

CS9840
Learning and Computer Vision
Prof. Olga Veksler

Lecture 3
Linear Machines
Information Theory (a little BIT)

Today

- Linear Classifier
- Mutual Information
- Next time:
 - paper: “Object Recognition with Informative Features and Linear Classification” by M. Naquet and S. Ullman
 - Ignore section of tree-augmented network

Last Time: Supervised Learning

- Training samples (or examples) X^1, X^2, \dots, X^n
- Each example is typically multi-dimensional
 - $X^i_1, X^i_2, \dots, X^i_d$ are typically called *features*, X^i is sometimes called a *feature vector*
 - **How many features and which features do we take?**
- Know desired output for each example (labeled samples) Y^1, Y^2, \dots, Y^n
 - This learning is supervised (“teacher” gives desired outputs).
 - Y^i are often one-dimensional, but can be multidimensional

Last Time: Supervised Learning

- Wish to design a *machine* $f(X, W)$ s.t.
 $f(X, W) = \text{true output value at } X$
 - In classification want $f(X, W) = \text{label of } X$
 - **How do we choose f ?**
 - when we choose a particular f , we are making implicit assumptions about our problem
 - W is typically multidimensional vector of weights (also called *parameters*) which enable the machine to “learn”
 - $W = [w_1, w_2, \dots, w_k]$

Training and Testing

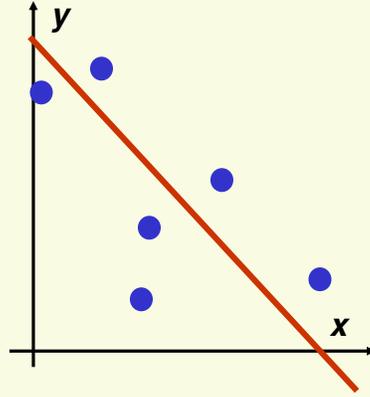
- There are 2 phases, training and testing
 - Divide all labeled samples X^1, X^2, \dots, X^n into 2 sets, *training* set and *testing* set
 - Training phase is for “teaching” our machine (finding optimal weights W)
 - Testing phase is for evaluating how well our machine works on unseen examples
- Training phase
 - Find the weights W s.t. $f(X^i, W) = Y^i$ “as much as possible” for the *training* samples X^i
 - “as much as possible” needs to be defined
 - Training can be quite complex and time-consuming

Loss Function

- How do we quantify what it means for the machine $f(X, W)$ do well in the training and testing phases?
- $f(X, W)$ has to be “close” to the true output on X
- Define Loss (or Error) function L
 - This is up to the designer (that is you)
- Typically first define per-sample loss $L(X^i, Y^i, W)$
 - Some examples:
 - for classification, $L(X^i, Y^i, W) = \mathbf{I}[f(X^i, W) \neq Y^i]$, where $\mathbf{I}[\text{true}] = 1$, $\mathbf{I}[\text{false}] = 0$
 - we just care if the sample has been classified correctly
 - For continuous Y , $L(X^i, Y^i, W) = \|f(X^i, W) - Y^i\|^2$,
 - how far is the estimated output from the correct one?
- Then loss function $L = \sum_i L(X^i, Y^i, W)$
 - Number of missclassified example for classification
 - Sum of distances from the estimated output to the correct output

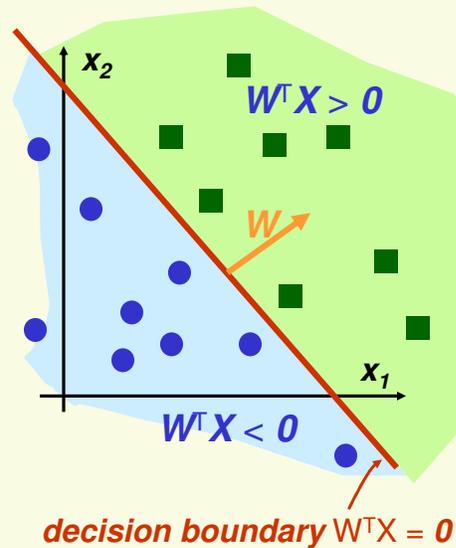
Linear Machine, Continuous Y

- $f(X,W) = w_0 + \sum_{i=1,2,\dots,d} w_i x_i$
 - w_0 is called bias
- In vector form, if we let $X = (1, x_1, x_2, \dots, x_d)$, then $f(X,W) = W^T X$
 - notice abuse of notation, I made $X = [1 \ X]$
- This is standard linear regression (line fitting)
 - assume $L(X^i, Y^i, W) = \|f(X^i, W) - Y^i\|^2$
 - optimal W can be found by solving linear system of equations $W^* = [\sum X^i (X^i)^T]^{-1} \sum Y^i X^i$



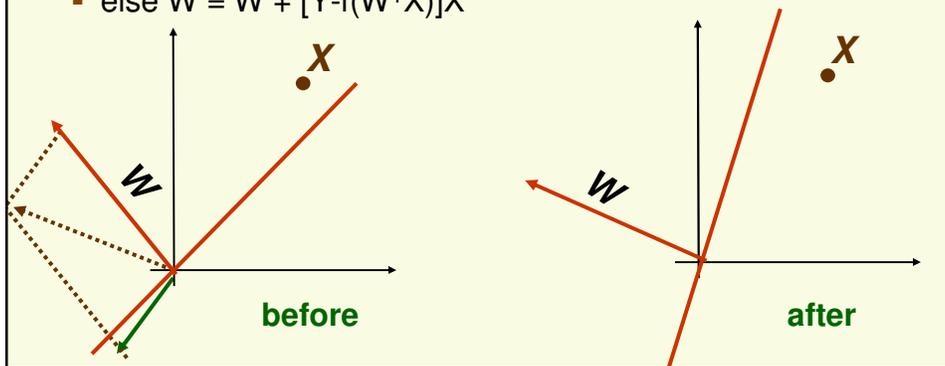
Linear Machine: binary Y

- $f(X,W) = \text{sign}(w_0 + \sum_{i=1,2,\dots,d} w_i x_i)$
 - $\text{sign}(\text{positive}) = 1$,
 $\text{sign}(\text{negative}) = -1$
 - w_0 is called bias
- In vector form, if we let $X = (1, x_1, x_2, \dots, x_d)$ then $f(X,W) = \text{sign}(W^T X)$



Perceptron Learning Procedure (Rosenblatt 1957)

- $f(X, W) = \text{sign}(w_0 + \sum_{i=1,2,\dots,d} w_i x_i)$
- Let $L(X^i, Y^i, W) = \mathbf{I}[f(X^i, W) \neq Y^i]$. How do we learn W ?
- A solution:
- Iterate over all training samples
 - if $f(X, W) = Y$ (correct label), do nothing
 - else $W = W + [Y - f(W^T X)] X$



Perceptron Learning Procedure (Rosenblatt 1957)

- Amazing fact: If the samples are linearly separable, the perceptron learning procedure will converge to a solution (separating hyperplane) in a finite amount of time
- Bad news: If the samples are not linearly separable, the perceptron procedure will not terminate, it will go on looking for a solution which does not exist!
- For most interesting problems the samples are not linearly separable
- Is there a way to learn W in non-separable case?
 - Remember, it's ok to have training error, so we don't have to have "perfect" classification

Optimization

- Need to minimize a function of many variables

$$J(\mathbf{x}) = J(x_1, \dots, x_d)$$

- We know how to minimize $J(\mathbf{x})$
 - Take partial derivatives and set them to zero

$$\begin{bmatrix} \frac{\partial}{\partial x_1} J(\mathbf{x}) \\ \vdots \\ \frac{\partial}{\partial x_d} J(\mathbf{x}) \end{bmatrix} = \nabla J(\mathbf{x}) = 0$$

gradient

- However solving analytically is not always easy

- Would you like to solve this system of nonlinear equations?

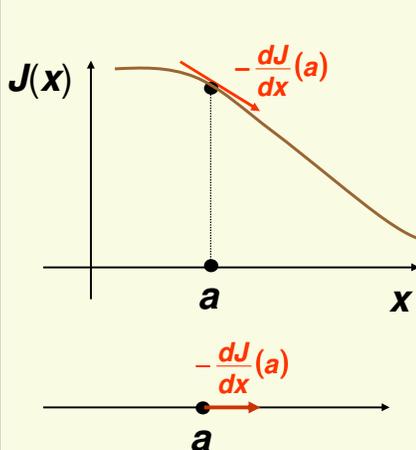
$$\begin{cases} \sin(x_1^2 + x_2^3) + e^{x_2} = 0 \\ \cos(x_1^2 + x_2^3) + \log(x_3^{x_2}) = 0 \end{cases}$$

- Sometimes it is not even possible to write down an analytical expression for the derivative, we will see an example later today

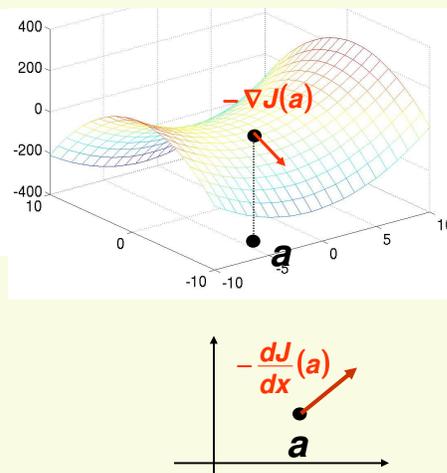
Optimization: Gradient Descent

- Gradient $\nabla J(\mathbf{x})$ points in direction of steepest increase of $J(\mathbf{x})$, and $-\nabla J(\mathbf{x})$ in direction of steepest decrease

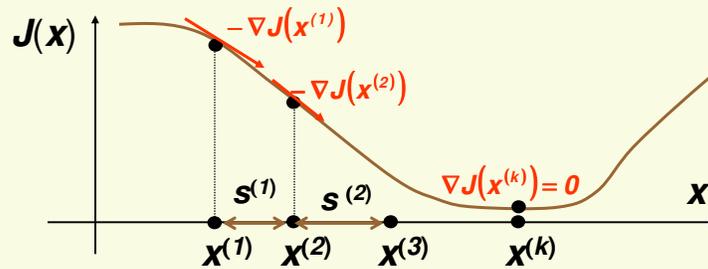
one dimension



two dimensions



Optimization: Gradient Descent



Gradient Descent for minimizing any function $J(x)$

set $k = 1$ and $x^{(1)}$ to some initial guess for the weight vector

while $\eta^{(k)} |\nabla J(x^{(k)})| > \epsilon$

choose **learning rate** $\eta^{(k)}$

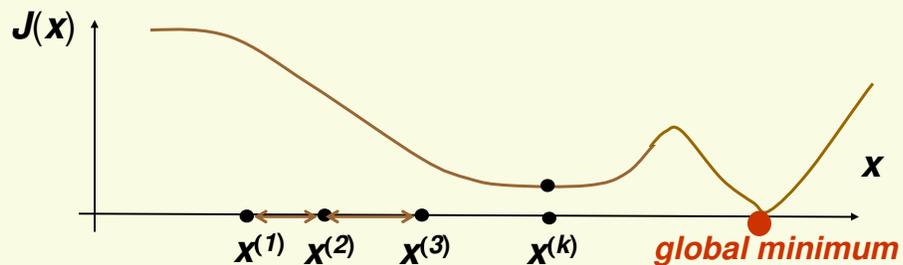
$$x^{(k+1)} = x^{(k)} - \eta^{(k)} \nabla J(x)$$

(update rule)

$$k = k + 1$$

Optimization: Gradient Descent

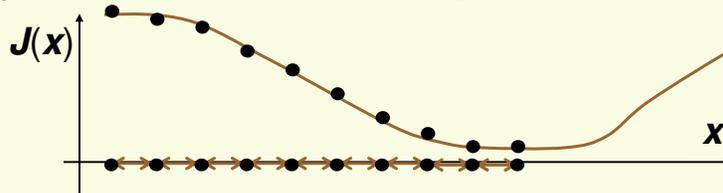
- Gradient descent is guaranteed to find only a local minimum



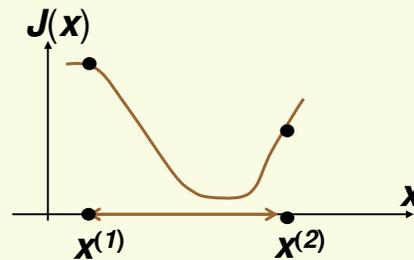
- Nevertheless gradient descent is very popular because it is simple and applicable to any differentiable function

Optimization: Gradient Descent

- Main issue: how to set parameter η (*learning rate*)
- If η is too small, need too many iterations



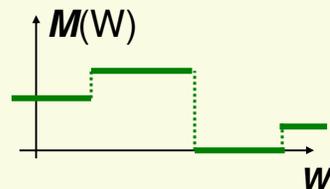
- If η is too large may overshoot the minimum and possibly never find it (if we keep overshooting)



“Optimal” W with Gradient Descent

- $f(X, W) = \text{sign}(w_0 + \sum_{i=1,2,\dots,d} w_i x_i)$
- If we let $L(X^i, Y^i, W) = \mathbf{I}[f(X^i, W) \neq Y^i]$, then $L(W)$ is the number of misclassified examples
- Let M be the set of examples misclassified by W

$$M(W) = \{ \text{sample } X^i \text{ s.t. } W^T X^i \neq Y^i \}$$
- Then $L(W) = |M(W)|$, the size of $M(W)$
- $L(W)$ is piecewise constant, gradient descent is useless

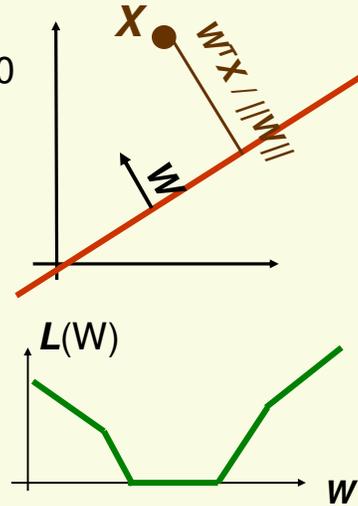


“Optimal” W with Gradient Descent

- Better choice:

$$L(W) = \sum_{X^i \in M} (-W^T X^i) Y^i$$

- If X^i is misclassified, $(W^T X^i) Y^i \leq 0$
- Thus $L(W, X^i, Y^i) \geq 0$
- $L(W, X^i, Y^i)$ is proportional to the distance of misclassified example to the decision boundary
- $L(W) = \sum L(W, X^i, Y^i)$ is piecewise linear and thus suitable for gradient descent



Batch Rule

$$L(W, X^i, Y^i) = \sum_{X \in M} (-W^T X) Y$$

- Gradient of L is $\nabla L(W) = \sum_{X \in M} (-X) Y$
 - M are samples misclassified by W
 - It is not possible to solve $\nabla L(W) = 0$ analytically
- Update rule for gradient descent: $x^{(k+1)} = x^{(k)} - \eta^{(k)} \nabla J(x)$
- Thus *gradient decent batch update rule* for $L(W)$ is:

$$W^{(k+1)} = W^{(k)} + \eta^{(k)} \sum_{Y \in M} X Y$$

- It is called **batch** rule because it is based on all misclassified examples

Single Sample Rule

- Thus *gradient decent single sample rule* for $L(W)$ is:

$$W^{(k+1)} = W^{(k)} + \eta^{(k)}(XY)$$

- apply for any sample X misclassified by $W^{(k)}$
- must have a consistent way of visiting samples

Convergence

- If classes are linearly separable, and $\eta^{(k)}$ is fixed to a constant, i.e. $\eta^{(1)} = \eta^{(2)} = \dots = \eta^{(k)} = c$ (*fixed learning rate*)
 - *both single sample and batch rules converge to a correct solution* (could be any W in the solution space)
- If classes are not linearly separable:
 - Single sample algorithm does not stop, it keeps looking for solution which does not exist
 - However by choosing appropriate learning rate, heuristically stop algorithm at hopefully good stopping point

$$\eta^{(k)} \rightarrow 0 \text{ as } k \rightarrow \infty$$

- for example, $\eta^{(k)} = \frac{\eta^{(1)}}{k}$
- for this learning rate convergence in the linearly separable case can also be proven

Learning by Gradient Descent

- Suppose we suspect that the machine has to have functional form $f(X,W)$, not necessarily linear
- Pick differentiable per-sample loss function $L(X^i, Y^i, W)$
- We need to find W that minimizes $L = \sum_i L(X^i, Y^i, W)$
- Use gradient-based minimization:
 - Batch rule: $W = W - \eta \nabla L(W)$
 - Or single sample rule: $W = W - \eta \nabla L(X^i, Y^i, W)$

Important Questions

- How do we choose the feature vector X ?
- How do we split labeled samples into training/testing sets?
- How do we choose the machine $f(X,W)$?
- How do we choose the loss function $L(X^i, Y^i, W)$?
- How do we find the optimal weights W ?

Information theory

- Information Theory regards information as only those symbols that are uncertain to the receiver
 - **only information essential to understand must be transmitted**
- Shannon made clear that uncertainty is the very commodity of communication
- The amount of information, or uncertainty, output by an information source is a measure of its entropy
- In turn, a source's entropy determines the amount of bits per symbol required to encode the source's information
- Messages are encoded with strings of 0 and 1 (bits)

Information theory

- Suppose we toss a **fair** die with 8 sides
 - need 3 bits to transmit the results of each toss
 - 1000 throws will need 3000 bits to transmit
- Suppose the die is biased
 - side A occurs with probability $1/2$, chances of throwing B are $1/4$, C are $1/8$, D are $1/16$, E are $1/32$, F $1/64$, G and H are $1/128$
 - Encode A = 0, B = 10, C = 110, D = 1110, ..., so on until G = 1111110, H = 1111111
 - We need, on average, $1/2 + 2/4 + 3/8 + 4/16 + 5/32 + 6/64 + 7/128 + 7/128 = 1.984$ bits to encode results of a toss
 - 1000 throws require 1984 bits to transmit
 - Less bits to send = less "information"
 - Biased die tosses contain less "information" than unbiased die tosses (know in advance biased sequence will have a lot of A's)
 - What's the number of bits in the best encoding?
- Extreme case: if a die always shows side A, a sequence of 1,000 tosses has no information, 0 bits to encode

Information theory

- if a die is fair (any side is equally likely, or uniform distribution), for any toss we need $\log(8) = 3$ bits
- Suppose any of n events is equally likely (uniform distribution)
 - $P(x) = 1/n$, therefore $-\log P = -\log(1/n) = \log n$
- In the “good” encoding strategy for our biased die example, every side x has $-\log p(x)$ bits in its code
- Expected number of bits is

$$-\sum_x p(x) \log p(x)$$

Shannon's Entropy

$$H[p(x)] = -\sum_x p(x) \log p(x) = \sum_x p(x) \log \frac{1}{p(x)}$$

- How much randomness (or uncertainty) is there in the value of signal x if it has distribution $p(x)$
 - For uniform distribution (every event is equally likely), $H[x]$ is maximum
 - If $p(x) = 1$ for some event x , then $H[x] = 0$
 - Systems with one very common event have less entropy than systems with many equally probable events
- Gives the expected length of optimal encoding (in binary bits) of a message following distribution $p(x)$
 - doesn't actually give this optimal encoding

Conditional Entropy of X given Y

$$H[x | y] = \sum_{x,y} p(x,y) \log \frac{1}{p(x|y)} = - \sum_{x,y} p(x,y) \log p(x|y)$$

- Measures average uncertainty about x when y is known
- Property:
 - $H[x] \geq H[x|y]$, which means after seeing new data (y), the uncertainty about x is not increased, on average

Mutual Information of X and Y

$$I[x, y] = H(x) - H(x | y)$$

- Measures the average reduction in uncertainty about x after y is known
- or, equivalently, it **measures the amount of information that y conveys about x**
- Properties
 - $I(x,y) = I(y,x)$
 - $I(x,y) \geq 0$
 - If x and y are independent, then $I(x,y) = 0$
 - $I(x,x) = H(x)$

MI for Feature Selection

$$I[x,c] = H(c) - H(c|x)$$

- Let x be a proposed feature and c be the class
- If $I[x,c]$ is high, we can expect feature x be good at predicting class c