## Recitation 5

#### Kernels

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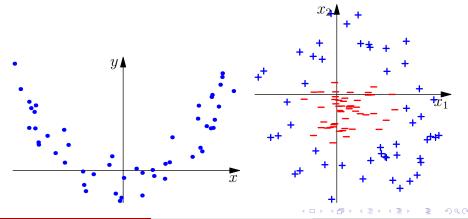
CDS at NYU

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## Intro Question

#### Question

Consider applying linear regression to the data set on the left, and an SVM to the data set on the right. What is the issue? Can it be improved?



### Intro Solution

#### Regression Solution

We want to allow for non-linear regression functions, but we would like to reuse the same fitting procedures we have already developed. To do this we will expand our feature set by adding non-linear functions of old features. We change our features from (1,x) to  $(1,x,x^2)$ . That is

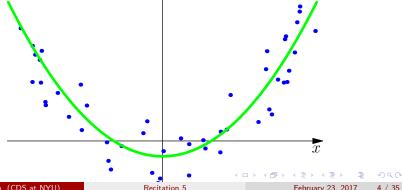
$$X = \begin{pmatrix} 1 & -1 \\ 1 & -.7 \\ \vdots & \vdots \\ 1 & 1 \end{pmatrix} \implies \Phi = \begin{pmatrix} 1 & -1 & (-1)^2 \\ 1 & -.7 & (-.7)^2 \\ \vdots & \vdots & \vdots \\ 1 & 1 & 1^2 \end{pmatrix}.$$

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### Intro Solution

#### Regression Solution

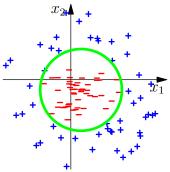
Using features  $(1, x, x^2)$  and w = (-.1, 0, 1) gives us  $f_w(x) = -.1 + 0x + 1x^2 = x^2 - .1$ . Our prediction function is quadratic but we obtained it through standard linear methods.



### Intro Solution

#### **SVM Solution**

For the SVM we expand our feature vector from  $(1, x_1, x_2)$  to  $(1, x_1, x_2, x_1x_2, x_1^2, x_2^2)$ . Using w = (-1.875, 2.5, -2.5, 0, 1, 1) gives  $-1.875 + 2.5x_1 - 2.5x_2 + x_1^2 + x_2^2 = (x_1 + 1.25)^2 + (x_2 - 1.25)^2 - 5 = 0$  as our decision boundary.



## Cost of Adding Features

#### Question

Suppose we begin with d features (and a bias)  $x=(1,x_1,\ldots,x_d)$ . We add all monomials up to degree M. More precisely, all terms of the form  $x_1^{p_1}\cdots x_d^{p_d}$  where  $p_i\geq 0$  and  $p_1+\cdots+p_d\leq M$ . How many features will we have in total?

# Cost of Adding Features

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#### Solution

There will be  $\binom{M+d}{M}$  terms total. If M is fixed and we let d grow, this behaves like  $\frac{d^M}{M!}$ . For example, if d=40 and M=8 we get  $\binom{40+8}{8}=377348994$ . If we are training or predicting with a linear model  $w^Tx$ , this product now takes  $O(d^M)$  operations to evaluate.



### Kernel Trick

Consider the polynomial kernel  $k(x,y)=(1+x^Ty)^M$  where  $x,y\in\mathbb{R}^d$ . This computes the inner product of all monomials up to degree M in time O(d). For example, if M=2 we have

$$(1 + x^T y)^2 = 1 + 2x^T y + x^T y x^T y = 1 + 2 \sum_{i=1}^d x_i y_i + \sum_{i,j=1}^d x_i y_i x_j y_j.$$

The resulting feature map is

$$\varphi(x) = (1, \sqrt{2}x_1, \dots, \sqrt{2}x_d, x_1^2, \dots, x_d^2, \sqrt{2}x_1x_2, \sqrt{2}x_1x_3, \dots, \sqrt{2}x_{d-1}x_d).$$

Then  $k(x, y) = \varphi(x)^T \varphi(y)$ .



# Kernel Ridge Regression

Recall that our ridge regression loss is given by

$$J(w) = \frac{1}{n} \sum_{i=1}^{n} (w^{T} x_{i} - y_{i})^{2} + \lambda ||w||_{2}^{2}$$

ullet If we map to a larger feature space  $arphi:\mathbb{R}^d o\mathbb{R}^D$  we get

$$J(\tilde{w}) = \frac{1}{n} \sum_{i=1}^{n} (\tilde{w}^{T} \varphi(x_i) - y_i)^2 + \lambda ||\tilde{w}||_2^2.$$

• Using the kernel trick we can try to write this (incorrectly) as

$$J(w) = \frac{1}{n} \sum_{i=1}^{n} (k(w, x_i) - y_i)^2 + \lambda k(w, w).$$

• What are the issues with this?



# Kernel Ridge Regression

Issues with

$$J(w) = \frac{1}{n} \sum_{i=1}^{n} (k(w, x_i) - y_i)^2 + \lambda k(w, w)$$

- Writing  $\tilde{w}^T \varphi(x_i)$  isn't the same as  $k(w, x_i) = \varphi(w)^T \varphi(x_i)$  since  $\varphi$  isn't onto. That is,  $\varphi(w)$  is a very specific type of element of  $\mathbb{R}^D$ , the larger feature space.
- The L(w) written above isn't a ridge regression problem any more, since  $k(w,x_i)$  and k(w,w) can be weird functions of w. Thus our previous code and theory for dealing with ridge regression doesn't immediately carry over.



# Math Thought Experiment

• Suppose we have a standard ridge regression problem for  $w \in \mathbb{R}^d$ :

$$J(w) = \frac{1}{n} \sum_{i=1}^{n} (w^{T} x_{i} - y_{i})^{2} + \lambda ||w||_{2}^{2}.$$

• Now suppose we add a new feature to x that is always zero:  $\tilde{x} = (x, 0)$ . For  $w \in \mathbb{R}^{d+1}$  we now have

$$J(w) = \frac{1}{n} \sum_{i=1}^{n} (w^{T} \tilde{x}_{i} - y_{i})^{2} + \lambda ||w||_{2}^{2}.$$

Does the answer change? In other words, will the new minimizer  $w_* \in \mathbb{R}^{d+1}$  have a non-zero value in its last coordinate?



# Math Thought Experiment

Does the answer change for  $w \in \mathbb{R}^{d+1}$  with

$$J(w) = \frac{1}{n} \sum_{i=1}^{n} (w^{T} \tilde{x}_{i} - y_{i})^{2} + \lambda ||w||_{2}^{2}$$

No!

- Suppose  $w_* = (v, a)$  where  $v \in \mathbb{R}^d$  and  $a \in \mathbb{R}$ .
- Recall that  $\tilde{x}_i = (x, 0)$ .
- Then  $w_*^T \tilde{x}_i = v^T x$ . In other words, a has **no** effect on the prediction  $w_*^T \tilde{x}_i$ .
- But if  $a \neq 0$  then  $\|(v, a)\|_2^2 > \|(v, 0)\|_2^2$ .
- Can you think of the more general version of this phenomenon?



# Representer Theorem (Baby Version)

### Theorem ((Baby) Representer Theorem)

Suppose you have a loss function of the form

$$J(w) = L(w^T \varphi(x_1), \dots, w^T \varphi(x_n)) + R(\|w\|_2)$$

where

- $w, \varphi(x_i) \in \mathbb{R}^D$ .
- $L: \mathbb{R}^n \to \mathbb{R}$  is an arbitrary function (loss term).
- $R: \mathbb{R}_{\geq 0} \to \mathbb{R}$  is increasing (regularization term).

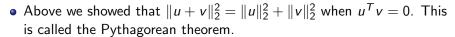
Assume J has at least one minimizer. Then J has a minimizer  $w^*$  of the form  $w^* = \sum_{i=1}^n \alpha_i \varphi(x_i)$  for some  $\alpha \in \mathbb{R}^n$ . If R is strictly increasing, then all minimizers have this form.



# Representer Theorem: Proof

#### Proof.

- Let  $w^* \in \mathbb{R}^D$  and let  $S = \operatorname{Span}(\varphi(x_1), \dots, \varphi(x_n))$ .
- Write  $w^* = u + v$  where  $u \in S$  and  $v \in S^{\perp}$ . Here u is the orthogonal projection of  $w^*$  onto S, and  $S^{\perp}$  is the subspace of all vectors orthogonal to S.
- Then  $(w^*)^T \varphi(x_i) = (u+v)^T \varphi(x_i) = u^T \varphi(x_i) + v^T \varphi(x_i) = u^T \varphi(x_i)$ .
- But  $||w^*||_2^2 = ||u + v||_2^2 = ||u||_2^2 + ||v||_2^2 + 2u^Tv = ||u||_2^2 + ||v||_2^2 \ge ||u||_2^2$ .
- Thus  $R(\|w^*\|_2) \ge R(\|u\|_2)$  showing  $J(w^*) \ge J(u)$ .





# Representer Theorem: Meaning

- If your loss function only depends on w via its inner products with the inputs, and the regularization is an increasing function of the  $\ell_2$  norm, then we can write  $w^*$  as a linear combination of the training data.
- This applies to ridge regression and SVM.

### Question

Suppose you have n=100 samples, d=40 features, and M=8 degree monomial terms giving 377348994 features. This implies  $w \in \mathbb{R}^{377348994}$  for ridge regression. What does the representer theorem say?

# Representer Theorem: Meaning

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#### Question

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#### Solution

As  $y \in \mathbb{R}^n$  varies, the solution w must lie in a 100-dimensional subspace of  $\mathbb{R}^{377348994}$ 



# Representer Theorem: Ridge Regression

By adding features to ridge regression we had

$$J(\tilde{w}) = \frac{1}{n} \sum_{i=1}^{n} (\tilde{w}^{T} \varphi(x_{i}) - y_{i})^{2} + \lambda ||\tilde{w}||_{2}^{2}$$
$$= \frac{1}{n} ||\Phi \tilde{w} - y||_{2}^{2} + \lambda \tilde{w}^{T} \tilde{w},$$

where  $\Phi \in \mathbb{R}^{n \times D}$  is the matrix with  $\varphi(x_i)^T$  as its *i*th row.

- Representer Theorem applies giving  $\tilde{w} = \sum_{i=1}^{n} \alpha_i \varphi(x_i) = \Phi^T \alpha$ .
- Plugging in gives

$$J(\alpha) = \frac{1}{n} \left\| \Phi \Phi^T \alpha - y \right\|_2^2 + \lambda \alpha^T \Phi \Phi^T \alpha.$$



# Representer Theorem: Ridge Regression

• Let  $K \in \mathbb{R}^{n \times n}$  be given by  $K = \Phi \Phi^T$ . This is called the **Gram** Matrix and satisfies  $K_{ij} = k(x_i, x_j) = \varphi(x_i)^T \varphi(x_j)$ :

$$K = \begin{pmatrix} \varphi(x_1)^T \varphi(x_1) & \cdots & \varphi(x_1)^T \varphi(x_n) \\ \vdots & \ddots & \vdots \\ \varphi(x_n)^T \varphi(x_1) & \cdots & \varphi(x_n)^T \varphi(x_n) \end{pmatrix}.$$

We can write ridge regression in the kernelized form by turning

$$J(\alpha) = \frac{1}{n} \left\| \Phi \Phi^{\mathsf{T}} \alpha - y \right\|_{2}^{2} + \lambda \alpha^{\mathsf{T}} \Phi \Phi^{\mathsf{T}} \alpha.$$

into

$$J(\alpha) = \frac{1}{n} ||K\alpha - y||_2^2 + \lambda \alpha^T K\alpha.$$

- Can derive the solution algebraically (see Homework 4).
- Prediction function is  $f_{\alpha}(x) = (w^*)^T \varphi(x) = \sum_{i=1}^n \alpha_i k(x_i, x)$ .

# Representer Theorem: Primal SVM

• For a general linear model, the same derivation above shows

$$J(w) = L(\Phi w) + R(\|w\|_2)$$

becomes

$$J(\alpha) = L(K\alpha) + R(\sqrt{\alpha^T K\alpha}).$$

Here  $\varphi(x_i)^T w$  became  $(K\alpha)_i$ .

• The primal SVM (bias in features) has loss function

$$J(w) = \frac{c}{n} \sum_{i=1}^{n} (1 - y_i(\varphi(x_i)^T w))_+ + ||w||_2^2.$$

This is kernelized to

$$J(\alpha) = \frac{c}{n} \sum_{i=1}^{n} (1 - y_i(K\alpha)_i)_+ + \alpha^T K\alpha.$$

• Positive decision made if  $(w^*)^T \varphi(x) = \sum_{i=1}^n \alpha_i k(x_i, x) > 0$ .

### Dual SVM

The dual SVM problem (with features) is given by

$$\begin{aligned} & \underset{i=1}{\operatorname{maximize}}_{\alpha} & & \sum_{i=1}^{n} \alpha_{i} - \frac{1}{2} \sum_{i,j=1}^{n} \alpha_{i} \alpha_{j} y_{i} y_{j} \varphi(x_{i})^{T} \varphi(x_{j}) \\ & \text{subject to} & & \sum_{i=1}^{n} \alpha_{i} y_{i} = 0 \\ & & \alpha_{i} \in \left[0, \frac{c}{n}\right] \quad \text{for } i = 1, \dots, n. \end{aligned}$$

- We can immediately kernelize (no representer theorem needed) by replacing  $\varphi(x_i)^T \varphi(x_i) = k(x_i, x_i)$ .
- Recall that we were able to derive the conclusion of the representer theorem using strong duality for SVMs.



### **RBF Kernel**

As we saw last time, the most frequently used kernel is the RBF kernel

$$k(w,x) = \exp\left(-\frac{\|w-x\|_2^2}{2\sigma^2}\right).$$

- Is there a corresponding feature map  $\varphi : \mathbb{R}^d \to \mathbb{R}^D$  so that  $k(w,x) = \varphi(w)^T \varphi(x)$ ?
- Unfortunately there is no finite D that will work.
- We will handle this by allowing infinite dimensional spaces called Hilbert spaces.

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# Rough Review of Abstract Linear Algebra

- Vector Spaces: Spaces where linear combinations (scaling and adding) make sense.
- A vector space has no sense of length or angle, but has concepts like subspace, basis, span, dimension, linear transformation, eigenvector.
- We can account for length and angle by looking at inner products.
- Recall that in 2d, there is a law of cosines that relates side lengths to the angles of a triangle. In other words, reasonable definitions of length and angle are not independent.

### Definition of Inner Product

### Definition (Inner Product)

Let V be a vector space. We say  $\langle \cdot, \cdot \rangle : V^2 \to \mathbb{R}$  is an **inner product** if it satisfies the following:

Symmetry:

$$\langle v, w \rangle = \langle w, v \rangle$$

for  $v, w \in V$ .

Bilinearity:

$$\langle \alpha \mathbf{v}_1 + \beta \mathbf{v}_2, \mathbf{w} \rangle = \alpha \langle \mathbf{v}_1, \mathbf{w} \rangle + \beta \langle \mathbf{v}_2, \mathbf{w} \rangle$$

for  $v_1, v_2, w \in V$  and  $\alpha, \beta \in \mathbb{R}$ . Same holds for second argument by symmetry.

**3** Positive-Definiteness:  $\langle v, v \rangle \geq 0$  with  $\langle v, v \rangle = 0$  if and only if v = 0.

If we associate an inner product with a vector space we call the pair an inner product space.

### More on Inner Products and Norms

- On  $\mathbb{R}^n$  the "dot-product"  $v^T w$  is an inner product.
- Using an inner product we can obtain the norm (length) of a vector:  $\|v\| = \sqrt{\langle v, v \rangle}$ . We obtain the angle via  $\cos(\angle(v, w)) := \frac{\langle v, w \rangle}{\|v\| \|w\|}$ .
- We say v, w are orthogonal if  $\langle v, w \rangle = 0$ .
- Not all norms we have seen have an associated inner product. For example,  $||v||_1$  has no associated inner product.
- Roughly, certain norms (like  $\ell_1$ ) do not lead to a consistent way of measuring angles. More precisely:

### Theorem (Parallelogram Law)

A norm  $\|\cdot\|$  has an associated inner product if and only if

$$2||v||^2 + 2||w||^2 = ||v - w||^2 + ||v + w||^2.$$



## Hilbert Space

- Once we have a norm, we can talk about the distance between two vectors: ||v w||. This allows one to introduce concepts such as convergence and continuity.
- Certain facts we need about projections are false in infinite dimensions unless we impose an extra constraint on our inner product spaces: completeness.
- A space is complete if all Cauchy sequences converge.

## Definition (Hilbert Space)

An inner product space V with inner product  $\langle \cdot, \cdot \rangle$  is called a **Hilbert** space if it is complete with respect to the associated norm.

All finite dimensional inner product spaces are Hilbert space.



## Projection Theorem

#### Theorem (Projection Theorem)

Let H be a Hilbert space and let S be a finite dimensional subspace. Then any vector  $w \in H$  can be written uniquely as w = u + v where  $u \in S$  and  $v \in S^{\perp}$ .

#### More can be said:

- u is the projection of w onto S:  $u = \arg\min_{x \in S} ||x w||$ .
- Can extend theorem to hold for any closed subspace (finite dimensional subspaces are examples of closed subspaces).
- A variant of the theorem holds for non-empty closed convex subsets.

# Representer Theorem (Adult Version)

### Theorem (Representer Theorem)

Suppose you have a loss function of the form

$$J(w) = L(\langle w, \varphi(x_1) \rangle, \dots, \langle w, \varphi(x_n) \rangle) + R(\|w\|)$$

#### where

- $w, \varphi(x_i) \in H$  for some Hilbert space H.
- $L: \mathbb{R}^n \to \mathbb{R}$  is an arbitrary function (loss term).
- $R: \mathbb{R}_{\geq 0} \to \mathbb{R}$  is increasing (regularization term).
- $\|\cdot\|$  is the norm associated with H.

Assume J has at least one minimizer. Then J has a minimizer  $w^*$  of the form  $w^* = \sum_{i=1}^n \alpha_i \varphi(x_i)$  for some  $\alpha \in \mathbb{R}^n$ . If R is strictly increasing, then all minimizers have this form.

# Representer Theorem (Adult Version)

- Same proof as before, but apply the projection theorem and use  $\langle v, w \rangle$  in place of  $v^T w$ .
- Now we have a representer theorem that works for infinite dimensional Hilbert spaces. Why do we care?
- We will show that kernels, such as the RBF Kernel, correspond to inner products in Hilbert spaces.

### Positive Semi-Definite

### Definition (Positive Semi-Definite)

A matrix  $A \in \mathbb{R}^{n \times n}$  is **positive semi-definite** if it is symmetric and

$$x^T A x \ge 0$$

for all  $x \in \mathbb{R}^n$ .

 Equivalent to saying the matrix is symmetric with non-negative eigenvalues.

### Mercer's Theorem

### Theorem (Mercer's Theorem)

Fix a kernel  $k: \mathcal{X} \times \mathcal{X} \to \mathbb{R}$ . There is a Hilbert space H and a feature map  $\varphi: \mathcal{X} \to H$  such that  $k(x,y) = \langle \varphi(x), \varphi(y) \rangle_H$  if and only if for any  $x_1, \ldots, x_n \in \mathcal{X}$  the associated matrix K is positive semi-definite:

$$K = \begin{pmatrix} k(x_1, x_1) & \cdots & k(x_1, x_n) \\ \vdots & \ddots & \vdots \\ k(x_n, x_1) & \cdots & k(x_n, x_n) \end{pmatrix}.$$

Such a kernel k is called **positive semi-definite**.

# Finding Your Own Kernels

Let  $k_1, k_2 : \mathcal{X} \times \mathcal{X} \to \mathbb{R}$  be positive semi-definite kernels. Then so are the following:

- $k_3(w,x) = k_1(w,x) + k_2(w,x)$
- $k_4(w,x) = \alpha k_1(w,x)$  for  $\alpha \ge 0$
- $k_5(w,x) = f(w)f(x)$  for any function  $f: \mathcal{X} \to \mathbb{R}$
- $k_6(w,x) = k_1(w,x)k_2(w,x)$

Furthermore, if  $k_1, k_2, ...$  is a sequence of positive semi-definite kernels then

$$k(w,x) = \lim_{n \to \infty} k_n(w,x)$$

is also positive semi-definite (assuming the limit exists for all w, x).

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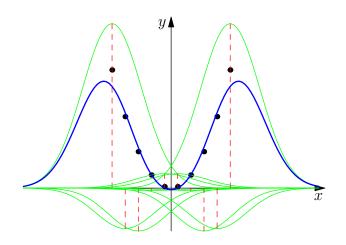
## Representer Theorem for RBF Kernels

- As we saw earlier for ridge regression and SVM classification, the decision function has the form  $f_{\alpha}(x) = \sum_{i=1}^{n} \alpha_{i} k(x_{i}, x)$ .
- For ridge regression, this means that using the RBF kernel amounts to approximating our data by a linear combination of Gaussian bumps.
- For SVM classification, each  $k(x_i, x) = \exp\left(-\|x_i x\|_2^2/(2\sigma^2)\right)$  represents a exponentially decaying distance between  $x_i$  and x. Thus our decisions depend on our proximities to data points.

# **RBF** Regression

- Below we use 10 uniformly spaced x-values between -2 and 2, with  $y_i = x_i^2$ . We fit kernelized ridge regression with the RBF kernel using  $\sigma = 1$  and  $\lambda = .1$ .
- Each green curve is  $g(x) = \alpha_i k(x_i, x)$ . The predicted function is drawn in blue.
- As you might expect, extrapolating outside of [-2, 2] can have poor results.
- People will often normalize the RBF kernel (see Hastie, Tibshirani, Friedman p. 213).

# **RBF** Regression

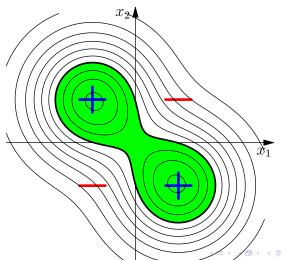


### **RBF** Classification

 Next we show 4 points placed on the corners of a square with positive and negative points on each diagonal.

### **RBF** Classification

Contours of  $f(x) = k(x_1, x) + k(x_2, x)$  where  $x_1, x_2$  are positive examples, and  $\sigma = 1$ .



### **RBF** Classification

Contours of  $f(x) = k(x_1, x) + k(x_2, x) - k(x_3, x) - k(x_4, x)$  where  $x_1, x_2$  are positive examples, and  $\sigma = 1$ .

