

## Module 5: Classification

# Mixture Models for Classification

STAT/BIOSTAT 527, University of Washington

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## Overview of Classification So Far

- Supervised methods

$(x_1, y_1), \dots, (x_n, y_n)$   
↑ input data      ↖ class label

"labeled data"  
as training data

Generative

Discriminative

LDA, QDA  
KDE for class.  
Naive Bayes  
⋮

logistic reg.  
CART  
⋮

- Objectives:

$\max P(Y, X)$

$\max P(Y|X)$

- Unsupervised methods (generative)

mixture models

$x_1, \dots, x_n$

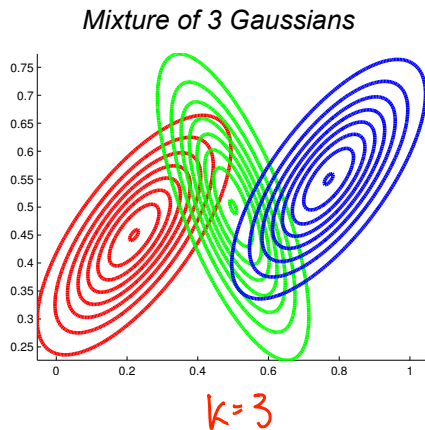
"unlabeled data"

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# Density as Mixture of Gaussians

- Approximate density with a mixture of Gaussians



$$p(x_i | \pi, \mu, \Sigma) = \sum_{k=1}^K \pi_k N(x_i | \mu_k, \Sigma_k)$$

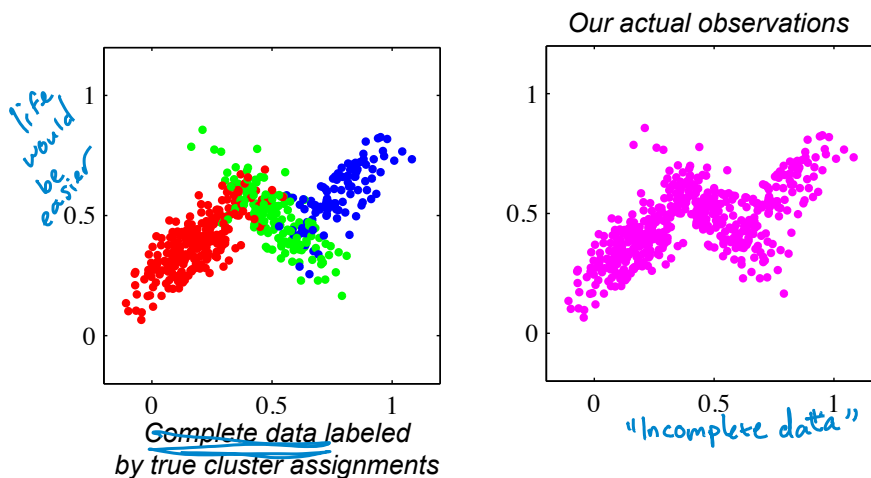
*Handwritten notes:*  
 Gauss. kernel, just like in KDE, but not centered at obs.  
 $\pi = [\pi_1, \dots, \pi_K]$   
 $\mu = \{\mu_1, \dots, \mu_K\}$   
 $\Sigma = \{\Sigma_1, \dots, \Sigma_K\}$   
 # of mix comp.  $\rightarrow K$   
 $\pi_k$  mix. weights  $\rightarrow$  shape params  
 In 1D:  $\leftarrow P = \text{target density}$   
 $\Sigma \pi_k = 1$

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# Clustering our Observations

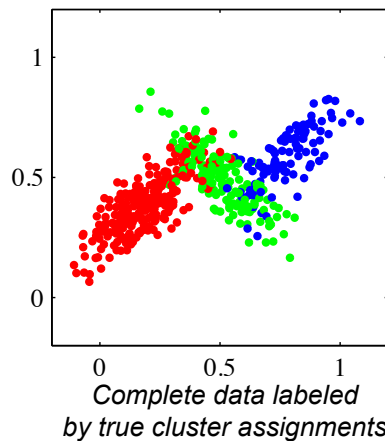
- Imagine we have an assignment of each  $x_i$  to a Gaussian



C. Bishop, Pattern Recognition & Machine Learning

# Clustering our Observations

- Imagine we have an assignment of each  $x_i$  to a Gaussian

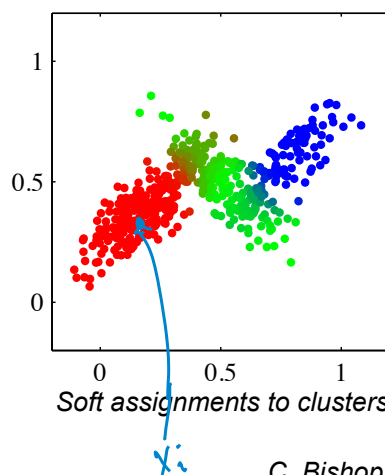


- Introduce latent cluster indicator variable  $z_i$   
 $z_i \in \{1, \dots, K\}$   
 $Pr(z_i = k) = \pi_k$
- Then we have  
 $p(x_i | z_i, \pi, \mu, \Sigma) = N(x_i | \mu_{z_i}, \Sigma_{z_i})$   
*think of as class indicator  $y_i$ , but no training data labels*  
*param. est. is easy if we have  $\{z_i\}$   $\Rightarrow$  decoupled into K Gauss. est.*

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# Clustering our Observations

- We must infer the cluster assignments from the observations



- Posterior probabilities of assignments to each cluster \*given\* model parameters:  
*"responsibilities"*  
 $r_{ik} = p(z_i = k | x_i, \pi, \theta) = \frac{\pi_k N(x_i | \mu_k, \Sigma_k)}{\sum_j \pi_j N(x_i | \mu_j, \Sigma_j)}$   
*motivates an iterative alg.*

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# Mixture Models for Classification

- Can use mixture models as a generative classifier in the unsupervised setting

✓ expectation maximization (EM)

- EM algorithm = iteratively: → point est.

- Estimate responsibilities given parameter estimates

$$\hat{r}_{ik} = \frac{\hat{\pi}_k N(x_i, \hat{\mu}_k, \hat{\Sigma}_k)}{\sum_{\ell} \hat{\pi}_{\ell} N(x_i, \hat{\mu}_{\ell}, \hat{\Sigma}_{\ell})}$$

- Maximize parameters given responsibilities

$\{\hat{\pi}_k, \hat{\mu}_k, \hat{\Sigma}_k\}$

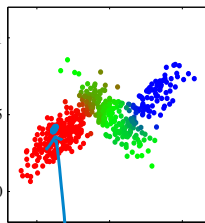
iterate

- For classification, threshold the estimated responsibilities

- E.g.,  $\hat{g}(x_i) = \arg \max_k \hat{r}_{ik}$

- Note: allows non-linear boundaries as in QDA

$\Sigma_k$



0-1 loss  
 $r_{ik}$  for  $k=1,2,3$

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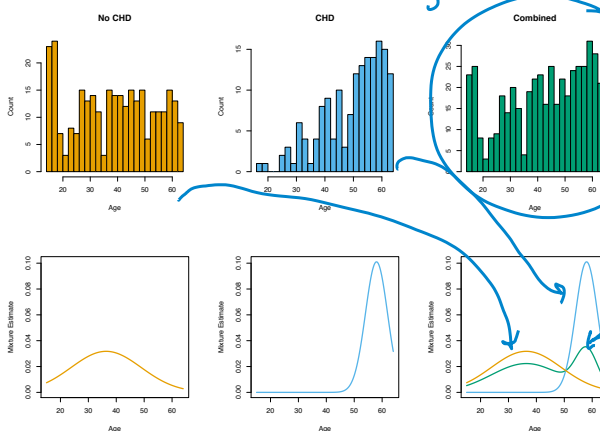
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# Example: Heart Disease Data

- Binary response = CHD (coronary heart disease)

- Predictor = ~~systolic blood pressure~~ age

using this



EM alg.

combined

From Hastie, Tibshirani, Friedman book

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# What you need to know

- Discriminative vs. Generative classifiers
- LDA and QDA assume Gaussian class-conditional densities
  - Results in linear and quadratic decision boundaries, respectively
- KDE for classification
  - Challenging in areas with little data or in high dimensions
  - Estimating class-conditionals is not optimizing classification objective
- Naïve Bayes assumes factored form
  - Results in log odds that have GAM form
- Mixture models allow for unsupervised generative approach

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# Readings

- Hastie, Tibshirani, Friedman – 4.3, 4.4.5, 6.6.2-6.6.3, 6.8

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## Module 5: Classification

# Online Learning Perceptron Algorithm

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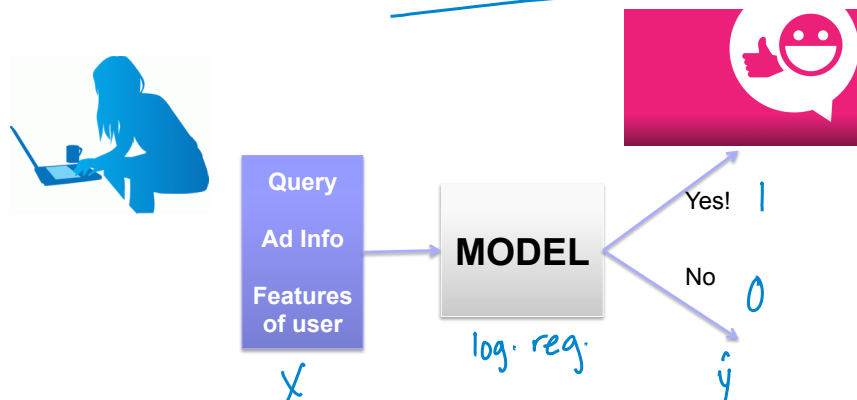
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## Estimating Click Probabilities

- **Goal:** Predict whether a person clicks on an ad
- **Basic approach:** Logistic regression



# Challenge 1: Complexity of Computing Gradients *in terms of n, d*

- What's the cost of a gradient update step for <sup>reg:</sup> LR???

$$\beta_j^{(t+1)} \leftarrow \beta_j^{(t)} + \eta \left\{ -\lambda \beta_j^{(t)} + \sum_i x_{ij} \left( y_i - \hat{p}(y=1 | x_i, \beta^{(t)}) \right) \right\}$$

*instead, talked about stochastic grad. ascent ... small change after each obs.*

*for each j*

*O(d)*

*O(nd)*

Naively,  $O(nd^2)$

but cache  $\hat{p}$  (same  $\forall j$ )  $\rightarrow O(nd)$

However, if n is huge (or streaming), this is very slow (infeasible) per little gradient step

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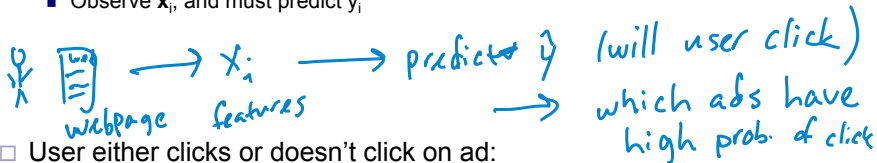
# Challenge 2: Data is streaming

- Assumption thus far: **Batch data**

*have data, learn model*

- But, e.g., click prediction for ads is a streaming data task:

- User enters query, and ad must be selected:
  - Observe  $x_i$ , and must predict  $y_i$



- User either clicks or doesn't click on ad:
  - Label  $y_i$  is revealed afterwards

- Google gets a reward if user clicks on ad

*reward is just like 0-1 loss in class. setting*

- Weights must be updated for next time:

$$\beta^{(t+1)} \leftarrow \beta^{(t)} + \Delta \leftarrow \mathbb{1}(y_i \neq \hat{y})$$

*.. what is this*

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# Online Learning Problem

- At each time step  $t$ :

- Observe features (covariates) of data point:

- Note: many assumptions are possible, e.g., data is iid, data is adversarially chosen... details beyond scope of course

- Make a prediction:

- Note: many models are possible, we focus on linear models

$$\beta_0^{(t)} + \sum_j \beta_j^{(t)} X_{tj} > 0 \Rightarrow \text{"click"} \quad \parallel \quad \beta^{(t)} \cdot X_t > 0$$

*vec.  $X_{t0}=1$*

- Observe true label:

- Note: other observation models are possible, e.g., we don't observe the label directly, but only a noisy version... Details beyond scope of course

$$y_t \begin{cases} \rightarrow \text{clicked} \\ \rightarrow \text{not clicked} \end{cases}$$

- Update model:

$$\beta^{(t+1)} \leftarrow \beta^{(t)} + \Delta^{(t)} \text{ something}$$

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# The Perceptron Algorithm [Rosenblatt '58, '62]

- Classification setting:  $y$  in  $\{-1, +1\}$

*(note diff. from  $\{0, 1\}$ ... just a practical change)*

- Linear model

- Prediction:

$$\hat{y} = \text{sign}(\beta \cdot x)$$

- Training:

- Initialize weight vector:

*e.g.  $\beta^{(0)} = 0$  (or smarter)*

- At each time step:

- Observe covariates:

$$X_t$$

- Make prediction:

$$\hat{y} = \text{sign}(\beta^{(t)} \cdot x)$$

- Observe true class:

$$y_t \leftarrow \text{true label}$$

- Update model:

- If prediction is not equal to truth

*... if we made mistake*

$$\text{if } \hat{y} \neq y_t, \beta^{(t+1)} \leftarrow \beta^{(t)} \\ \text{else } \beta^{(t+1)} \leftarrow \beta^{(t)} + y_t X_t$$

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## Intuition

If  $\hat{y} = y_t$ ,

$$\beta^{(t+1)} \leftarrow \beta^{(t)}$$

else

$$\underline{\beta^{(t+1)}} \leftarrow \beta^{(t)} + \underline{y_t x_t}$$

$$\hat{y} = \text{sign}(\beta^{(t)} \cdot x_t)$$

- Why is this a reasonable update rule?

if mistake e.g.  $y_t = +1$  but  $\beta^{(t)} \cdot x_t < 0$  ... wanted  $\beta^{(t)} \cdot x_t > 0$

What  $u$  max  $u \cdot x_t \rightarrow u^* = x_t$

by adding  $x_t$  to  $\beta^{(t)}$ , we increase  $\beta^{(t+1)} \cdot x_t$  the most.  
similarly for  $y_t = -1$

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## Which weight vector to report?

- Practical problem for all online learning methods
- Suppose you run online learning method and want to sell your learned weight vector... Which one do you sell???

- Last one?  $\beta^{(T)}$ ? no... very noisy. Influenced by last mistake

- Random  $\beta^{(\tau)}$  random no

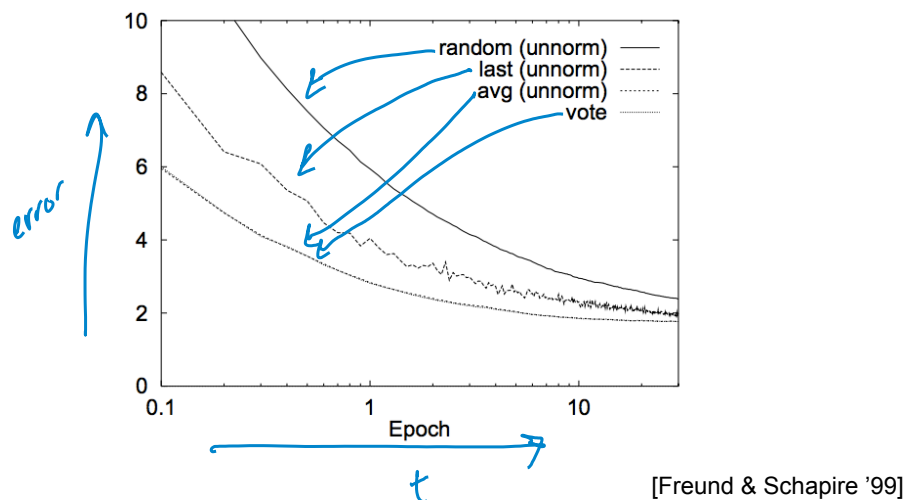
- Average  $\hat{\beta} = \frac{1}{T} \sum_{t=1}^T \beta^{(t)}$  easy to maintain

- Voting + more advanced : how long has this param been ground?

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## Choice can make a huge difference!!



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## Mistake Bounds

Why does it work?

- Algorithm “pays” every time it makes a mistake:

Loss fcn in online setting:  
# mistakes up to time  $T$

⇒ Google pays for its mistakes

- How many mistakes is it going to make?

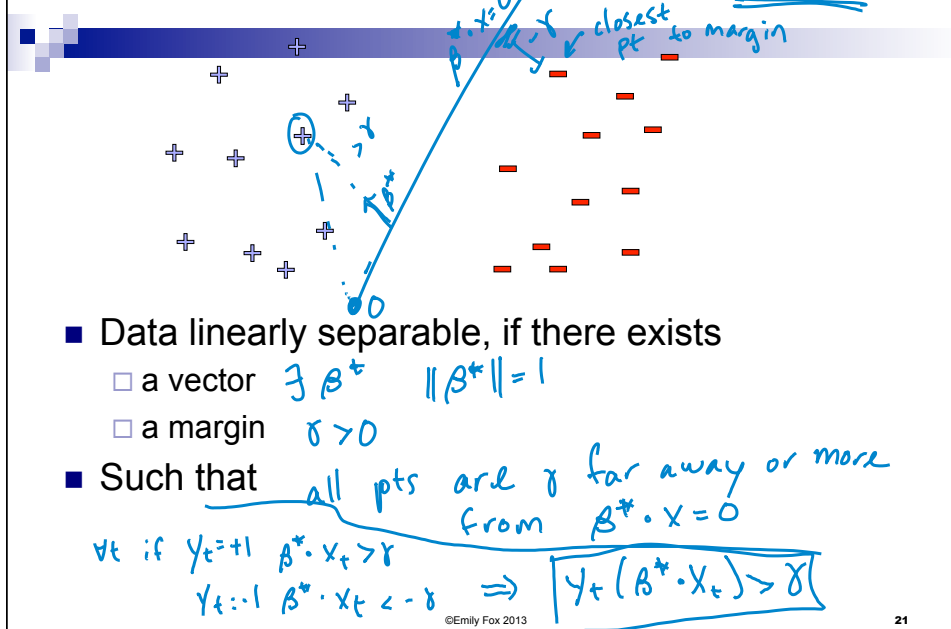
“mistake bound”

in its  
lifetime

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## Linear Separability: More formally, Using Margin



## Perceptron Analysis: Linearly Separable Case

### ■ Theorem [Block, Novikoff]:

- Given a sequence of labeled examples:  $(x_1, y_1), \dots, (x_n, y_n)$   
*examples need not be iid nor random*
- Each covariate vector has bounded norm:

$$\forall t \quad \|x_t\| \leq R$$

- If dataset is linearly separable:

$$\exists \beta^* \quad \|\beta^*\|=1 \quad \text{s.t.} \quad y_t (\beta^* \cdot x_t) \geq \delta \quad \text{for some } \delta > 0$$

- Then the number of mistakes made by the online perceptron on this sequence is bounded by

$$\left(\frac{R}{\delta}\right)^2$$

*crazy!*

*constant... doesn't depend on T or dim X*

## Perceptron Proof for Linearly Separable case

- Every time we make a mistake, we get  $\gamma$  closer to  $\beta^*$ :

- Mistake at time  $t$ :  $\beta^{(t+1)} \leftarrow \beta^{(t)} + y_t x_t$

- Taking dot product with  $\beta^*$ :

- Thus after  $m$  mistakes:

by induction,  $\beta^* \cdot \beta^{(t+1)} \geq m\gamma$

- Similarly, norm of  $\beta^{(t+1)}$  doesn't grow too fast:

- $\|\beta^{(t+1)}\|^2 = \|\beta^{(t)}\|^2 + 2y_t(\beta^{(t)} \cdot x_t) + \|x_t\|^2$

- Thus, after  $m$  mistakes:

$$\|\beta^{(t+1)}\|^2 \leq mR^2$$

- Putting all together:

$$m\gamma \leq \beta^* \cdot \beta^{(t+1)} \leq \|\beta^*\| \|\beta^{(t+1)}\| \leq \sqrt{m} R$$

$$\Rightarrow m\gamma \leq \sqrt{m} R \Rightarrow m \leq \left(\frac{R}{\gamma}\right)^2$$

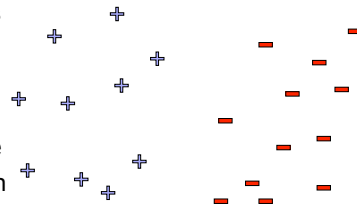
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## Beyond Linearly Separable Case

- Perceptron algorithm is super cool!

- No assumption about data distribution!
  - Could be generated by an oblivious adversary, no need to be iid
- Makes a fixed number of mistakes, and it's done for ever!
  - Even if you see infinite data



- However, real world not linearly separable
  - Can't expect never to make mistakes again
  - Analysis extends to non-linearly separable case
  - Very similar bound, see Freund & Schapire
  - Converges, but ultimately may not give good accuracy (make many many many mistakes)

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# What is the Perceptron Doing???

- When we discussed logistic regression:
  - Started from maximizing conditional log-likelihood

$$\max_{\beta} P(Y|X, \beta)$$

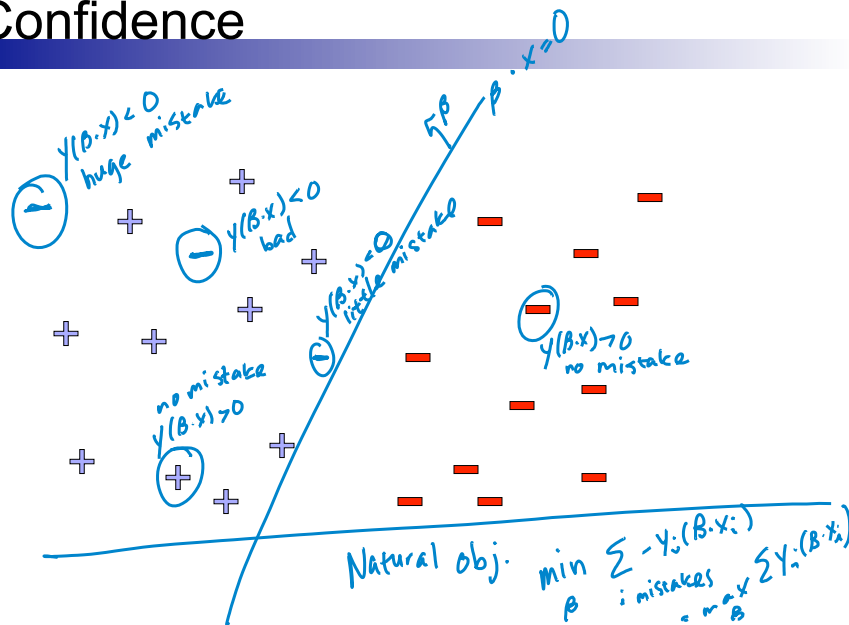
- When we discussed the perceptron:
  - Started from description of an algorithm

- What is the perceptron optimizing????

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# Perceptron Prediction: Margin of Confidence

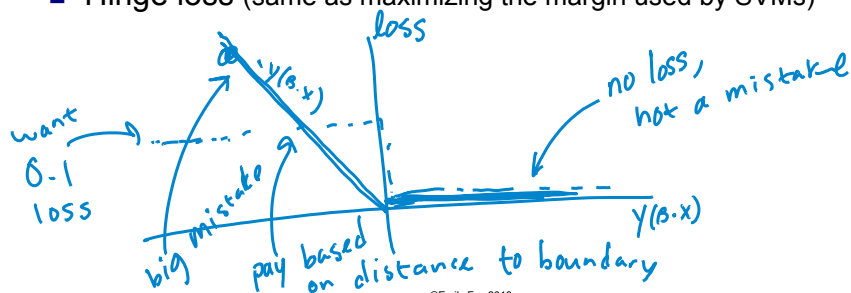


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# Hinge Loss

- Perceptron prediction:  $\text{sign}(\beta \cdot x)$
- Makes a mistake when:  $y(\beta \cdot x) < 0 \Rightarrow l(\beta, x) = \begin{cases} 0 & y(\beta \cdot x) > 0 \\ -y(\beta \cdot x) & \text{else} \end{cases} \Rightarrow (-y(\beta \cdot x))_+$
- Hinge loss (same as maximizing the margin used by SVMs)

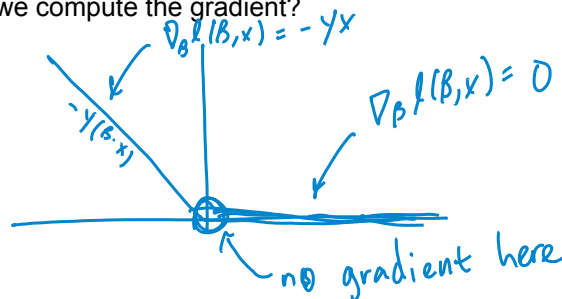


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# Minimizing Hinge Loss in Batch Setting

- Given a dataset:  $(x_1, y_1), \dots, (x_n, y_n)$
- Minimize average hinge loss:  $\min_{\beta} \frac{1}{n} \sum_{i=1}^n l(\beta, x_i)$   $(-y_i(\beta \cdot x_i))_+$
- How do we compute the gradient?

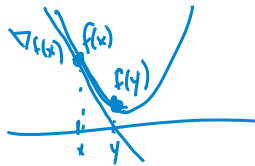


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# Subgradients of Convex Functions

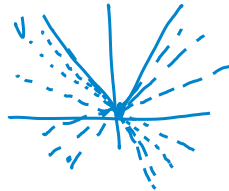
- Gradients lower bound convex functions:



$$f(y) \geq f(x) + \nabla f(x)(y-x)$$

- Gradients are unique at  $x$  if function differentiable at  $x$
- Subgradients: Generalize gradients to non-differentiable points:

- Any plane that lower bounds function:



For  $|\beta_j| \leq 1$   
 $\forall \beta \in [-1, 1]$

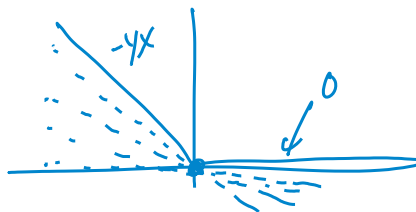
$\forall \beta \in \partial f(x)$  subgradient  
 if  
 $f(y) \geq f(x) + \beta(y-x)$

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# Subgradient of Hinge

- Hinge loss:



- Subgradient of hinge loss:

- If  $y_t(\beta \cdot \mathbf{x}_t) > 0$ :  $\partial \ell(\beta, \mathbf{x}) = 0$
- If  $y_t(\beta \cdot \mathbf{x}_t) < 0$ :  $\partial \ell(\beta, \mathbf{x}) = -y \mathbf{x}$
- If  $y_t(\beta \cdot \mathbf{x}_t) = 0$ :  $\partial \ell(\beta, \mathbf{x}) = [-y \mathbf{x}, 0]$  e.g.  $-y \mathbf{x}$
- In one line:

$$\partial \ell(\beta, \mathbf{x}) = \mathbb{1}(y(\beta \cdot \mathbf{x}) \leq 0) (-y \mathbf{x})$$

mistake indicator

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## Subgradient Descent for Hinge Minimization

- Given data:  $(x_1, y_1), \dots, (x_n, y_n)$

- Want to minimize:

$$\frac{1}{n} \sum_{i=1}^n \ell(\beta, x_i) = \frac{1}{n} \sum_{i=1}^n (-y_i(\beta \cdot x_i))_+$$

- Subgradient descent works the same as gradient descent:
  - But if there are multiple subgradients at a point, just pick (any) one:

$$\beta^{(t+1)} \leftarrow \beta^{(t)} - \eta \sum_{i=1}^n \underbrace{\partial \ell(\beta, x_i)}_{\mathbb{I}(y_i(\beta \cdot x_i) \leq 0) (-y_i x_i)}$$

↑  
step size

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## Perceptron Revisited

- Perceptron update:

$$\beta^{(t+1)} \leftarrow \beta^{(t)} + \eta \mathbb{I}[y_t(\beta^{(t)} \cdot x_t) \leq 0] y_t x_t$$

↑  $\eta=1$  if mistake neg. subgrad.

- Batch hinge minimization update:

$$\beta^{(t+1)} \leftarrow \beta^{(t)} + \eta \frac{1}{n} \sum_{i=1}^n \left\{ \mathbb{I}[y_i(\beta^{(t)} \cdot x_i) \leq 0] y_i x_i \right\}$$

↑ if mistake neg. subgrad.

param, hopefully  $\eta > 1$

- Difference? Perceptron algorithm = SGD for hinge loss minimization using  $\eta=1$

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# What you need to know

- Notion of online learning
- Perceptron algorithm
- Mistake bounds and proof
- In online learning, report averaged weights at the end
- Perceptron is optimizing hinge loss
- Subgradients and hinge loss
- (Sub)gradient decent for hinge objective