

CSCI567 Machine Learning (Fall 2025)

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Sep 12, 2025

Administration

- HW 1 is due on Wed, Sep 17th.

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- recall the late day policy: 3 in total, at most 1 for each homework

Outline

- 1 Review of Last Lecture
- 2 Linear Classifiers and Surrogate Losses
- 3 A Detour of Numerical Optimization Methods
- 4 Perceptron
- 5 Logistic Regression

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Regression

Predicting a continuous outcome variable using past observations

- temperature, amount of rainfall, house price, etc.

Key difference from classification

- continuous vs discrete
- measure *prediction errors* differently.
- lead to quite different learning algorithms.

Linear Regression: regression with linear models: $f(x) = w^T x$

Least square solution

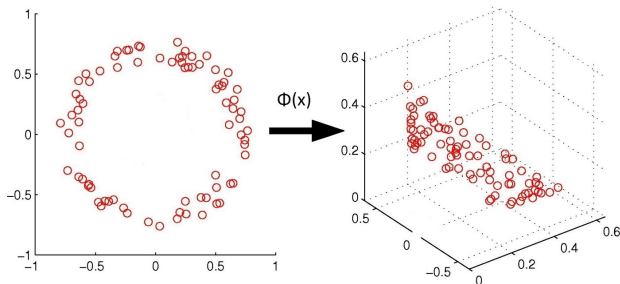
$$\begin{aligned}
 \mathbf{w}^* &= \underset{\mathbf{w}}{\operatorname{argmin}} \operatorname{RSS}(\mathbf{w}) \\
 &= \underset{\mathbf{w}}{\operatorname{argmin}} \|\mathbf{X}\mathbf{w} - \mathbf{y}\|_2^2 \\
 &= (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{y}
 \end{aligned}$$

$$\mathbf{X} = \begin{pmatrix} \mathbf{x}_1^T \\ \mathbf{x}_2^T \\ \vdots \\ \mathbf{x}_N^T \end{pmatrix}, \quad \mathbf{y} = \begin{pmatrix} y_1 \\ y_2 \\ \vdots \\ y_N \end{pmatrix}$$

Two approaches to find the minimum:

- find **stationary points** by setting gradient = 0
- “**complete the square**”

Regression with nonlinear basis



Model: $f(x) = w^T \phi(x)$ where $w \in \mathbb{R}^M$

Similar least square solution: $w^* = (\Phi^T \Phi)^{-1} \Phi^T y$

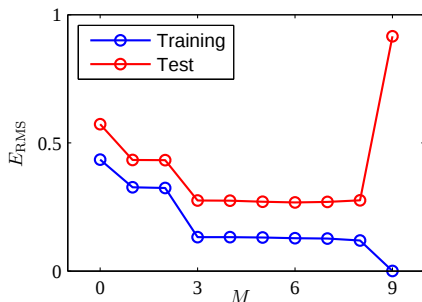
Underfitting and Overfitting

$M \leq 2$ is *underfitting* the data

- large training error
- large test error

$M \geq 9$ is *overfitting* the data

- small training error
- **large test error**



How to prevent overfitting? more data + regularization

$$\mathbf{w}^* = \underset{\mathbf{w}}{\operatorname{argmin}} \left(\text{RSS}(\mathbf{w}) + \lambda \|\mathbf{w}\|_2^2 \right) = (\Phi^T \Phi + \lambda \mathbf{I})^{-1} \Phi^T \mathbf{y}$$

General idea to derive ML algorithms

Step 1. Pick a set of **models** $\mathcal{F}$

- e.g. $\mathcal{F} = \{f(\mathbf{x}) = \mathbf{w}^T \mathbf{x} \mid \mathbf{w} \in \mathbb{R}^D\}$
- e.g. $\mathcal{F} = \{f(\mathbf{x}) = \mathbf{w}^T \Phi(\mathbf{x}) \mid \mathbf{w} \in \mathbb{R}^M\}$

Step 2. Define **error/loss** $L(y', y)$

Step 3. Find **(regularized) empirical risk minimizer (ERM)**:

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Today: another exercise of this recipe + a closer look at Step 3

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Classification

Recall the setup:

- input (feature vector): $\mathbf{x} \in \mathbb{R}^D$
- output (label): $y \in [C] = \{1, 2, \dots, C\}$
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This lecture: **binary classification**

- Number of classes: $C = 2$
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We have discussed **nearest neighbor classifier**:

- require carrying the training set
- intuitive but more like a heuristic

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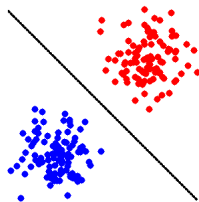
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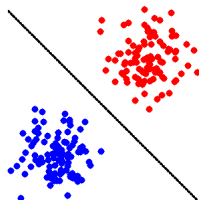
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Sign of $w^T x$ predicts the label:

$$\text{sign}(w^T x) = \begin{cases} +1 & \text{if } w^T x > 0 \\ -1 & \text{if } w^T x \leq 0 \end{cases}$$

(Sometimes use sgn for sign too.)



The models

The set of **(separating) hyperplanes**:

$$\mathcal{F} = \{f(\mathbf{x}) = \text{sgn}(\mathbf{w}^T \mathbf{x}) \mid \mathbf{w} \in \mathbb{R}^D\}$$

The models

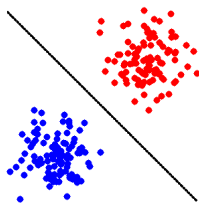
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Good choice for *linearly separable* data, i.e., $\exists w$ s.t.

$$\text{sgn}(w^T x_n) = y_n$$

for all $n \in [N]$.



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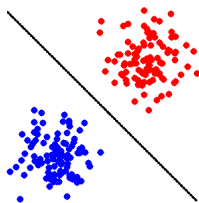
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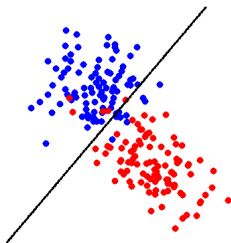
$$\text{sgn}(w^T x_n) = y_n \quad \text{or} \quad y_n w^T x_n > 0$$

for all $n \in [N]$.



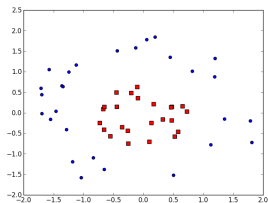
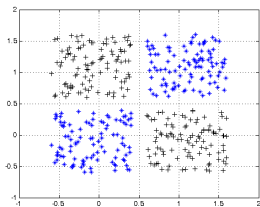
The models

Still makes sense for “almost” linearly separable data



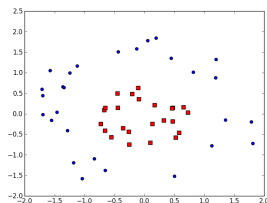
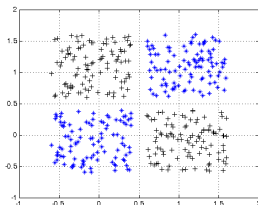
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Again can apply a **nonlinear mapping** Φ :

$$\mathcal{F} = \{f(x) = \text{sgn}(w^T \Phi(x)) \mid w \in \mathbb{R}^M\}$$

More discussions in future lectures.

0-1 Loss

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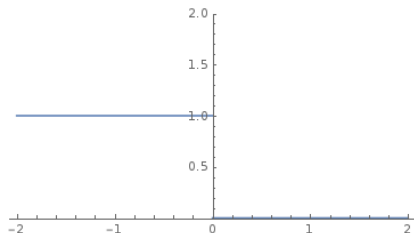
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For classification, more convenient to look at the loss **as a function of** $yw^T x$. That is, with

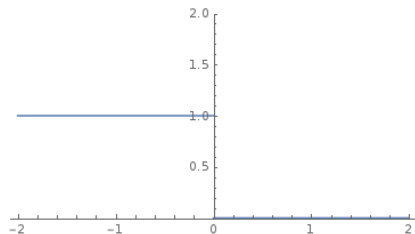
$$\ell_{0-1}(z) = \mathbb{I}[z \leq 0]$$



the loss for hyperplane w on example (x, y) is $\ell_{0-1}(yw^T x)$

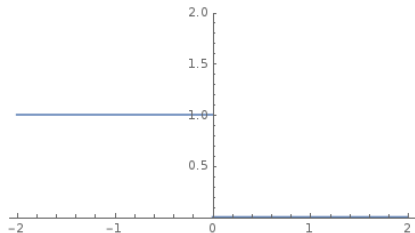
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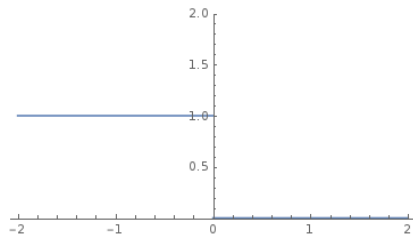
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Even worse, minimizing 0-1 loss is *NP-hard in general*.

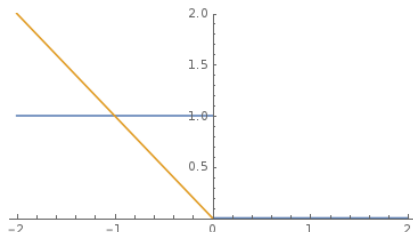
Surrogate Losses

Solution: find a **convex surrogate loss**



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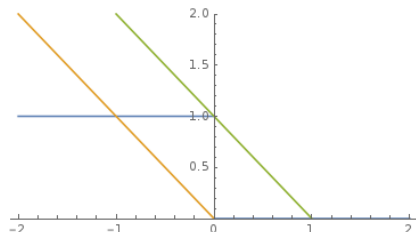
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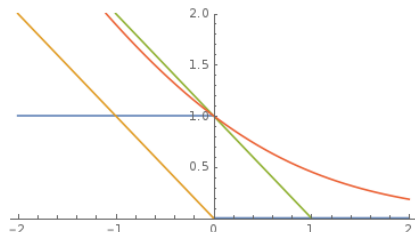
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- **logistic loss** $\ell_{\text{logistic}}(z) = \log(1 + \exp(-z))$ (used in logistic regression; the base of $\log$ doesn't matter)

ML becomes convex optimization

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where $\ell(\cdot)$ can be perceptron/hinge/logistic loss

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Note: minimizing perceptron loss *does not really make sense* (try $\mathbf{w} = \mathbf{0}$), but the algorithm derived from this perspective does.

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 - First-order methods
 - Second-order methods
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Numerical optimization

Problem setup

- Given: a function $F(\boldsymbol{w})$
- Goal: minimize $F(\boldsymbol{w})$ (approximately)

First-order optimization methods

Two simple yet extremely popular methods

- **Gradient Descent (GD)**: simple and fundamental
- **Stochastic Gradient Descent (SGD)**: faster, effective for large-scale problems

First-order optimization methods

Two simple yet extremely popular methods

- **Gradient Descent (GD)**: simple and fundamental
- **Stochastic Gradient Descent (SGD)**: faster, effective for large-scale problems

Gradient is sometimes referred to as *first-order* information of a function. Therefore, these methods are called *first-order methods*.

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GD: keep moving in the *negative gradient direction*

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Start from some (random) $\mathbf{w}^{(0)}$. For $t = 0, 1, 2, \dots$

$$\mathbf{w}^{(t+1)} \leftarrow \mathbf{w}^{(t)} - \eta \nabla F(\mathbf{w}^{(t)})$$

where $\eta > 0$ is called step size or learning rate

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- in practice we often try different small values

Stop when $F(\mathbf{w}^{(t)})$ **does not change much** or t **reaches a fixed number**

Why GD?

Intuition: by first-order **Taylor approximation**

$$F(\mathbf{w}) \approx F(\mathbf{w}^{(t)}) + \nabla F(\mathbf{w}^{(t)})^T (\mathbf{w} - \mathbf{w}^{(t)})$$

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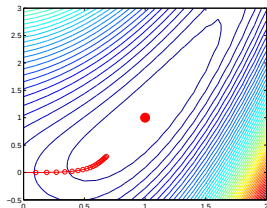
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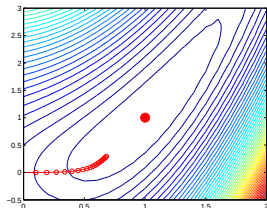
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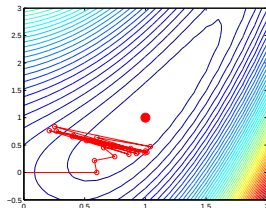
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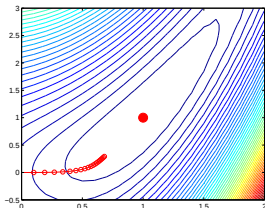
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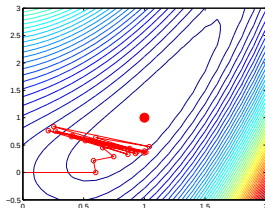
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See Colab Example 1



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More on learning rate

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Adaptive and automatic step size tuning is an active research area

- notable examples: AdaGrad, Adam, etc.
- ideas: tune η based on past gradient information

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More examples coming soon. Key point: it could be *much faster to obtain a stochastic gradient!*

Convergence guarantees — convex objectives

Many for both GD and SGD on **convex objectives**.

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Many for both GD and SGD on **convex objectives**.

They tell you how many iterations t (in terms of ϵ) needed to achieve

$$F(\mathbf{w}^{(t)}) - F(\mathbf{w}^*) \leq \epsilon$$

Convergence guarantees — convex objectives

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They tell you how many iterations t (in terms of ϵ) needed to achieve

$$F(\mathbf{w}^{(t)}) - F(\mathbf{w}^*) \leq \epsilon$$

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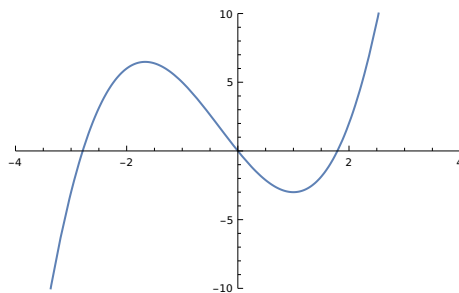
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Convergence guarantees — nonconvex objectives

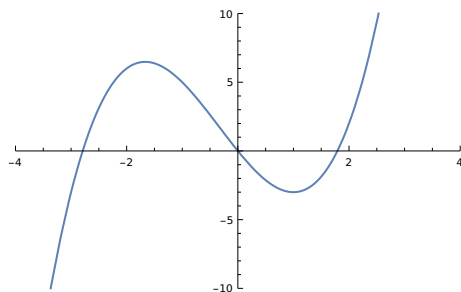
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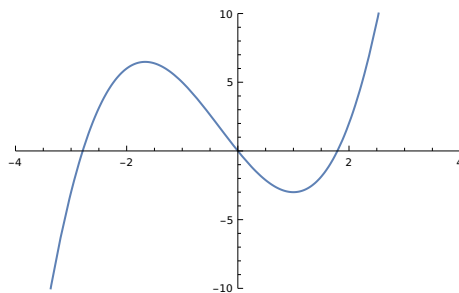
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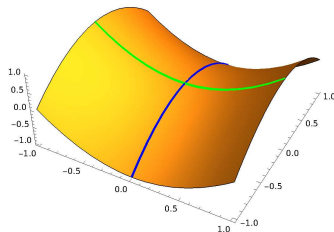
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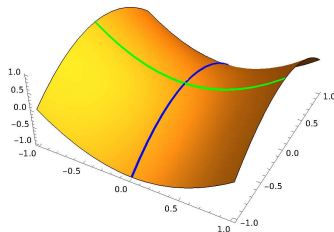
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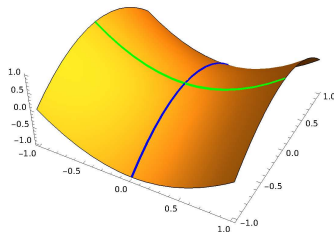
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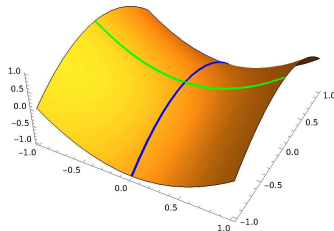
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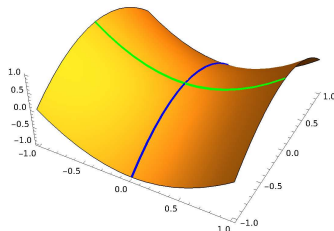
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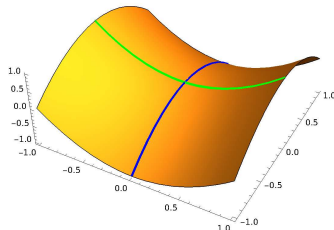
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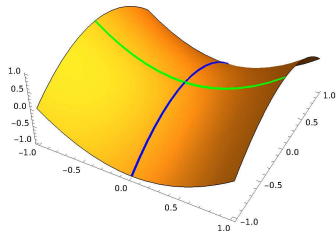
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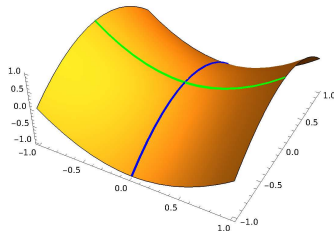
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- so not a real issue especially *when initialized randomly* (Colab Example 2)



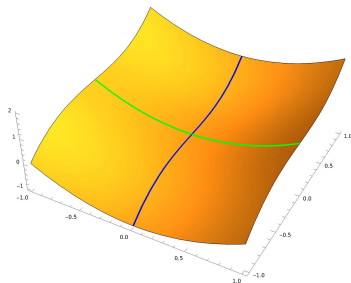
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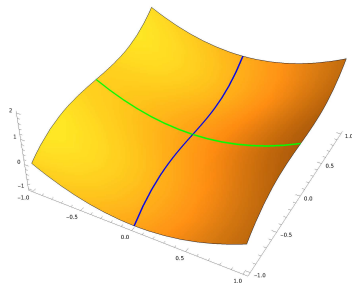
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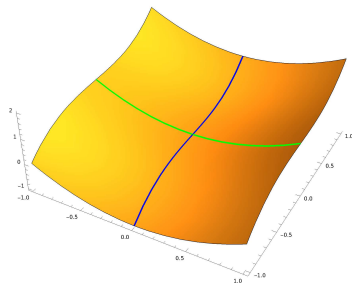
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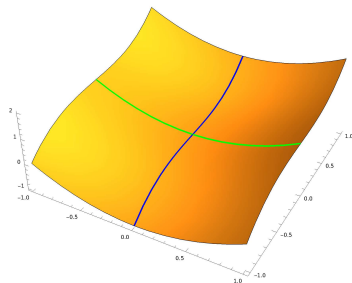
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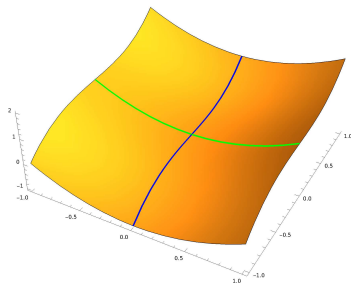
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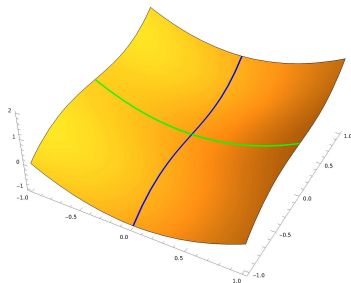
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Even worse, distinguishing local min and saddle point is generally **NP-hard**.

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- recent research shows that *many problems have no “bad” saddle points or even “bad” local minimizers*
- justify the practical effectiveness of GD/SGD (default method to try)

Second-order methods

Recall the intuition of GD: we look at first-order **Taylor approximation**

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where $\mathbf{H}_t = \nabla^2 F(\mathbf{w}^{(t)}) \in \mathbb{R}^{D \times D}$ is the *Hessian* of F at $\mathbf{w}^{(t)}$, i.e.,

$$H_{t,ij} = \left. \frac{\partial^2 F(\mathbf{w})}{\partial w_i \partial w_j} \right|_{\mathbf{w}=\mathbf{w}^{(t)}}$$

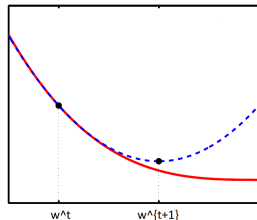
(think “second derivative” when $D = 1$)

Newton method

If we minimize the second-order approximation (via “complete the square”)

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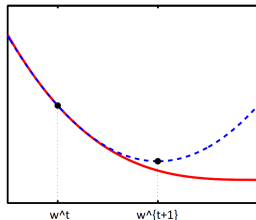
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for strictly convex F (so H_t is *positive definite*), we obtain **Newton method**:

$$\mathbf{w}^{(t+1)} \leftarrow \mathbf{w}^{(t)} - \mathbf{H}_t^{-1} \nabla F(\mathbf{w}^{(t)})$$



Comparing GD and Newton

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- does not really make sense for *nonconvex objectives*

Outline

- 1 Review of Last Lecture
- 2 Linear Classifiers and Surrogate Losses
- 3 A Detour of Numerical Optimization Methods
- 4 Perceptron**
- 5 Logistic Regression

Recall the perceptron loss

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Let's approximately minimize it with GD/SGD.

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Slow: each update makes one pass of the entire training set!

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Applying SGD to perceptron loss

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One common trick: pick one example $n \in [N]$ uniformly at random, let

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Conveniently, objective of most ML tasks is a *finite sum* (over each training point) and the above trick applies!

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- why $\eta = 1$? Does not really matter in terms of prediction of $\mathbf{w}$

Why does it make sense?

If the current weight $\mathbf{w}$ makes a mistake

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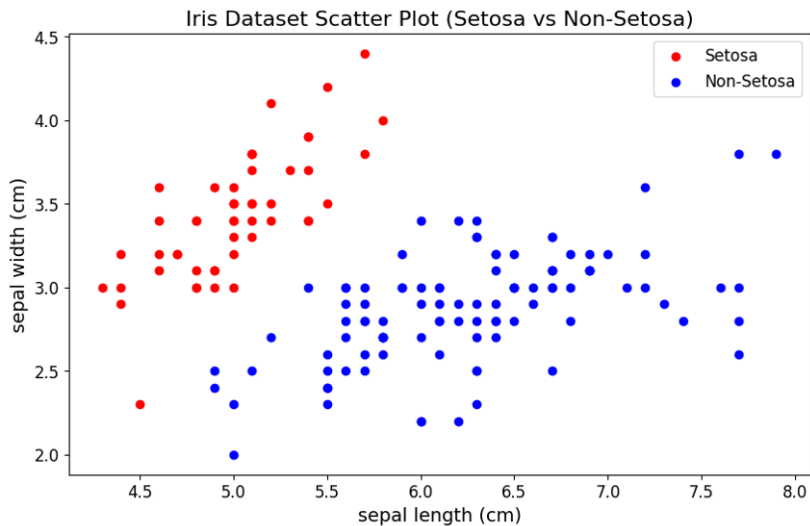
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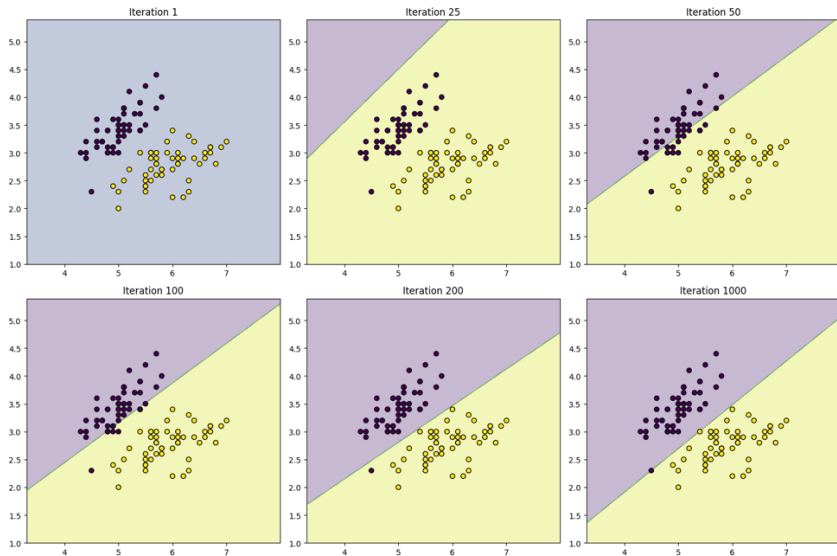
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Thus it is more likely to get it right after the update.

Example: Iris Dataset



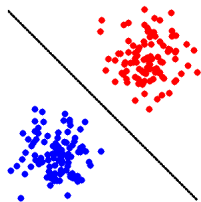
Example: Perceptron for Iris Dataset



Any theory?

If training set is linearly separable

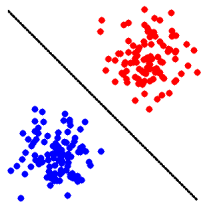
- Perceptron *converges in a finite number of steps*
- training error is 0



Any theory?

If training set is linearly separable

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There are also guarantees when the data are not linearly separable.

Outline

- 1 Review of Last Lecture
- 2 Linear Classifiers and Surrogate Losses
- 3 A Detour of Numerical Optimization Methods
- 4 Perceptron
- 5 Logistic Regression**
 - A probabilistic view

A simple view

In one sentence: find the minimizer of

$$\begin{aligned} F(\mathbf{w}) &= \frac{1}{N} \sum_{n=1}^N \ell_{\text{logistic}}(y_n \mathbf{w}^T \mathbf{x}_n) \\ &= \frac{1}{N} \sum_{n=1}^N \ln(1 + e^{-y_n \mathbf{w}^T \mathbf{x}_n}) \end{aligned}$$

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Before optimizing it: *why logistic loss? and why "regression"?*

Predicting probability

Instead of predicting a discrete label, can we *predict the probability of each label?* i.e. regress the probabilities

Predicting probability

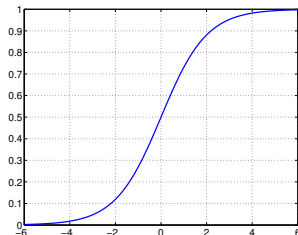
Instead of predicting a discrete label, can we *predict the probability of each label?* i.e. regress the probabilities

One way: **sigmoid function + linear model**

$$\mathbb{P}(y = +1 \mid \mathbf{x}; \mathbf{w}) = \sigma(\mathbf{w}^T \mathbf{x})$$

where σ is the sigmoid function:

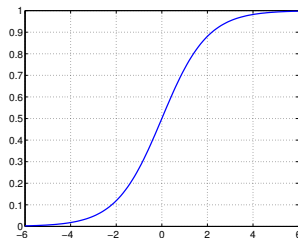
$$\sigma(z) = \frac{1}{1 + e^{-z}}$$



Properties

Properties of sigmoid $\sigma(z) = \frac{1}{1+e^{-z}}$

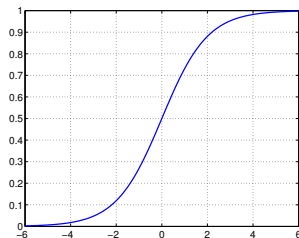
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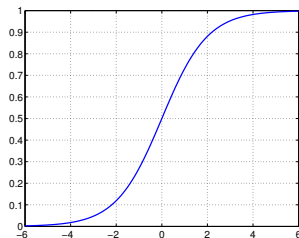
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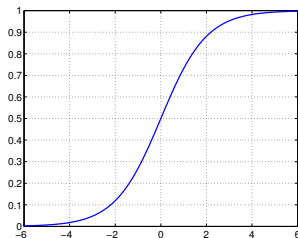
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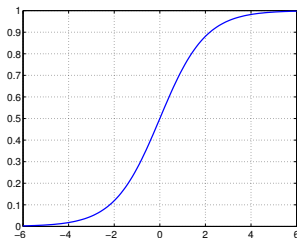
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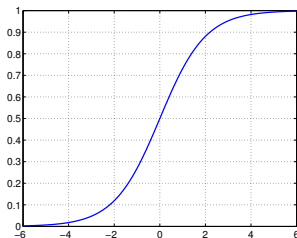
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and thus

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Take a **probabilistic view**

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Specifically, what is the probability of seeing label $y_1, \dots, y_n$ given $x_1, \dots, x_n$, as a function of some w ?

$$P(w) = \prod_{n=1}^N \mathbb{P}(y_n \mid x_n; w)$$

MLE: find w^* that **maximizes the probability** $P(w)$

The MLE solution

$$\boldsymbol{w}^* = \operatorname{argmax}_{\boldsymbol{w}} P(\boldsymbol{w}) = \operatorname{argmax}_{\boldsymbol{w}} \prod_{n=1}^N \mathbb{P}(y_n \mid \boldsymbol{x}_n; \boldsymbol{w})$$

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i.e. *minimizing logistic loss is exactly doing MLE for the sigmoid model!*

Back to algorithms: apply SGD again

$$\boldsymbol{w} \leftarrow \boldsymbol{w} - \eta \tilde{\nabla} F(\boldsymbol{w})$$

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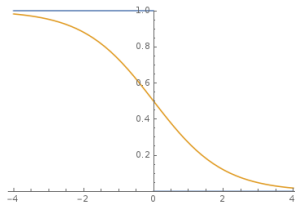
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 \end{aligned}$$

This is a *soft version of Perceptron!*

$\mathbb{P}(-y_n \mid \mathbf{x}_n; \mathbf{w})$ versus $\mathbb{I}[y_n \neq \text{sgn}(\mathbf{w}^T \mathbf{x}_n)]$



Applying Newton to logistic loss

$$\nabla_{\mathbf{w}} \ell_{\text{logistic}}(y_n \mathbf{w}^T \mathbf{x}_n) = -\sigma(-y_n \mathbf{w}^T \mathbf{x}_n) y_n \mathbf{x}_n$$

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Exercises:

- why is the Hessian of logistic loss positive semidefinite?

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Exercises:

- why is the Hessian of logistic loss positive semidefinite?
- can we apply Newton method to perceptron/hinge loss?

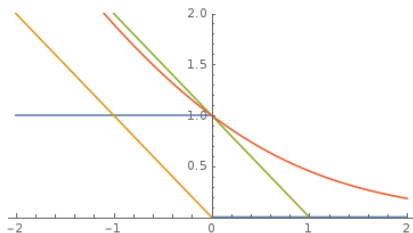
Summary

Linear models for classification:

Step 1. Model is the set of **separating hyperplanes**

$$\mathcal{F} = \{f(\mathbf{x}) = \text{sgn}(\mathbf{w}^T \mathbf{x}) \mid \mathbf{w} \in \mathbb{R}^D\}$$

Step 2. Pick the **surrogate loss**



- **perceptron loss** $\ell_{\text{perceptron}}(z) = \max\{0, -z\}$ (used in Perceptron)
- **hinge loss** $\ell_{\text{hinge}}(z) = \max\{0, 1 - z\}$ (used in SVM and many others)
- **logistic loss** $\ell_{\text{logistic}}(z) = \log(1 + \exp(-z))$ (used in logistic regression)

Step 3. Find empirical risk minimizer (ERM):

$$\mathbf{w}^* = \operatorname{argmin}_{\mathbf{w} \in \mathbb{R}^D} \frac{1}{N} \sum_{n=1}^N \ell(y_n \mathbf{w}^T \mathbf{x}_n)$$

using

- **GD:** $\mathbf{w} \leftarrow \mathbf{w} - \eta \nabla F(\mathbf{w})$
- **SGD:** $\mathbf{w} \leftarrow \mathbf{w} - \eta \tilde{\nabla} F(\mathbf{w})$ $(\mathbb{E}[\tilde{\nabla} F(\mathbf{w})] = \nabla F(\mathbf{w}))$
- **Newton:** $\mathbf{w} \leftarrow \mathbf{w} - (\nabla^2 F(\mathbf{w}))^{-1} \nabla F(\mathbf{w})$