

Controlling Complexity: Feature Selection and Regularization

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Complexity of Hypothesis Spaces

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To control the “size” of \mathcal{F} , we need some measure of its **complexity**:

- Number of variables / features
- Degree of polynomial

General Approach to Control Complexity

1. Learn a sequence of models varying in complexity from the training data

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2. Select one of these models based on a score (e.g. validation error)

Feature Selection in Linear Regression

Nested sequence of hypothesis spaces: $\mathcal{F}_1 \subset \mathcal{F}_2 \subset \mathcal{F}_n \cdots \subset \mathcal{F}$

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Best subset selection:

- Choose the subset of features that is best according to the score (e.g. validation error)
 - Example with two features: Train models using $\{\}, \{X_1\}, \{X_2\}, \{X_1, X_2\}$, respectively
- **Not an efficient search algorithm**; iterating over all subsets becomes very expensive with a large number of features

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Backward Selection:

- Start with all features; in each iteration, remove the worst feature

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 - Find the subset of features that maximizes the score
- Forward & backward selection do not guarantee to find the best solution.
- Forward & backward selection do not in general result in the same subset.

ℓ_2 and ℓ_1 Regularization

Complexity Penalty

An objective that balances number of features and prediction performance:

$$\text{score}(S) = \text{training_loss}(S) + \lambda|S| \quad (1)$$

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λ balances the training loss and the number of features used:

- Adding an extra feature must be justified by at least λ improvement in training loss
- Larger $\lambda \rightarrow$ complex models are penalized more heavily

Complexity Penalty

Goal: Balance the complexity of the hypothesis space \mathcal{F} and the training loss

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For complexity measure $\Omega : \mathcal{F} \rightarrow [0, \infty)$ and fixed $\lambda \geq 0$,

$$\min_{f \in \mathcal{F}} \frac{1}{n} \sum_{i=1}^n \ell(f(x_i), y_i) + \lambda \Omega(f)$$

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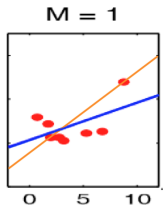
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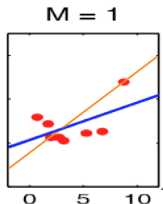
Number of features as complexity measure is hard to optimize—other measures?

Weight Shrinkage: Intuition



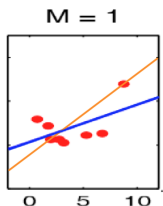
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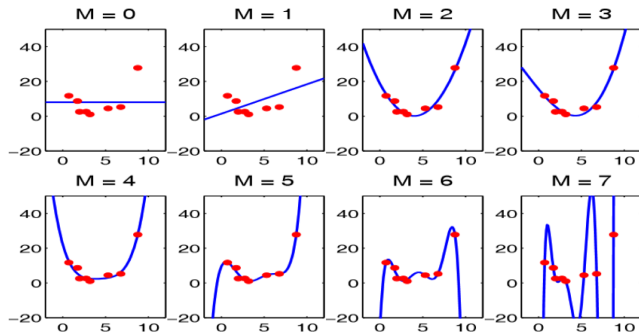
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- Why would we prefer a regression line with **smaller slope** (unless the data strongly supports a larger slope)?
- More conservative: small change in the input does not cause large change in the output
- If we push the estimated weights to be small, re-estimating them on a new dataset wouldn't cause the prediction function to change dramatically (**less sensitive to noise in data**)

Weight Shrinkage: Polynomial Regression



- Large weights are needed to make the curve wiggle sufficiently to overfit the data
- $\hat{y} = 0.001x^7 + 0.003x^3 + 1$ less likely to overfit than $\hat{y} = 1000x^7 + 500x^3 + 1$

(Adapted from Mark Schmidt's slide)

Linear Regression with ℓ_2 Regularization

- We have a linear model

$$\mathcal{F} = \{f : \mathbb{R}^d \rightarrow \mathbb{R} \mid f(x) = w^T x \text{ for } w \in \mathbb{R}^d\}$$

- Square loss: $\ell(\hat{y}, y) = (y - \hat{y})^2$
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- Training data $\mathcal{D}_n = ((x_1, y_1), \dots, (x_n, y_n))$
- Linear least squares regression is ERM for square loss over \mathcal{F} :

$$\hat{w} = \arg \min_{w \in \mathbb{R}^d} \frac{1}{n} \sum_{i=1}^n (w^T x_i - y_i)^2$$

- This often overfits, especially when d is large compared to n (e.g. in NLP one can have 1M features for 10K documents).

Linear Regression with L2 Regularization

Penalizes large weights:

$$\hat{w} = \arg \min_{w \in \mathbb{R}^d} \frac{1}{n} \sum_{i=1}^n \{w^T x_i - y_i\}^2 + \lambda \|w\|_2^2,$$

where $\|w\|_2^2 = w_1^2 + \dots + w_d^2$ is the square of the ℓ_2 -norm.

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- Also known as **ridge regression**.
- Equivalent to linear least square regression when $\lambda = 0$.
- ℓ_2 regularization can be used for other models too (e.g. neural networks).

ℓ_2 regularization reduces sensitivity to changes in input

- $\hat{f}(x) = \hat{w}^T x$ is **Lipschitz continuous** with Lipschitz constant $L = \|\hat{w}\|_2$: when moving from x to $x + h$, \hat{f} changes no more than $L\|h\|$.

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- Other norms also provide a bound on L due to the equivalence of norms:
 $\exists C > 0$ s.t. $\|\hat{w}\|_2 \leq C \|\hat{w}\|_p$

Linear Regression vs. Ridge Regression

Objective:

- Linear: $L(w) = \frac{1}{2} \|Xw - y\|_2^2$
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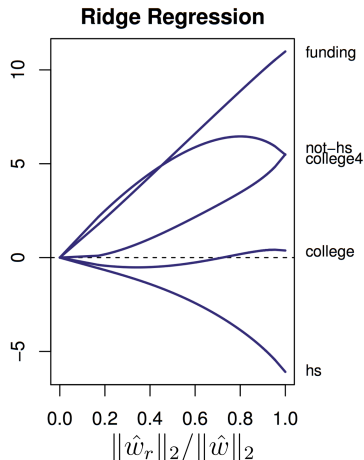
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Closed-form solution:

- Linear: $X^T X w = X^T y$
- Ridge: $(X^T X + \lambda I) w = X^T y$
 - $(X^T X + \lambda I)$ is always invertible

Ridge Regression: Regularization Path



$$\hat{w}_r = \arg \min_{\|w\|_2^2 \leq r^2} \frac{1}{n} \sum_{i=1}^n (w^T x_i - y_i)^2$$
$$\hat{w} = \hat{w}_\infty = \text{Unconstrained ERM}$$

- For $r = 0$, $\|\hat{w}_r\|_2 / \|\hat{w}\|_2 = 0$.
- For $r = \infty$, $\|\hat{w}_r\|_2 / \|\hat{w}\|_2 = 1$

Modified from Hastie, Tibshirani, and Wainwright's *Statistical Learning with Sparsity*, Fig 2.1. About predicting crime in 50 US cities.

Lasso Regression

Penalize the ℓ_1 norm of the weights:

Lasso Regression (Tikhonov Form, soft penalty)

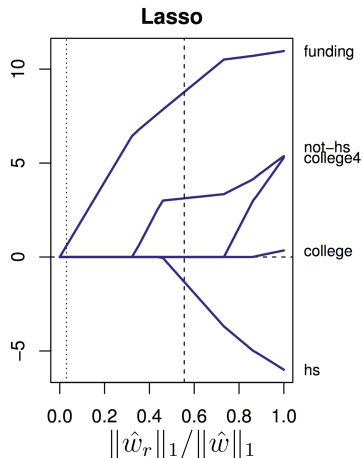
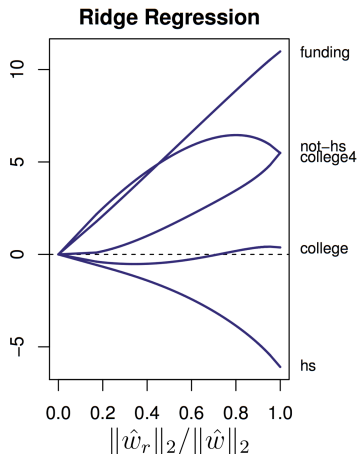
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where $\|w\|_1 = |w_1| + \dots + |w_d|$ is the ℓ_1 -norm.

(“Least Absolute Shrinkage and Selection Operator”)

Ridge vs. Lasso: Regularization Paths

Lasso yields sparse weights:



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- Prediction function may generalize better (model is less complex)

Why does ℓ_1 Regularization Lead to Sparsity?

Regularization as Constrained Empirical Risk Minimization

Constrained ERM (Ivanov regularization)

For complexity measure $\Omega : \mathcal{F} \rightarrow [0, \infty)$ and fixed $r \geq 0$,

$$\begin{aligned} \min_{f \in \mathcal{F}} \frac{1}{n} \sum_{i=1}^n \ell(f(x_i), y_i) \\ \text{s.t. } \Omega(f) \leq r \end{aligned}$$

Lasso Regression (Ivanov Form, hard constraint)

The lasso regression solution for complexity parameter $r \geq 0$ is

$$\hat{w} = \arg \min_{\|w\|_1 \leq r} \frac{1}{n} \sum_{i=1}^n \{w^T x_i - y_i\}^2.$$

r has the same role as λ in penalized ERM (Tikhonov).

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- In practice, both approaches are effective: we will use whichever one is more convenient for training or analysis.

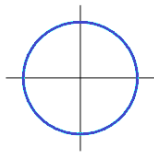
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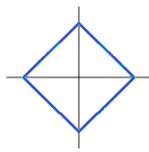
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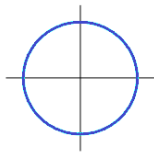
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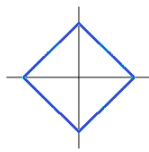
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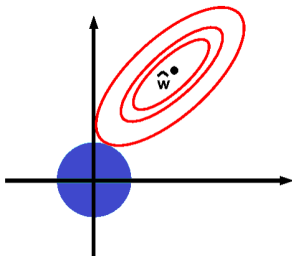
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- Where are the sparse solutions?

Visualizing Regularization

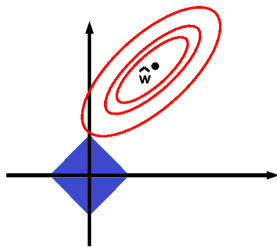
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- Blue region: Area satisfying complexity constraint: $w_1^2 + w_2^2 \leq r$
- Red lines: contours of the empirical risk $\hat{R}_n(w) = \sum_{i=1}^n (w^T x_i - y_i)^2$.

Why Does ℓ_1 Regularization Encourage Sparse Solutions?

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- ℓ_1 solution tends to touch the **corners**.

Why Does ℓ_1 Regularization Encourage Sparse Solutions?

Geometric intuition: Projection onto diamond encourages solutions at corners.

- \hat{w} in red/green regions are closest to corners in the ℓ_1 “ball”.

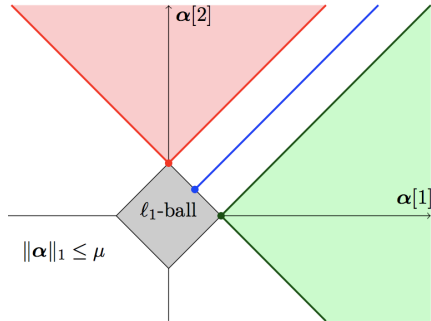


Fig from Mairal et al.'s Sparse Modeling for Image and Vision Processing Fig 1.6

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Geometric intuition: Projection onto ℓ_2 sphere favors all directions equally.

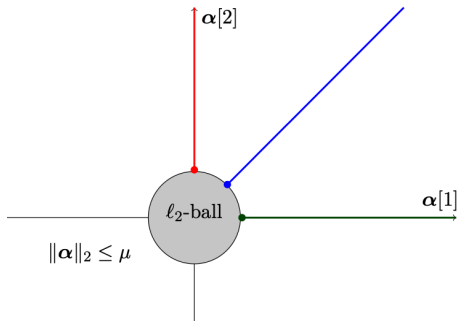


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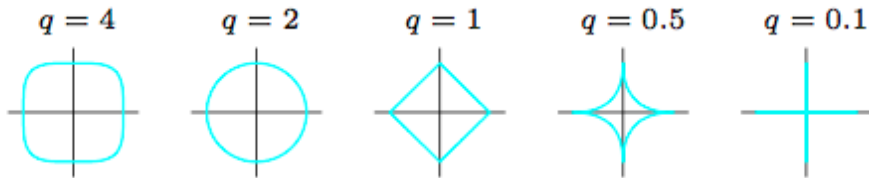
- The gradient stays the same as the weights approach zero
- This pushes the weights to be exactly zero even if they are already small

(ℓ_q) Regularization

- We can generalize to ℓ_q : $(\|w\|_q)^q = |w_1|^q + |w_2|^q$.

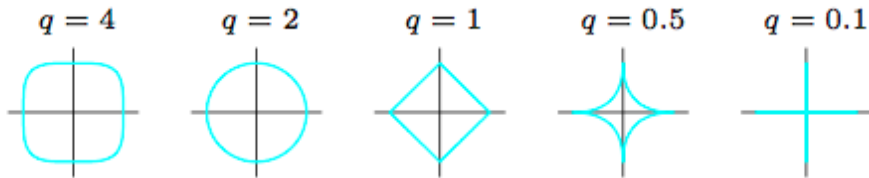
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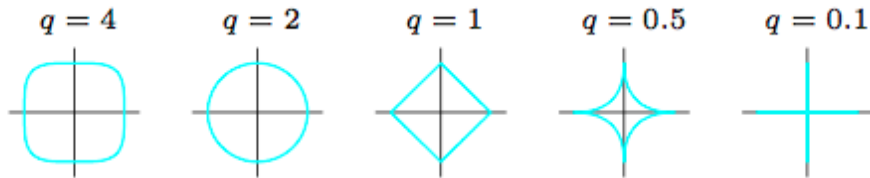
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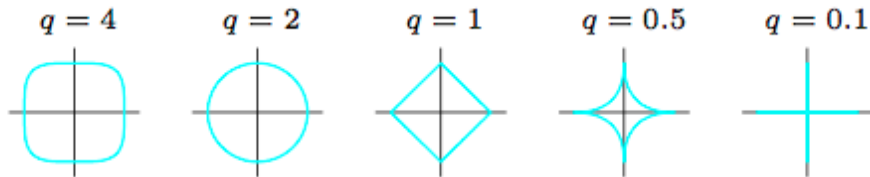
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- When $q < 1$, the ℓ_q constraint is non-convex, so it is hard to optimize; lasso is good enough in practice
- ℓ_0 ($\|w\|_0$) is defined as the number of non-zero weights, i.e. subset selection

Minimizing the lasso objective

Minimizing the lasso objective

- The ridge regression objective is differentiable (and there is a closed form solution)
- Lasso objective function:

$$\min_{w \in \mathbb{R}^d} \sum_{i=1}^n (w^T x_i - y_i)^2 + \lambda \|w\|_1$$

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- $\|w\|_1 = |w_1| + \dots + |w_d|$ is not differentiable!
- We will briefly review three approaches for finding the minimum:
 - Quadratic programming
 - Projected SGD
 - Coordinate descent

Rewriting the Absolute Value

- Consider any number $a \in \mathbb{R}$.
- Let the **positive part** of a be

$$a^+ = a1(a \geq 0).$$

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- Is it always the case that $a^+ \geq 0$ and $a^- \geq 0$?
- How do you write a in terms of a^+ and a^- ?
- How do you write $|a|$ in terms of a^+ and a^- ?

The Lasso as a Quadratic Program

Substituting $w = w^+ - w^-$ and $|w| = w^+ + w^-$ results in an **equivalent** problem:

$$\begin{aligned} \min_{w^+, w^-} \quad & \sum_{i=1}^n \left((w^+ - w^-)^T x_i - y_i \right)^2 + \lambda \mathbf{1}^T (w^+ + w^-) \\ \text{subject to} \quad & w_i^+ \geq 0 \text{ for all } i \quad \text{and} \quad w_i^- \geq 0 \text{ for all } i, \end{aligned}$$

- This objective is **differentiable** (in fact, **convex and quadratic**)
- How many variables does the new objective have?
- This is a **quadratic program**: a convex quadratic objective with linear constraints.
- Quadratic programming is a very well understood problem; we can plug this into a generic QP solver.

Are we missing some constraints?

We have claimed that the following objective is equivalent to the lasso problem:

$$\begin{aligned} \min_{w^+, w^-} \quad & \sum_{i=1}^n \left((w^+ - w^-)^T x_i - y_i \right)^2 + \lambda \mathbf{1}^T (w^+ + w^-) \\ \text{subject to} \quad & w_i^+ \geq 0 \text{ for all } i \quad w_i^- \geq 0 \text{ for all } i, \end{aligned}$$

- When we plug this optimization problem into a QP solver,
 - it just sees $2d$ variables and $2d$ constraints.
 - Doesn't know we want w_i^+ and w_i^- to be positive and negative parts of w_i .
- Turns out that these constraints will be satisfied anyway!
- To make it clear that the solver isn't aware of the constraints of w_i^+ and w_i^- , let's denote them a_i and b_i

The Lasso as a Quadratic Program

(Trivially) reformulating the lasso problem:

$$\begin{aligned} \min_w \min_{a,b} \quad & \sum_{i=1}^n \left((a-b)^T x_i - y_i \right)^2 + \lambda \mathbf{1}^T (a+b) \\ \text{subject to} \quad & a_i \geq 0 \text{ for all } i \quad b_i \geq 0 \text{ for all } i, \\ & a - b = w \\ & a + b = |w| \end{aligned}$$

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Claim: Don't need the constraint $a + b = |w|$.

Exercise: Prove by showing that the optimal solutions a^* and b^* satisfies $\min(a^*, b^*) = 0$, hence $a^* + b^* = |w|$.

The Lasso as a Quadratic Program

$$\begin{aligned} \min_w \min_{a,b} \quad & \sum_{i=1}^n \left((a-b)^T x_i - y_i \right)^2 + \lambda \mathbf{1}^T (a+b) \\ \text{subject to} \quad & a_i \geq 0 \text{ for all } i \quad b_i \geq 0 \text{ for all } i, \\ & a - b = w \end{aligned}$$

Claim: Can remove \min_w and the constraint $a - b = w$.

Exercise: Prove by switching the order of the minimization.

Projected SGD

- Now that we have a differentiable objective, we could also use gradient descent
- But how do we handle the **constraints**?

$$\min_{w^+, w^- \in \mathbb{R}^d} \sum_{i=1}^n \left((w^+ - w^-)^T x_i - y_i \right)^2 + \lambda \mathbf{1}^T (w^+ + w^-)$$

subject to $w_i^+ \geq 0$ for all i
 $w_i^- \geq 0$ for all i

- Projected SGD is just like SGD, but after each step
 - We project w^+ and w^- into the constraint set.
 - In other words, if any component of w^+ or w^- becomes negative, we set it back to 0.

Coordinate Descent Method

Goal: Minimize $L(w) = L(w_1, \dots, w_d)$ over $w = (w_1, \dots, w_d) \in \mathbb{R}^d$.

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Coordinate Descent Method

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- In gradient descent or SGD, each step potentially changes **all entries** of w .
- In **coordinate descent**, each step adjusts only a **single coordinate** w_i .

$$w_i^{\text{new}} = \arg \min_{w_i} L(w_1, \dots, w_{i-1}, w_i, w_{i+1}, \dots, w_d)$$

- Solving the argmin for a particular coordinate may itself be an iterative process.
- Coordinate descent is an effective method when it's easy (or easier) to minimize w.r.t. one coordinate at a time

Coordinate Descent Method

Goal: Minimize $L(w) = L(w_1, \dots, w_d)$ over $w = (w_1, \dots, w_d) \in \mathbb{R}^d$.

- **Initialize** $w^{(0)} = 0$
- **while** not converged:
 - Choose a coordinate $j \in \{1, \dots, d\}$
 - $w_j^{\text{new}} \leftarrow \arg \min_{w_j} L(w_1^{(t)}, \dots, w_{j-1}^{(t)}, w_j, w_{j+1}^{(t)}, \dots, w_d^{(t)})$
 - $w_j^{(t+1)} \leftarrow w_j^{\text{new}}$ and $w^{(t+1)} \leftarrow w^{(t)}$
 - $t \leftarrow t + 1$
- Random coordinate choice \implies **stochastic coordinate descent**
- Cyclic coordinate choice \implies **cyclic coordinate descent**

Coordinate Descent Method for Lasso

The lasso objective coordinate minimization has a closed form! If

$$\hat{w}_j = \arg \min_{w_j \in \mathbb{R}} \sum_{i=1}^n (w^T x_i - y_i)^2 + \lambda |w|_1$$

Then

$$\hat{w}_j = \begin{cases} (c_j + \lambda)/a_j & \text{if } c_j < -\lambda \\ 0 & \text{if } c_j \in [-\lambda, \lambda] \\ (c_j - \lambda)/a_j & \text{if } c_j > \lambda \end{cases}$$

$$a_j = 2 \sum_{i=1}^n x_{i,j}^2$$

$$c_j = 2 \sum_{i=1}^n x_{i,j} (y_i - w_{-j}^T x_{i,-j})$$

where w_{-j} is w without the j -th component, and $x_{i,-j}$ is x_i without the j -th component.

Coordinate Descent in General

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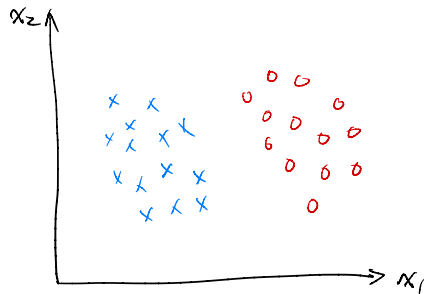
Coordinate Descent in General

- In general, coordinate descent is not competitive with gradient descent: its convergence rate is slower and the iteration cost is similar
- But it works very well for certain problems
- Very simple and easy to implement
- Example applications: lasso regression, SVMs

Maximum Margin Classifier

Linearly Separable Data

Consider a linearly separable dataset \mathcal{D} :



Find a separating hyperplane such that

- $w^T x_i > 0$ for all x_i where $y_i = +1$
- $w^T x_i < 0$ for all x_i where $y_i = -1$

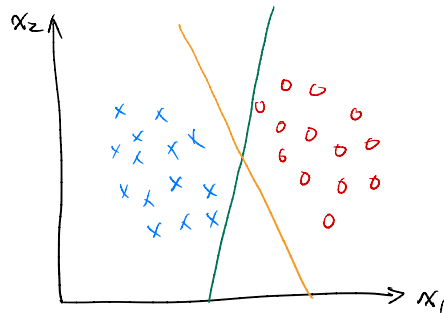
The Perceptron Algorithm

- Initialize $w \leftarrow 0$
- While not converged (exists misclassified examples)
 - For $(x_i, y_i) \in \mathcal{D}$
 - If $y_i w^T x_i < 0$ (wrong prediction)
 - Update $w \leftarrow w + y_i x_i$
- Intuition: move towards misclassified positive examples and away from negative examples
- Guarantees to find a zero-error classifier (if one exists) in finite steps
- What is the loss function if we consider this as a SGD algorithm?

Maximum-Margin Separating Hyperplane

For separable data, there are infinitely many zero-error classifiers.

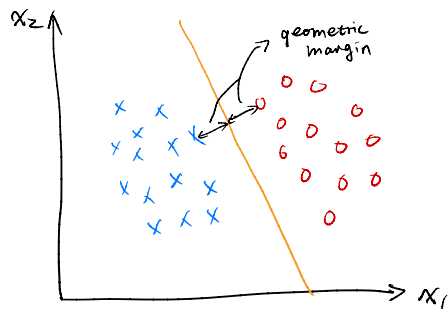
Which one do we pick?



(Perceptron does not return a unique solution.)

Maximum-Margin Separating Hyperplane

We prefer the classifier that is farthest from both classes of points



- Geometric margin: smallest distance between the hyperplane and the points
- Maximum margin: *largest* distance to the closest points

Geometric Margin

We want to maximize the distance between the **separating hyperplane** and the **closest** points.

Let's formalize the problem.

Definition (separating hyperplane)

We say (x_i, y_i) for $i = 1, \dots, n$ are **linearly separable** if there is a $w \in \mathbb{R}^d$ and $b \in \mathbb{R}$ such that $y_i(w^T x_i + b) > 0$ for all i . The set $\{v \in \mathbb{R}^d \mid w^T v + b = 0\}$ is called a **separating hyperplane**.

Definition (geometric margin)

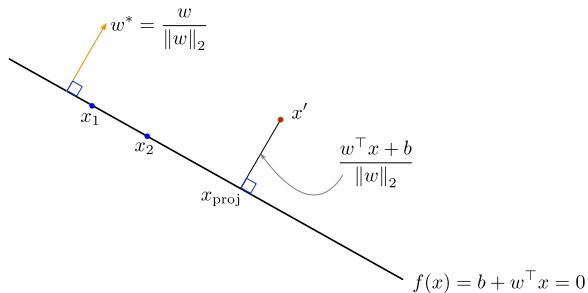
Let H be a hyperplane that separates the data (x_i, y_i) for $i = 1, \dots, n$. The **geometric margin** of this hyperplane is

$$\min_i d(x_i, H),$$

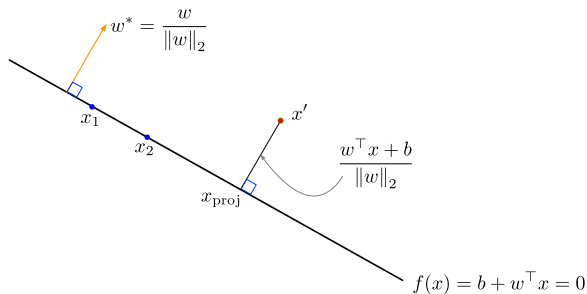
the distance from the hyperplane to the closest data point.

Distance between a Point and a Hyperplane

- Any point on the plane p , and normal vector $w/\|w\|_2$

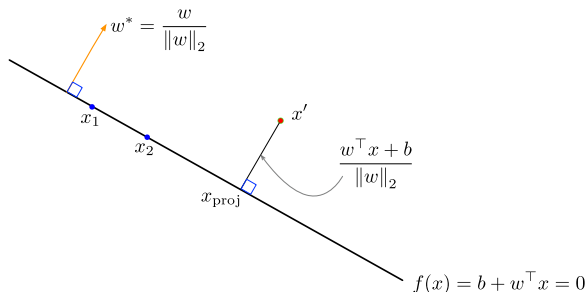


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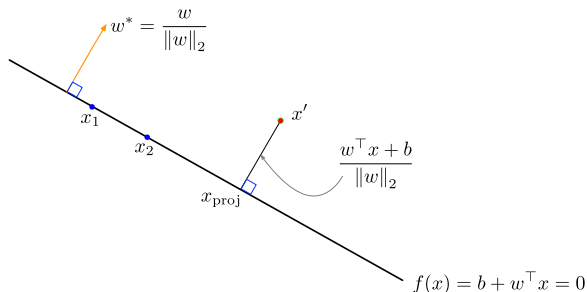
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- Projection of x onto the normal: $\frac{(x'-p)^T w}{\|w\|_2}$

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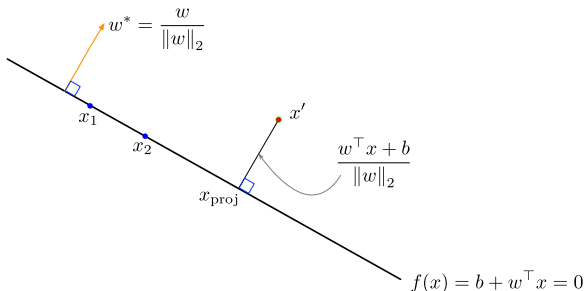
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- Signed distance between x' and Hyperplane H : $\frac{w^T x' + b}{\|w\|_2}$
- Taking into account of the label y :
$$d(x', H) = \frac{y(w^T x' + b)}{\|w\|_2}$$

Maximize the Margin

We want to maximize the geometric margin:

$$\text{maximize } \min_i d(x_i, H).$$

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Let's remove the inner minimization problem by

$$\begin{aligned} &\text{maximize} && M \\ &\text{subject to} && \frac{y_i(w^T x_i + b)}{\|w\|_2} \geq M \quad \text{for all } i \end{aligned}$$

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Note that the solution is not unique (why?).

Maximize the Margin

Let's fix the norm $\|w\|_2$ to $1/M$ to obtain:

$$\begin{array}{ll}\text{maximize} & \frac{1}{\|w\|_2} \\ \text{subject to} & y_i(w^T x_i + b) \geq 1 \quad \text{for all } i\end{array}$$

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It's equivalent to solving the minimization problem

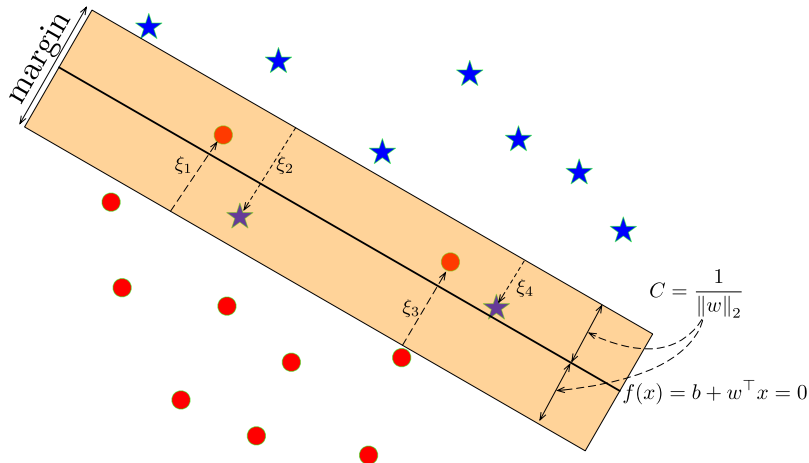
$$\begin{array}{ll}\text{minimize} & \frac{1}{2} \|w\|_2^2 \\ \text{subject to} & y_i(w^T x_i + b) \geq 1 \quad \text{for all } i\end{array}$$

Note that $y_i(w^T x_i + b)$ is the (functional) margin. The optimization finds the minimum norm solution which has a margin of at least 1 on all examples.

Not linearly separable

What if the data is *not* linearly separable?

For any w , there will be points with a negative margin.



Introduce **slack variables** ξ 's to penalize small margin:

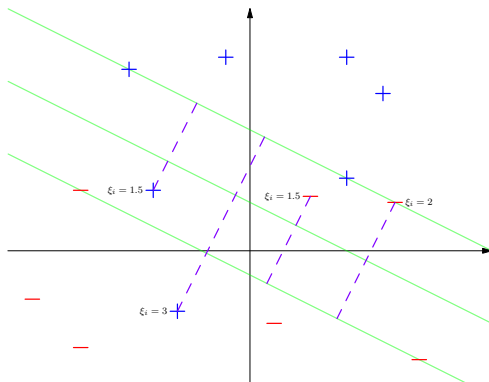
$$\begin{aligned} & \text{minimize} && \frac{1}{2} \|w\|_2^2 + \frac{C}{n} \sum_{i=1}^n \xi_i \\ & \text{subject to} && y_i (w^T x_i + b) \geq 1 - \xi_i \quad \text{for all } i \\ & && \xi_i \geq 0 \quad \text{for all } i \end{aligned}$$

- If $\xi_i = 0 \forall i$, it's reduced to hard SVM.
- What does $\xi_i > 0$ mean?
- What does C control?

Slack Variables

$d(x_i, H) = \frac{y_i(w^T x_i + b)}{\|w\|_2} \geq \frac{1 - \xi_i}{\|w\|_2}$, thus ξ_i measures the violation by multiples of the geometric margin:

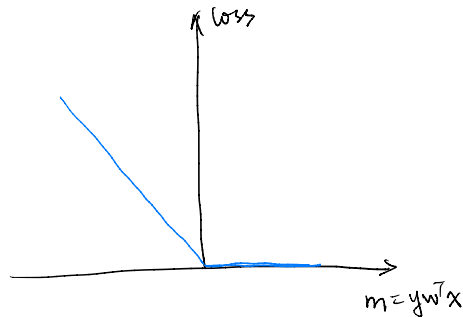
- $\xi_i = 1$: x_i lies on the hyperplane
- $\xi_i = 3$: x_i is past 2 margin width beyond the decision hyperplane



Minimize the Hinge Loss

Perceptron Loss

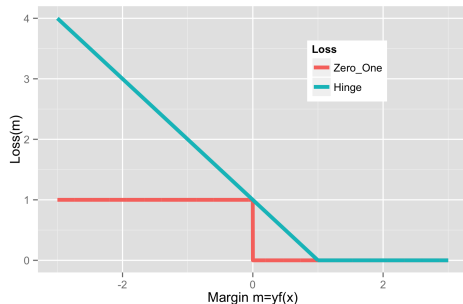
$$\ell(x, y, w) = \max(0, -yw^T x)$$



If we do ERM with this loss function, what happens?

Hinge Loss

- SVM/Hinge loss: $\ell_{\text{Hinge}} = \max\{1 - m, 0\} = (1 - m)_+$
- Margin $m = yf(x)$; “Positive part” $(x)_+ = x1(x \geq 0)$.



Hinge is a **convex, upper bound** on 0–1 loss. Not differentiable at $m = 1$.
We have a “margin error” when $m < 1$.

SVM as an Optimization Problem

- The SVM optimization problem is equivalent to

$$\begin{aligned} \text{minimize} \quad & \frac{1}{2} \|w\|^2 + \frac{c}{n} \sum_{i=1}^n \xi_i \\ \text{subject to} \quad & \xi_i \geq (1 - y_i [w^T x_i + b]) \text{ for } i = 1, \dots, n \\ & \xi_i \geq 0 \text{ for } i = 1, \dots, n \end{aligned}$$

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which is equivalent to

$$\begin{aligned} \text{minimize} \quad & \frac{1}{2} \|w\|^2 + \frac{c}{n} \sum_{i=1}^n \xi_i \\ \text{subject to} \quad & \xi_i \geq \max(0, 1 - y_i [w^T x_i + b]) \text{ for } i = 1, \dots, n. \end{aligned}$$

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Move the constraint into the objective:

$$\min_{w \in \mathbb{R}^d, b \in \mathbb{R}} \frac{1}{2} \|w\|^2 + \frac{c}{n} \sum_{i=1}^n \max(0, 1 - y_i [w^T x_i + b]).$$

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- The first term is the L2 regularizer.
- The second term is the Hinge loss.

Support Vector Machine

Using ERM:

- Hypothesis space $\mathcal{F} = \{f(x) = w^T x + b \mid w \in \mathbb{R}^d, b \in \mathbb{R}\}$.
- ℓ_2 regularization (Tikhonov style)
- Hinge loss $\ell(m) = \max\{1 - m, 0\} = (1 - m)_+$
- The SVM prediction function is the solution to

$$\min_{w \in \mathbb{R}^d, b \in \mathbb{R}} \frac{1}{2} \|w\|^2 + \frac{c}{n} \sum_{i=1}^n \max(0, 1 - y_i [w^T x_i + b]).$$

- **Not differentiable** because of the max

Two ways to derive the SVM optimization problem:

- Maximize the margin
- Minimize the hinge loss with ℓ_2 regularization

Both leads to the minimum norm solution satisfying certain margin constraints.

- **Hard-margin SVM:** all points must be correctly classified with the margin constraints
- **Soft-margin SVM:** allow for margin constraint violation with some penalty