## EE445 Mod2-Lec4: Kernel Regression

[Lecturer: L.J. Ratliff]

## What is Kernel Regression?

- We have been talking about supervised ML in the context of data fitting with least squares
- Cost function paradigm for supervised machine learning
  - ► Features x
  - Output/response y
  - Goal: Find f(x) such that  $f(x^{(i)}) \approx y^{(i)}$
  - objective/cost function  $F(\theta) = ||A\theta y||^2$

## Kernel Motivation

- But what we really want are flexible non-linear classifers/predictors!
- We can get this via a linear model using the kernel trick
- Note that feature maps are already all non-linear

$$x \mapsto 1, x, x^2, \dots$$

- Yet, we want something a little more automatic that implicitly captures nonlinearities without expanding out data to many times the original size
- Kernels give us this

## Kernel Trick: Starting Point

• [A2]:

$$heta = \sum_{i=1}^m lpha_i x^{(i)}$$
 for some  $lpha_1, \dots, lpha_m \in \mathbb{R}$ 

i.e.,  $\boldsymbol{\theta}$  is in the span of the feature vectors

• We will see shortly how to find these  $\alpha_i$ 's

## Kernel Trick: Linear Regression

• [A2]: 
$$\theta = \sum_{i=1}^{m} \alpha_i x^{(i)}$$
 for some  $\alpha_1, \dots, \alpha_m \in \mathbb{R}$   
$$f(x) = \theta^\top x = \left(\sum_{i=1}^{m} \alpha_i x^{(i)}\right)^\top x = \sum_{i=1}^{m} \alpha_i (x^{(i)})^\top x = \sum_{i=1}^{m} \alpha_i K(x^{(i)}, x)$$

- Kernel function:  $K(x, z) = x^{\top} z$
- Predictions only depend on training data through kernel function which is just a dot product.

## Linear Regression: Objective Function

- [A2]:  $\theta = \sum_{i=1}^m \alpha_i x^{(i)}$  for some  $\alpha_1, \ldots, \alpha_m \in \mathbb{R}$
- The predictor has the form

$$f(x) = \sum_{i=1}^{m} \alpha_i K(x^{(i)}, x)$$

• The objective function has the form

$$\frac{1}{2}\sum_{i=1}^{m} (f(x^{(i)}) - y^{(i)})^2 = \frac{1}{2}\sum_{i=1}^{m} \left(\sum_{j=1}^{m} \alpha_j K(x^{(j)}, x^{(i)}) - y^{(i)}\right)^2 =: F(\alpha)$$

- Objective function only depends on training data through kernel function which is just dot products
- Choose  $\alpha$  by minimizing  $F(\alpha)$

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## Kernel Trick: Take Aways

- Predictor and objective only depend on training data through the kernel which is itself just dot products
- Hence, if we only have the ability to do dot product operations, then we can still suprisingly train a model (i.e., find a prediction of y)

#### Kernelized Linear Regression

• Rewrite linear regression as a different linear regression model:

$$f(x) = \sum_{i=1}^{m} \alpha_i K(x^{(i)}, x) = \alpha^\top k(x)$$

#### where

$$\alpha^{\top} = \begin{bmatrix} \alpha_1 & \cdots & \alpha_m \end{bmatrix}$$
 and  $k(x) = \begin{bmatrix} K(x^{(1)}, x) \\ \vdots \\ K(x^{(m)}, x) \end{bmatrix}$ 

• i.e., we map x to a new "feature vector" k(x) (= kernel evaluation between x and each training feature vector).

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# What happens to original data matrix X under this mapping?

- Recall: *i*-th row of X is *i*-th feature vector  $x^{(i)}$
- Kernel Matrix: new "data matrix" K such that the *i*-th row contains dot products between  $x^{(i)}$  and every other training point:

$$K_{ij} = K(x^{(i)}, x^{(j)}) = (x^{(i)})^{\top} x^{(j)}$$

- Sometimes this is called the Kernel Trick.
- Take-Away: you can learn an equivalent linear model using the kernel matrix in place of the original data matrix.
- this equivalence is only exact without regularization (I will talk about this shortly)

## Nonlinear Feature Maps

• Suppose we want to do feature mappings before learning such as

$$f(x) = \theta^{\top} \phi(x), \quad \phi : \mathbb{R}^n \to \mathbb{R}^p$$

• Kernel corresponding to *φ*: To solve the learning problem and make predictions, we only need to be able to compute

$$K(x,z) = \phi(x)^{\top} \phi(z)$$

## Examples: Polynomial Kernel

Note: we can often compute kernel without actually doing the expansion

- Consider  $K(x,z) = (x^{\top}z)^2$
- What is  $\phi: \mathbb{R}^2 \to \mathbb{R}^4$ ?

$$\phi(x) = (x_1^2, x_1 x_2, x_2 x_1, x_2^2)$$

• Check:

$$K(x,z) = (x^{\top}z)^2 = \left( \begin{bmatrix} x_1 \\ x_2 \end{bmatrix}^{\top} \begin{bmatrix} z_1 \\ z_2 \end{bmatrix} \right)^2 = (z_1x_1 + z_2x_2)^2 = x_1^2z_1^2 + 2z_1x_1 \cdot x_2z_2 + z_2^2x_2^2$$

and

$$\phi(x)^{\top}\phi(z) = \begin{bmatrix} x_1^2 \\ x_1 x_2 \\ x_2 x_1 \\ x_2^2 \end{bmatrix}^{\top} \begin{bmatrix} z_1^2 \\ z_1 z_2 \\ z_2 z_1 \\ z_2^2 \end{bmatrix} = x_1^2 z_1^2 + 2z_1 x_1 \cdot x_2 z_2 + z_2^2 x_2^2$$

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## Examples: Polynomial Kernel

Note: computational complexity is lower

- Consider  $K(x,z) = (x^{\top}z + 1)^2$
- What is  $\phi : \mathbb{R}^2 \to \mathbb{R}^4$ ?

$$\phi(x) = (1, \sqrt{2}x_1, \sqrt{2}x_2, x_1^2, x_1x_2, x_2x_1, x_2^2)$$

- complexity of  $\phi(x)^{\top}\phi(z)$ :  $O(n^2)$
- complexity of  $x^{\top}z$ : O(n)
- complexity of  $(x^{\top}z+1)^2$ : O(n)
- If using kernel trick, can implement a non-linear feature expansion at no additional cost
- More general:  $K(x,z) = (x^{\top}z + 1)^d$ 
  - complexity of computing corresponding features with  $\phi$ :  $O(n^d)$
  - complexity of computing K: O(n)

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## Example: Gaussian Kernel

$$K(x, z) = \exp(-\gamma ||x - z||^2)$$

Some observations:

- non-linear kernel with a lot of flexibility
- corresponds to an infinite dimensional  $\phi$ —i.e., cannot implement the corresponding feature mapping  $\phi$ .

### In Practice: Regularization

• We often introduce a regularization term in practice:

$$F(\theta) = \sum_{k=1}^{m} (\theta^{\top} \phi(x^{(k)}) - y^{(k)})^2 + \frac{\lambda}{2} \|\theta\|^2$$

- why?: Regularization improves the conditioning of the problem and reduces the variance of the estimates.
- Taking derivatives and setting them to zero we have

$$\sum_{k=1}^{m} (\theta^{\top} \phi(x^{(k)}) - y^{(k)}) \phi(x^{(k)}) = \lambda \theta$$
  
$$\implies \hat{\theta} = \left(\lambda I + \sum_{k=1}^{m} \phi(x^{(k)}) \phi(x^{(k)})^{\top}\right)^{-1} \sum_{j=1}^{m} \phi(x^{(j)}) y^{(j)}$$

i.e.,  $\lambda$  helps stabilize the inverse.

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#### Deriving the $\alpha$ -dependent regularization term

Recall that we converted  $F(\theta)$  to a cost in terms of  $\alpha.$  We will do the same thing for the regularized cost.

• [A2]: 
$$\theta = \sum_{i=1}^{m} \alpha_i x^{(i)}$$
 for some  $\alpha_1, \ldots, \alpha_m \in \mathbb{R}$ 

$$\|\theta\|^{2} = \theta^{\top}\theta = \left(\sum_{k=1}^{m} \alpha_{k}\phi(x^{(k)})\right)^{\top} \left(\sum_{k=1}^{m} \alpha_{k}\phi(x^{(k)})\right)$$
$$= \sum_{k=1}^{m} \sum_{j=1}^{m} \alpha_{k}\alpha_{j}\phi(x^{(k)})^{\top}\phi(x^{(j)})$$
$$= \sum_{k=1}^{m} \sum_{j=1}^{m} \alpha_{k}\alpha_{j}K(x^{(k)}, x^{(j)})$$
$$= \alpha^{\top}K\alpha$$

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## Kernelized Regression Regularized Cost (Ridge Regression)

$$F(\alpha) = \frac{1}{2} \|K\alpha - y\|^2 + \frac{\lambda}{2} \alpha^\top K \alpha$$

so that taking derivatives, we have

$$K(K\alpha - y) + \lambda K\alpha = 0 \iff K(K + \lambda I)\alpha = Ky$$

One solution is

$$\alpha = (K + \lambda I)^{-1} y$$

- This turns out to be the only solution we care about due to the form of our predictor  $f(\cdot)$ ; any other solution won't affect the final form of our predictor.
- Choose  $\lambda$  via cross validation!

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