### Regression recap

Recall that in regression we are given training data

$$(\mathbf{x}_1,y_1),\ldots,(\mathbf{x}_n,y_n)$$

where  $\mathbf{x}_i \in \mathbb{R}^d$  and  $y_i \in \mathbb{R}$ 

In linear regression we assume that we are trying to estimate a function of the form

$$f(\mathbf{x}) = \boldsymbol{\beta}^T \mathbf{x} + \beta_0$$

where  $oldsymbol{eta} \in \mathbb{R}^d$  ,  $eta_0 \in \mathbb{R}$ 

**Least squares regression:** Select  $\beta$ ,  $\beta_0$  to minimize

$$SSE(\beta, \beta_0) := \sum_{i=1}^n (y_i - \beta^T \mathbf{x}_i - \beta_0)^2$$

### Regularization and regression

Overfitting occurs as the number of features  $\,d\,$  begins to approach the number of observations  $\,n\,$ 

In this regime, we have too many degrees of freedom

Idea: penalize candidate solutions for using too many features

One candidate regularizer:  $r(\theta) = \|\theta\|_2^2$ 

$$\widehat{\boldsymbol{\theta}} = \operatorname*{arg\,min}_{\boldsymbol{\theta}} \| \boldsymbol{y} - \boldsymbol{A} \boldsymbol{\theta} \|_2^2 + \lambda \| \boldsymbol{\theta} \|_2^2$$

 $\lambda > 0$  is a "tuning parameter" that controls the tradeoff between fit and complexity

### Least squares

$$y = \begin{bmatrix} y_1 \\ y_2 \\ \vdots \\ y_n \end{bmatrix} A = \begin{bmatrix} 1 & x_1(1) & \cdots & x_1(d) \\ 1 & x_2(1) & \cdots & x_2(d) \\ \vdots & \vdots & \ddots & \vdots \\ 1 & x_n(1) & \cdots & x_n(d) \end{bmatrix} \theta = \begin{bmatrix} \beta_0 \\ \beta(1) \\ \vdots \\ \beta(d) \end{bmatrix}$$

$$SSE(\boldsymbol{\theta}) = \sum_{i=1}^{n} (y_i - \boldsymbol{\beta}^T \mathbf{x}_i - \beta_0)^2 = \|\boldsymbol{y} - \boldsymbol{A}\boldsymbol{\theta}\|_2^2$$

Minimizer given by

$$\widehat{m{ heta}} = \left(m{A}^Tm{A}
ight)^{-1}m{A}^Tm{y}$$

provided that  $oldsymbol{A}^Toldsymbol{A}$  is **nonsingular** 

### Tikhonov regularization

This is one example of a more general technique called *Tikhonov regularization* 

$$\widehat{m{ heta}} = rg\min_{m{ heta}} \|m{y} - m{A}m{ heta}\|_2^2 + \|\Gammam{ heta}\|_2^2$$

(Note that  $\lambda$  has been replaced by the matrix  $\Gamma$ )



#### **Solution:**

$$\widehat{\theta} = (A^T A + \Gamma^T \Gamma)^{-1} A^T y$$

always well-conditioned

$$\Gamma = \sqrt{\lambda} I$$
  $\widehat{\boldsymbol{\theta}} = \left( \underline{\boldsymbol{A}}^T \boldsymbol{A} + \lambda \boldsymbol{I} \right)^{-1} \boldsymbol{A}^T \boldsymbol{y}$  for suitable choice of  $\lambda$  .

### Ridge regression

In the context of regression, Tikhonov regularization has a special name: *ridge regression* 

Ridge regression is essentially exactly what we have been talking about, but in the special case where

$$\Gamma = \begin{bmatrix} 0 & 0 & 0 & \cdots & 0 \\ 0 & \sqrt{\lambda} & 0 & \cdots & 0 \\ 0 & 0 & \sqrt{\lambda} & \cdots & 0 \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & 0 & \cdots & \sqrt{\lambda} \end{bmatrix}$$

We are penalizing all coefficients in  $\beta$  equally, but not penalizing the offset  $\beta_0$ 

### The LASSO

LASSO

$$\widehat{m{ heta}} = rg \min_{m{ heta}} \|m{y} - m{A}m{ heta}\|_2^2 + \lambda \|m{ heta}\|_1$$

Can also be stated in a constrained form

$$\begin{split} \widehat{\theta} &= \mathop{\arg\min}_{\theta} \| \boldsymbol{y} - \boldsymbol{A}\boldsymbol{\theta} \|_2^2 \qquad \widehat{\boldsymbol{\theta}} = \mathop{\arg\min}_{\theta} \| \boldsymbol{\theta} \|_1 \\ \text{s.t.} \quad \| \boldsymbol{\theta} \|_1 &\leq \tau \qquad \qquad \text{s.t.} \quad \| \boldsymbol{y} - \boldsymbol{A}\boldsymbol{\theta} \|_2^2 \leq \sigma \end{split}$$

For Tikhonov, we have a closed form solution, but LASSO requires solving an optimization problem

**Note:** Just like in ridge regression, in practice we may just want to penalize the elements of  $\beta$  (not  $\beta_0$ )

### Alternative regularizers

- Akaike information criterion (AIC)
- Bayesian information criterion (BIC)

$$r(\theta) \approx \|\theta\|_0 := |\mathsf{supp}(\theta)|$$

• Least absolute shrinkage and selection operator (LASSO)

$$r(\theta) = \|\theta\|_1 = \sum_j |\theta(j)|$$

- also results in shrinkage, but where all coordinates are shrunk by the same amount (in case of an orthobasis)
- promotes sparsity
- can think of  $\| \boldsymbol{\theta} \|_1$  as a more computationally tractable replacement for  $\| \boldsymbol{\theta} \|_0$

### Sparsity and the LASSO

One can show (see supplemental notes) that if we have a data set of size n, then the solution to the LASSO  $\widehat{\theta}$  will have at most n nonzeros (for any possible dataset /A)

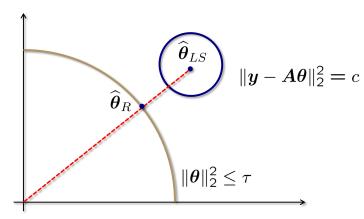
This is a nice property when  $n \ll d$  , since in this setting we are very susceptible to overfitting

- fewer observations than unknowns
- $oldsymbol{A}$  has nontrivial nullspace
- we can achieve  $y=A\theta$ , with infinitely many different choices of  $\theta$  and no obvious way to know which one is best
- limiting the number of nonzeros addresses this problem

In practice, the number of nonzeros is usually much smaller than  $\boldsymbol{n}$ 

### Tikhonov versus least squares

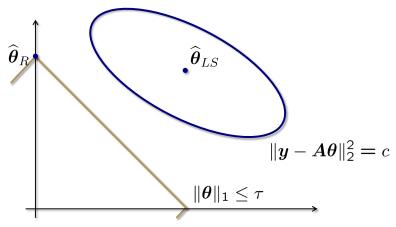
Assume  $\Gamma=I$  and that A has orthonormal columns



Tikhonov regularization is equivalent to shrinking the least squares solution towards the origin

### Lasso versus least squares

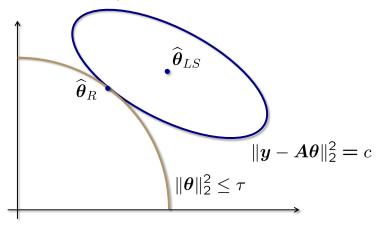
For the LASSO we get something like this...



LASSO still shrinking the least squares solution towards the origin, but now in a way that promotes sparsity

### Tikhonov versus least squares

In general, we have this picture



Tikhonov regularization still shrinking the least squares solution towards the origin

### A general approach to regression

Least squares, ridge regression, and the LASSO call all be viewed as particular instances of the following general approach to regression

$$\hat{\theta} = \underset{\theta}{\operatorname{arg\,min}} L(\theta) + \lambda r(\theta)$$

- $L(m{ heta})$ , often called the *loss function*, enforces data fidelity  $f_{m{ heta}}(\mathbf{x}_i) pprox y_i$
- $r(\theta)$  is a **regularizer** which serves to quantify the "complexity" of  $\theta$

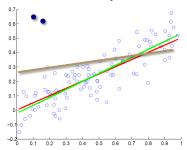
We have seen some examples of regularizers, what about other loss functions?

### Outliers in regression

The squared error loss function is sensitive to *outliers* 

If  $f(\mathbf{x}_i) - y_i$  is small, then  $(f(\mathbf{x}_i) - y_i)^2$  is not too large But if  $f(\mathbf{x}_i) - y_i$  is big, then  $(f(\mathbf{x}_i) - y_i)^2$  is **really** big

Normally this is not a bad property - we want to penalize big errors - but this can make us very sensitive to large outliers



# Regularized robust regression

Suppose we combine this loss with an  $\ell_2$  regularizer

$$\widehat{\boldsymbol{\beta}}, eta_0 = \operatorname*{arg\,min}_{(oldsymbol{eta}, eta_0)} \ \sum_{i=1}^n L_{\epsilon}(y_i - (oldsymbol{eta}^T \mathbf{x}_i + eta_0)) + rac{\lambda}{2} \|oldsymbol{eta}\|_2^2$$

Note that the  $\epsilon\text{-insensitive}$  loss has no penalty as long as your prediction is within a "margin" of  $\epsilon$ 

This looks like an SVM...

### Robust regression

What else could we do aside from least squares?

Mean absolute error

$$L_{AE}(r) = |r|$$

**Huber loss** 

$$L_H(r) = egin{cases} rac{1}{2}r^2 & ext{if } |r| \leq c \ c|r| - rac{c^2}{2} & ext{if } |r| > c \end{cases}$$

 $\epsilon$  -insensitive loss

$$L_{\epsilon}(r) = \begin{cases} 0 & \text{if } |r| \le c \\ |r| - \epsilon & \text{if } |r| > \epsilon \end{cases}$$

### Support vector regression

The previous problem can also be cast in the following dual form

$$\min_{\alpha,\alpha^*} \sum_{i} ((\epsilon - y_i)\alpha_i^* + (\epsilon + y_i)\alpha_i) + \frac{1}{2} \sum_{i,j} (\alpha_i^* - \alpha_i)(\alpha_j^* - \alpha_j) \mathbf{x}_j^T \mathbf{x}_i$$

subject to 
$$0 \le \alpha_i^*, \alpha_i \le \frac{1}{\lambda}$$

$$\sum_{i} (\alpha_i^* - \alpha_i) = 0$$

$$\alpha_i^*\alpha_i=0$$

The solution has the form  $\hat{f}(\mathbf{x}) = \sum_{i} (\hat{\alpha}_{i}^{*} - \hat{\alpha}_{i}) \mathbf{x}_{i}^{T} \mathbf{x} + \beta_{0}$ 

### Can we kernelize regression?

In order to "kernelize" an algorithm, the general approach consists of three main steps:

- 1. Show that the training process only involves the training data via inner products (i.e.,  $\mathbf{x}_i^T \mathbf{x}_j$ )
- 2. Show that applying the decision rule to a new x only involves computing inner products (i.e.,  $\mathbf{w}^T \mathbf{x}$ )
- 3. Replace all inner products with evaluations of the kernel function  $k(\cdot, \cdot)$

This approach extends well beyond SVMs

### Kernelized LASSO?

Can we kernelize this?

$$\widehat{ heta} = rg \min_{ heta} \|y - A heta\|_2^2 + \lambda \| heta\|_1$$

Not exactly...

Nevertheless, we can assert (with no justification) that

$$\widehat{\boldsymbol{\theta}} = \sum_{i} \alpha_{i} \mathbf{x}_{i}$$

and then replace  $\| heta\|_1$  with  $\|lpha\|_1$ 

This yields an algorithm that can be easily kernelized, although it is really something different than the LASSO

- promotes sparsity in  $\alpha$ , not  $\theta$ 

### Kernel support vector regression

$$\begin{aligned} \min_{\alpha,\alpha^*} & \sum_{i} \left( (\epsilon - y_i) \alpha_i^* + (\epsilon + y_i) \alpha_i \right) + \frac{1}{2} \sum_{i,j} (\alpha_i^* - \alpha_i) (\alpha_j^* - \alpha_j) \mathbf{x}_j^T \mathbf{x}_i \\ \text{subject to} & 0 \leq \alpha_i^*, \alpha_i \leq \frac{1}{\lambda} \\ & \sum_{i} (\alpha_i^* - \alpha_i) = 0 \\ & \alpha_i^* \alpha_i = 0 \end{aligned}$$

$$\hat{f}(\mathbf{x}) = \sum_{i} (\hat{\alpha}_i^* - \hat{\alpha}_i) \mathbf{x}_i^T \mathbf{x} + \beta_0$$

Straightforward to kernelize

### Ridge regression revisited

Ridge regression: Given  $\{(\mathbf{x}_i,y_i)\}_{i=1}^n$  ,  $\mathbf{x}_i \in \mathbb{R}^d, y_i \in \mathbb{R}$ 

$$(\widehat{\boldsymbol{\beta}}, \widehat{\beta}_0) = \underset{\boldsymbol{\beta}, \beta_0}{\arg\min} \sum_{i=1}^n (y_i - \boldsymbol{\beta}^T \mathbf{x}_i - \beta_0)^2 + \lambda \|\boldsymbol{\beta}\|_2^2$$
Solution: 
$$\frac{\partial}{\partial \beta_0} = -2 \sum_{i=1}^n (y_i - \boldsymbol{\beta}^T \mathbf{x}_i - \beta_0) = 0$$

$$\widehat{\beta}_0 = \frac{1}{n} \sum_{i=1}^n y_i - \widehat{\boldsymbol{\beta}}^T \mathbf{x}_i$$

$$= \bar{y} - \widehat{\boldsymbol{\beta}}^T \bar{\mathbf{x}}$$

$$\bar{\mathbf{x}} = \frac{1}{n} \sum_{i=1}^n y_i$$

$$\bar{\mathbf{x}} = \frac{1}{n} \sum_{i=1}^n y_i$$

# Ridge regression revisited

Plugging this back in we are left to minimize

$$\sum_{i=1}^{n} (y_i - \bar{y} - \beta^T (\mathbf{x}_i - \bar{\mathbf{x}}))^2 + \lambda \|\beta\|_2^2$$

with respect to  $\beta$ 

$$\widehat{\boldsymbol{\beta}} = (\mathbf{A}^T \mathbf{A} + \lambda \mathbf{I})^{-1} \mathbf{A}^T \widetilde{\mathbf{y}}$$

$$\mathbf{A} = egin{bmatrix} (\mathbf{x}_1 - \mathbf{ar{x}})^T \ dots \ (\mathbf{x}_n - \mathbf{ar{x}})^T \end{bmatrix} \qquad \widetilde{\mathbf{y}} = egin{bmatrix} y_1 - ar{y} \ dots \ y_n - ar{y} \end{bmatrix}$$

### Woodbury matrix inversion identity

$$(\mathbf{P} + \mathbf{Q}\mathbf{R}\mathbf{S})^{-1} = \mathbf{P}^{-1} - \mathbf{P}^{-1}\mathbf{Q}(\mathbf{R}^{-1} + \mathbf{S}\mathbf{P}^{-1}\mathbf{Q})^{-1}\mathbf{S}\mathbf{P}^{-1}$$

$$\mathbf{P} = \lambda \mathbf{I} \quad \mathbf{Q} = \mathbf{A}^T \quad \mathbf{R} = \mathbf{I} \quad \mathbf{S} = \mathbf{A}$$

$$(\lambda \mathbf{I} + \mathbf{A}^T\mathbf{A})^{-1} = \frac{1}{\lambda}\mathbf{I} - \frac{1}{\lambda}\mathbf{I}\mathbf{A}^T\left(\mathbf{I} + \frac{1}{\lambda}\mathbf{A}\mathbf{A}^T\right)^{-1}\mathbf{A}\frac{1}{\lambda}$$

$$= \frac{1}{\lambda}\left[\mathbf{I} - \mathbf{A}^T(\lambda \mathbf{I} + \mathbf{A}\mathbf{A}^T)^{-1}\mathbf{A}\right]$$

$$(\lambda \mathbf{I} + \mathbf{A}^T\mathbf{A})^{-1}\mathbf{A}^T\widetilde{\mathbf{y}} = \frac{1}{\lambda}\left[\mathbf{A}^T - \mathbf{A}^T(\lambda \mathbf{I} + \mathbf{A}\mathbf{A}^T)^{-1}\mathbf{A}\mathbf{A}^T\right]\widetilde{\mathbf{y}}$$

### Kernel ridge regression

$$\widehat{\boldsymbol{\beta}} = (\mathbf{A}^T \mathbf{A} + \lambda \mathbf{I})^{-1} \mathbf{A}^T \widetilde{\mathbf{y}}$$

$$\mathbf{A} = \begin{bmatrix} (\mathbf{x}_1 - \bar{\mathbf{x}})^T \\ \vdots \\ (\mathbf{x}_n - \bar{\mathbf{x}})^T \end{bmatrix} \qquad \widetilde{\mathbf{y}} = \begin{bmatrix} y_1 - \bar{y} \\ \vdots \\ y_n - \bar{y} \end{bmatrix}$$

$$\widehat{f}(\mathbf{x}) = \bar{y} + \widehat{\boldsymbol{\beta}}^T (\mathbf{x} - \bar{\mathbf{x}})$$

Can we express ridge regression in terms of inner products  $\langle \mathbf{x}_i, \mathbf{x}_j \rangle$  and  $\langle \mathbf{x}_i, \mathbf{x} \rangle$ ?

Not immediately.  $[A^T A](i,j) \neq \mathbf{x}_i^T \mathbf{x}_j$ 

### Kernelizing ridge regression

$$(\lambda \mathbf{I} + \mathbf{A}^T \mathbf{A})^{-1} \mathbf{A}^T \widetilde{\mathbf{y}} = \frac{1}{\lambda} \left[ \mathbf{A}^T - \mathbf{A}^T (\lambda \mathbf{I} + \mathbf{A} \mathbf{A}^T)^{-1} \mathbf{A} \mathbf{A}^T \right] \widetilde{\mathbf{y}}$$

$$K(i, j) = (\mathbf{x}_i - \overline{\mathbf{x}})^T (\mathbf{x}_i - \overline{\mathbf{x}})$$

$$\mathbf{X}(i,j) = (\mathbf{x}_i - \bar{\mathbf{x}})^T (\mathbf{x}_j - \bar{\mathbf{x}})$$

$$= \mathbf{x}_i^T \mathbf{x}_j - \frac{1}{n} \sum_{r=1}^n \mathbf{x}_i^T \mathbf{x}_r - \frac{1}{n} \sum_{s=1}^n \mathbf{x}_s^T \mathbf{x}_j + \frac{1}{n^2} \sum_{r,s=1}^n \mathbf{x}_r^T \mathbf{x}_s$$

$$(\lambda \mathbf{I} + \mathbf{A}^T \mathbf{A})^{-1} \mathbf{A}^T \widetilde{\mathbf{y}} = \frac{1}{\lambda} \left[ \mathbf{A}^T - \mathbf{A}^T (\lambda \mathbf{I} + \mathbf{K})^{-1} \mathbf{K} \right] \widetilde{\mathbf{y}}$$

### Kernelizing ridge regression

What about the remaining  $A^T$ ?

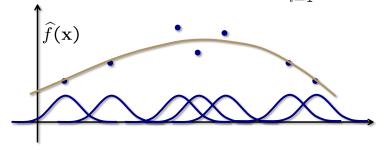
$$\begin{split} \widehat{f}(\mathbf{x}) &= \overline{y} + \widehat{\boldsymbol{\beta}}^T (\mathbf{x} - \overline{\mathbf{x}}) \\ &= \overline{y} + \frac{1}{\lambda} \widetilde{\mathbf{y}}^T \left[ \mathbf{A} - \mathbf{K} (\lambda \mathbf{I} + \mathbf{K})^{-1} \mathbf{A} \right] (\mathbf{x} - \overline{\mathbf{x}}) \\ &= \overline{y} + \frac{1}{\lambda} \widetilde{\mathbf{y}}^T \left[ \mathbf{I} - \mathbf{K} (\lambda \mathbf{I} + \mathbf{K})^{-1} \mathbf{I} \right] \mathbf{k}(\mathbf{x}) \\ \text{where } \mathbf{k}(\mathbf{x}) &= \begin{bmatrix} (\mathbf{x}_1 - \overline{\mathbf{x}})^T (\mathbf{x} - \overline{\mathbf{x}}) \\ \vdots \\ (\mathbf{x}_n - \overline{\mathbf{x}})^T (\mathbf{x} - \overline{\mathbf{x}}) \end{bmatrix} \end{split}$$

$$[k(\mathbf{x})](i) = \mathbf{x}_i^T \mathbf{x} - \frac{1}{n} \sum_{j=1}^n \mathbf{x}_i^T \mathbf{x}_j - \frac{1}{n} \sum_{j=1}^n \mathbf{x}^T \mathbf{x}_j + \frac{1}{n^2} \sum_{j,k=1}^n \mathbf{x}_j^T \mathbf{x}_k$$

# Example: Gaussian kernel

$$k(\mathbf{u}, \mathbf{v}) = \exp\left(-\frac{\|\mathbf{u} - \mathbf{v}\|_2^2}{2\sigma^2}\right)$$

If we omit 
$$\beta_0$$
, then  $\widehat{f}(\mathbf{x}) = \mathbf{y}^T (\mathbf{K} + \lambda \mathbf{I})^{-1} \mathbf{k}(\mathbf{x})$ 
$$= \alpha^T \mathbf{k}(\mathbf{x}) = \sum_{i=1}^n \alpha(i) k(\mathbf{x}, \mathbf{x}_i)$$



### Homogenous kernel ridge regression

For many kernels,  $\Phi(x)$  already contains a constant component, in which case we often omit  $\beta_0$ 

- inhomogenous polynomial kernel
- Gaussian kernel does not seem to require a constant

In this case, the kernel ridge regression solution becomes

$$\widehat{f}(\mathbf{x}) = \mathbf{y}^T (\mathbf{K} + \lambda \mathbf{I})^{-1} \mathbf{k}(\mathbf{x})$$
where  $\mathbf{y} = \begin{bmatrix} y_1, \cdots, y_n \end{bmatrix}^T$ 

$$\mathbf{k}(\mathbf{x}) = \begin{bmatrix} k(\mathbf{x}_1, \mathbf{x}), \cdots, k(\mathbf{x}_n, \mathbf{x}) \end{bmatrix}^T$$

$$\mathbf{K} = \begin{bmatrix} k(\mathbf{x}_1, \mathbf{x}_1) & \cdots & k(\mathbf{x}_1, \mathbf{x}_n) \\ \vdots & \ddots & \vdots \\ k(\mathbf{x}_n, \mathbf{x}_1) & \cdots & k(\mathbf{x}_n, \mathbf{x}_n) \end{bmatrix}$$

### Loss functions and regularization

We have talked about a whole range of algorithms for regression that can be viewed through the lens of minimizing a loss function plus a regularization term

$$\widehat{\theta} = \underset{\theta}{\operatorname{arg\,min}} L(\theta) + \lambda r(\theta)$$

Does this viewpoint also apply to classification?

### Regularized logistic regression

Everything we have said so far about least squares regression can be extended to many classification problems

For example, in logistic regression we can replace

$$\min_{\boldsymbol{\theta}} - \ell(\boldsymbol{\theta})$$

with

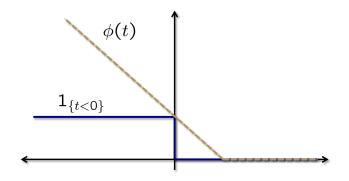
$$\min_{\boldsymbol{\theta}} - \ell(\boldsymbol{\theta}) + \lambda \|\boldsymbol{\theta}\|_2^2$$

Has a similar interpretation to least squares regularization

- makes the Hessian matrix well conditioned
- super useful when the number of observations is small
- · also helpful when data is separable

### Hinge loss

Let's take 
$$\phi(t) = \max\{0, 1 - t\} =: (1 - t)_+$$



### Regularization for linear classification

The kinds of regularization we have talked about can also give us a new way to think about designing linear classifiers

Goal (ideal): Find( $\mathbf{w}, b$ ) minimizing

$$\frac{1}{n}\sum_{i=1}^n \mathbb{1}_{\{y_i(\mathbf{w}^T\mathbf{x}_i+b)<0\}}$$

This is actually much harder than it sounds and is not computationally tractable for large problems

Instead, we can consider replacing this with

$$\frac{1}{n} \sum_{i=1}^{n} \phi(y_i(\mathbf{w}^T \mathbf{x}_i + b))$$

where  $\phi(t)$  is some upper bound on  $\mathbf{1}_{\{t<0\}}$ 

### Adding regularization

Let's try to minimize

$$\frac{1}{n}\sum_{i=1}^{n}(1-y_i(\mathbf{w}^T\mathbf{x}_i+b))_+$$

but to prevent overfitting, let's add a regularization penalty on  $\mathbf{W}$ 

$$\min_{\mathbf{w},b} \ \frac{1}{n} \sum_{i=1}^{n} (1 - y_i(\mathbf{w}^T \mathbf{x}_i + b))_+ + \frac{\lambda}{2} \|\mathbf{w}\|_2^2$$

# Soft-margin hyperplane

Compare this to the optimization problem we considered previously for the optimal soft-margin hyperplane:

$$\min_{\mathbf{w},b} \ \frac{1}{n} \sum_{i=1}^{n} (1 - y_i(\mathbf{w}^T \mathbf{x}_i + b))_+ + \frac{\lambda}{2} ||\mathbf{w}||_2^2$$

VS

$$\min_{\mathbf{w},b,\xi} \frac{1}{2} \|\mathbf{w}\|_2^2 + \frac{C}{n} \sum_{i=1}^n \xi_i$$
s.t.  $y_i(\mathbf{w}^T \mathbf{x}_i + b) \ge 1 - \xi_i$   $i = 1, \dots, n$ 

$$\xi_i \ge 0 \quad i = 1, \dots, n$$

As you all just showed, these are equivalent!