think of latent factor analysis as a sort of clustering that permits clusters to overlap. Another way in which it differs from traditional clustering is that the u -vectors contain real numbers, and so some points have stronger cluster membership than others. One book might be just a bit romance, another a lot.]

Application in market research:

identifying consumer types (hipster, suburban mom) & items bought together.

[For applications like this, the first few singular vectors are the most useful. Most of the singular vectors are mostly noise, and they have small singular values to tell you so. This motivates approximating a matrix by using only some of its singular vectors.]

Truncated sum $X' = \sum_{i=1}^r \delta_i u_i v_i^\top$ is a low-rank approximation of X , of rank r . [Assuming $\delta_r > 0$.]

[We choose the singular vectors with the largest singular values, because they carry the most information.]

$$\begin{array}{c|c}
 \boxed{X'} & = \\
 \hline
 n \times d & \begin{array}{c|c}
 \boxed{u_1} & \boxed{v_1} \\
 \vdots & \vdots \\
 \boxed{u_r} & \boxed{v_r} \\
 \hline
 n \times r & r \times d
 \end{array}
 \end{array}$$

[Draw this by hand. [truncate.pdf](#)]

X' is the rank- r matrix that minimizes the [squared] Frobenius norm

$$\|X - X'\|_F^2 = \sum_{i,j} (X_{ij} - X'_{ij})^2$$

Applications:

- Fuzzy search. [Suppose you want to find a document about gasoline prices, but the document you want doesn't have the word "gasoline"; it has the word "petrol." One cool thing about the reduced-rank matrix X' is that it will probably associate that document with "gasoline," because the SVD tends to group synonyms together.]
- Denoising. [The idea is to assume that X is a noisy measurement of some unknown matrix that probably has low rank. If that assumption is partly true, then the reduced-rank matrix X' might be better than the input X .]
- Matrix compression. [As you can see above, if we use a low-rank approximation with a small rank r , we can express the approximate matrix as an SVD that takes up much less space than the original matrix. Often this row-rank approximation supports faster matrix computations.]
- Collaborative filtering: fills in unknown values, e.g. user ratings.
[Suppose the rows of X represents Netflix users and the columns represent movies. The entry X_{ij} is the review score that user i gave to movie j . But most users haven't reviewed most movies. We want to fill in the missing values. Just as the rank reduction will associate "petrol" with "gasoline," it will tend to associate users with similar tastes in movies, so the reduced-rank matrix X' can predict ratings for users who didn't supply any.]

24 Boosting

ADABOOST (Yoav Freund and Robert Schapire, 1997)

[We're done with unsupervised learning. This week I'm going back to classifiers.]

AdaBoost ("adaptive boosting") is an ensemble method for classification (or regression) that

- trains multiple learners on weighted sample points [like bagging];
- uses different weights for each learner;
- increases weights of misclassified sample points;
- gives bigger votes to more accurate learners.

Input: $n \times d$ design matrix X , vector of labels $y \in \mathbb{R}^n$ with $y_i = \pm 1$.

Ideas:

- Train T classifiers G_1, \dots, G_T .
- Weight for sample point X_i in G_t grows according to how many of G_1, \dots, G_{t-1} misclassified it.
[Moreover, if X_i is misclassified by very accurate learners, its weight grows even more.]
[And, the weight shrinks every time X_i is correctly classified.]
- Train G_t to try harder to correctly classify sample pts with larger weights.
- Metalearner is a linear combination of learners. For test point z , $M(z) = \sum_{t=1}^T \beta_t G_t(z)$.
Each G_t is ± 1 , but M is continuous. Return sign of $M(z)$.

[In the previous lecture on ensemble methods, I talked briefly about how to assign different weights to sample points. It varies for different learning algorithms. For example, in regression we usually modify the risk function by multiplying each point's loss function by its weight. In a soft-margin support vector machine, we modify the objective function by multiplying each point's slack by its weight.]

[Boosting works with most learning algorithms, but it was originally developed for decision trees, and boosted decision trees are still very popular and successful. For decision trees, we use a weighted entropy where instead of computing the proportion of points in each class, we compute the proportion of weight in each class.]

In iteration T , what classifier G_T and coefficient β_T should we choose? Pick a loss fn L (prediction, label).

Find G_T & β_T that minimize

$$\text{Risk} = \frac{1}{n} \sum_{i=1}^n L(M(X_i), y_i), \quad M(X_i) = \sum_{t=1}^T \beta_t G_t(X_i).$$

AdaBoost metalearner uses exponential loss function

$$L(\rho, \ell) = e^{-\rho\ell} = \begin{cases} e^{-\rho} & \ell = +1 \\ e^\rho & \ell = -1 \end{cases}$$

[This loss function is for the metalearner only. The individual learners G_t usually use other loss functions, if they use a loss function at all.]

Important: label ℓ is binary, G_t is binary, but $\rho = M(X_i)$ is continuous!

[The exponential loss function has the advantage that it pushes hard against badly misclassified points. That's one reason why it's usually better than the squared error loss function for classification in a metalearner. It's similar to why in neural networks we prefer the cross-entropy loss function to the squared error.]

$$\begin{aligned}
n \cdot \text{Risk} &= \sum_{i=1}^n L(M(X_i), y_i) = \sum_{i=1}^n e^{-y_i M(X_i)} \\
&= \sum_{i=1}^n \exp\left(-y_i \sum_{t=1}^T \beta_t G_t(X_i)\right) = \sum_{i=1}^n \prod_{t=1}^T e^{-\beta_t y_i G_t(X_i)} \quad \begin{matrix} y_i G_t(X_i) = \pm 1 \\ -1 \rightarrow G_t \text{ misclassifies } X_i \end{matrix} \\
&= \sum_{i=1}^n w_i^{(T)} e^{-\beta_T y_i G_T(X_i)}, \quad \text{where } w_i^{(T)} = \prod_{t=1}^{T-1} e^{-\beta_t y_i G_t(X_i)} \\
&= e^{-\beta_T} \sum_{y_i = G_T(X_i)} w_i^{(T)} + e^{\beta_T} \sum_{y_i \neq G_T(X_i)} w_i^{(T)} \quad [\text{correctly classified and misclassified}] \\
&= e^{-\beta_T} \sum_{i=1}^n w_i^{(T)} + (e^{\beta_T} - e^{-\beta_T}) \sum_{y_i \neq G_T(X_i)} w_i^{(T)}.
\end{aligned}$$

What G_T minimizes the risk? The learner that minimizes the sum of $w_i^{(T)}$ over all misclassified pts X_i !

[This is interesting. By manipulating the formula for the risk, we've discovered what weight we should assign to each sample point. If we want to minimize the risk, we should find the classifier that minimizes the total weight of the misclassified points for this weight function $w_i^{(T)}$. It's a complicated function, but we can compute it. A useful observation is that each learner's weights are related to the previous learner's weights:]

Recursive definition of weights:

$$w_i^{(T+1)} = w_i^{(T)} e^{-\beta_T y_i G_T(X_i)} = \begin{cases} w_i^{(T)} e^{-\beta_T} & y_i = G_T(X_i), \\ w_i^{(T)} e^{\beta_T} & y_i \neq G_T(X_i). \end{cases}$$

[This recursive formulation is a nice benefit of the exponential loss. Notice that a weight shrinks if the point was classified correctly, and grows if the point was misclassified.]

[Now, you might wonder if we should just pick a learner that classifies all the training points correctly. But that's not always possible. If we're using a linear classifier on data that's not linearly separable, some points must be classified wrongly. Moreover, it's NP-hard to find the optimal linear classifier, so in practice G_T will be an approximate best learner, not the true minimizer of training error. But that's okay.]

[You might ask, if we use decision trees, can't we get 100% training accuracy? Usually we can. But interestingly, boosting is usually used with short, imperfect decision trees instead of tall, pure decision trees, for reasons I'll explain later.]

[Now, let's derive the optimal value of β_T .]

To choose β_T , set $\frac{d}{d\beta_T} \text{Risk} = 0$:

$$\begin{aligned}
0 &= -e^{-\beta_T} \sum_{i=1}^n w_i^{(T)} + (e^{\beta_T} + e^{-\beta_T}) \sum_{y_i \neq G_T(X_i)} w_i^{(T)}; \quad [\text{now divide both sides by the first term}] \\
0 &= -1 + (e^{2\beta_T} + 1) \text{err}_T, \quad \text{where } \text{err}_T = \frac{\sum_{y_i \neq G_T(X_i)} w_i^{(T)}}{\sum_{i=1}^n w_i^{(T)}}; \quad [G_T \text{'s weighted error rate}] \\
\beta_T &= \frac{1}{2} \ln\left(\frac{1 - \text{err}_T}{\text{err}_T}\right).
\end{aligned}$$

[So now we have derived the optimal metalearner!]

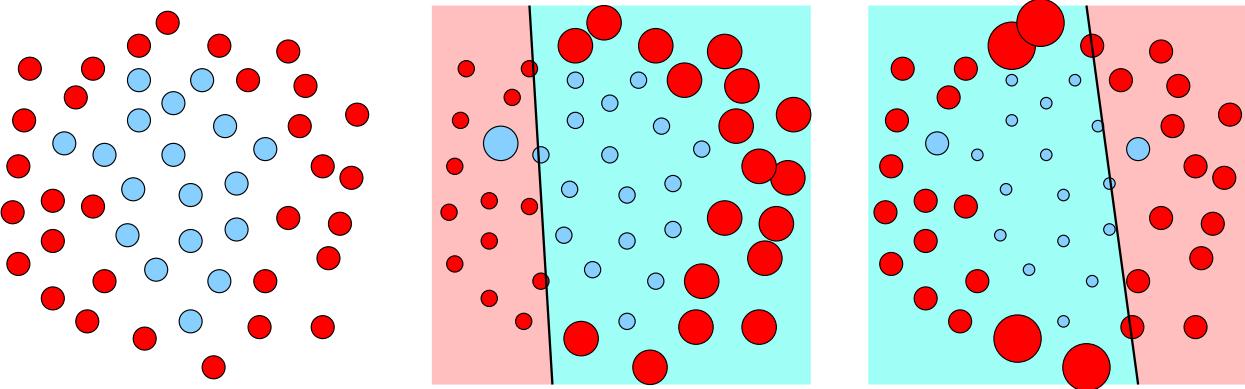
- If $\text{err}_T = 0, \beta_T = \infty$. [So a perfect learner gets an infinite vote.]
- If $\text{err}_T = 1/2, \beta_T = 0$. [So a learner with 50% training accuracy gets no vote at all.]

[More accurate learners get bigger votes in the metalearner. Interestingly, a learner with worse than 50% training accuracy gets a negative vote. A learner with 40% accuracy is just as useful as a learner with 60% accuracy; the metalearner just reverses the sign of its votes.]

[Now we can state the AdaBoost algorithm.]

AdaBoost alg:

1. Initialize weights $w_i \leftarrow \frac{1}{n}, \forall i \in [1, n]$.
2. for $t \leftarrow 1$ to T
 - a. Train G_t with weights w_i
 - b. Compute weighted error rate $\text{err} \leftarrow \frac{\sum_{\text{misclassified}} w_i}{\sum_{\text{all}} w_i}$; coefficient $\beta_t \leftarrow \frac{1}{2} \ln \left(\frac{1-\text{err}}{\text{err}} \right)$.
 - c. Reweight pts: $w_i \leftarrow w_i \cdot \begin{cases} e^{\beta_t}, & G_t \text{ misclassifies } X_i \\ e^{-\beta_t}, & \text{otherwise} \end{cases} = w_i \cdot \begin{cases} \sqrt{\frac{1-\text{err}}{\text{err}}}, & \\ \sqrt{\frac{\text{err}}{1-\text{err}}}. & \end{cases}$
3. return metalearner $h(z) = \text{sign} \left(\sum_{t=1}^T \beta_t G_t(z) \right)$.



`boost0.png, boost2.png, boost4.png` [At left, all the training points have equal weight. After choosing a first linear classifier, we increase the weights of the misclassified points and decrease the weights of the correctly classified points (center). We train a second classifier with these weighted points, then again adjust the weights of the points according to whether they are misclassified by the second classifier.]

Why boost decision trees? [As opposed to other learning algorithms?] Why **short** trees?

- Fast. [We're training many learners, and running many learners at classification time too. Short decision trees that only look at a few features are very fast at both training and testing.]
- No hyperparameter search needed. [Unlike SVMs, neural nets, etc.] [UC Berkeley's Leo Breiman called AdaBoost with decision trees "the best off-the-shelf classifier in the world."]
- Easy to make a tree beat 55% training accuracy [or other threshold] consistently.
- Easy bias-variance control. Boosting can overfit. AdaBoost trees are usually short, to reduce overfitting!

[As you train more learners, the bias decreases. The variance is more complicated: it often decreases at first, because successive trees focus on different features, but often it later increases. Sometimes boosting overfits after many iterations, and sometimes it doesn't; it's hard to predict when it will and when it won't.]

- AdaBoost + short trees is a form of subset selection.

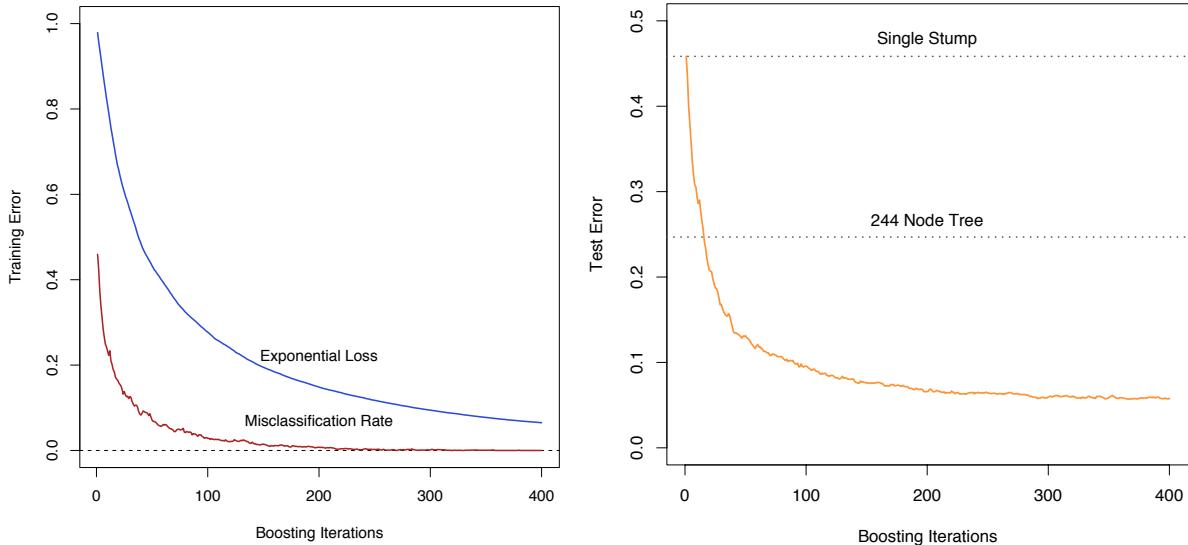
[Features that don't improve the metalearner's predictive power enough aren't used at all. This helps reduce overfitting and running time, especially if there are a lot of irrelevant features.]

- Linear decision boundaries don't boost well.

[Boosting linear classifiers gives you an approximately linear classifier, so SVMs aren't a great choice. Methods with nonlinear decision boundaries benefit more from boosting, because they allow boosting to reduce the bias faster. Sometimes you'll see examples where people do AdaBoost with depth-one decision trees with just one decision each. But that's not ideal, because depth-one decision trees are linear. Even depth-two decision trees boost substantially better.]

More about AdaBoost:

- Posterior prob. can be approximated: $P(Y = 1|x) \approx 1/(1 + e^{-2M(x)})$.
- Exponential loss is vulnerable to outliers; for corrupted data, use other loss.
[Better loss functions have been derived for dealing with outliers. Unfortunately, they have more complicated weight computations.]
- If every learner beats training accuracy μ for $\mu > 50\%$, metalearner accuracy will eventually be 100%.
- [The AdaBoost paper and its authors, Freund and Schapire, won the 2003 Gödel Prize, a prize for outstanding papers in theoretical computer science.]



`trainboost.pdf, testboost.pdf` [Training and testing errors for AdaBoost with stumps, depth-one decision trees that make only one decision each. At left, observe that the training error eventually drops to zero, and even after that the average loss (which is continuous, not binary) continues to decay exponentially. At right, the test error drops to 5.8% after 400 iterations, even though each learner has an error rate of about 46%. AdaBoost with more than 25 stumps outperforms a single 244-node decision tree. In this example no overfitting is observed, but there are other datasets for which overfitting is a problem.]

NEAREST NEIGHBOR CLASSIFICATION

[I saved the simplest classifier for the end of the semester.]

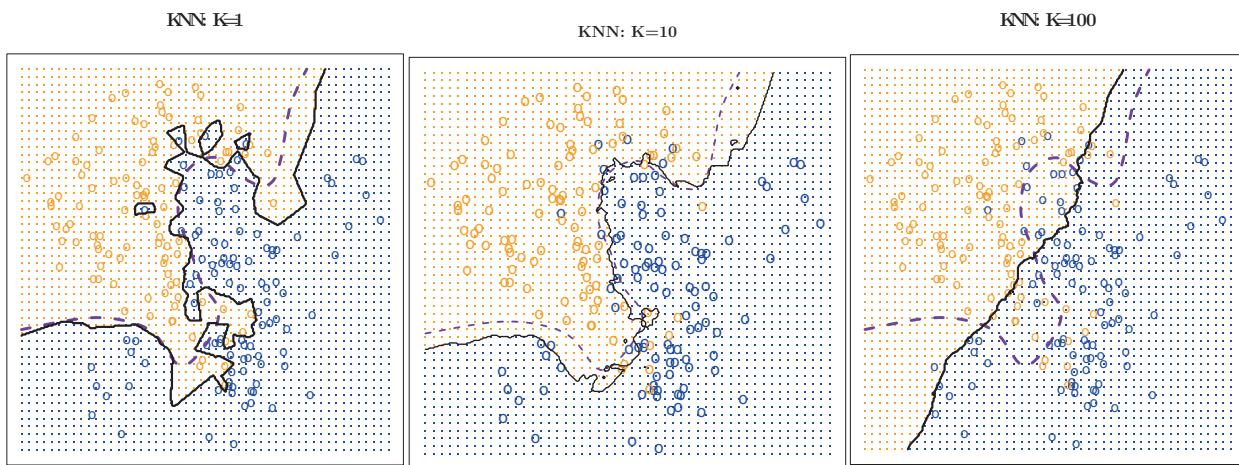
Idea: Given query point q , find the k sample pts nearest q .

Distance metric of your choice.

Regression: Return average label of the k pts.

Classification: Return class with the most votes from the k pts OR
return histogram of class probabilities.

[The histogram of class probabilities tries to estimate the posterior probabilities of the classes. Obviously, the histogram has limited precision. If $k = 3$, then the only probabilities you'll ever return are 0, 1/3, 2/3, or 1. You can improve the precision by making k larger, but you might underfit. The histogram works best when you have a huge amount of data.]



allnn.pdf (ISL, Figures 2.15, 2.16) [Examples of 1-NN, 10-NN, and 100-NN. A larger k smooths out the boundary. In this example, the 1-NN classifier is badly overfitting the data, and the 100-NN classifier is badly underfitting. The 10-NN classifier does well: it's reasonably close to the Bayes decision boundary (purple). Generally, the ideal k depends on how dense your data is. As your data gets denser, the best k increases.]

[There are theorems showing that if you have a lot of data, nearest neighbors can work quite well.]

Theorem (Cover & Hart, 1967):

As $n \rightarrow \infty$, the 1-NN error rate is $< B(2 - B)$ where B = Bayes risk.
if only 2 classes, $\leq 2B(1 - B)$

[There are a few technical requirements of this theorem. The most important is that the training points and the test points all have to be drawn independently from the same probability distribution. The theorem applies to any separable metric space, so it's not just for the Euclidean metric.]

Theorem (Fix & Hodges, 1951):

As $n \rightarrow \infty$, $k \rightarrow \infty$, $k/n \rightarrow 0$, k -NN error rate converges to B . [Which means Bayes optimal.]

25 Nearest Neighbor Algorithms: Voronoi Diagrams and k-d Trees

NEAREST NEIGHBOR ALGORITHMS

Exhaustive k -NN Alg.

Given query point q :

- Scan through all n sample pts, computing (squared) distances to q .
- Maintain a max-heap with the k shortest distances seen so far.

[Whenever you encounter a sample point closer to q than the point at the top of the heap, you remove the heap-top point and insert the better point. Obviously you don't need a heap if $k = 1$ or even 5, but if $k = 101$ a heap will substantially speed up keeping track of the k th-shortest distance.]

Time to construct classifier: $O(0)$ [This is the only $O(0)$ -time algorithm we'll learn this semester.]

Query time: $O(nd + n \log k)$

expected $O(nd + k \log n \log k)$ if random pt order

[It's a cute theoretical observation that you can slightly improve the expected running time by randomizing the point order so that only expected $O(k \log n)$ heap operations occur. But in practice I don't recommend it; you'll probably lose more from cache misses than you'll gain from fewer heap operations.]

Can we preprocess training pts to obtain sublinear query time?

2–5 dimensions: Voronoi diagrams

Medium dim (up to ~ 30): k -d trees

Large dim: exhaustive k -NN, but can use PCA or random projection [or another dimensionality reduction method]

locality sensitive hashing [still researchy, not widely adopted]

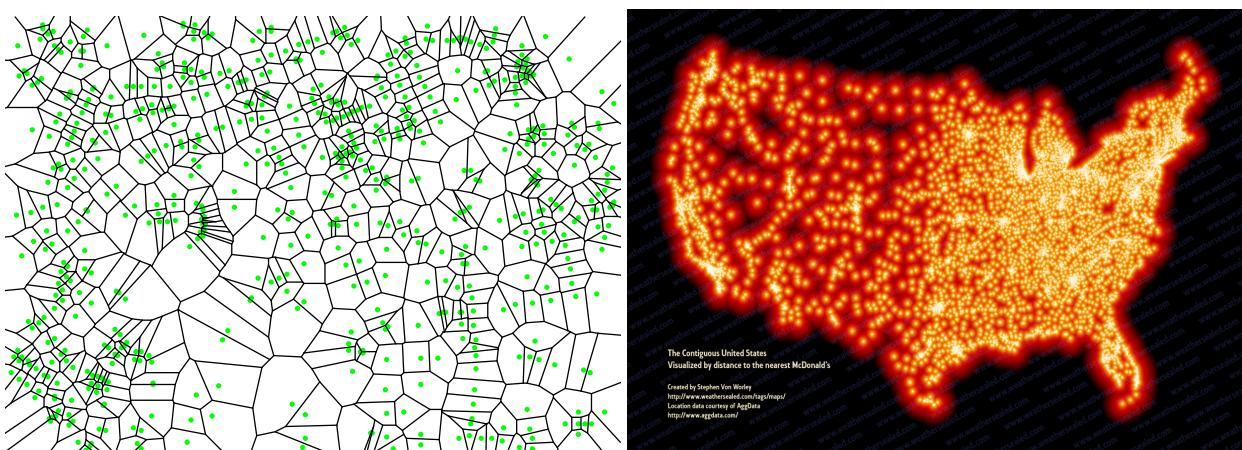
Voronoi Diagrams

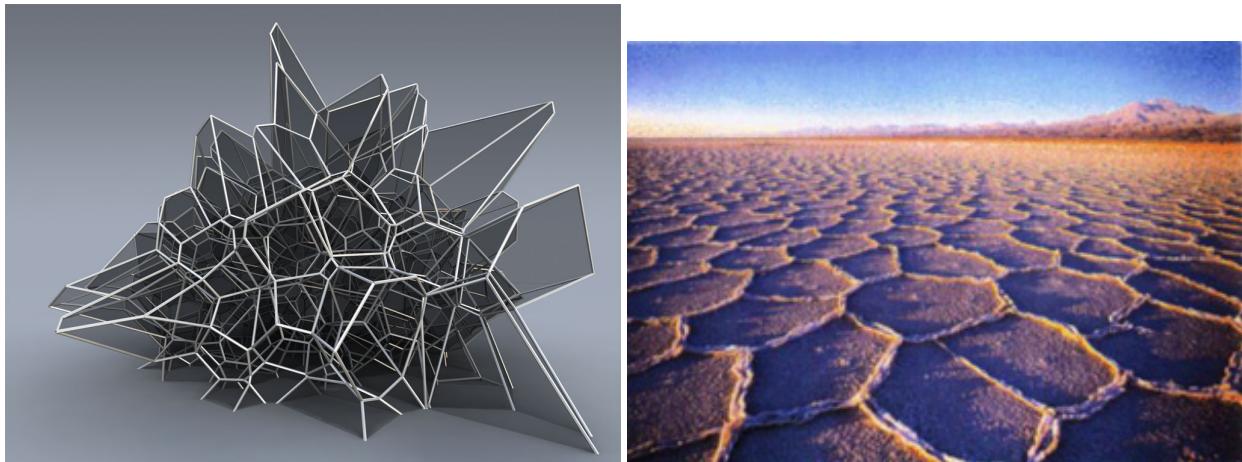
Let X be a point set. The Voronoi cell of $w \in X$ is

$$\text{Vor } w = \{p \in \mathbb{R}^d : |pw| \leq |pv| \quad \forall v \in X\}$$

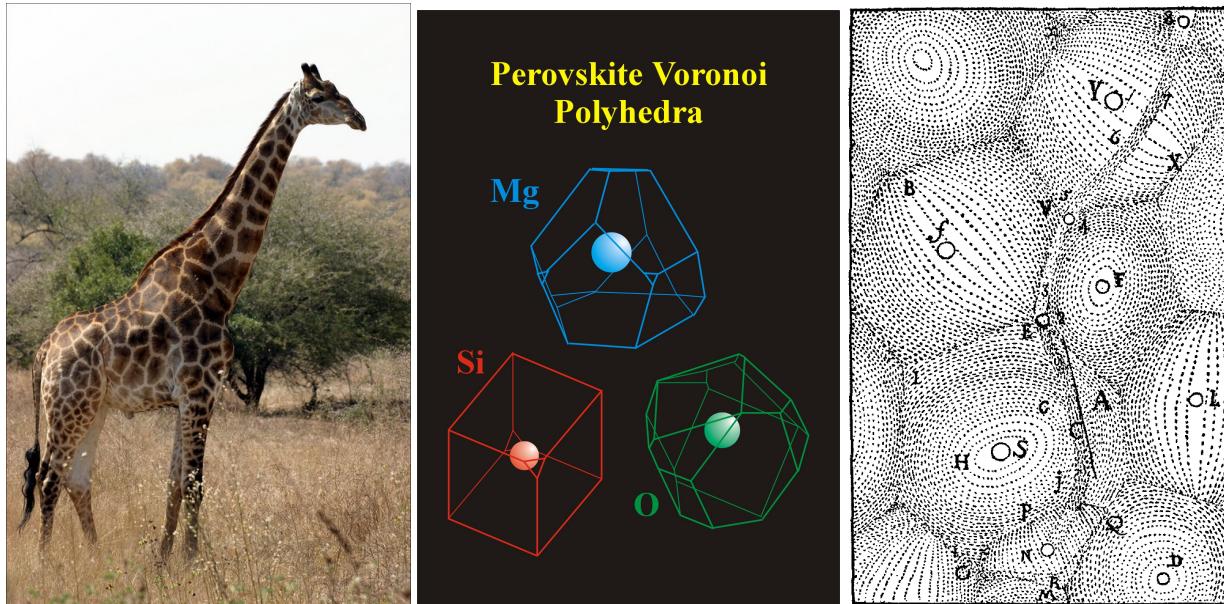
[A Voronoi cell is always a convex polyhedron or polytope.]

The Voronoi diagram of X is the set of X 's Voronoi cells.





[voronoi.pdf, vormcdonalds.jpg, voronoiGregorEichinger.jpg, saltflat-1.jpg]
 [Voronoi diagrams sometimes arise in nature (salt flats, giraffe, crystallography).]



[giraffe-1.jpg, perovskite.jpg, vortex.pdf]

[Believe it or not, the first published Voronoi diagram dates back to 1644, in the book “Principia Philosophiae” by the famous mathematician and philosopher René Descartes. He claimed that the solar system consists of vortices. In each region, matter is revolving around one of the fixed stars (vortex.pdf). His physics was wrong, but his idea of dividing space into polyhedral regions has survived.]

Size (e.g., # of vertices) $\in O(n^{\lceil d/2 \rceil})$

[This upper bound is tight when d is a small constant. As d grows, the tightest asymptotic upper bound is somewhat smaller than this, but the complexity still grows exponentially with d .]

...but often in practice it is $O(n)$.

[Here I'm leaving out a constant that may grow exponentially with d .]

Point location: Given query point $q \in \mathbb{R}^d$, find the point $w \in X$ for which $q \in \text{Vor } w$.

[We need a second data structure that can perform this search on a Voronoi diagram efficiently.]

2D: $O(n \log n)$ time to compute V.d. and a trapezoidal map for pt location

$O(\log n)$ query time [because of the trapezoidal map]

[That's a pretty great running time compared to the linear query time of exhaustive search.]

dD: Use binary space partition tree (BSP tree) for pt location

[Unfortunately, it's difficult to characterize the running time of this strategy, although it is likely to be reasonably fast in 3–5 dimensions.]

1-NN only!

[A standard Voronoi diagram supports only 1-nearest neighbor queries. If you want the k nearest neighbors, there is something called an order- k Voronoi diagram that has a cell for each possible k nearest neighbors. But nobody uses those, for two reasons. First, the size of an order- k Voronoi diagram is $O(k^2n)$ in 2D, and worse in higher dimensions. Second, there's no software available to compute one.]

[There are also Voronoi diagrams for other distance metrics, like the L_1 and L_∞ norms.]

[Voronoi diagrams are good for 1-nearest neighbor in 2 or 3 dimensions, maybe 4 or 5, but k -d trees are much simpler and probably faster in 6 or more dimensions.]

k -d Trees

“Decision trees” for NN search. Differences: [compared to decision trees]

- Choose splitting feature w/greatest width: feature i in $\max_{i,j,k}(X_{ji} - X_{ki})$.

[With nearest neighbor search, we don't care about the entropy. Instead, what we want is that if we draw a sphere around the query point, it won't intersect very many boxes of the decision tree. So it helps if the boxes are nearly cubical, rather than long and thin.]

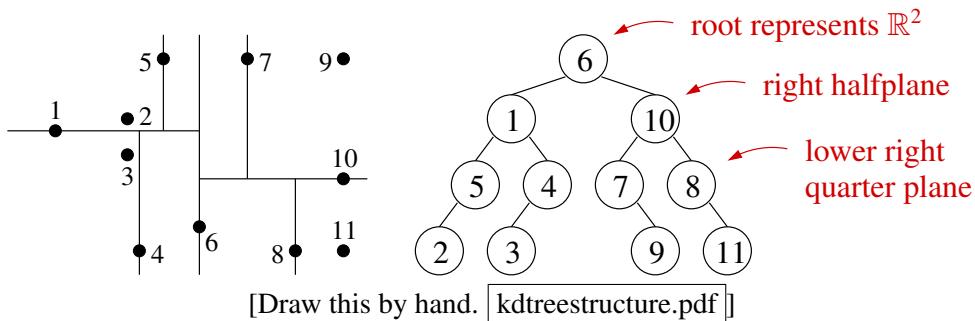
Cheap alternative: rotate through the features. [We split on the first feature at depth 1, the second feature at depth 2, and so on. This builds the tree faster, by a factor of $O(d)$.]

- Choose splitting value: median point for feature i , or $\frac{X_{ji} + X_{ki}}{2}$.

Median guarantees $\lceil \log_2 n \rceil$ tree depth; $O(nd \log n)$ tree-building time.

[... or just $O(n \log n)$ time if you rotate through the features. An alternative to the median is splitting at the box center, which improves the aspect ratios of the boxes, but it could unbalance your tree. You can alternate between medians at odd depths and centers at even depths, which also guarantees an $O(\log n)$ depth.]

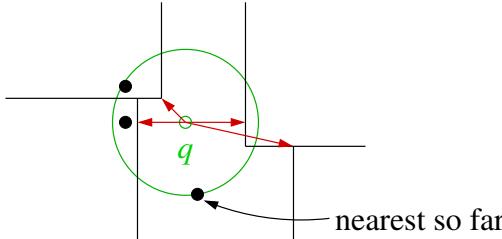
- Each internal node stores a sample point. [... that lies in the node's box. Usually the splitting point.] [Some k -d tree implementations have points only at the leaves, but it's better to have points in internal nodes too, so when we search the tree, we often stop searching earlier.]



Goal: given query pt q , find a sample pt w such that $|qw| \leq (1 + \epsilon) |qs|$, where s is the closest sample pt.
 $\epsilon = 0 \Rightarrow$ exact NN; $\epsilon > 0 \Rightarrow$ approximate NN.

Query alg. maintains:

- Nearest neighbor found so far (or k nearest). goes down ↓
- Binary heap of unexplored subtrees, keyed by distance from q . goes up ↑



[Draw this by hand. [kdtreequery.pdf](#)] [A query in progress.]

[Each subtree represents an axis-aligned box. The query tries to avoid searching most of the subtrees by searching the boxes close to q first. We measure the distance from q to a box and use it as a key for the subtree in the heap. The search stops when the distance to the k th-nearest neighbor found so far \leq the distance to the nearest unexplored box (times $1 + \epsilon$). For example, in the figure above, the query never visits the box at far lower right, because it doesn't intersect the circle.]

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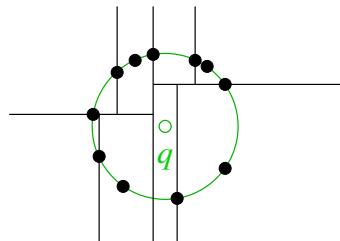
 $Q \leftarrow$  heap containing root node with key zero
 $r \leftarrow \infty$ 
while  $Q$  not empty and  $(1 + \epsilon) \cdot \text{minkey}(Q) < r$ 
     $B \leftarrow \text{removemin}(Q)$ 
     $w \leftarrow B$ 's sample point
     $r \leftarrow \min\{r, \text{dist}(q, w)\}$  [For speed, store square of  $r$  instead.]
     $B', B'' \leftarrow$  child boxes of  $B$ 
    if  $(1 + \epsilon) \cdot \text{dist}(q, B') < r$  then insert( $Q, B', \text{dist}(q, B')$ ) [The key for  $B'$  is  $\text{dist}(q, B')$ ]
    if  $(1 + \epsilon) \cdot \text{dist}(q, B'') < r$  then insert( $Q, B'', \text{dist}(q, B'')$ )
return point that determined  $r$ 

```

For k -NN, replace “ r ” with a max-heap holding the k nearest neighbors
[... just like in the exhaustive search algorithm.]

Works with any L_p norm for $p \in [1, \infty]$.
[k -d trees are not limited to the Euclidean (L_2) norm.]

Why ϵ -approximate NN?



[Draw this by hand. [kdtreeproblem.pdf](#)] [A worst-case exact NN query.]

[In the worst case, we may have to visit every node in the k -d tree to find the exact nearest neighbor. In that case, the k -d tree is slower than simple exhaustive search. This is an example where an *approximate* nearest neighbor search can be much faster. In practice, settling for an approximate nearest neighbor sometimes improves the speed by a factor of 10 or even 100, because you don't need to look at most of the tree to do a query. This is especially true in high dimensions—remember that in high-dimensional space, the nearest point often isn't much closer than a *lot* of other points.]

Software: ANN (U. Maryland), FLANN (U. British Columbia), GeRaF (U. Athens) [random forests!]

Example: im2gps

[I want to emphasize the fact that exhaustive nearest neighbor search really is one of the first classifiers you should try in practice, even if it seems too simple. So here's an example of a modern research paper that uses 1-NN and 120-NN search to solve a problem.]

Paper by James Hays and [our own] Prof. Alexei Efros.

[Goal: given a query photograph, determine where on the planet the photo was taken. Called geolocation. They evaluated both 1-NN and 120-NN. What they did not do, however, is treat each photograph as one long vector. That's okay for tiny digits, but too expensive for millions of travel photographs. Instead, they reduced each photo to a small descriptor made up of a variety of features that extract the essence of each photo.]
[Show slides (im2gps.pdf). Sorry, images not included here. <http://graphics.cs.cmu.edu/projects/im2gps/>]

[Bottom line: With 120-NN, their most sophisticated implementation came within 64 km of the correct location about 50% of the time.]

RELATED CLASSES [if you like machine learning, consider taking these courses]

CS 182 (spring): Deep Neural Networks

CS C281A (fall): Statistical Learning Theory [C281A is the most direct continuation of CS 189/289A.]

EECS 127 (both), EE 227BT (fall), EE C227C (spring): Numerical Optimization [a core part of ML]

[It's hard to overemphasize the importance of numerical optimization to machine learning, as well as other CS fields like graphics, theory, and scientific computing.]

EECS 126 (both): Random Processes [Markov chains, expectation maximization, PageRank]

EE C106A/B (fall): Intro to Robotics [dynamics, control, sensing]

Math 110: Linear Algebra [but the real gold is in Math 221]

Math 221: Matrix Computations [how to solve linear systems, compute SVDs, eigenvectors, etc.]

CS C267 (spring): Scientific Computing [parallelization, practical matrix algebra, some graph partitioning]

CS C280 (spring): Computer Vision

CS 285 (fall): Deep Reinforcement Learning (Levine)

CS 287 (fall): Advanced Robotics (Abbeel, Canny)

CS 287H (spring): Algorithmic Human-Robot Interaction (Dragan)

CS 288 (spring): Natural Language Processing (DeNero, Klein)

CS 294-43 (spring): Visual Object and Activity Recognition (Darrell)

CS 294-131 (both): Deep Learning Seminar (Darrell)

CS 294-150 (spring): ML & Biology (Listgarten)

CS 294-158 (spring): Deep Unsupervised Learning (Abbeel)

CS 294-162 (fall): ML Systems (Gonzalez)

VS 265 (fall): Neural Computation

A Bonus Lecture: Learning Theory

LEARNING THEORY: WHAT IS GENERALIZATION?

[One thing humans do well is generalize. When you were a young child, you only had to see a few examples of cows before you learned to recognize cows, including cows you had never seen before. You didn't have to see every cow. You didn't even have to see $\log n$ of the cows.]

[Learning theory tries to explain how machine learning algorithms generalize, so they can classify data they've never seen before. It also tries to derive mathematically how much training data we need to generalize well. Learning theory starts with the observation that if we want to generalize, we must constrain what hypotheses we allow our learner to consider.]

A range space (aka set system) is a pair (P, H) , where

P is set of all possible test/training points (can be infinite)

H is hypothesis class, a set of hypotheses (aka ranges, aka classifiers):

each hypothesis is a subset $h \subseteq P$ that specifies which points h predicts are in class C.

[So each hypothesis h is a 2-class classifier, and H is a set of sets of points.]

Examples:

1. Power set classifier: P is a set of k numbers; H is the power set of P , containing all 2^k subsets of P .
e.g. $P = \{1, 2\}$, $H = \{\emptyset, \{1\}, \{2\}, \{1, 2\}\}$
2. Linear classifier: $P = \mathbb{R}^d$; H is the set of all halfspaces; each halfspace has the form $\{x : w \cdot x \geq -\alpha\}$.
[In this example, both P and H are infinite. In particular, H contains every possible halfspace—that is, every possible linear classifier in d dimensions.]

[The power set classifier sounds very powerful, because it can learn every possible hypothesis. But the reality is that it can't generalize at all. Imagine we have three training points and three test points in a row.]

? C C ? N ?

[The power set classifier can classify these three test points any way you like. Unfortunately, that means it has learned nothing about the test points from the training points. By contrast, the linear classifier can learn only two hypotheses that fit this training data. The leftmost test point must be classified class C, and the rightmost test point must be classified class Not-C. Only the test point in the middle can swing either way. So the linear classifier has a big advantage: it can generalize rapidly. That's also a big disadvantage if the data isn't close to linearly separable, but that's another story.]

[Now we will investigate how well the training error predicts the test error, and how that differs for these two classifiers.]

Suppose all training pts & test pts are drawn independently from same prob. distribution \mathcal{D} defined on domain P . [\mathcal{D} also determines each point's label. Classes C and Not-C may have overlapping distributions.]

Let $h \in H$ be a hypothesis [a classifier]. h predicts a pt x is in class C if $x \in h$.

The risk aka generalization error $R(h)$ of h is the probability that h misclassifies a random pt x drawn from \mathcal{D} —i.e. the prob. that $x \in C$ but $x \notin h$ or vice versa.

[Risk is almost the same as the test error. To be precise, the risk is the average test error for test points drawn from \mathcal{D} . Sometimes the test error is higher, sometimes lower, but on average it is $R(h)$. If you had an infinite amount of test data, the risk and the test error would be the same.]

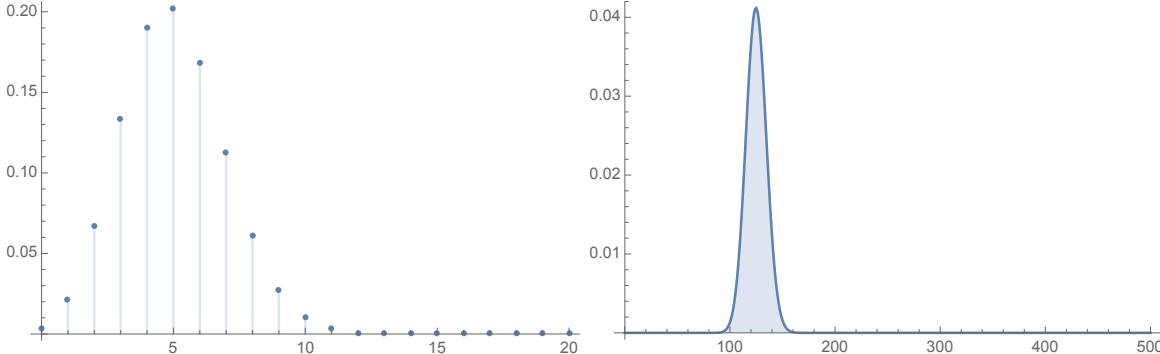
Let $X \subseteq P$ be a set of n training pts drawn from \mathcal{D}

The empirical risk aka training error $\hat{R}(h)$ is % of X misclassified by h .

[This matches the definition of empirical risk I gave you in Lecture 12, if you use the 0-1 loss function.]

h misclassifies each training pt w/prob. $R(h)$, so total misclassified has a binomial distribution.

As $n \rightarrow \infty$, $\hat{R}(h)$ better approximates $R(h)$.



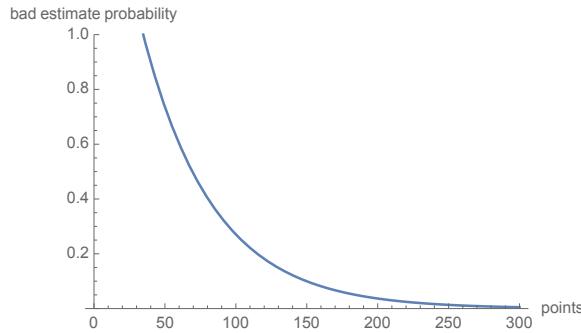
binom20.pdf, binom500.pdf [Consider a hypothesis whose risk of misclassification is 25%. Plotted are distributions of the number of misclassified training points for 20 points and 500 points, respectively. For 20 points, the training error is not a reliable estimate of the risk: the hypothesis might get “lucky” with misleadingly low training error.]

[If we had infinite training data, this distribution would become infinitely narrow and the training error would always be equal to the risk, just like the test error when we have infinite test data. But we can't have infinite training data. So, how well does the training error approximate the risk?]

Hoeffding's inequality tells us prob. of bad estimate:

$$\Pr(|\hat{R}(h) - R(h)| > \epsilon) \leq 2e^{-2\epsilon^2 n}.$$

[Hoeffding's inequality is a standard result about how likely it is that a number drawn from a binomial distribution will be far from its mean. If n is big enough, then it's very unlikely.]



hoeffding.pdf [Hoeffding's bound for the unambitious $\epsilon = 0.1$. It takes at least 200 training points to have some reasonable confidence of attaining that error bound.]

[One reason this matters is because we will try to choose the best hypothesis. If the training error is a bad estimate of the test error, we might choose a hypothesis we think is good but really isn't. So we are happy to see that the likelihood of that decays exponentially in the amount of training data.]

Idea for learning alg: choose $\hat{h} \in H$ that minimizes $\hat{R}(\hat{h})$! Empirical risk minimization.

[None of the classification algorithms we've studied actually do this, but only because it's computationally infeasible to pick the best hypothesis. Support vector machines can find a linear classifier with zero training error when the training data is linearly separable. When it isn't, SVMs try to find a linear classifier with low training error, but they don't generally find the one with minimum training error. That's NP-hard.]

[Nevertheless, for the sake of understanding learning theory, we're going to pretend that we have the computational power to try every hypothesis and pick the one with the lowest training error.]

Problem: if too many hypotheses, some h with high $R(h)$ will get lucky and have very low $\hat{R}(h)$!

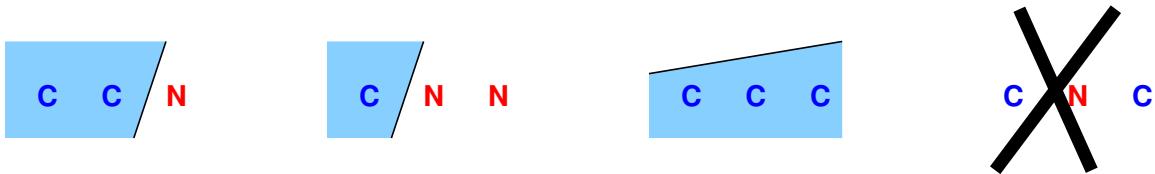
[This brings us to a central idea of learning theory. You might think that the ideal learning algorithm would have the largest class of hypotheses, so it could find the perfect one to fit the data. But the reality is that you can have so many hypotheses that some of them just get lucky and score far lower training error than their actual risk. That's another way to understand what "overfitting" is.]

[More precisely, the problem isn't too many *hypotheses*. Usually we have infinitely many hypotheses, and that's okay. The problem is too many dichotomies.]

Dichotomies

A dichotomy of X is $X \cap h$, where $h \in H$.

[A dichotomy picks out the training points that h predicts are in class C. Think of each dichotomy as assigning each training point to class C or class Not-C.]



[Draw this by hand. [dichotomies.pdf](#)] [Three examples of dichotomies for three points in a hypothesis class of linear classifiers, and one example (right) that is not a dichotomy.]

[For n training points, there could be up to 2^n dichotomies. The more dichotomies there are, the more likely it is that one of them will get lucky and have misleadingly low empirical risk.]

Extreme case: if H allows all 2^n possible dichotomies, $\hat{R}(\hat{h}) = 0$ even if every $h \in H$ has high risk.

[If our hypothesis class permits all 2^n possible assignments of the n training points to classes, then one of them will have zero training error. But that's true even if all of the hypotheses are terrible and have a large risk. Because the hypothesis class imposes no structure, we overfit the training points.]

Given Π dichotomies, $\Pr(\text{at least one dichotomy has } |\hat{R} - R| > \epsilon) \leq \delta$, where $\delta = 2\Pi e^{-2\epsilon^2 n}$.

[Let's fix a value of δ and solve for ϵ .] Hence with prob. $\geq 1 - \delta$, for every $h \in H$,

$$|\hat{R}(h) - R(h)| \leq \epsilon = \sqrt{\frac{1}{2n} \ln \frac{2\Pi}{\delta}}$$

[This tells us that the smaller we make Π , the number of possible dichotomies, and the larger we make n , the number of training points, the more likely it is that the training error will be an accurate measure of how well the classifier actually performs on test data.]

smaller Π or larger $n \Rightarrow$ training error probably closer to true risk (& test error).

[Smaller Π means we're less likely to overfit. We have less variance, but more bias. This doesn't necessarily mean the risk will be small. If our hypothesis class H doesn't fit the data well, both the training error and the test error will be large. In an ideal world, we want a hypothesis class that fits the data well, yet doesn't have many hypotheses.]

Let $h^* \in H$ minimize $R(h^*)$; "best" classifier.

[Remember we picked the classifier \hat{h} that minimizes the empirical risk. We really want the classifier h^* that minimizes the actual risk, but we can't know what h^* is. But if Π is small and n is large, the hypothesis \hat{h} we have chosen is probably nearly as good as h^* .]

With prob. $\geq 1 - \delta$, our chosen \hat{h} has nearly optimal risk:

$$R(\hat{h}) \leq \hat{R}(\hat{h}) + \epsilon \leq \hat{R}(h^*) + \epsilon \leq R(h^*) + 2\epsilon, \quad \epsilon = \sqrt{\frac{1}{2n} \ln \frac{2\Pi}{\delta}}.$$

[This is excellent news! It means that with enough training data and a limit on the number of dichotomies, empirical risk minimization usually chooses a classifier close to the best one in the hypothesis class.]

Choose a δ and an ϵ .

The sample complexity is the # of training pts needed to achieve this ϵ with prob. $\geq 1 - \delta$:

$$n \geq \frac{1}{2\epsilon^2} \ln \frac{2\Pi}{\delta}.$$

[If Π is small, we won't need too many samples to choose a good classifier. Unfortunately, if $\Pi = 2^n$ we lose, because this inequality says that n has to be bigger than n . So the power set classifier can't learn much or generalize at all. We need to severely reduce Π , the number of possible dichotomies. One way to do that is to use a linear classifier.]

The Shatter Function & Linear Classifiers

[How many ways can you divide n points into two classes with a hyperplane?]

$$\begin{aligned} \text{\# of dichotomies: } \Pi_H(X) &= |\{X \cap h : h \in H\}| && \in [1, 2^n] \text{ where } n = |X| \\ \text{shatter function: } \Pi_H(n) &= \max_{|X|=n, X \subseteq P} \Pi_H(X) && [\text{The most dichotomies of any point set of size } n] \end{aligned}$$

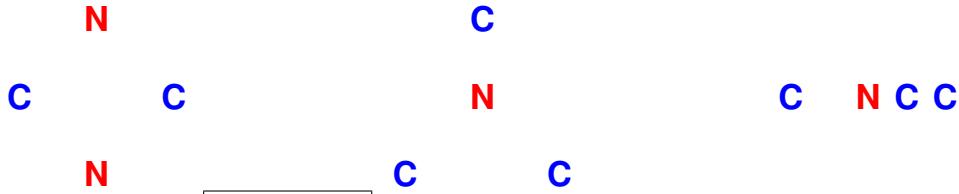
Example: Linear classifiers in plane. H = set of all halfplanes. $\Pi_H(3) = 8$:



[Draw this by hand. [shatter.pdf](#)] [Linear classifiers can induce all eight dichotomies of these three points. The other four dichotomies are the complements of these four.]

$\Pi_H(4) = 14$:

[Instead of showing you all 14 dichotomies, let me show you dichotomies that halfplanes *cannot* learn, which illustrate why no four points have 16 dichotomies.]



[Draw this by hand. [unshatter.pdf](#)] [Examples of dichotomies of four points in the plane that no linear classifier can induce.]

[This isn't a proof that 14 is the maximum, because we have to show that 16 is not possible for *any* four points in the plane. The standard proof uses a result called Radon's Theorem.]

Fact: for all range spaces, either $\Pi_H(n)$ is polynomial in n , or $\Pi_H(n) = 2^n \ \forall n \geq 0$.

[This is a surprising fact with deep implications. Imagine that you have m points, some of them training points and some of them test points. Either a range space permits every possible dichotomy of the points, and the training points don't help you classify the test points at all; or the range space permits only a polynomial subset of the 2^m possible dichotomies, so once you have labeled the training points, you have usually cut down the number of ways you can classify the test points dramatically. No shatter function ever occupies the no-man's-land between polynomial and 2^m .]

[For linear classifiers, we know exactly how many dichotomies there can be.]

Cover's Theorem [1965]: linear classifiers in \mathbb{R}^d allow up to $\Pi_H(n) = 2 \sum_{i=0}^d \binom{n-1}{i}$ dichotomies of n pts.

For $n \leq d + 1$, $\Pi_H(n) = 2^n$.

For $n \geq d + 1$, $\Pi_H(n) \leq 2 \left(\frac{e(n-1)}{d}\right)^d$ [Observe that this is polynomial in n ! With exponent d .] and the sample complexity needed to achieve $R(\hat{h}) \leq \hat{R}(\hat{h}) + \epsilon \leq R(h^*) + 2\epsilon$ with prob. $\geq 1 - \delta$ is

$$n \geq \frac{1}{2\epsilon^2} \left(d \ln \frac{n-1}{d} + d + \ln \frac{4}{\delta} \right). \quad [\text{Observe that the logarithm turned the exponent } d \text{ into a factor!}]$$

Corollary: linear classifiers need only $n \in O(d)$ training pts
for training error to accurately predict risk/test error.

[In a d -dimensional feature space, we need more than d training points to train an accurate linear classifier. But it's reassuring to know that the number we need is linear in d . By contrast, if we have a classifier that permits all 2^n possible dichotomies however large n is, then no amount of training data will guarantee that the training error of the hypothesis we choose approximates the true risk.]

[The constant hidden in that big-O notation can be quite large. For comforting choices of ϵ and δ , it's at least 100. If you want a lot of confidence that you've chosen one of the best hypotheses, you have to pay for it with a large sample size.]

[This sample complexity applies even if you add polynomial features or other features, but you have to count the extra features in d . So the number of training points you need increases with the number of polynomial terms.]

VC Dimension

The Vapnik–Chervonenkis dimension of (P, H) is

$$\text{VC}(H) = \max\{n : \Pi_H(n) = 2^n\}. \quad \Leftarrow \text{Can be } \infty.$$

Say that H shatters a set X of n pts if $\Pi_H(X) = 2^n$.

$\text{VC}(H)$ is size of largest X that H can shatter.

[This means that X is a point set for which all 2^n dichotomies are possible.]

[I told you earlier that if the shatter function isn't 2^n for all n , then it's a polynomial in n . The VC dimension is motivated by an observation that sometimes makes it easy to bound that polynomial.]

$$\text{Theorem: } \Pi_H(n) \leq \sum_{i=0}^{\text{VC}(H)} \binom{n}{i}. \quad \text{Hence for } n \geq \text{VC}(H), \quad \Pi_H(n) \leq \left(\frac{en}{\text{VC}(H)}\right)^{\text{VC}(H)}.$$

[So the VC dimension is an upper bound on the exponent of the polynomial. This theorem is useful because often we can find an easy upper bound on the VC dimension. You just need to show that for some number n , no set of n points can have all 2^n dichotomies.]

Corollary: $O(\text{VC}(H))$ training pts suffice for accuracy. [Again, the hidden constant is big.]

[If the VC dimension is finite, it tells us how the sample complexity grows with the number of features. If the VC dimension is infinite, no amount of training data will make the classifier generalize well.]

Example: Linear classifiers in plane.

Recall $\Pi_H(3) = 8$: there exist 3 pts shattered by halfplanes.

But $\Pi_H(4) = 14$: no 4 pts are shattered.

Hence:

- $\text{VC}(H) = 3$ [The VC dimension of halplanes is 3.]
- $\Pi_H(n) \leq \frac{e^3}{27} n^3$ [The shatter function is polynomial.]
- $O(1)$ sample complexity.

[The VC dimension doesn't always give us the tightest bound. In this example, the VC dimension promises that the number of ways halfplanes can classify the points is at worst cubic in n ; but in reality, it's quadratic in n . In general, linear classifiers in d dimensions have VC dimension $d + 1$, which is one dimension looser than the exponent Cover proved. That's not a big deal, though, as the sample complexity and the accuracy bound are both based on the *logarithm* of the shatter function. So if we get the exponent wrong, it only changes a constant in the sample complexity.]

[The important thing is simply to show that there is some polynomial bound on the shatter function at all. VC dimension is not the only way to do that, but often it's the easiest.]

[The main point you should take from this lecture is that if you want to have generalization, you need to limit the expressiveness of your hypothesis class so that you limit the number of possible dichotomies of a point set. This may or may not increase the bias, but if you don't limit the number of dichotomies at all, the overfitting will be very bad. If you limit the hypothesis class, your artificial child will only need to look at $O(d)$ cows to learn the concept of cows. If you don't, your artificial child will need to look at every cow in the world, and every non-cow too.]