

CPSC 340: Machine Learning and Data Mining

Linear Classifiers
Spring 2022 (2021W2)

Admin

- A4 released on Feb 25 (Fri); Due Friday Mar 11

Last Time: L1-Regularization

- We discussed L1-regularization:

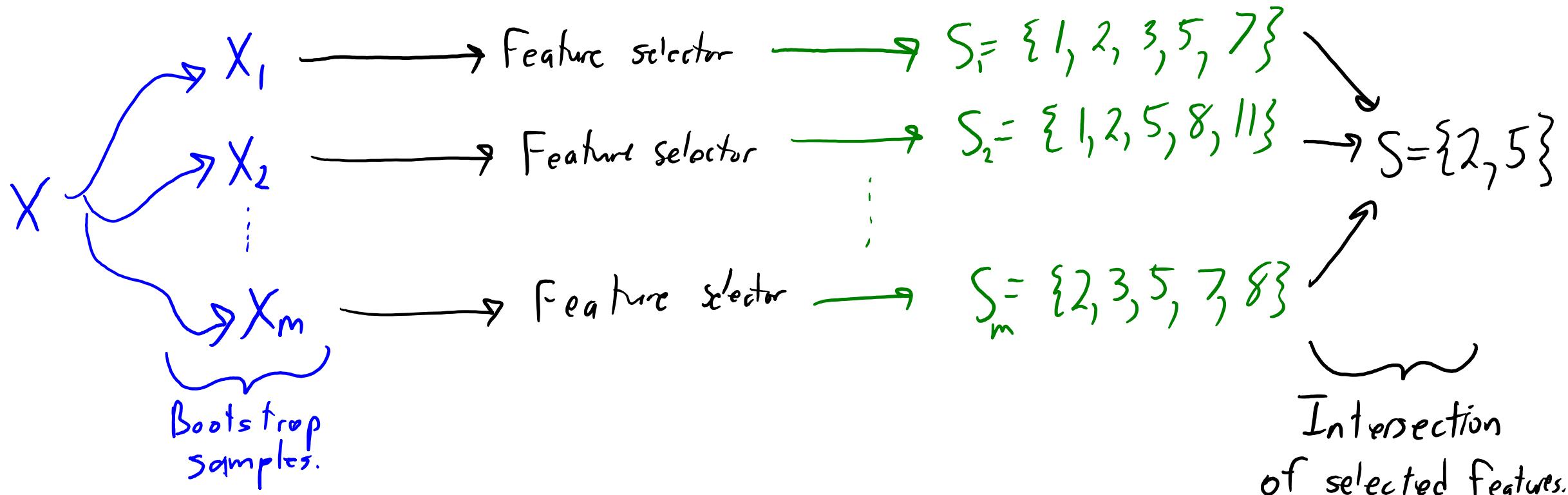
$$f(w) = \frac{1}{2} \|Xw - y\|^2 + \gamma \|w\|_1$$

- Also known as “LASSO” and “basis pursuit denoising”.
- Regularizes ‘w’ so we decrease our test error (like L2-regularization).
- Yields sparse ‘w’ so it selects features (like L0-regularization).
- Properties:
 - It’s convex and fast to minimize (with “proximal-gradient” methods).
 - Solution is not unique (sometimes people do L2- and L1-regularization).
 - Usually includes “correct” variables but tends to yield false positives.

Ensemble Feature Selection

- We can also use **ensemble methods** for feature selection.
 - Usually designed to **reduce false positives** or **reduce false negatives**.
- In this case of L1-regularization, we **want to reduce false positives**.
 - Unlike L0-regularization, the **non-zero w_j** are still “shrunken”.
 - “Irrelevant” variables can be included before “relevant” w_j reach best value.
- A **bootstrap** approach to reducing false positives:
 - Apply the method to bootstrap samples of the training data.
 - Only take the **features selected** in all bootstrap samples.

Ensemble Feature Selection



- Example: LASSO on Bootstrapped samples [Bolasso: Model Consistent Lasso Estimation through the Bootstrap, *F. Bach 2008*].
 - Reduces false positives.
 - It's possible to show it recovers “correct” variables.
 - Can replace “intersection” with “selected frequency” if has false negatives too.

(pause)

Motivation: Identifying Important E-mails

- How can we automatically identify ‘important’ e-mails?



- A **binary classification** problem (“important” vs. “not important”).
 - Labels are approximated by whether you took an “action” based on mail.
 - High-dimensional feature set (that we’ll discuss later).
- Gmail uses **regression** for this binary classification problem.

Binary Classification Using Regression?

- Can we apply linear models for **binary classification**?
 - Set $y_i = +1$ for one class (“important”).
 - Set $y_i = -1$ for the other class (“not important”).
- At training time, **fit a linear regression model**:

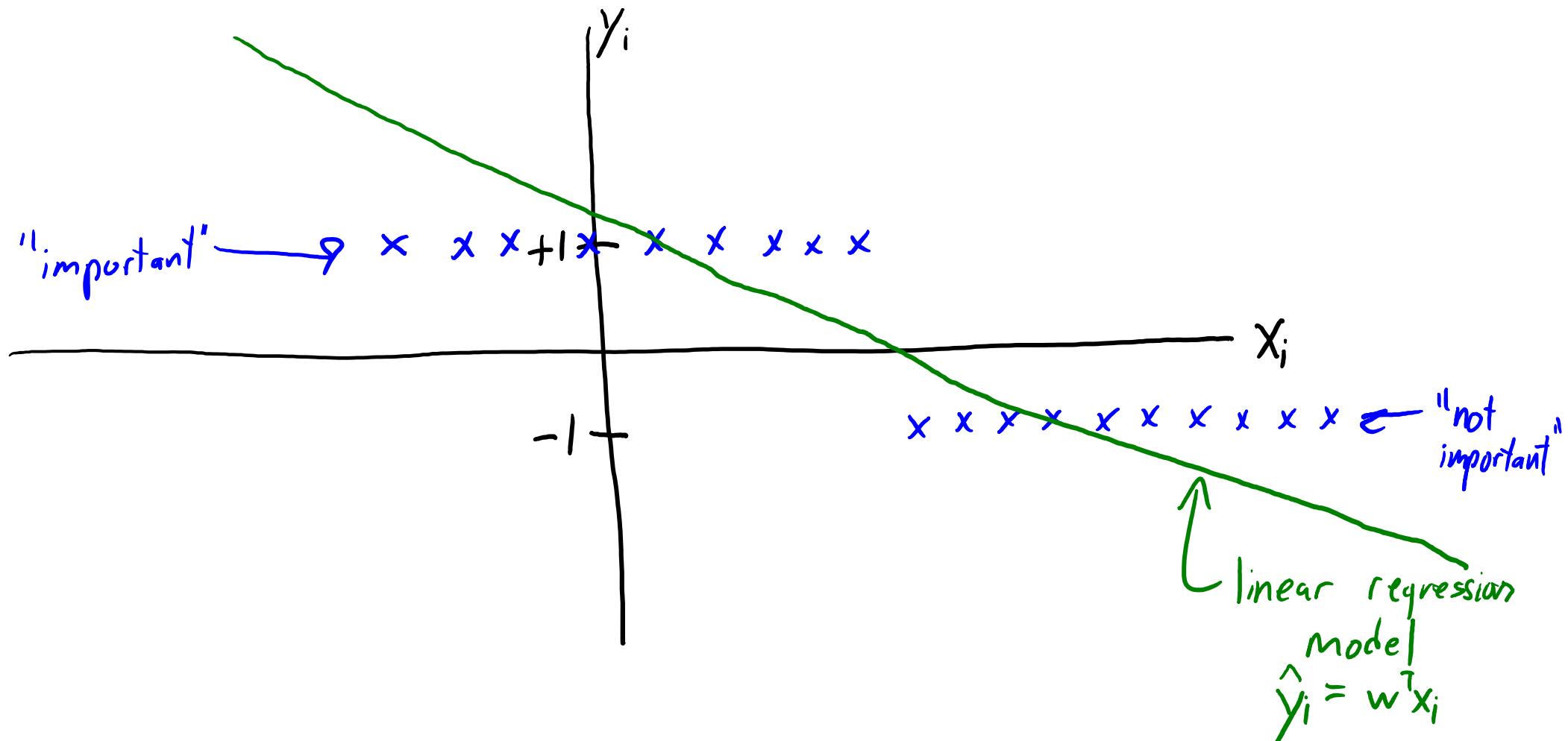
$$\begin{aligned}\hat{y}_i &= w_1 x_{i1} + w_2 x_{i2} + \dots + w_d x_{id} \\ &= w^T x_i\end{aligned}$$

- The model will try to make $w^T x_i = +1$ for “important” e-mails, and $w^T x_i = -1$ for “not important” e-mails.

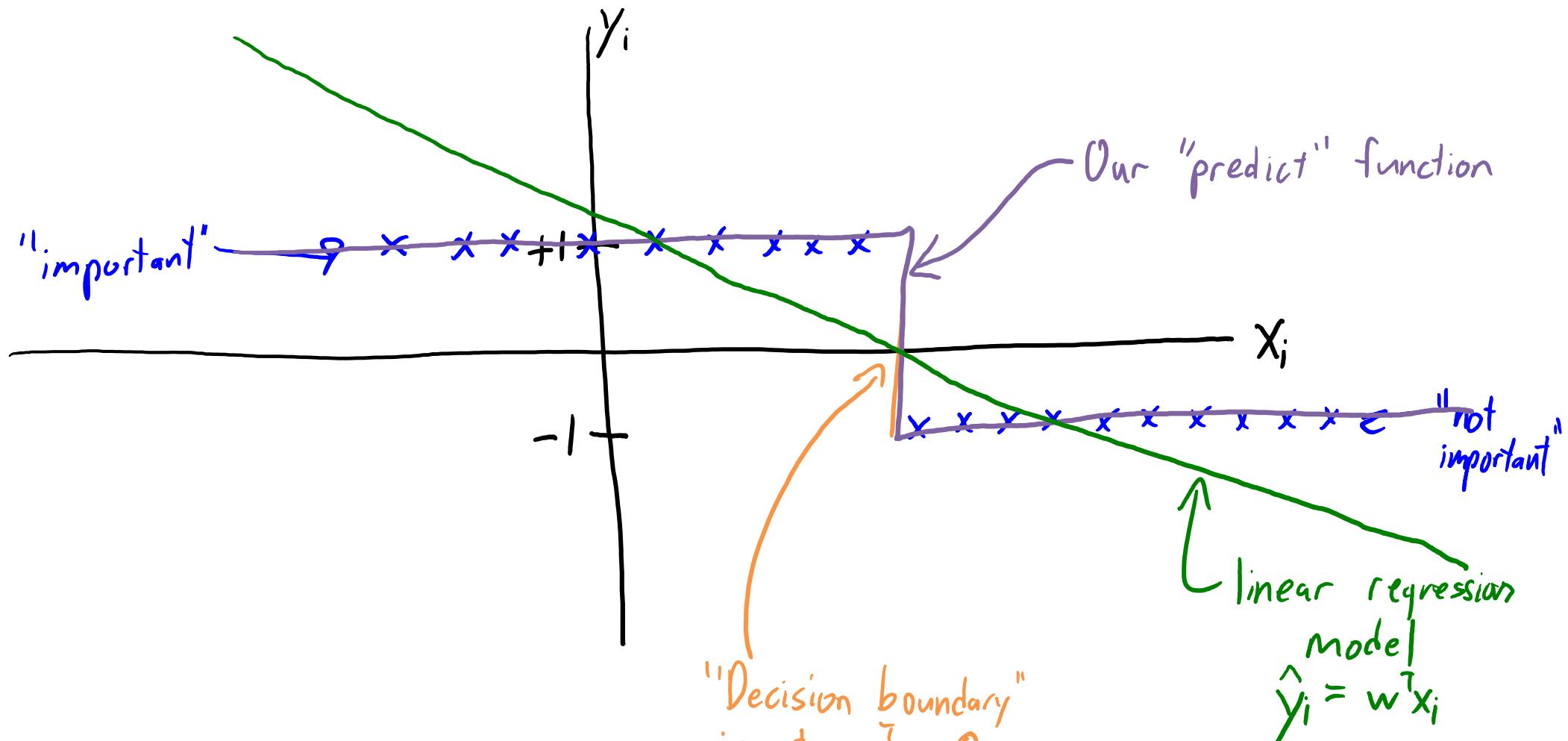
Binary Classification Using Regression?

- Can we apply linear models for **binary classification**?
 - Set $y_i = +1$ for one class (“important”).
 - Set $y_i = -1$ for the other class (“not important”).
- Linear model gives **real numbers** like 0.9, -1.1, and so on.
- So to predict, we look at whether $w^T x_i$ is closer to +1 or -1.
 - If $w^T x_i = 0.9$, predict $\hat{y}_i = +1$.
 - If $w^T x_i = -1.1$, predict $\hat{y}_i = -1$.
 - If $w^T x_i = 0.1$, predict $\hat{y}_i = +1$.
 - If $w^T x_i = -100$, predict $\hat{y}_i = -1$.
 - We write this operation (rounding to +1 or -1) as $\hat{y}_i = \text{sign}(w^T x_i)$.

Decision Boundary in 1D



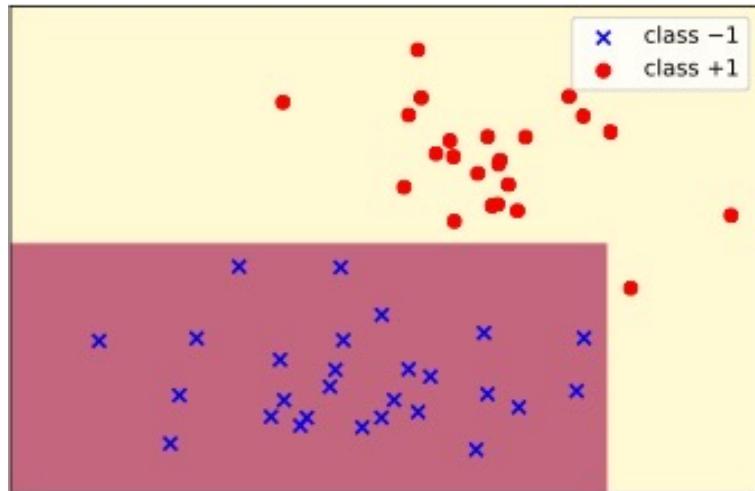
Decision Boundary in 1D



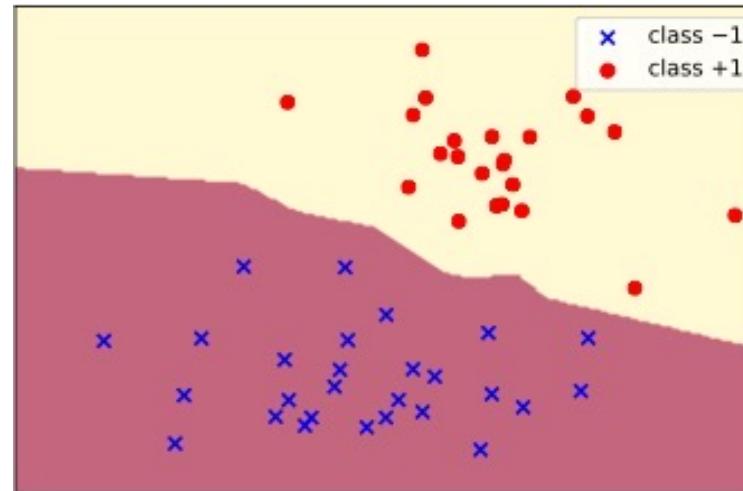
- We can interpret 'w' as a hyperplane separating x into sets:
 - Set where $w^T x_i > 0$ and set where $w^T x_i < 0$.

Decision Boundary in 2D

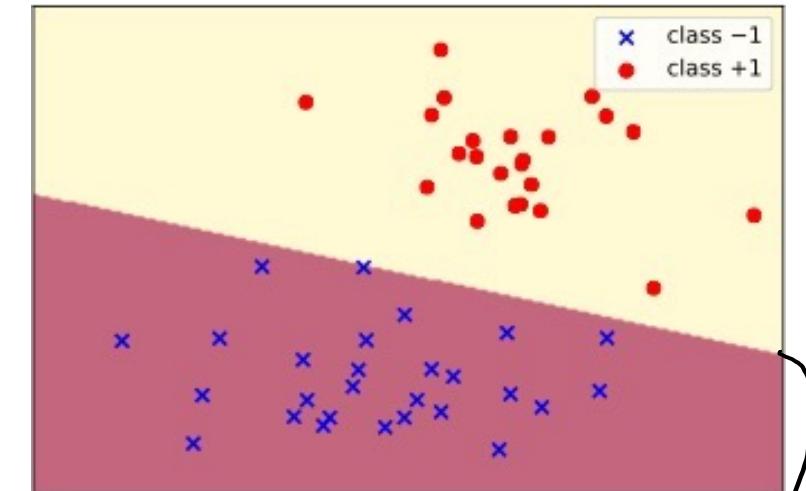
decision tree



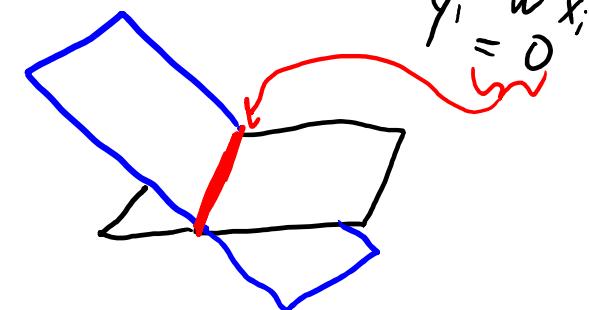
KNN



linear classifier



- Linear classifier would be a $\hat{y}_i = w^T x_i$ function coming out of screen:
– The boundary is at $\hat{y}_i = 0$.



Should we use least squares for classification?

- Consider training by minimizing squared error with y_i that are +1 or -1:

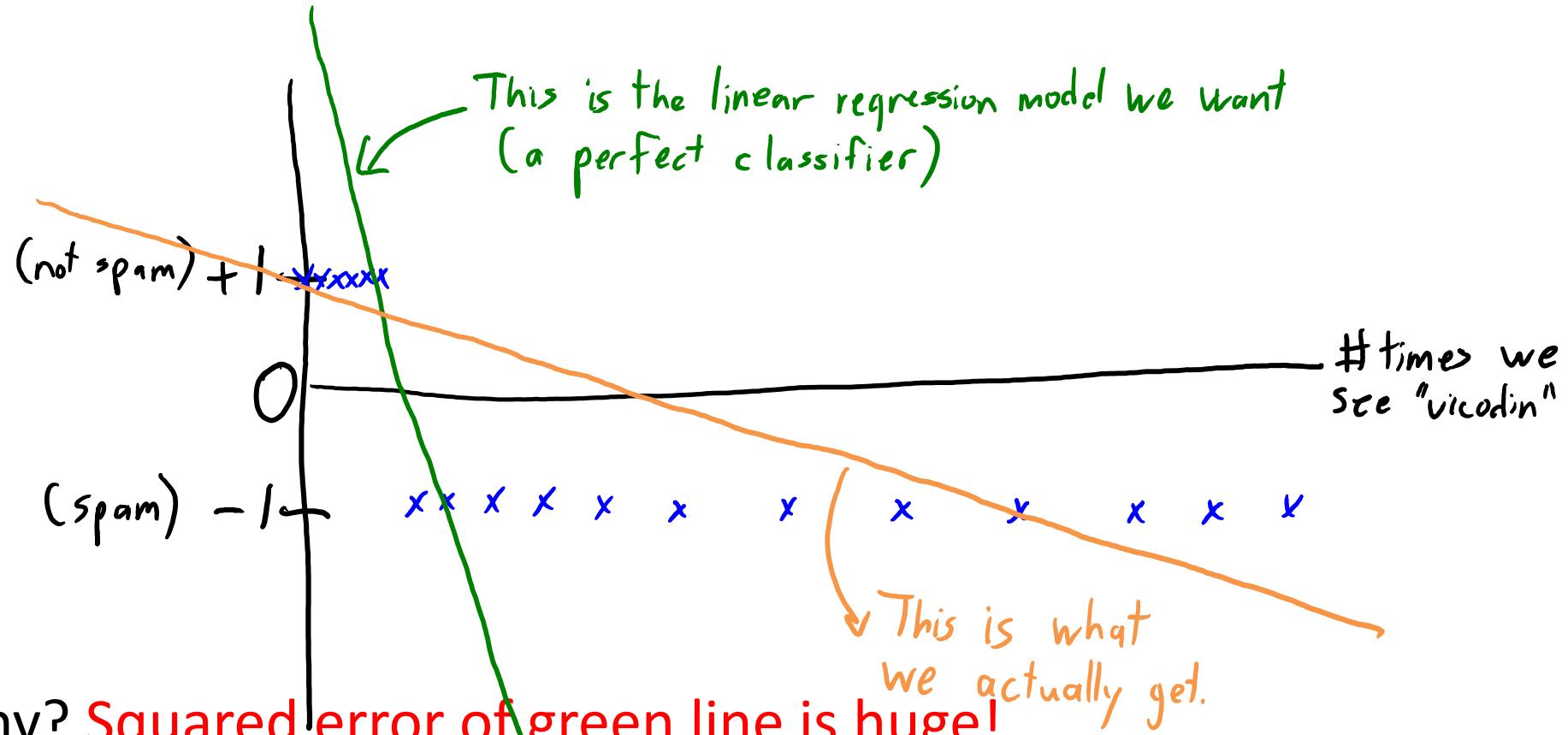
$$f(w) = \frac{1}{2} \|Xw - y\|^2$$

A handwritten-style equation $f(w) = \frac{1}{2} \|Xw - y\|^2$ is shown. To its right is a vertical vector labeled $y =$ with five entries: +1, +1, -1, +1, -1. A curly brace underlines the first two entries, and another curly brace underlines the last three entries.

- If we predict $w^T x_i = +0.9$ and $y_i = +1$, error is small: $(0.9 - 1)^2 = 0.01$.
- If we predict $w^T x_i = -0.8$ and $y_i = +1$, error is bigger: $(-0.8 - 1)^2 = 3.24$.
- If we predict $w^T x_i = +100$ and $y_i = +1$, error is huge: $(100 - 1)^2 = 9801$.
 - But it shouldn't be, the prediction is correct.
- Least squares penalized for being “too right”.
 - +100 has the right sign, so the error should not be large.

Should we use least squares for classification?

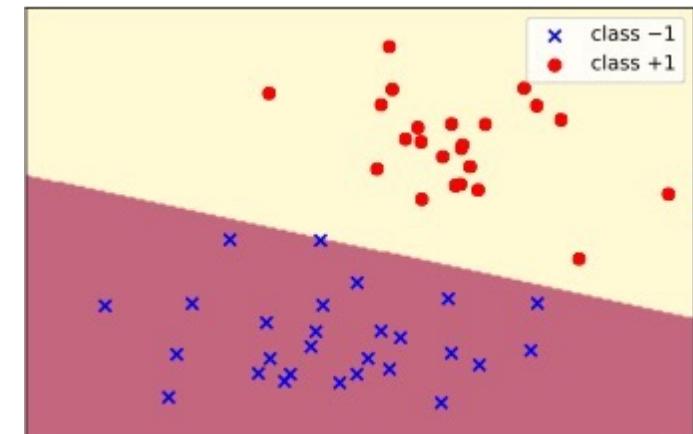
- Least squares can behave weirdly when applied to classification:



- Why? Squared error of green line is huge!
 - Make sure you understand why the green line achieves 0 training error.

“0-1 Loss” Function: Minimizing Classification Errors

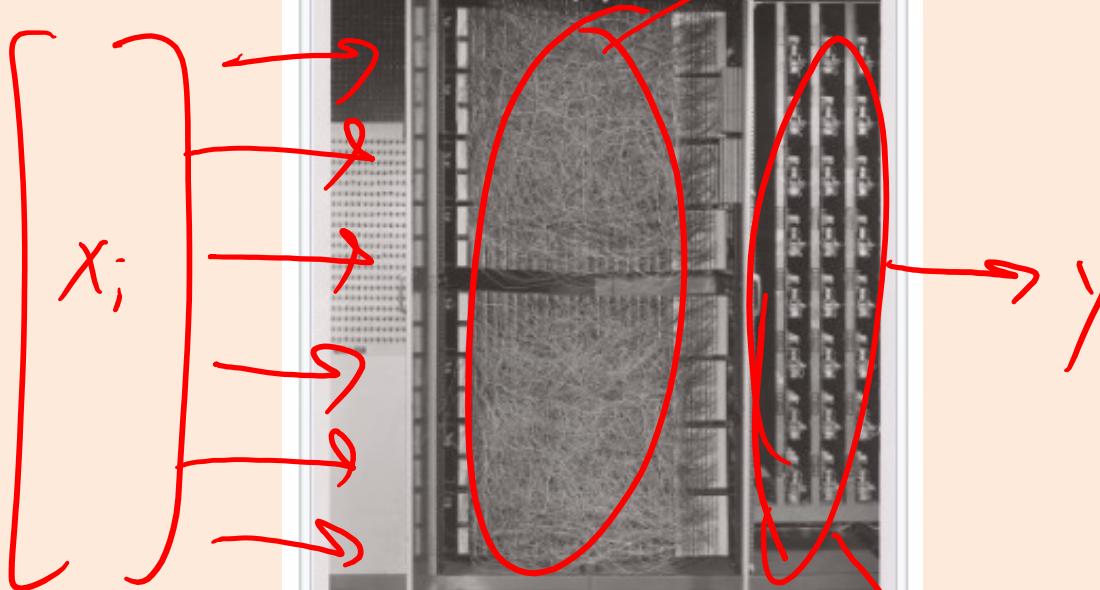
- Could we instead minimize **number of classification errors**?
 - This is called the **0-1 loss** function:
 - You either get the classification wrong (1) or right (0).
 - We can write using the L0-norm as $\|\hat{y} - y\|_0$.
 - Unlike regression, in classification it's reasonable that $\hat{y}_i = y_i$ (it's either +1 or -1).
- Important special case: “**linearly separable**” data.
 - Classes **can be “separated” by a hyper-plane**.
 - So a perfect linear classifier exists.



Perceptron Algorithm for Linearly-Separable Data

- One of the first “learning” algorithms was the “perceptron” (1957).
 - Searches for a ‘w’ such that $\text{sign}(w^T x_i) = y_i$ for all i.
- Perceptron algorithm:
 - Start with $w^0 = 0$.
 - Go through examples in any order until you make a mistake predicting y_i .
 - Set $w^{t+1} = w^t + y_i x_i$.
 - Keep going through examples until you make no errors on training data.
- If a perfect classifier exists, this algorithm finds one in finite number of steps.
- Intuition:
 - Consider a case where $w^T x_i < 0$ but $y_i = +1$.
 - In this case the update “adds more of x_i to w ” so that $w^T x_i$ is larger.
$$(w^{t+1})^T x_i = (w^t + x_i)^T x_i = (w^t)^T x_i + x_i^T x_i = (\text{old prediction}) + \|x_i\|^2$$
 - If $y_i = -1$, you would be subtracting the squared norm.

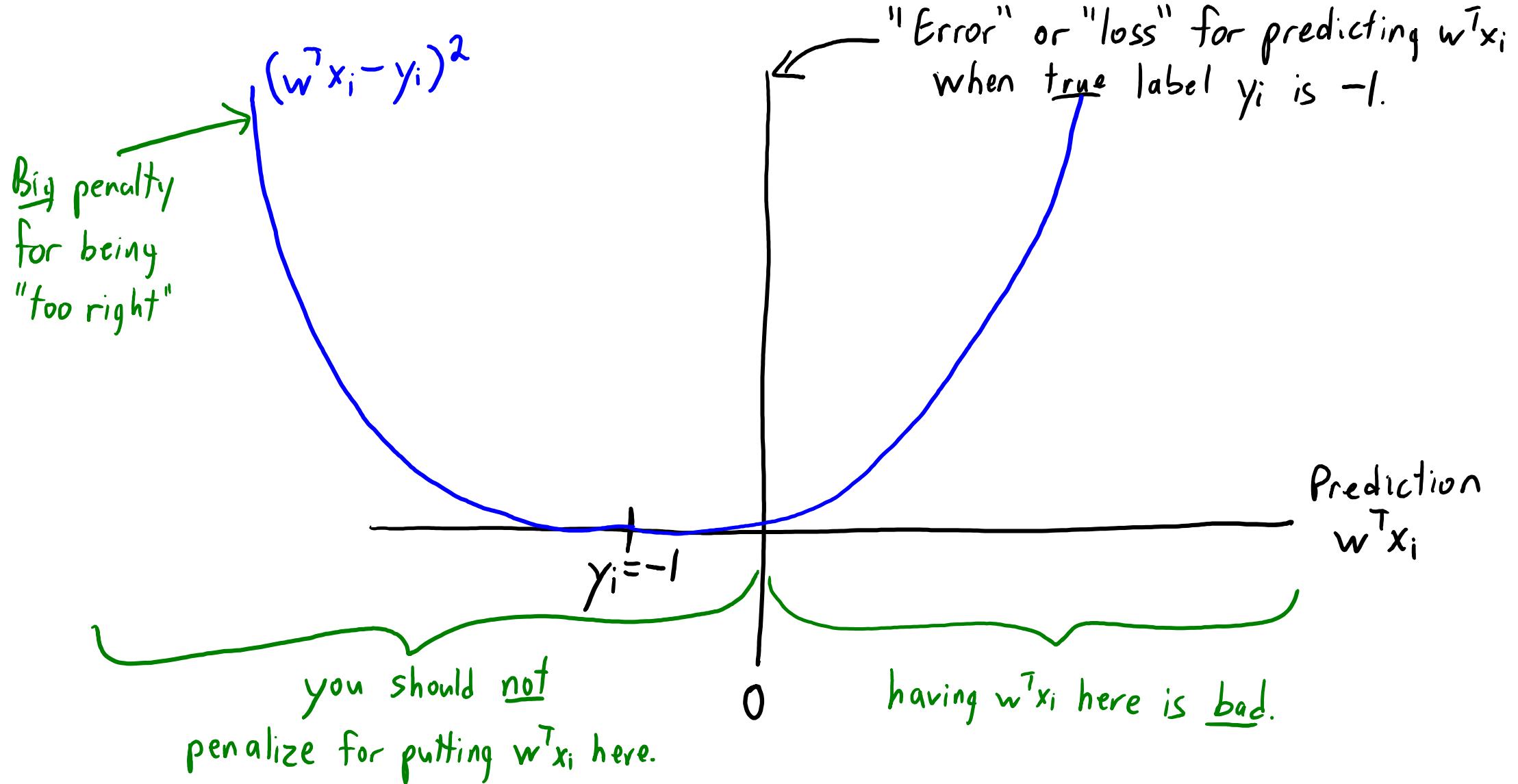
History [edit]



The Mark I Perceptron machine was the first implementation of the perceptron algorithm. The machine was connected to a camera that used 20×20 cadmium sulfide photocells to produce a 400-pixel image. The main visible feature is a patchboard that allowed experimentation with different combinations of input features. To the right of that are arrays of potentiometers that implemented the adaptive weights.^{[2]:213}

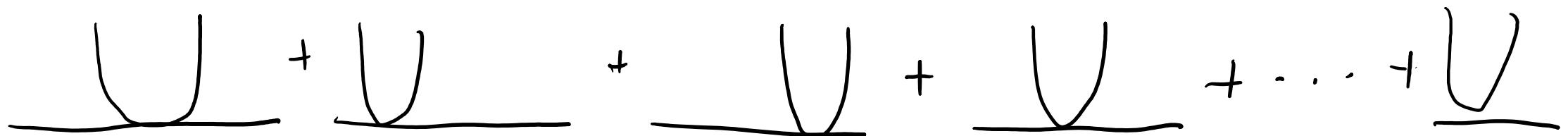
$$z_i = [x_i^2 \ x_i x_2 \ x_3]^{\text{bonus!}}$$

Geometry of why we want the 0-1 loss

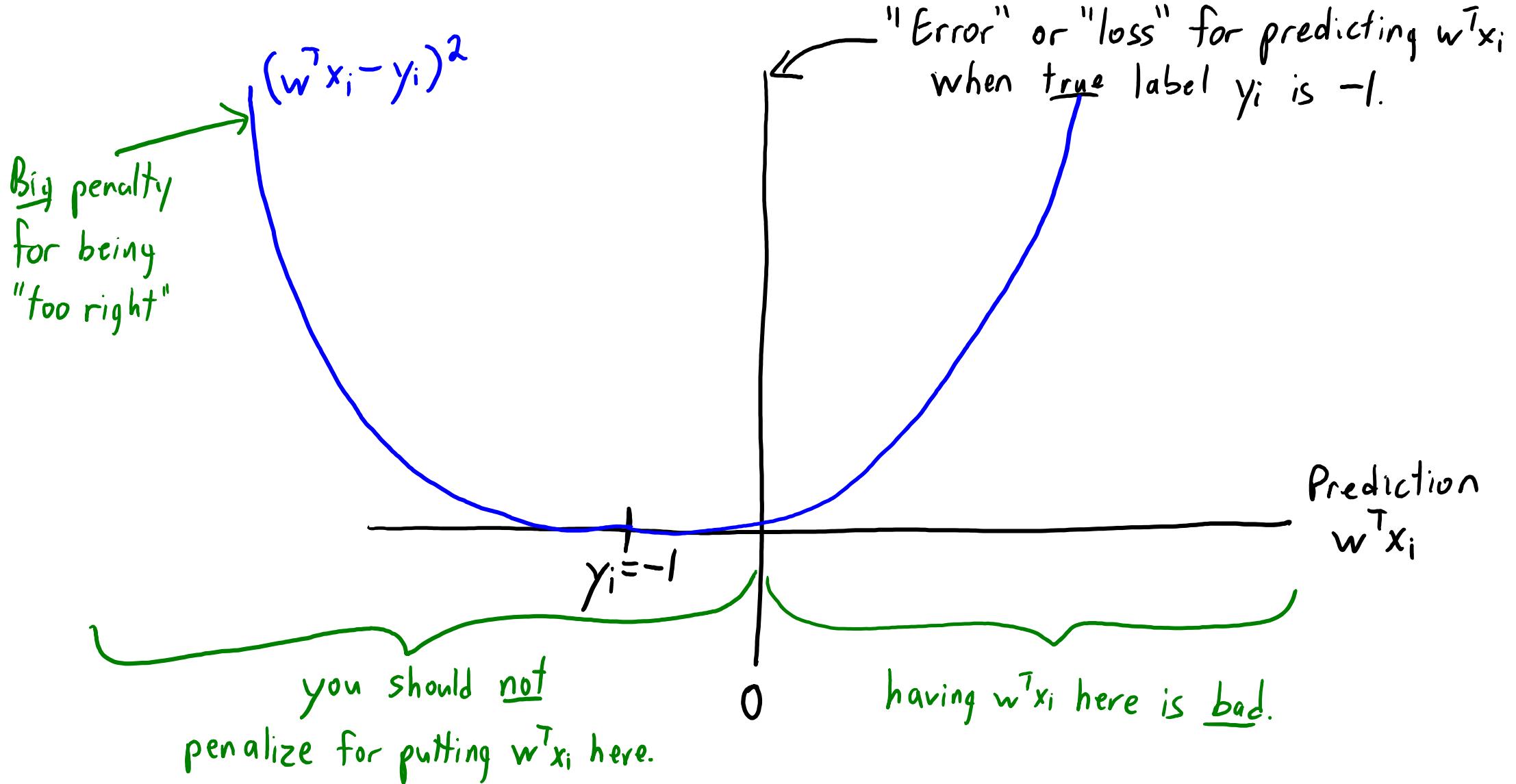


Wait, what does that picture mean?

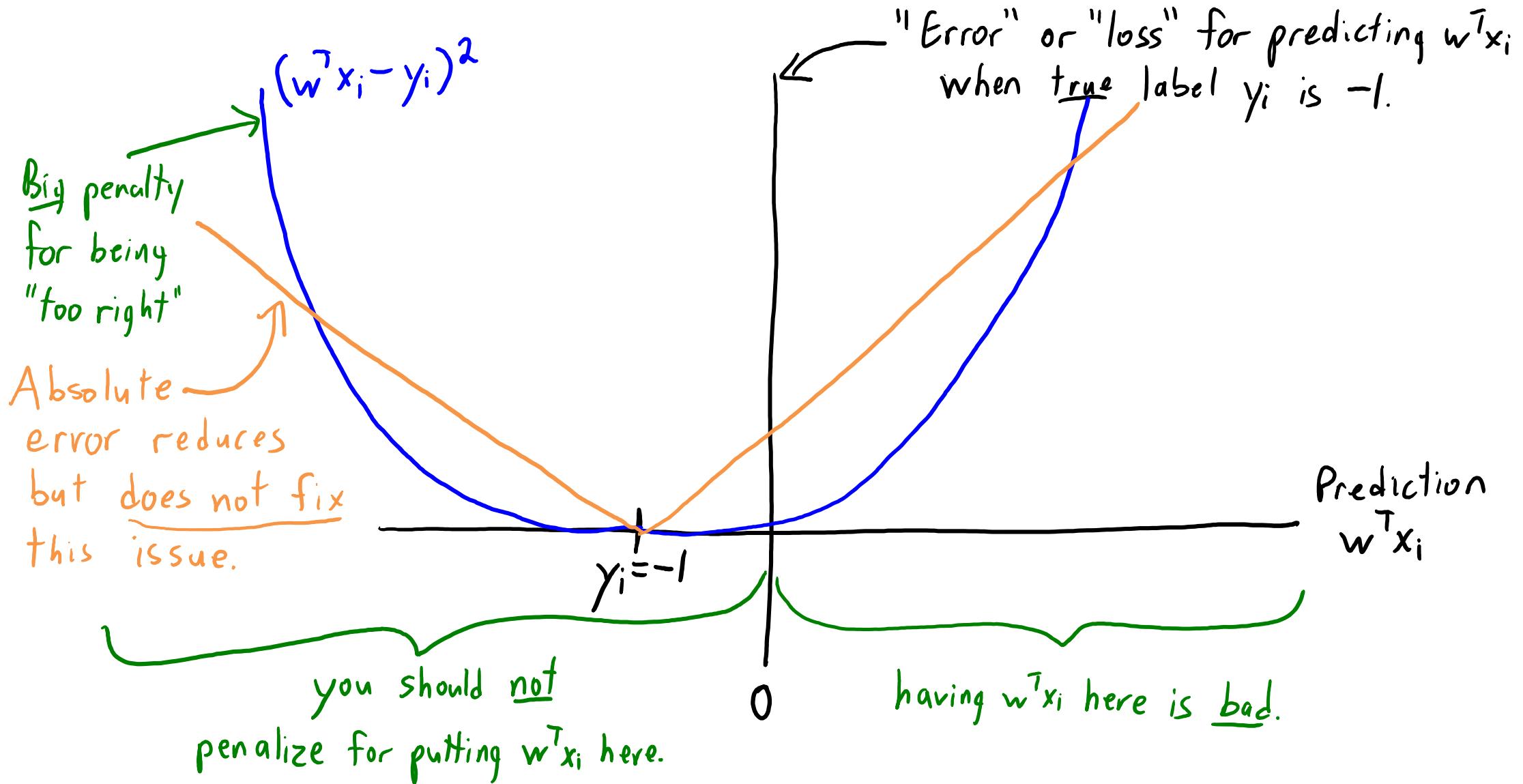
- We are now plotting the loss vs. the predicted $w^T x_i$.
 - “Loss space”, which is different than parameter space or data space.
 - Close (but not quite same) as when we were plotting losses for robust regression
- We're plotting the individual loss for a particular training example.
 - In the figure the label is $y_i = -1$ (so loss is centered at -1).
 - It will be centered at +1 when $y_i = +1$.
 - The objective in least squares regression is a sum of ‘n’ of these losses:



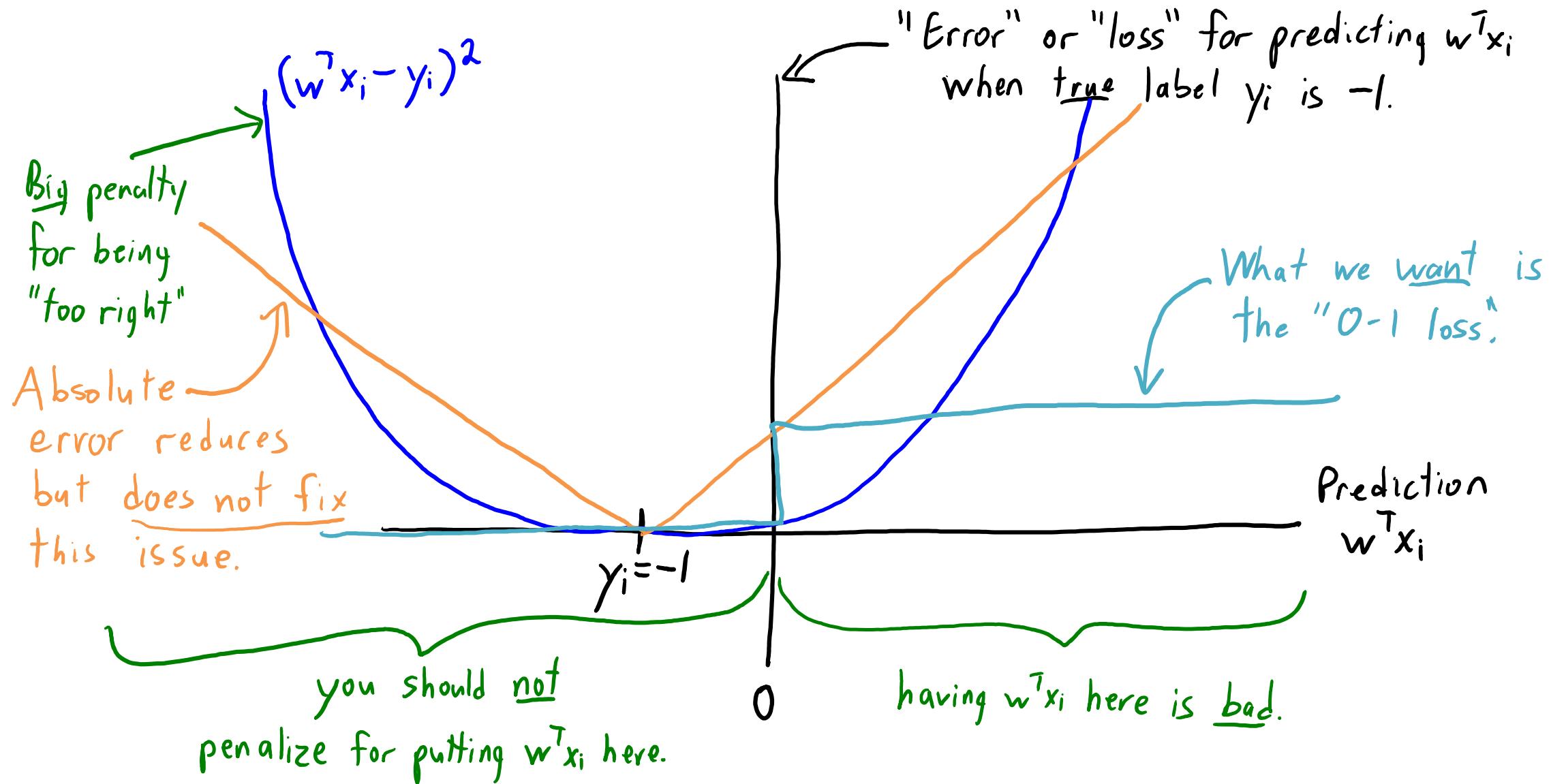
Geometry of why we want the 0-1 loss



Geometry of why we want the 0-1 loss



Geometry of why we want the 0-1 loss



0-1 Loss Function

- Unfortunately the **0-1 loss is non-convex** in ‘w’.
 - It’s easy to minimize if a perfect classifier exists (perceptron).
 - Otherwise, finding the ‘w’ **minimizing 0-1 loss is (NP-)hard**.
 - Gradient is zero everywhere: don’t even know “which way to go”.
 - NOT the same type of problem we had with using the squared loss.
 - We can minimize the squared error, but it might give a bad model for classification.
- Motivates **convex approximations to 0-1 loss...**

A Convex Approximation to 0-1 Loss

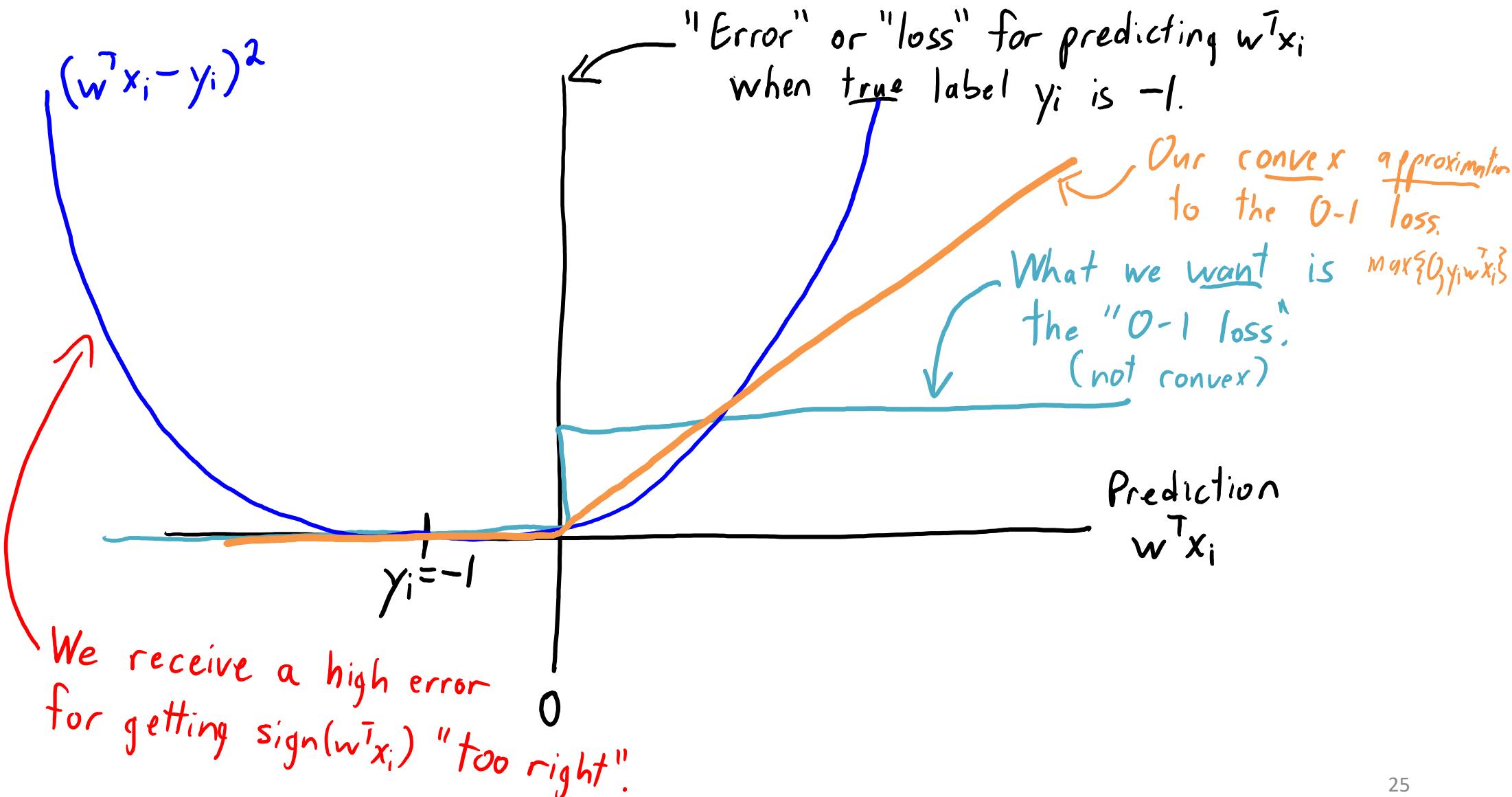
- If $y_i = +1$, we get the label right if $w^T x_i > 0$.
- If $y_i = -1$, we get the label right if $w^T x_i < 0$, or equivalently $-w^T x_i > 0$.
- So “classifying ‘i’ correctly” is equivalent to having $y_i w^T x_i > 0$.
- One possible convex approximation to 0-1 loss:
 - Minimize how much this constraint is violated.

If $y_i w^T x_i > 0$ then you get an “error” of 0.

If $y_i w^T x_i < 0$ then you get an “error” of $-y_i w^T x_i$

→ So the “error” is given by $\max\{0, -y_i w^T x_i\}$
 $\max\{\text{constant}, \text{linear}\} \Rightarrow \text{convex}$

A Convex Approximation to 0-1 Loss



A (Bad) Convex Approximation to 0-1 Loss

- Our convex approximation of the error for one example is:

$$\max\{0, -y_i w^\top x_i\}$$

- We could train by minimizing sum over all examples:

$$f(w) = \sum_{i=1}^n \max\{0, -y_i w^\top x_i\}$$

- But this has a **degenerate solution**:
 - We have $f(0) = 0$, and this is the lowest possible value of 'f'.
- There are two standard fixes: **hinge loss** and **logistic loss**.

Hinge Loss

- We saw that we classify examples ‘i’ correctly if $y_i w^T x_i > 0$.
 - Our convex approximation is the amount this inequality is violated.
- Consider replacing $y_i w^T x_i > 0$ with $y_i w^T x_i \geq 1$.
(the “1” is arbitrary: we could make $\|w\|$ bigger/smaller to use any positive constant)
- The violation of this constraint is now given by:

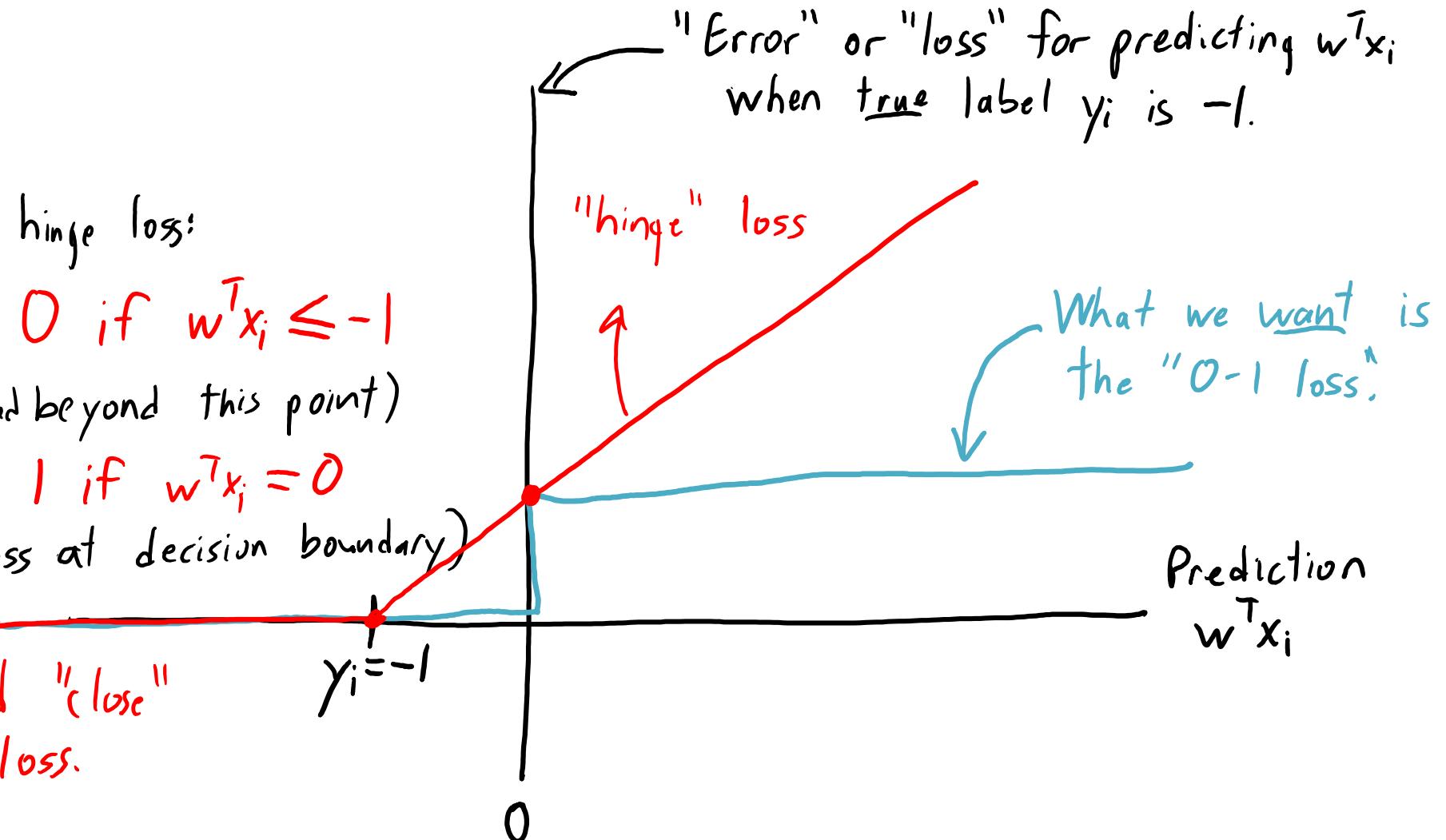
$$\max \{0, 1 - y_i w^T x_i\}$$

- This is called hinge loss.
 - It’s convex: $\max(\text{constant}, \text{linear})$.
 - It’s not degenerate: $w=0$ now gives an error of 1 instead of 0.

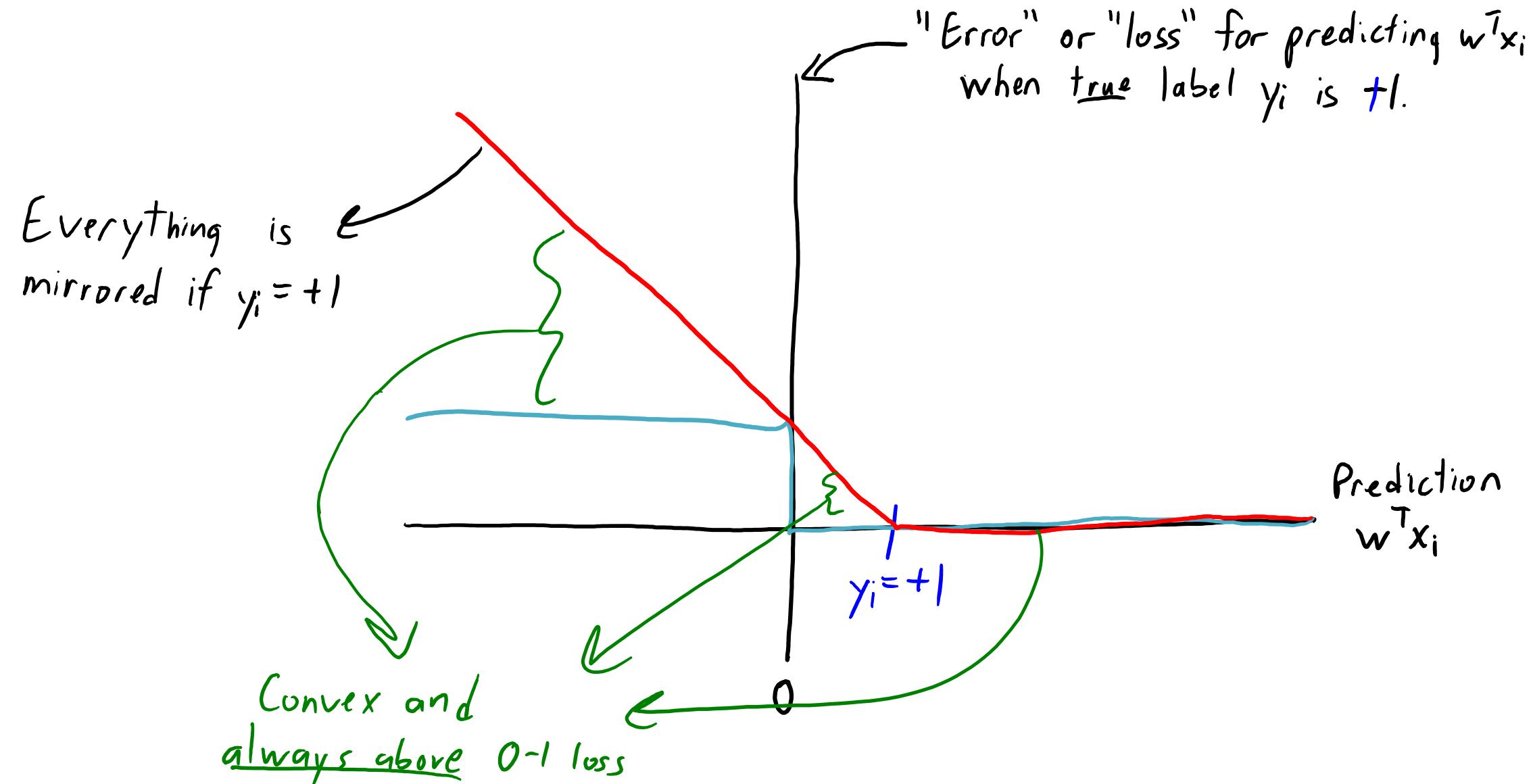
Hinge Loss: Convex Approximation to 0-1 Loss

Properties of the hinge loss:

1. Has error of 0 if $w^T x_i \leq -1$
(no penalty applied beyond this point)
2. Has a loss of 1 if $w^T x_i = 0$
(matches 0-1 loss at decision boundary)
3. Is convex and "close" to 0-1 loss.



Hinge Loss: Convex Approximation to 0-1 Loss



Hinge Loss

- Hinge loss for all 'n' training examples is given by:

$$f(w) = \sum_{i=1}^n \max\{0, 1 - y_i w^T x_i\}$$

- Convex upper bound on 0-1 loss.
 - If the hinge loss is 18.3, then number of training errors is at most 18.
 - So minimizing hinge loss indirectly tries to minimize training error.
 - Like perceptron, finds a perfect linear classifier if one exists.
- Support vector machine (SVM) is hinge loss with L2-regularization.

$$f(w) = \sum_{i=1}^n \max\{0, 1 - y_i w^T x_i\} + \frac{1}{2} \|w\|^2$$

- There exist specialized optimization algorithm for this problems.
- SVMs can also be viewed as “maximizing the margin” (later).

Summary

- Ensemble feature selection reduces false positives or negatives.
- Binary classification using regression:
 - Encode using y_i in $\{-1,1\}$.
 - Use $\text{sign}(w^T x_i)$ as prediction.
 - “Linear classifier” (a hyperplane splitting the space in half).
- Least squares is a weird error for classification.
- Perceptron algorithm: finds a perfect classifier (if one exists).
- 0-1 loss is the ideal loss, but is non-smooth and non-convex.
- Hinge loss is a convex upper bound on 0-1 loss.
 - SVMs add L2-regularization.
- Next time: one of the best “out of the box” classifiers.

bonus!

L1-Regularization as a Feature Selection Method

- Advantages:
 - Deals with conditional independence (if linear).
 - Sort of deals with collinearity:
 - Picks at least one of “mom” and “mom_again”.
 - Very fast with specialized algorithms.
- Disadvantages:
 - Tends to give **false positives** (selects too many variables).
- Neither good nor bad:
 - Does not take small effects.
 - Says “sex” is relevant if we know “baby”.
 - **Good for prediction if we want fast training and don’t care about having some irrelevant variables included.**

bonus!

“Elastic Net”: L2- and L1-Regularization

- To address **non-uniqueness**, sometimes **use both L2- and L1-**:

$$f(w) = \frac{1}{2} \|Xw - y\|^2 + \frac{\lambda_2}{2} \|w\|^2 + \lambda_1 \|w\|_1$$

- Called “**elastic net**” regularization.
 - Solution is **sparse and unique**.
 - Slightly better with feature dependence:
 - Selects both “mom” and “mom_again”.
- Optimization is easier, though still non-differentiable.

bonus!

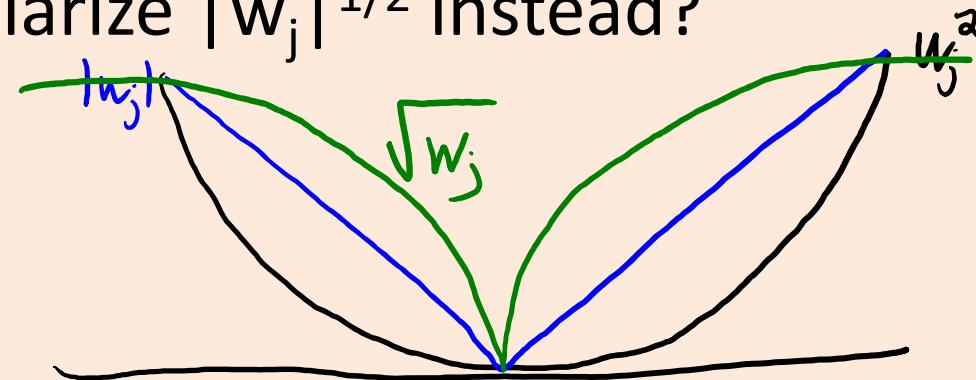
L1-Regularization Debiasing and Filtering

- To remove **false positives**, some authors add a **debiasing step**:
 - Fit ‘w’ using L1-regularization.
 - Grab the non-zero values of ‘w’ as the “relevant” variables.
 - Re-fit relevant ‘w’ using least squares or L2-regularized least squares.
- A related use of L1-regularization is as a **filtering method**:
 - Fit ‘w’ using L1-regularization.
 - Grab the non-zero values of ‘w’ as the “relevant” variables.
 - Run standard (slow) variable selection restricted to relevant variables.
 - Forward selection, exhaustive search, stochastic local search, etc.

bonus!

Non-Convex Regularizers

- Regularizing $|w_j|^2$ selects **all features**.
- Regularizing $|w_j|$ selects fewer, but still has many **false positives**.
- What if we regularize $|w_j|^{1/2}$ instead?

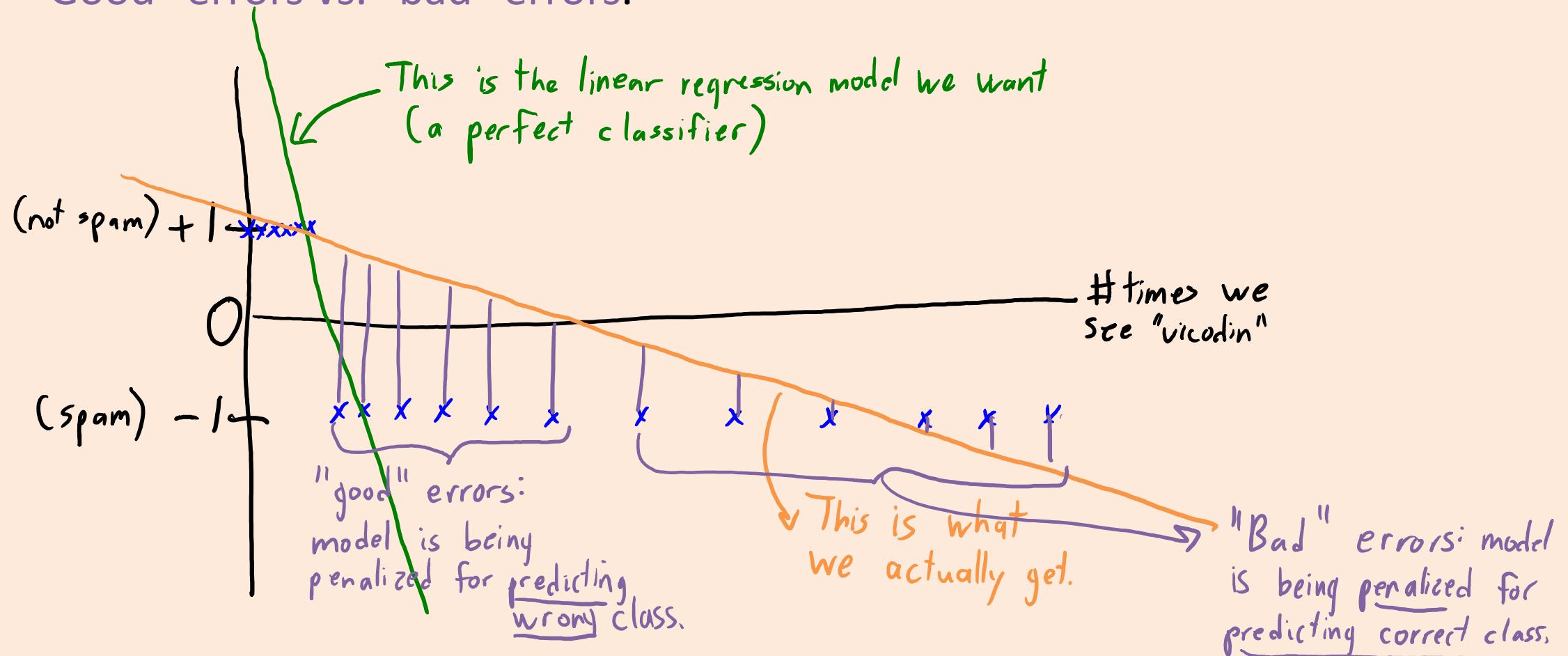


- Minimizing this objective would lead to **fewer false positives**.
 - Less need for debiasing, but it's not convex and **hard to minimize**.
- There are many non-convex regularizers with similar properties.
 - L1-regularization is (basically) the “most sparse” convex regularizer.

bonus!

Can we just use least squares??

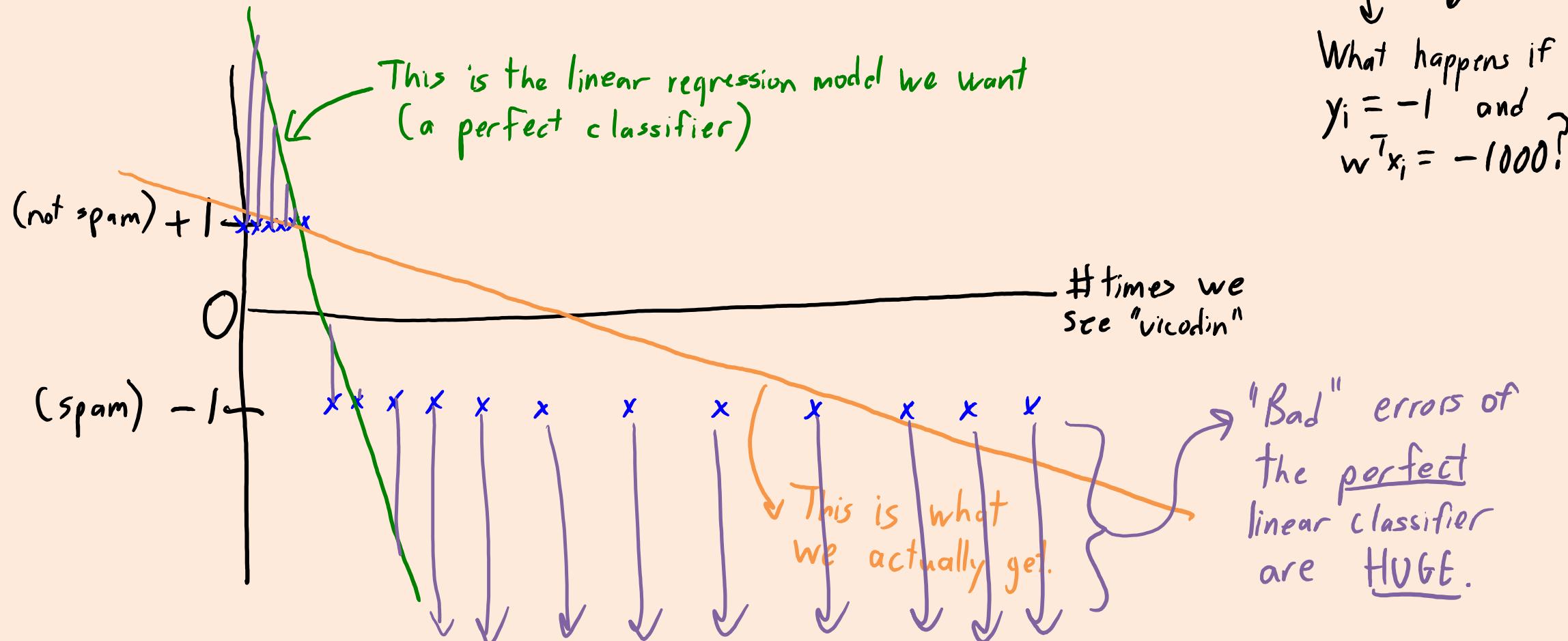
- What went wrong?
 - “Good” errors vs. “bad” errors.



bonus!

Can we just use least squares??

- What went wrong?
 - “Good” errors vs. “bad” errors.



bonus!

Online Classification with Perceptron

- Perceptron for online linear binary classification [Rosenblatt, 1957]
 - Start with $w_0 = 0$.
 - At time ‘t’ we receive features x_t .
 - We predict $\hat{y}_t = \text{sign}(w_t^T x_t)$.
 - If $\hat{y}_t \neq y_t$, then set $w_{t+1} = w_t + y_t x_t$.
 - Otherwise, set $w_{t+1} = w_t$.

(Slides are old so above I'm using subscripts of ‘t’ instead of superscripts.)

- Perceptron mistake bound [Novikoff, 1962]:
 - Assume data is linearly-separable with a “margin”:
 - There exists w^* with $\|w^*\|=1$ such that $\text{sign}(x_t^T w^*) = \text{sign}(y_t)$ for all ‘t’ and $|x^T w^*| \geq \gamma > 0$.
 - Then the number of total mistakes is bounded.
 - No requirement that data is IID.

bonus!

Perceptron Mistake Bound

- Let's normalize each x_t so that $\|x_t\| = 1$.
 - Length doesn't change label.
- Whenever we make a mistake, we have $\text{sign}(y_t) \neq \text{sign}(w_t^T x_t)$ and

$$\begin{aligned}\|w_{t+1}\|^2 &= \|w_t + yx_t\|^2 \\ &= \|w_t\|^2 + 2 \underbrace{y_t w_t^T x_t}_{<0} + 1 \\ &\leq \|w_t\|^2 + 1 \\ &\leq \|w_{t-1}\|^2 + 2 \\ &\leq \|w_{t-2}\|^2 + 3.\end{aligned}$$

- So after 'k' errors we have $\|w_t\|^2 \leq k$.

bonus!

Perceptron Mistake Bound

- Let's consider a solution w^* , so $\text{sign}(y_t) = \text{sign}(x_t^T w^*)$.
 - And let's choose a w^* with $\|w^*\| = 1$,
- Whenever we make a mistake, we have:

$$\begin{aligned}\|w_{t+1}\| &= \|w_{t+1}\| \|w^*\| \\ &\geq w_{t+1}^T w^* \\ &= (w_t + y_t x_t)^T w^* \\ &= w_t^T w^* + y_t x_t^T w^* \\ &= w_t^T w^* + |x_t^T w^*| \\ &\geq w_t^T w^* + \gamma.\end{aligned}$$

- Note: $w_t^T w^* \geq 0$ by induction (starts at 0, then at least as big as old value plus γ).
- So after 'k' mistakes we have $\|w_t\| \geq \gamma k$.

bonus!

Perceptron Mistake Bound

- So our two bounds are $\|w_t\| \leq \sqrt{k}$ and $\|w_t\| \geq \gamma k$.
- This gives $\gamma k \leq \sqrt{k}$, or a maximum of $1/\gamma^2$ mistakes.
 - Note that $\gamma > 0$ by assumption and is upper-bounded by one by $\|x\| \leq 1$.
 - After this ‘k’, under our assumptions we’re guaranteed to have a perfect classifier.

bonus!

Hinge-Loss Perceptron

- A **perceptron-like algorithm for minimizing the hinge loss:**
 - Start with any w^0 .
 - Go through examples until you **find an example with $y_i w^T x_i > 1$.**
 - Set $w^{t+1} = w^t + \frac{1 - y_i (w^t)^T x_i}{x_i^T x_i} y_i x_i$ (minimum change to w_t that satisfies constraint).
- **If a classifier with hinge loss of 0 exists,** this converges to one.
 - Looks like perceptron, but with a step size added to update (**green term**).
 - Get perceptron algorithm if you replace green term with '1'.
 - A special case of the "projection onto convex sets" (POCS) algorithm.