CS485/685 Lecture 9: Jan 31, 2012

Kernel methods
[B] Sections 6.1, 6.2

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Non-linear Models Recap

- Generalized linear models:
- Neural networks:

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Kernel Methods

- Idea: use large (possibly infinite) set of fixed nonlinear basis functions
- Normally, complexity depends on number of basis functions, but by a "dual trick", complexity depends on the amount of data
- Examples:
 - Gaussian Processes (next class)
 - Support Vector Machines (next week)
 - Kernel Perceptron
 - Kernel Principal Component Analysis

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Kernel Function

- Let $\phi(x)$ be a set of basis functions that map inputs x to a feature space.
- In many algorithms, this feature space only appears in the dot product $\phi(x)^T \phi(x')$ of pairs inputs x, x'.
- Define the kernel function $k(x, x') = \phi(x)^T \phi(x')$ to be the dot product of any pair x, x' in feature space.
 - We only need to know k(x, x'), not $\phi(x)$

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Dual Representations

• Recall linear regression objective

$$E(\mathbf{w}) = \frac{1}{2} \sum_{n=1}^{N} [\mathbf{w}^{T} \phi(\mathbf{x}_{n}) - y_{n}]^{2} + \frac{\lambda}{2} \mathbf{w}^{T} \mathbf{w}$$

Solution: set gradient to 0

$$\nabla E(\mathbf{w}) = \sum_{n} (\mathbf{w}^{T} \phi(\mathbf{x}_{n}) - y_{n}) \phi(\mathbf{x}_{n}) + \lambda \mathbf{w} = 0$$
$$\mathbf{w} = -\frac{1}{\lambda} \sum_{n} (\mathbf{w}^{T} \phi(\mathbf{x}_{n}) - y_{n}) \phi(\mathbf{x}_{n})$$

 $\dot{}$ w is a linear combination of inputs in feature space

$$\{\phi(x_n)|1 \le n \le N\}$$

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Dual Representations

- Substitute $\mathbf{w} = \mathbf{\Phi} \mathbf{a}$
- Where $\Phi = [\phi(x_1) \phi(x_2) \dots \phi(x_N)]$

$$a = \begin{bmatrix} a_1 \\ a_2 \\ \vdots \\ a_N \end{bmatrix}$$
 and $a_n = -\frac{1}{\lambda} (\mathbf{w}^T \phi(\mathbf{x}_n) - y_n)$

• Dual objective: minimize $\it E$ with respect to $\it a$

$$E(\boldsymbol{a}) = \frac{1}{2}\boldsymbol{a}^T\boldsymbol{\Phi}^T\boldsymbol{\Phi}\boldsymbol{\Phi}^T\boldsymbol{\Phi}\boldsymbol{a} - \boldsymbol{a}^T\boldsymbol{\Phi}^T\boldsymbol{\Phi}\boldsymbol{y} + \frac{\boldsymbol{y}^T\boldsymbol{y}}{2} + \frac{\lambda}{2}\boldsymbol{a}^T\boldsymbol{\Phi}^T\boldsymbol{\Phi}\boldsymbol{a}$$

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Gram Matrix

- Let $K = \Phi^T \Phi$ be the Gram matrix
- Substitute in objective:

$$E(\boldsymbol{a}) = \frac{1}{2}\boldsymbol{a}^{T}\boldsymbol{K}\boldsymbol{K}\boldsymbol{a} - \boldsymbol{a}^{T}\boldsymbol{K}\boldsymbol{y} + \frac{\boldsymbol{y}^{T}\boldsymbol{y}}{2} + \frac{\lambda}{2}\boldsymbol{a}^{T}\boldsymbol{K}\boldsymbol{a}$$

• Solution: set gradient to 0

$$\nabla E(\mathbf{a}) = \mathbf{K}\mathbf{K}\mathbf{a} - \mathbf{K}\mathbf{y} + \lambda \mathbf{K}\mathbf{a} = 0$$
$$\mathbf{K}(\mathbf{K} + \lambda \mathbf{I})\mathbf{a} = \mathbf{K}\mathbf{y}$$
$$\mathbf{a} = (\mathbf{K} + \lambda \mathbf{I})^{-1}\mathbf{y}$$

• Prediction:

$$y_n = \mathbf{w}^T \phi(\mathbf{x}_n) = \mathbf{a}^T \mathbf{\Phi} \phi(\mathbf{x}_n) = k(\cdot, \mathbf{x}_n)^T (\mathbf{K} + \lambda \mathbf{I})^{-1} \mathbf{y}$$

where $k(\cdot, \mathbf{x}_n) = K(\mathbf{x}, \mathbf{x}_n) \ \forall \mathbf{x}$

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Dual Linear Regression

- Prediction: $y_n = \boldsymbol{a}^T \boldsymbol{\Phi} \phi(\boldsymbol{x}_n)$ = $k(\cdot, \boldsymbol{x}_n)^T (\boldsymbol{K} + \lambda \boldsymbol{I})^{-1} \boldsymbol{y}$
- Linear regression where we find dual solution a instead of primal solution w.
- Complexity:
 - Primal solution: # of basis functions
 - Dual solution: amount of data
 - Advantage: can use very large # of basis functions
 - Just need to know kernel k

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Constructing Kernels

- Two possibilities:
 - Find mapping ϕ to feature space and let $K = \phi^T \phi$
 - Directly specify K
- Can any function that takes two argument serve as a kernel?
- No, a valid kernel must be positive semi-definite
 - In other words, k must factor into the product of a transposed matrix by itself (e.g., $K = \phi^T \phi$)
 - Or, all eigenvalues must be greater than or equal to 0.

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Example

• Let $k(\mathbf{x}, \mathbf{z}) = (\mathbf{x}^T \mathbf{z})^2$

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Constructing Kernels

- Can we construct k directly without knowing ϕ ?
- Yes, any positive semi-definite k is fine since there is a corresponding implicit feature space. But positive semi-definiteness is not always easy to verify.
- Alternative, construct kernels from other kernels using rules that preserve positive semi-definiteness

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Rules to construct Kernels

- Let $k_1(x, x')$ and $k_2(x, x')$ be valid kernels
- The following kernels are also valid:
 - 1. $k(x, x') = ck_1(x, x') \quad \forall c > 0$
 - 2. $k(x, x') = f(x)k_1(x, x')f(x') \quad \forall f$
 - 3. $k(x, x') = q(k_1(x, x'))$ q is polynomial with coeffs ≥ 0
 - 4. $k(x, x') = \exp(k_1(x, x'))$
 - 5. $k(\mathbf{x}, \mathbf{x}') = k_1(\mathbf{x}, \mathbf{x}') + k_2(\mathbf{x}, \mathbf{x}')$
 - 6. $k(x, x') = k_1(x, x')k_2(x, x')$
 - 7. $k(x, x') = k_3(\phi(x), \phi(x'))$
 - 8. $k(x, x') = x^T A x'$ A is symmetric positive semi-definite
 - 9. $k(x, x') = k_a(x_a, x'_a) + k_b(x_b, x'_b)$
 - 10. $k(\mathbf{x}, \mathbf{x}') = k_a(\mathbf{x}_a, \mathbf{x}'_a)k_b(\mathbf{x}_b, \mathbf{x}'_b)$

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Common Kernels

- Polynomial kernel: $k(x, x') = (x^T x')^M$
 - − *M* is the degree
 - Feature space: all degree M products of entries in x
 - Example: Let x and x' be two images, then feature space could be all products of M pixel intensities
- More general polynomial kernel:

$$k(x, x') = (x^T x' + c)^M$$
 with $c > 0$

- Feature space: all products of up to M entries in $oldsymbol{x}$

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Common Kernels

- Gaussian Kernel: $k(x, x') = \exp\left(-\frac{\left||x-x'|\right|^2}{2\sigma^2}\right)$
- Valid Kernel because:

• Implicit feature space is infinite!

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Non-vectorial Kernels

- Kernels can be defined with respect to other things than vectors such as sets, strings or graphs
- Example for sets: $k(A_1, A_2) = 2^{|A_1 \cap A_2|}$

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