# CSE 417T Introduction to Machine Learning

Lecture 8

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## Logistics

- HW1: Due Sep 23
  - Reserve time if you have never used Gradescope
  - Check that submission is readable (if you scan your handwriting)
  - Correctly assign pages to each problem (you won't get points otherwise)
- HW2: Will be announce later today or tomorrow
  - Expect roughly two weeks to work on it
- Exam dates
  - Exam 1: October 27
    - We expect to finish the content for exam1 several lectures before the exam.
  - Exam 2: December 8

# Recap

### Linear Models

This is why it's called linear models

• H contains hypothesis  $h(\vec{x})$  as some function of  $\vec{w}^T\vec{x}$ 

	Domain	Model
Linear Classification	$y \in \{-1, +1\}$	$H = \{h(\vec{x}) = sign(\vec{w}^T \vec{x})\}$
Linear Regression	$y \in \mathbb{R}$	$H = \{h(\vec{x}) = \vec{w}^T \vec{x}\}$
Logistic Regression	$y \in [0,1]$	$H = \{h(\vec{x}) = \theta(\vec{w}^T \vec{x})\}$

#### Credit Card Example

Approve or not

Credit line

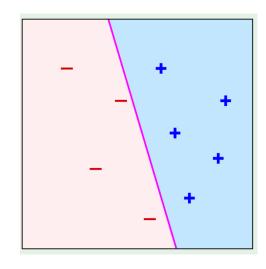
Prob. of default

$$\theta(s) = \frac{e^s}{1 + e^s}$$

- Algorithm:
  - Focus on  $g = argmin_{h \in H} E_{in}(h)$

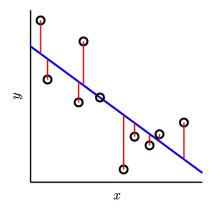
# Linear Classification (Perceptron)

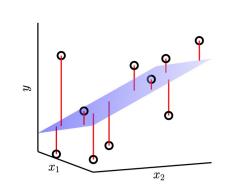
- Formulation
  - Hypothesis set  $H = \{h(\vec{x}) = sign(\vec{w}^T\vec{x})\}$
  - Error measure: binary error  $e(h(\vec{x}), y) = \mathbb{I}[h(\vec{x}) \neq y]$
- Data is linearly separable
  - Run PLA =>  $E_{in} = 0$  => Low  $E_{out}$
- Data is not linearly separable
  - Minimizing  $E_{in}$  is NP hard
  - Pocket algorithm
  - Engineering the features (e.g., handwritten digits)
  - More discussion later in the semester



## Linear Regression

- Formulation
  - Hypothesis set  $H = \{h(\vec{x}) = \vec{w}^T \vec{x}\}$
  - Squared error  $e(h(\vec{x}), y) = (h(\vec{x}) y)^2$





- Given dataset  $D = \{(\vec{x}_1, y_1), ..., (\vec{x}_N, y_N)\}$ 
  - $E_{in}(\vec{w}) = \frac{1}{N} \sum_{n=1}^{N} (\vec{w}^T \vec{x}_n y_n)^2$
- Goal: find  $\overrightarrow{w}_{lin} = argmin_{\overrightarrow{w}} E_{in}(\overrightarrow{w})$

# Linear Regression "Algorithm"

- There is a closed-form solution for minimizing  $E_{in}$ 
  - Closed-form solution for  $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}) = 0$ 
    - ( $E_{in}$  is convex; you can check the second derivate of  $E_{in}$ )
- One-step algorithm
  - Given  $D = \{(\vec{x}_1, y_1), ..., (\vec{x}_N, y_N)\}$

• Construct 
$$X = \begin{bmatrix} \vec{x}_1^T \\ \vdots \\ \vec{x}_N^T \end{bmatrix} = \begin{bmatrix} x_{1,0} & x_{1,1} & \cdots & x_{1,d} \\ x_{2,0} & x_{2,1} & \cdots & x_{2,d} \\ \vdots & \vdots & \ddots & \vdots \\ x_{2,0} & x_{N,1} & \cdots & x_{N,d} \end{bmatrix}$$
 and  $\vec{y} = \begin{bmatrix} y_1 \\ y_2 \\ \vdots \\ y_n \end{bmatrix}$ 

• Output  $\overrightarrow{w}_{lin} = (X^T X)^{-1} X^T \overrightarrow{y}$  (Assume  $X^T X$  is invertible)

# Today's Lecture

The notes are not intended to be comprehensive. They should be accompanied by lectures and/or textbook. Let me know if you spot errors.

# Logistic Regression

	Domain	Model
Linear Classification	$y \in \{-1, +1\}$	$H = \{h(\vec{x}) = sign(\vec{w}^T \vec{x})\}\$
Linear Regression	$y \in \mathbb{R}$	$H = \{h(\vec{x}) = \vec{w}^T \vec{x}\}$
Logistic Regression	$y \in [0,1]$	$H = \{h(\vec{x}) = \theta(\vec{w}^T \vec{x})\}\$

# Logistic Regression: Predicting a Probability

Will this patient have a heart attack within the next year?

age	62 years
gender	male
blood sugar	120 mg/dL40,000
HDL	50
LDL	120
Mass	190 lbs
Height	5' 10"
	• • •

Classification: Yes/No

Logistic regression: Probability of Yes

- A hypothesis  $h(\vec{x})$  outputs a value in [0,1]
  - Interpreting it as the probability of yes

# Logistic Regression: Predicting a Probability

- Hypothesis set  $H = \{h(\vec{x}) = \theta(\vec{w}^T\vec{x})\}$ 
  - Want  $\theta$  to map from  $(-\infty, \infty)$  to [0,1]

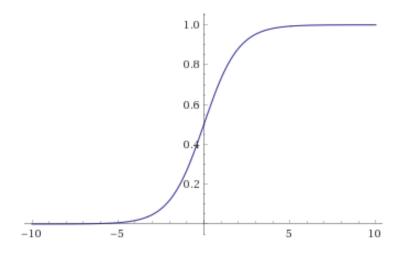
• 
$$\theta(s) = \frac{e^s}{1+e^s} = \frac{1}{1+e^{-s}}$$

A sigmoid function ("S"-shaped function)

• 
$$\theta(s) = \begin{cases} 1 & \text{when } s \to \infty \\ 0.5 & \text{when } s = 0 \\ 0 & \text{when } s \to -\infty \end{cases}$$

Useful property

• 
$$1 - \theta(s) = \frac{1 + e^s}{1 + e^s} - \frac{e^s}{1 + e^s} = \frac{1}{1 + e^s} = \theta(-s)$$



#### What Kind of Dataset do We Get?

• Dataset  $D = \{(\vec{x}_1, y_1), ..., (\vec{x}_N, y_N)\}$ 

age	62 years
gender	male
blood sugar	120 mg/dL40,000
HDL	50
LDL	120
Mass	190 lbs
Height	5' 10"
	• • •

- What are the values of  $y_n$ ?
  - Ideally, we want to have  $y_n$  to be the probability value
  - In practice, we cannot measure a probability
  - We can only see the occurrence of an event and infer the probability
  - (We often only observe whether the person had heart attack, we don't observe the "probability")
- Need to address the case when  $y_n \in \{-1, +1\}$  in the given dataset D

# Error Measure: Quantifying $g \approx f$

• Target function  $f(\vec{x}) = \Pr(y = +1|\vec{x})$ 

Side note:

You probably can guess why the property  $1 - \theta(s) = \theta(-s)$  might be helpful

- Another way to write it:  $\Pr(y|\vec{x}) = \begin{cases} f(\vec{x}) & \text{for } y = +1 \\ 1 f(\vec{x}) & \text{for } y = -1 \end{cases}$
- How do we define the error measure to quantify  $g \approx f$ 
  - Ideally, we want it to be meaningful
    - Binary error for classification: tell us the number of mistakes we make
    - Squared error for regression: the error minimizer is the "mean (average)"
  - We also want it to be easy to optimize

## Cross Entropy Error

$$E_{in}(\vec{w}) = \frac{1}{N} \sum_{n=1}^{N} \ln(1 + e^{-y_n \vec{w}^T \vec{x}_n})$$

- It looks complicated, but
  - It has nice interpretations (min error => max likelihood)
  - It is easy to optimize (continuous, differentiable, convex)

# Minimizing Cross Entropy Error



Maximizing Likelihood

#### Maximum Likelihood Estimation

- Likelihood Pr(D|h)
  - The probability of seeing dataset D if D is generated according to h (i.e., if h is the target function)
  - $Pr(D|h) = Pr((\vec{x}_1, y_1), ..., (\vec{x}_N, y_N)|h)$
  - Maximum likelihood estimation (MLE)
    - $g = argmax_{h \in H} Pr(D|h)$
- Sidenote: Two different concepts in ML
  - Likelihood: Pr(D|h) [Focus of this course]
  - Posterior: Pr(h|D) [Focus of Bayesian machine learning: More in 515T]
  - Connection:  $Pr(h|D) = \frac{Pr(h)Pr(D|h)}{Pr(D)}$ 
    - Prior Pr(h): the additional assumption Bayesian ML makes

#### Write Down the Likelihood

- How are  $D = \{(\vec{x}_1, y_1), ..., (\vec{x}_N, y_N)\}$  generated?
  - $(\vec{x}_1, ..., \vec{x}_N)$  are i.i.d. drawn from a distribution
  - $(y_1, ..., y_N)$  are labeled according to target function  $f(\vec{x})$
- UNKNOWN TARGET DISTRIBUTION  $P(y \mid \mathbf{x})$ UNKNOWN
  INPUT DISTRIBUTION  $P(\mathbf{x})$ TRAINING EXAMPLES  $(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_N, y_N)$ ERROR
  MEASURE  $\mathbf{x}_1, \mathbf{y}_2, \dots, \mathbf{y}_N$ HYPOTHESIS SET  $\mathcal{H}$

- Likelihood Pr(D|h)
  - The probability of seeing dataset D if D is generated according to h

• 
$$\Pr(D|h) = \Pr((\vec{x}_1, y_1), ..., (\vec{x}_N, y_N)|h)$$
  
 $= \Pr(\vec{x}_1, ..., \vec{x}_N) \Pr((y_1, ..., y_N)|(\vec{x}_1, ..., \vec{x}_N), h)$   
 $= \prod_{n=1}^{N} \Pr(\vec{x}_n) \prod_{n=1}^{N} \Pr(y_n|\vec{x}_n, h)$  (Assumption of independent data)

#### Maximum Likelihood

Choosing the hypothesis that maximizes the likelihood

```
• g = argmax_{h \in H} \Pr(D|h)

= argmax_{h \in H} \prod_{n=1}^{N} \Pr(\vec{x}_n) \prod_{n=1}^{N} \Pr(y_n|\vec{x}_n, h)

= argmax_{h \in H} \prod_{n=1}^{N} \Pr(y_n|\vec{x}_n, h)
```

 $\prod_{n=1}^{N} \Pr(\vec{x}_n)$  doesn't depend on h

• We interpret  $h(\vec{x})$  as the probability of y=+1

• 
$$\Pr(y|\vec{x},h) = \begin{cases} h(\vec{x}) = \theta(\vec{w}^T \vec{x}) & \text{for } y = +1 \\ 1 - h(\vec{x}) = 1 - \theta(\vec{w}^T \vec{x}) & \text{for } y = -1 \end{cases}$$

- Since  $1 \theta(s) = \theta(-s)$ 
  - $Pr(y|\vec{x}, h) = \theta(y \vec{w}^T \vec{x})$

#### Maximum Likelihood

Choosing the hypothesis that maximizes the likelihood

```
• g = argmax_{h \in H} \Pr(D|h)
= argmax_{h \in H} \prod_{n=1}^{N} \Pr(y_n | \vec{x}_n, h)
```

• 
$$\overrightarrow{w}^* = argmax_{\overrightarrow{w}} \prod_{n=1}^{N} \theta(y_n \overrightarrow{w}^T \overrightarrow{x}_n)$$
  

$$= argmax_{\overrightarrow{w}} \ln(\prod_{n=1}^{N} \theta(y_n \overrightarrow{w}^T \overrightarrow{x}_n))$$
  

$$= argmax_{\overrightarrow{w}} \sum_{n=1}^{N} \ln(\theta(y_n \overrightarrow{w}^T \overrightarrow{x}_n))$$
  

$$= argmin_{\overrightarrow{w}} - \sum_{n=1}^{N} \ln(\theta(y_n \overrightarrow{w}^T \overrightarrow{x}_n))$$
  

$$= argmin_{\overrightarrow{w}} \sum_{n=1}^{N} \ln \frac{1}{\theta(y_n \overrightarrow{w}^T \overrightarrow{x}_n)}$$
  

$$= argmin_{\overrightarrow{w}} \sum_{n=1}^{N} \ln(1 + e^{-y_n \overrightarrow{w}^T \overrightarrow{x}_n})$$
  

$$= argmin_{\overrightarrow{w}} \sum_{n=1}^{N} \ln(1 + e^{-y_n \overrightarrow{w}^T \overrightarrow{x}_n})$$

$$\theta(s) = \frac{e^s}{1 + e^s} = \frac{1}{1 + e^{-s}}$$

$$E_{in}(\vec{w}) = \frac{1}{N} \sum_{n=1}^{N} \ln(1 + e^{-y_n \vec{w}^T \vec{x}_n})$$

## Cross Entropy Error

$$E_{in}(\vec{w}) = \frac{1}{N} \sum_{n=1}^{N} \ln(1 + e^{-y_n \vec{w}^T \vec{x}_n})$$

- Minimizing  $E_{in}(\vec{w})$  is the same as maximizing likelihood
- Next question: How to solve  $\vec{w}^* = argmin_{\vec{w}} E_{in}(\vec{w})$ 
  - Answer: Solve for  $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}) = 0$
  - No single-step solution like we have in linear regression

# Using Logistic Regression for Classification

• Let  $\overrightarrow{w}^*$  or g be the learned logistic regression model, how can we make classification predictions using  $\overrightarrow{w}^*$ ?

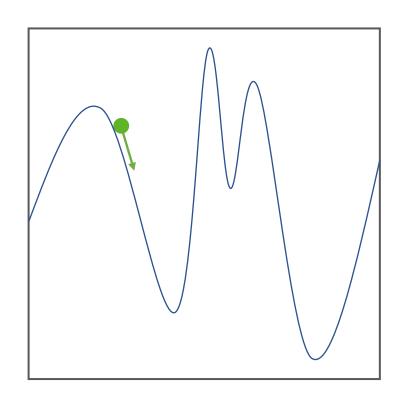
- Set a cutoff probability C% (e.g., 50%).
  - Classify +1 if  $g(\vec{x}) = \theta(\vec{w}^* \vec{x}) > C\%$
  - Classify -1 if  $g(\vec{x}) = \theta(\vec{w}^* \vec{x}) < C\%$
- When C is 50 (a common choice)
  - $\theta(\vec{w}^{*T}\vec{x}) > 50\% = \vec{w}^{*T}\vec{x} > 0$
  - Equivalent to using  $\vec{w}^*$  as the linear classification hypothesis, i.e.,  $g(\vec{x}) = sign(\vec{w}^{*T}\vec{x})$

# Gradient Descent

A general optimization technique

#### **Gradient Descent**

• A technique for optimizing functions that gradients exist everywhere.



• An iterative method that converges to local optimum.

 Converge to global optimum if the function is convex (since there is only one local optimum).

# Gradient Descent: Minimizing $E_{in}(\vec{w})$

An iterative method of the form:

$$\vec{w}(t+1) \leftarrow \vec{w}(t) + \eta_t \vec{v}_t$$

- $\vec{v}_t$ : a unit vector, determining the direction of the update
- $\eta_t$ : a scalar, determining how much to update
- How to choose  $\vec{v}_t$  and  $\eta_t$ ?

Choosing 
$$\vec{v}_t$$
 in  $\vec{w}(t+1) \leftarrow \vec{w}(t) + \eta_t \vec{v}_t$ 

- Intuition: Choose  $\vec{v}_t$  that moves towards the "steepest" direction
  - Approaching the minimum faster
- Taylor's approximation:

• 
$$E_{in}(\overrightarrow{w}(t) + \eta_t \overrightarrow{v}_t) = E_{in}(\overrightarrow{w}(t)) + \eta_t \nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}(t))^T \overrightarrow{v}_t + O(\eta_t^2)$$
  
•  $E_{in}(\overrightarrow{w}(t+1)) - E_{in}(\overrightarrow{w}(t)) \approx \eta_t \nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}(t))^T \overrightarrow{v}_t$ 

• 
$$E_{in}(\vec{w}(t+1)) - E_{in}(\vec{w}(t)) \approx \eta_t \nabla_{\vec{w}} E_{in}(\vec{w}(t))^T \vec{v}_t$$

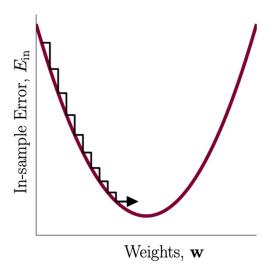
 $\eta_t$  is usually small, so ignore this term

- Choose unit vector  $\vec{v}_t$  that minimizes  $\nabla_{\vec{w}} E_{in}(\vec{w}(t))^T \vec{v}_t$ 
  - $\vec{v}_t$  should be in the opposite direction of  $\nabla_{\vec{w}} E_{in}(\vec{w}(t))$

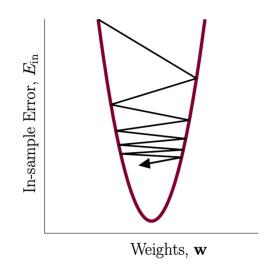
• 
$$\vec{v}_t = \frac{-\nabla_{\vec{w}} E_{in}(\vec{w}(t))}{\|\nabla_{\vec{w}} E_{in}(\vec{w}(t))\|}$$

# Choosing $\eta_t$ in $\vec{w}(t+1) \leftarrow \vec{w}(t) + \eta_t \vec{v}_t$

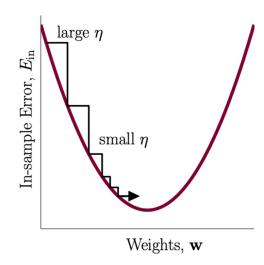
 $\eta$  too small



 $\eta$  too large

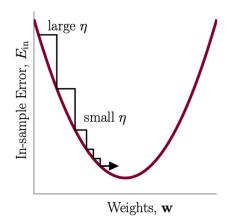


variable  $\eta_t$  – just right



Choosing 
$$\eta_t$$
 in  $\vec{w}(t+1) \leftarrow \vec{w}(t) + \eta_t \vec{v}_t$ 

- Intuition (for convex  $E_{in}$ )
  - When  $E_{in}$  is closer to the minimum,
    - $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}(t))$  is smaller
    - We should set  $\eta_t$  smaller



• Therefore, set  $\eta_t = \eta \|\nabla_{\vec{w}} E_{in}(\vec{w}(t))\|$ 

# Putting Them Together

• Iterative update rule:  $\vec{w}(t+1) \leftarrow \vec{w}(t) + \eta_t \vec{v}_t$ 

• 
$$\vec{w}(t+1) \leftarrow \vec{w}(t) - \eta \nabla_{\vec{w}} E_{in}(\vec{w}(t))$$

$$\vec{v}_t = \frac{-\nabla_{\vec{w}} E_{in}(\vec{w}(t))}{\|\nabla_{\vec{w}} E_{in}(\vec{w}(t))\|}$$

$$\eta_t = \eta \| \nabla_{\vec{w}} E_{in}(\vec{w}(t)) \|$$

Gradient calculations

• 
$$E_{in}(\vec{w}) = \frac{1}{N} \sum_{n=1}^{N} \ln(1 + e^{-y_n \vec{w}^T \vec{x}_n})$$

• 
$$\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}) = \frac{1}{N} \sum_{n=1}^{N} \frac{-y_n \overrightarrow{x} e^{-y_n \overrightarrow{w}^T \overrightarrow{x}_n}}{1 + e^{-y_n \overrightarrow{w}^T \overrightarrow{x}_n}} = -\frac{1}{N} \sum_{n=1}^{N} \frac{y_n \overrightarrow{x}_n}{1 + e^{y_n \overrightarrow{w}^T \overrightarrow{x}_n}}$$

## Gradient Descent for Logistic Regression

- Initialize  $\vec{w}(0)$
- For t = 0, ...
  - Compute  $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}(t)) = -\frac{1}{N} \sum_{n=1}^{N} \frac{y_n \overrightarrow{x}_n}{1 + e^{y_n \overrightarrow{w}(t)} \overrightarrow{T} \overrightarrow{x}_n}$
  - $\vec{w}(t+1) \leftarrow \vec{w}(t) \eta \nabla_{\vec{w}} E_{in}(\vec{w}(t))$
  - Terminate if the stop conditions are met
- Return the final weights

 $\eta$ : learning rate. A parameter the learner can choose.

# Gradient Descent for Logistic Regression

- Initialization
  - Random initialization is a good idea and a common approach
  - (we specify the initialization in HW2 mostly for grading purposes)
- Stop conditions (a mix of the following criteria)
  - When the number of iteration exceeds the pre-set threshold
  - When the improvement on  $E_{in}$  (e.g., check  $\nabla_{\overrightarrow{w}}E_{in}$ ) is too small
  - When  $E_{in}$  is small enough

## Computation of Gradient Descent

- Gradient Descent for Logistic Regression
  - Initialize  $\vec{w}(0)$
  - For t = 0, ...
    - Compute  $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}(t)) = -\frac{1}{N} \sum_{n=1}^{N} \frac{y_n \overrightarrow{x}_n}{1 + e^{y_n \overrightarrow{w}(t)^T \overrightarrow{x}_n}}$
    - $\vec{w}(t+1) \leftarrow \vec{w}(t) \eta \nabla_{\vec{w}} E_{in}(\vec{w}(t))$
    - Terminate if the stop conditions are met
  - Return the final weights
- Which step requires the most computation?
  - Calculate  $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}) = -\frac{1}{N} \sum_{n=1}^{N} \frac{y_n \vec{x}_n}{1 + e^{y_n \overrightarrow{w}^T \vec{x}_n}}$
  - The time complexity is O(N)
    - *N* is large for big datasets

# Stochastic Gradient Descent

# Deal with Heavy Computation of $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w})$

- Speed up the implementation of  $\nabla_{\vec{w}} E_{in}(\vec{w}) = -\frac{1}{N} \sum_{n=1}^{N} \frac{y_n \vec{x}_n}{1 + e^{y_n \vec{w}^T \vec{x}_n}}$ 
  - Vectorization can make your HW2 running time in several order of magnitudes faster
  - Example:
    - Given  $[x_1, ..., x_N]$ , want to calculate  $[e^{x_1}, ..., e^{x_N}]$
    - Using for loop:
      - Loop from n=1 to N, calculate  $e^{x_n}$
    - Vectorized method:
      - Using numpy library: np.exp( $[x_1, ..., x_N]$ )
  - Why? Matrix operations are optimized in a low level using numpy operations (or other scientific computing libraries).
    - Try to replace loops with numpy matrix operations in your HW2

# Deal with Heavy Computation of $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w})$

- Speed up the implementation of  $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}) = -\frac{1}{N} \sum_{n=1}^{N} \frac{y_n \vec{x}_n}{1 + e^{y_n \overrightarrow{w}} T_{\overrightarrow{x}_n}}$ 
  - Vectorization can make your HW2 running time in several order of magnitudes faster
- Solve  $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w})$  "in expectation"
  - Define  $e_n(\vec{w}) = \ln(1 + e^{-y_n \vec{w}^T \vec{x}_n})$ , the point-wise error caused by  $(\vec{x}_n, y_n)$
  - Observe that
    - $E_{in}(\overrightarrow{w}) = \frac{1}{N} \sum_{n=1}^{N} e_n(\overrightarrow{w})$
    - $\nabla_{\overrightarrow{w}} E_{in}(\overrightarrow{w}) = \frac{1}{N} \sum_{n=1}^{N} \nabla_{\overrightarrow{w}} e_n(\overrightarrow{w})$
    - Draw a point  $\vec{x}_n$  from  $\{\vec{x}_1, \dots, \vec{x}_N\}$  uniformly at random
      - $E_{\vec{x}_n}[\nabla_{\vec{w}}e_n(\vec{w})] = \nabla_{\vec{w}}E_{in}(\vec{w})$

# Stochastic Gradient Descent (SGD)

- Algorithm
  - Initialize  $\vec{w}(0)$
  - For t = 0, ...
    - Randomly choose n from  $\{1, ..., N\}$
    - $\vec{w}(t+1) \leftarrow \vec{w}(t) \eta \nabla_{\vec{w}} e_n(\vec{w}(t))$
    - Terminate if the stop conditions are met
  - Return the final weights
- $\mathbb{E}[\nabla_{\overrightarrow{w}}e_n(\overrightarrow{w})] = \nabla_{\overrightarrow{w}}E_{in}(\overrightarrow{w})$ 
  - SGD is doing the same thing as GD in expectation
    - More efficient (scale to large dataset), suitable for online data, helps escaping local min, etc.
    - Noisier, harder to define stop criteria

#### Mini-Batch Gradient Descent

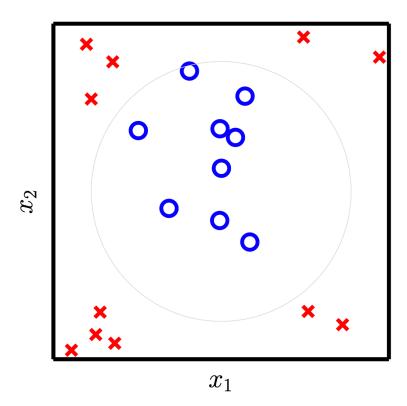
- GD: Computationally heavy, stable updates
- SGD: Computationally light, noisy updates
- Middle ground: Mini-Batch Gradient Descent
  - In each iteration, randomly choose k points  $\{n(1), ..., n(k)\}$
  - Update rule

• 
$$\overrightarrow{w}(t+1) \leftarrow \overrightarrow{w}(t) - \eta \frac{1}{k} \sum_{i=1}^{k} \nabla_{\overrightarrow{w}} e_{n(i)}(\overrightarrow{w}(t))$$

- Side note about HW2
  - Please report your results on GD (non-stochastic version)
    - You should feel free to play around with SGD or mini-batch on your own

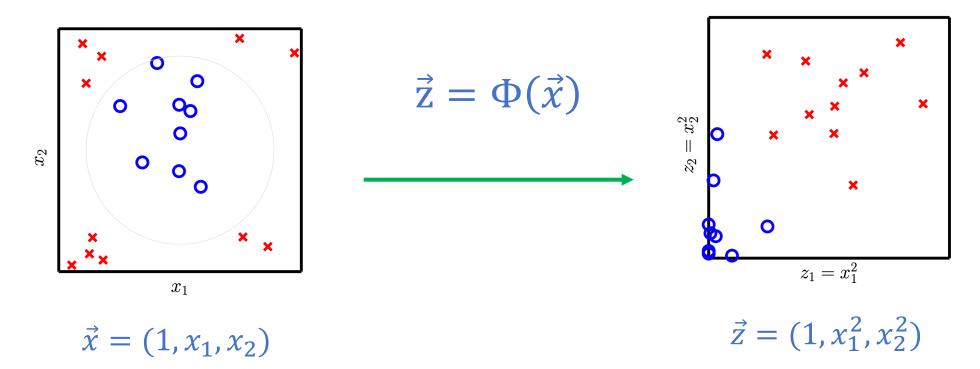
# Non-Linear Transformation

## Limitations of Linear Models



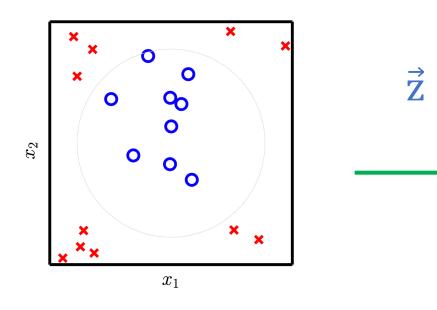
# Using Non-Linear Transformations

• Find a feature transform  $\Phi$  that maps data from  $\vec{x}$  space to  $\vec{z}$  space

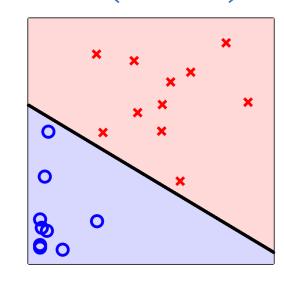


# Using Non-Linear Transformations

• Learn a linear classifier in  $\vec{z}$  space:  $g^{(z)}(\vec{z}) = sign(\vec{w}^{(z)}\vec{z})$ 



$$\vec{x} = (1, x_1, x_2)$$



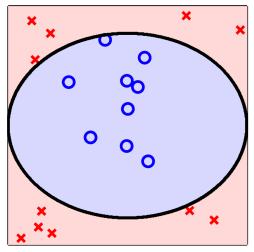
$$\vec{z} = (1, x_1^2, x_2^2)$$

$$g^{(z)}(\vec{z}) = sign(-0.6 + z_1 + z_2)$$

# Using Non-Linear Transformations

• Transform the learned hypothesis back to  $\vec{x}$  space

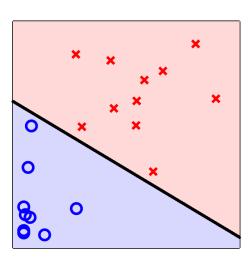
• 
$$g(\vec{x}) = g^{(z)}(\Phi(\vec{x})) = sign(\vec{w}^{(z)}\Phi(\vec{x}))$$



$$\vec{x} = (1, x_1, x_2)$$



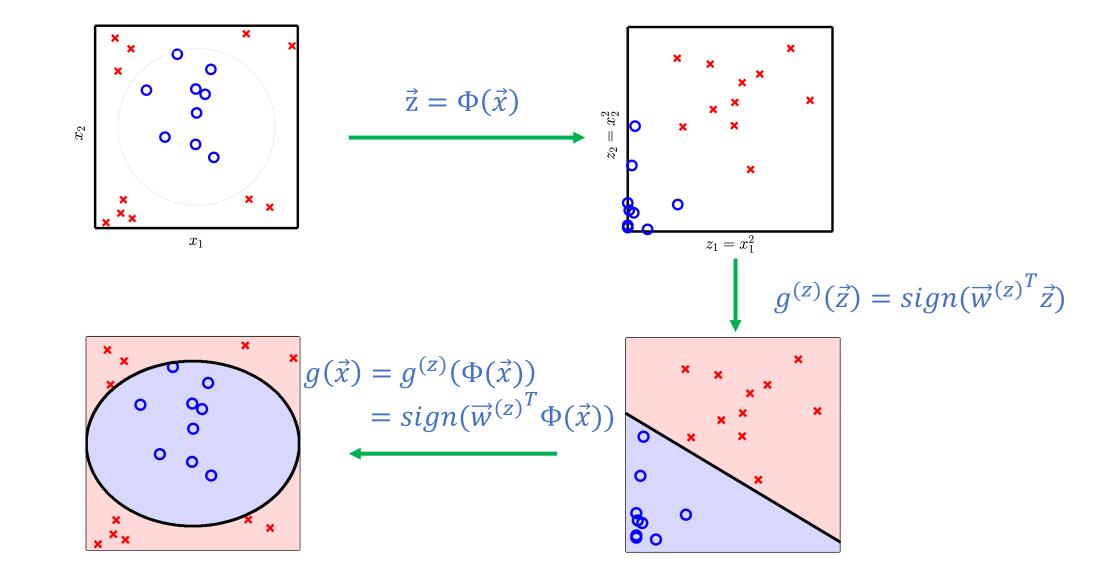
$$g(\vec{x}) = sign(-0.6 + x_1^2 + x_2^2)$$



$$\vec{z} = (1, x_1^2, x_2^2)$$

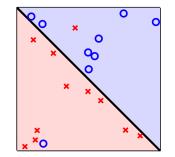
$$g^{(z)}(\vec{z}) = sign(-0.6 + z_1 + z_2)$$

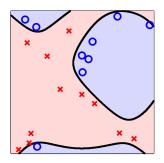
## Nonlinear Transformation



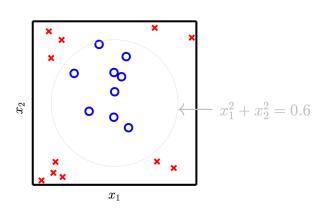
#### Generalization of Nonlinear Transformation

- Fact (We'll prove this later)
  - The VC Dimension of d-dim perceptron is d+1
- VC dimension of perceptron on input space  $\vec{x} = (x_0, ..., x_d)$ 
  - d+1
- VC dimension of perceptron on input space  $\vec{z} = (z_0, ..., z_{d'})$ 
  - $\leq d' + 1$  (usually treated as  $\approx d' + 1$ )
- Careful: Non-linear transform might lead to "nonsense" behavior

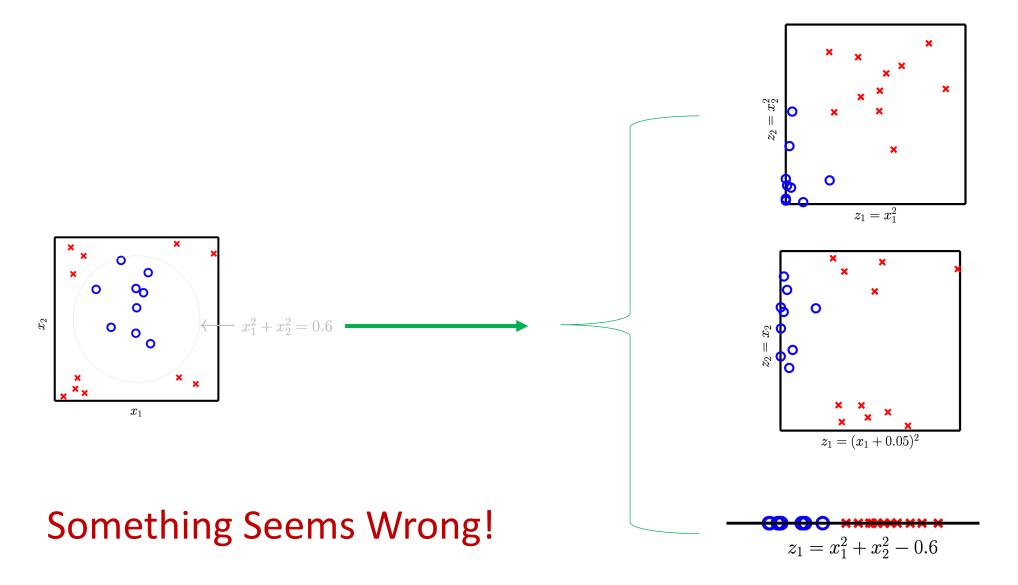




## How to Choose Feature Transform $\Phi$



### How to Choose Feature Transform $\Phi$



# Must choose Φ BEFORE looking at the data

Otherwise, you are doing "data snooping"

The hypothesis set H is as large as anything your brain can think of

# Choose Φ Before Seeing Data

- Rely on domain knowledge (feature engineering)
  - Handwriting digit recognition example
- Use common sets of feature transformation
  - Polynomial transformation
  - 2nd order Polynomial transformation
    - $\vec{x} = (1, x_1, x_2)$
    - $\Phi_2(\vec{x}) = (1, x_1, x_2, x_1 x_2, x_1^2, x_2^2)$
    - Pros: more powerful (contains circle, ellipse, hyperbola, etc)
    - Cons: 2-d => 5-d
      - More computation/storage
      - Worse generalization error

The VC dimension of d-dim perceptron is d+1