Statistical Machine Learning

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Learning from Data

In the real world, p(x, y) is unknown, but we have a training set \mathcal{D} . There's at least 3 approaches:

Definition

Given a training set \mathcal{D} , we call it

• a generative probabilistic approach: if we use \mathcal{D} to build a model $\hat{p}(x, y)$ of p(x, y), and then define

$$c(x) := \operatorname*{argmax}_{y \in \mathcal{Y}} \hat{p}(x, y) \quad \text{or} \quad c_{\ell}(x) := \operatorname*{argmin}_{y \in \mathcal{Y}} \mathbb{E}_{\bar{y} \sim \hat{p}(x, \bar{y})} \ell(\bar{y}, y).$$

• a discriminative probabilistic approach: if we use \mathcal{D} to build a model $\hat{p}(y|x)$ of p(y|x) and define

$$c(x) := \operatorname*{argmax}_{y \in \mathcal{Y}} \hat{p}(y|x) \quad \text{or} \quad c_{\ell}(x) := \operatorname*{argmin}_{y \in \mathcal{Y}} \mathbb{E}_{\bar{y} \sim \hat{p}(\bar{y}|x)} \ell(\bar{y}, y).$$

• a **decision theoretic approach**: if we use \mathcal{D} to directly seach for a classifier c in a hypothesis class \mathcal{H} .

Task: spam classification, $\mathcal{X} = \{\text{all possible emails}\}, \mathcal{Y} = \{\text{spam}, \text{ham}\}.$ What's, e.g., p(x|ham)? For every possible email, a value how likely it is to see that email, including:

- all possible languages,
- all possbile topics,
- an arbitrary length,
- all possible spelling mistakes, etc.

This is much more general (and much harder) than just deciding if an email is spam or not!

"When solving a problem, do not solve a more general problem as an intermediate step." (Vladimir Vapnik, 1998)

Instead of p(x, y) = p(x|y)p(y), we can also use p(x, y) = p(y|x)p(x). Since $\operatorname{argmax}_y p(x, y) = \operatorname{argmax}_y p(y|x)$, we don't need to model p(x), only p(y|x).

Let's use \mathcal{D} to estimate p(y|x).

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Let's use \mathcal{D} to estimate p(y|x).

Visual intuition:



class conditional densities = likelihood p(x|y)

joint density likelihood*prior: p(x|y)p(y)

class posteriors p(y|x)=p(x|y)p(y)/p(x)

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Let's use \mathcal{D} to estimate p(y|x).

Example (Spam Classification)

Is p(y|x) really easier than, e.g., p(x|y)?

- p("v1agra"|spam) is some positive value (not every spam is viagra)
- p(spam | "v1agra") is almost surely 1.

For p(y|x) we treat x as given, we don't need to know its probability.

Nonparametric Discriminative Model

Idea: split \mathcal{X} into regions, for each region store an estimate $\hat{p}(y|x)$.

p(1 x)=0.7 p(2 x)=0.2 p(3 x)=0.1	p(1 x)=0.9 p(2 x)=0.0 p(3 x)=0.1	X
	p(1 x)=0.1 p(2 x)=0.8 p(3 x)=0.1	
	p(1 x)=0.01 $p(2 x)=0.98p(3 x)=0.01$	

Idea: split \mathcal{X} into regions, for each region store an estimate $\hat{p}(y|x)$.

For example, using a decision tree:

- training: build a tree
- prediction: for new example x, find its leaf
- output $\hat{p}(y|x) = \frac{n_y}{n}$, where
 - n is the number of examples in the leaf,
 - n_y is the number of example of label y in the leaf.

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, where

- n is the number of examples in the leaf,
- n_y is the number of example of label y in the leaf.

Note: prediction rule

$$c(x) = \operatorname*{argmax}_{y} \hat{p}(y|x)$$

is predicts the most frequent label in each leaf (same as in first lecture).

Setting. We assume
$$\mathcal{X} \subseteq \mathbb{R}^d$$
 and $\mathcal{Y} = \{-1, +1\}$.

Definition (Logistic Regression (LogReg) Model)

Modeling

$$\hat{p}(y|x;w) = rac{1}{1 + \exp(-y\langle w, x \rangle)},$$

with parameter vector $w \in \mathbb{R}^d$ is called a *logistic regression* model.

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Lemma

 $\hat{p}(y|x; w)$ is a well defined probability density w.r.t. y for any $w \in \mathbb{R}^d$.

Proof. elementary.

Logistic Regression Training

Given a training set $\mathcal{D} = \{(x^1, y^1), \dots, (x^n, y^n)\}$, logistic regression training sets the free parameter vector as

$$w_{LR} = \operatorname*{\mathbf{argmin}}_{w \in \mathbb{R}^d} \; \sum_{i=1}^n \log \left(1 + \exp(-y^i \langle w, x^i \rangle)
ight)$$

Lemma (Conditional Likelihood Maximization)

 w_{LR} from Logistic Regression training maximizes the conditional data likelihood w.r.t. the LogReg model,

$$w_{LR} = \operatorname*{argmax}_{w \in \mathbb{R}^d} \hat{p}(y^1, \dots, y^n | x^1, \dots, x^n, w)$$

Maximizing

$$\hat{p}(\mathcal{D}^{Y}|\mathcal{D}^{X},w) \stackrel{i.i.d.}{=} \prod_{i=1}^{n} \hat{p}(y^{i}|x^{i},w)$$

is equivalent to minimizing its negative logarithm

$$-\log \hat{p}(\mathcal{D}^{Y}|\mathcal{D}^{X}, w) = -\log \prod_{i=1}^{n} \hat{p}(y^{i}|x^{i}, w) = -\sum_{i=1}^{n} \log \hat{p}(y^{i}|x^{i}, w)$$
$$= -\sum_{i=1}^{n} \log \frac{1}{1 + \exp(-y^{i}\langle w, x^{i} \rangle)},$$
$$= -\sum_{i=1}^{n} [\log 1 - \log(1 + \exp(-y^{i}\langle w, x^{i} \rangle)],$$
$$= \sum_{i=1}^{n} \log(1 + \exp(-y^{i}\langle w, x^{i} \rangle)).$$

Definition (Kullback-Leibler (KL) divergence)

Let p and q be two probability distributions (for discrete Z) or probability densities with respect to a measure $d\lambda$ (for continuous Z). The **Kullbach-Leibler (KL)-divergence** between p and q is defined as

$$\operatorname{KL}(p \| q) = \sum_{z \in \mathcal{Z}} p(z) \log \frac{p(z)}{q(z)}, \quad \text{or} \quad \operatorname{KL}(p \| q) = \int_{z \in \mathcal{Z}} p(z) \log \frac{p(z)}{q(z)} \, \mathrm{d}\lambda(\mathbf{z}),$$

(with convention $0 \log 0 = 0$, and $a \log \frac{a}{0} = \infty$ for a > 0).

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 KL is a similarity measure between probability distributions. It fulfills

$$0 \le KL(p \, \| \, q) \le \infty,$$
 and $KL(p \, \| \, q) = 0 \iff p = q.$

However, KL is not a metric.

- it is in general not symmetric, $KL(q \| p) \neq KL(p \| q)$,
- it does not fulfill the triangle inequality.

Definition (Expected Kullback-Leibler (KL) divergence)

Let p(x, y) be a probability distribution over $(x, y) \in \mathcal{X} \times \mathcal{Y}$ and let $\hat{p}(y|x)$ be an approximation of p(y|x). We measure the approximation quality by the **expected KL-divergence**

between p and q over all $x \in \mathcal{X}$:

$$\mathrm{KL}_{exp}(p \| q) = \mathbb{E}_{x \sim p(x)} \{ \mathrm{KL}(p(\cdot | x) \| q(\cdot | x)) \}$$

Theorem

The parameter w_{LR} obtained by logistic regression training approximately minimizes the KL divergence between $\hat{p}(y|x;w)$ and p(y|x).

We show how maximimzing the conditional likelihood relates to KL_{exp} :

$$\begin{split} \mathrm{KL}_{exp}(p \| \hat{p}) &= \mathbb{E}_{x \sim p(x)} \sum_{y \in \mathcal{Y}} p(y|x) \log \frac{p(y|x)}{\hat{p}(y|x,w)} \\ &= \underbrace{\mathbb{E}_{(x,y) \sim p(x,y)} \log p(y|x)}_{\text{indep. of } w} - \mathbb{E}_{(x,y) \sim p(x,y)} \log \hat{p}(y|x,w) \end{split}$$

We can't maximize $\mathbb{E}_{(x,y)\sim p(x,y)}\log \hat{p}(y|x,w)$ directly, because p(x,y) is unknown. But we can maximize its empirical estimate based on \mathcal{D} :

$$\mathbb{E}_{(x,y)\sim p(x,y)}\log \hat{p}(y|x,w) \approx \underbrace{\sum_{(x^i,y^i)\in\mathcal{D}}\log \hat{p}(y^i|x^i,w)}_{\log \text{ of conditional data likelihood}}$$

The approximation will get better the more data we have.

Theorem

Logistic Regression training,

$$w_{LR} = \operatorname*{argmin}_{w \in \mathbb{R}^d} \mathcal{L}(w) \quad \textit{for} \quad \mathcal{L}(w) = \sum_{i=1}^n \log\left(1 + \exp(-y^i \langle w, x^i \rangle)\right),$$

is a C^{∞} -smooth, unconstrained, convex optimization problem.

Proof.

- 1. it's an optimization problem,
- 2. it's unconstrained,
- 3. it's smooth (the objective function is C^{∞} differentiable),
- remains to show: the objective function is a *convex* function. Since L is smooth, it's enough to show that its *Hessian matrix* (the matrix of 2nd partial derivatives) is everywhere *positive definite*.

We compute first the gradient and then the Hessian of

$$\mathcal{L}(w) = \sum_{i=1}^{n} \log(1 + \exp(-y^{i} \langle w, x^{i} \rangle)).$$
$$\nabla_{w} \mathcal{L}(w) = \sum_{i=1}^{n} \nabla \log(1 + \exp(-y^{i} \langle w, x^{i} \rangle)).$$

use the chain rule, $\nabla f(g(w)) = \frac{d}{dt}(g(w)\nabla g(w), \text{ and } \frac{d}{dt}\log(t) = \frac{1}{t}$

$$=\sum_{i=1}^{n} \frac{\nabla[1 + \exp(-y^{i}\langle w, x^{i}\rangle]}{1 + \exp(-y^{i}\langle w, x^{i}\rangle)}$$
$$=\sum_{i=1}^{n} \underbrace{\frac{\exp(-y^{i}\langle w, x^{i}\rangle)}{1 + \exp(-y^{i}\langle w, x^{i}\rangle)}}_{=\hat{p}(-y^{i}|x^{i},w)} \nabla(-y^{i}\langle w, x^{i}\rangle)$$

use the chain rule again, $\frac{d}{dt}\exp(t)=\exp(t),$ and $\nabla_{\!w}\langle w,x^i\rangle=x^i$

$$= -\sum_{i=1}^{n} [\hat{p}(-y^{i}|x^{i}, w)] \ y^{i}x^{i}$$

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$$\begin{split} H_w \mathcal{L}(w) &= \nabla \nabla^\top \mathcal{L}(w) = -\sum_{i=1}^n [\nabla \hat{p}(-y^i | x^i, w)] \ y^i x^i \\ \nabla \hat{p}(-y^i | x^i, w) &= \nabla \frac{1}{1 + \exp(y^i \langle w, x^i \rangle)} \\ &= -\frac{\nabla [1 + \exp(y^i \langle w, x^i \rangle)]}{[1 + \exp(y^i \langle w, x^i \rangle)]^2} \\ \text{use quotient rule, } \nabla \frac{1}{f(w)} &= -\frac{\nabla f(w)}{f^2(w)}, \text{ and chain rule,} \\ &= -\frac{\exp(y^i \langle w, x^i \rangle)}{[1 + \exp(y^i \langle w, x^i \rangle)]^2} \nabla y^i \langle w, x^i \rangle \\ &= -(\hat{p}(-y^i | x^i)) \hat{p}(y^i | x^i, w) y^i x^i \end{split}$$

insert into above expression for $H_w \mathcal{L}(w)$

$$H = \sum_{i=1}^{n} \underbrace{\hat{p}(-y^{i}|x^{i})\hat{p}(y^{i}|x^{i},w)}_{>0} \underbrace{\underbrace{x^{i}x^{i\top}}_{sym.pos.\,def.}$$

A positively weighted linear combination of pos.def. matrices is pos.def. $_{\rm ff_{4/39}}$



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Convex optimization is a well understood field. We can use, e.g., *gradient descent* will converge to the globally optimal solution!

Steepest Descent Minimization with Line Search

 $\begin{array}{ll} \mbox{input} & \epsilon > 0 \mbox{ tolerance (for stopping criterion)} \\ 1: & w \leftarrow 0 \\ 2: \mbox{ repeat} \\ 3: & v \leftarrow -\nabla_w \, \mathcal{L}(w) & \{\mbox{descent direction}\} \\ 4: & \eta \leftarrow \mbox{argmin}_{\eta > 0} \, \mathcal{L}(w + \eta v) & \{\mbox{1D line search}\} \\ 5: & w \leftarrow w + \eta d \\ 6: \mbox{ until } \|v\| < \epsilon \\ \mbox{output } w \in \mathbb{R}^d \mbox{ learned weight vector} \\ \end{array}$

Faster conference from methods that use second-order information, e.g., conjugate gradients or (L-)BFGS \rightarrow convex optimization lecture

Binary classification with a LogReg Models

A discriminative probability model, $\hat{p}(y|x)$, is enough to make decisions:

$$c(x) = \operatorname*{argmax}_{y \in \mathcal{Y}} \hat{p}(y|x) \quad \text{or} \quad c(x) = \operatorname*{argmin}_{y \in \mathcal{Y}} \mathbb{E}_{\bar{y} \sim \hat{p}(y|x)} \ell(\bar{y}, y).$$

For Logistic Regression, this is particularly simple:

Lemma

The LogReg classification rule for 0/1-loss is

$$c(x) = \operatorname{sign} \langle w, x \rangle.$$
For a loss function $\ell = \begin{pmatrix} a & b \\ c & d \end{pmatrix}$ the rule is
$$c_{\ell}(x) = \operatorname{sign}[\langle w, x \rangle + \log \frac{c-d}{b-a}],$$

In particular, the decision boundaries is linear (or affine).

Proof. Elementary, since $\log \frac{\hat{p}(+1|x;w)}{p(-1|x;w)} = \langle w, x \rangle$

Multiclass Logistic Regression

For $\mathcal{Y} = \{1, \dots, M\}$, we can do two things:

• Parametrize $\hat{p}(y|x; \vec{w})$ using M-1 vectors, $w_1, \ldots, w_{M-1} \in \mathbb{R}^d$, as

$$\hat{p}(y|x,w) = \frac{\exp(\langle w_y, x \rangle)}{1 + \sum_{j=1}^{M-1} \exp(\langle w_j, x \rangle)} \quad \text{for } y = 1, \dots, M-1,$$
$$\hat{p}(M|x,w) = \frac{1}{1 + \sum_{j=1}^{M-1} \exp(\langle w_j, x \rangle)}.$$

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$$\hat{p}(y|x,w) = \frac{\exp(\langle w_y, x \rangle)}{\sum_{j=1}^{M} \exp(\langle w_j, x \rangle)} \quad \text{for } y = 1, \dots, M,$$

Second is more popular, since it's easier to implement and analyze. Decision boundaries are still *piecewise linear*, $c(x) = \operatorname{argmax}_{u} \langle w_{y}, x \rangle$. Discriminative models treats the input data, x, as fixed and only model the distribution of the output labels p(y|x).

Discriminative models, in particular LogReg, are popular, because

- they often need less training data than generative models,
- they provide an estimate of the uncertainty of a decision p(c(x)|x),
- training them is often efficient,
 e.g. Yahoo trains LogReg models routinely from billions of examples.
- But: they also have drawbacks
 - often $\hat{p}_{LR}(y|x) \not\rightarrow p(y|x)$, even for $n \rightarrow \infty$,
 - they usually are good for *prediction*, but they do not reflect the actual *mechanism*.

Note: there are much more complex discriminative models than LogReg, e.g. Conditional Random Fields (maybe later).

Maximum Margin Classifiers

Let's use \mathcal{D} to estimate a classifier $c: \mathcal{X} \to \mathcal{Y}$ directly.

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For a start, we fix

- $\mathcal{D} = \{(x^1, y^1), \dots, (x^n, y^n)\},\$
- $\mathcal{Y} = \{+1, -1\},\$
- we look for classifiers with linear decision boundary.

Several of the classifiers we saw had *linear* decision boundaries:

- Perceptron
- Generative classifiers for Gaussian class-conditional densities with shared covariance matrix
- Logistic Regression

What's the best linear classifier?

Definition

Let

$$\mathcal{F} = \{ f : \mathbb{R}^d \to \mathbb{R} \text{ with } f(x) = b + a_1 x_1 + \dots + a_d x_d = b + \langle w, x \rangle \}$$

be the set of linear (affine) function from $\mathbb{R}^d \to \mathbb{R}$.

A classifier $g: \mathcal{X} \to \mathcal{Y}$ is called **linear**, if it can be written as

 $g(x) = \operatorname{sign} f(x)$

for some $f \in \mathcal{F}$.

We write \mathcal{G} for the set of all linear classifiers.

A linear classifier, $g(x) = \operatorname{sign} \langle w, x \rangle$, with b = 0



A linear classifier $g(x) = \operatorname{sign}(\langle w, x \rangle + b)$, with b > 0



The bias term is good for intuition, but annoying in analysis:

Useful trick: feature augmentation

Adding a constant feature allows us to avoid models with explicit bias term:

- instead of
$$x=(x^1,\ldots,x^d)\in \mathbb{R}^d$$
, use $\tilde{x}=(x^1,\ldots,x^d,1)\in \mathbb{R}^{d+1}$

• for any $\tilde{w} \in \mathbb{R}^{d+1}$, think $\tilde{w} = (w, b)$ with $w \in \mathbb{R}^d$ and $b \in \mathbb{R}$

Linear function in \mathbb{R}^{d+1} :

$$f(\tilde{x}) = \langle \tilde{w}, \tilde{x} \rangle = \sum_{i=1}^{d+1} \tilde{w}_i \tilde{x}_i = \sum_{i=1}^d \tilde{w}_i \tilde{x}_i + \tilde{w}_{d+1} \tilde{x}_{d+1} = \langle w, x \rangle + b$$

Linear classifier with bias in $\mathbb{R}^d \equiv$ linear classifier with no bias in \mathbb{R}^{d+1}

Augmenting with other (larger) values than 1 can make sense, see later... $^{24/39}$

Definition (Ad hoc)

We call a classifier, g, **correct** (for a training set D), if it predicts the correct labels for all training examples:

$$g(x^i) = y^i$$
 for $i = 1, ..., n$.

Example (Perceptron)

- if the *Perceptron* converges, the result is an *correct* classifier.
- any classifier with zero training error is *correct*.

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Definition (Linear Separability)

A training set \mathcal{D} is called **linearly separable**, if it allows a correct linear classifier (i.e. the classes can be separated by a hyperplane).

A linearly separable dataset and a correct classifier



A linearly separable dataset and a correct classifier



A linearly separable dataset and a correct classifier



An incorrect classifier



Definition (Ad hoc)

The **robustness** of a classifier g (with respect to \mathcal{D}) is the largest amount, ρ , by which we can perturb the training samples without changing the predictions of g.

$$g(x^i + \epsilon) = g(x^i),$$
 for all $i = 1, \dots, n$.

for any $\epsilon \in \mathbb{R}^d$ with $\|\epsilon\| < \rho$.

Example

- constant classifier, e.g. c(x) ≡ 1: very robust (ρ = ∞), (but it is not *correct*, in the sense of the previous definition)
- robustness of the *Perceptron*: can be arbitrarily small (see Exercise...)

Robustness, $\rho\text{, of a linear classifier}$



Definition (Margin)

Let $f(x) = \langle w, x \rangle + b$ define a *correct* linear classifier. Then the smallest (Euclidean) distance of any training example from the decision hyperplane is called the **margin** of f (with respect to \mathcal{D}).

Lemma

We can compute the margin of a linear classifier $f = \langle w, x \rangle + b$ as

$$\rho = \min_{i=1,\dots,n} \left| \left\langle \frac{w}{\|w\|}, x^i \right\rangle + b \right|.$$

Proof.

High school maths: distance between a points and a hyperplane in *Hessian normal form.*

Margin, ρ , of a linear classifier



Theorem

The robustness of a linear classifier function $g(x) = \operatorname{sign} f(x)$ with $f(x) = \langle w, x \rangle$ is identical to the margin of f.

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Proof (blackboard). For any i = 1, ..., n and any $\epsilon \in \mathbb{R}^d$

$$f(x^{i} + \epsilon) = \langle w, x^{i} + \epsilon \rangle = \langle w, x^{i} \rangle + \langle w, \epsilon \rangle = f(x^{i}) + \langle w, \epsilon \rangle,$$

so it follows (Cauchy-Schwarz inequality) that

$$f(x^{i}) - ||w|| ||\epsilon|| \le f(x^{i} + \epsilon) \le f(x^{i}) + ||w|| ||\epsilon||.$$

Checking the cases $\epsilon = \pm \frac{\|\epsilon\|}{\|w\|} w$, we see that these inequalities are sharp. To ensure $g(x^i + \epsilon) = g(x^i)$ for all training samples, $f(x^i)$ and $f(x^i + \epsilon)$ have the same sign for all ϵ , i.e. $|f(x^i)| \ge \|w\| \|\epsilon\|$ for i = 1, ..., n.

This inequality holds for all samples, so in particular it holds for the one of minimal score, and $\min_i |f(x^i)| = \min_i |\langle w, x^i \rangle| = \rho$.

Theorem

Let \mathcal{D} be a linearly separable training set. Then the **most robust**, correct linear classifier (without bias term) is given by $g(x) = \operatorname{sign} \langle w^*, x \rangle$ where w^* are the solution to

$$\min_{w\in\mathbb{R}^d} \; rac{1}{2} \|w\|^2$$

subject to

$$y^i(\langle w, x^i \rangle) \ge 1$$
, for $i = 1, \dots, n$.

Remark

- The classifier defined above is call Maximum (Hard) Margin Classifier, or Hard-Margin Support Vector Machine (SVM)
- It is unique (follows from strictly convex optimization problem).

- **1.** All *w* that fulfill the inequalities yield *correct* classifiers.
- **2.** Since \mathcal{D} is linearly separable, there exists some v with

$$\operatorname{sign}\langle v, x^i \rangle = y_i, \quad \text{i.e.} \quad y_i \langle v, x^i \rangle \ge \gamma > 0.$$

for $\gamma = \min_i y_i \langle v, x^i \rangle$. So $\tilde{v} = v/\gamma$, fulfills the inequalities and we see that the constraint set is at least not empty.

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3. Now we check (with
$$i = 1, ..., n$$
):

$$\min_{w \in \mathbb{R}^d} \frac{1}{2} ||w||^2 \text{ sb.t. } y^i \langle w, x^i \rangle \ge 1$$

$$\Leftrightarrow \max_{w \in \mathbb{R}^d} \frac{1}{||w||} \quad \text{ sb.t. } y^i \langle w, x^i \rangle \ge 1$$

$$\Leftrightarrow \max_{\{w':|w'||=1\}, \rho \in \mathbb{R}} \quad \rho \quad \text{ sb.t. } y^i \langle \frac{w'}{\rho}, x^i \rangle \ge 1$$

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$$\Leftrightarrow \max_{\{w':|w'||=1\}, \rho \in \mathbb{R}} \quad \rho \quad \text{ sb.t. } |\langle w', x^i \rangle| \ge \rho \text{ and } \operatorname{sign}\langle w', x^i \rangle = y_i$$

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$$\min_{w \in \mathbb{R}^{d}} \frac{1}{2} ||w||^{2} \text{ sb.t. } y^{i} \langle w, x^{i} \rangle \geq 1$$

$$\Leftrightarrow \max_{w \in \mathbb{R}^{d}} \frac{1}{||w||} \quad \text{ sb.t. } y^{i} \langle w, x^{i} \rangle \geq 1$$

$$\Leftrightarrow \max_{\{w':|w'||=1\}, \rho \in \mathbb{R}} \rho \quad \text{ sb.t. } y^{i} \langle w', x^{i} \rangle \geq 1$$

$$\Leftrightarrow \max_{\{w':|w'||=1\}, \rho \in \mathbb{R}} \rho \quad \text{ sb.t. } y^{i} \langle w', x^{i} \rangle \geq \rho$$

$$\Leftrightarrow \underbrace{\max_{\{w':|w'||=1\}, \rho \in \mathbb{R}} \rho \quad \text{ sb.t. } |\langle w', x^{i} \rangle| \geq \rho}_{\text{maximal robustness}} \text{ and } \underbrace{\operatorname{sign}\langle w', x^{i} \rangle = y_{i}}_{\text{ and correct}}$$

Non-Separable Training Sets

Observation (Not all training sets are linearly separable.)



Definition (Maximum Soft-Margin Classifier)

Let \mathcal{D} be a training set, not necessarily linearly separable. Let C > 0. Then the classifier $g(x) = \operatorname{sign} \langle w^*, x \rangle$ where (w^*, b^*) are the solution to

$$\min_{w \in \mathbb{R}^d, \xi \in \mathbb{R}^n} \quad \frac{1}{2} \|w\|^2 + C \sum_{i=1}^n \xi^i$$

subject to

$$y^{i}(\langle w, x^{i} \rangle + b) \ge 1 - \xi^{i}, \quad \text{for } i = 1, \dots, n.$$

$$\xi^{i} \ge 0, \quad \text{for } i = 1, \dots, n.$$

is called Maximum (Soft-)Margin Classifier or Linear Support Vector Machine.

Theorem

The maximum soft-margin classifier exists and is unique for any C > 0.

Proof. optimization problem is strictly convex

Remark

The constant C > 0 is called **regularization** parameter.

It balances the wishes for robustness and for correctness

- $C \rightarrow 0$: mistakes don't matter much, emphasis on short w
- $C \rightarrow \infty$: as few errors as possible, might not be robust

We'll see more about this in the next lecture.

Remark

Sometimes, a soft margin is better even for linearly separable datasets!

