CSCE 990 Lecture 9:	Introduction • We are now very aware of the importance and
Designing Kernels*	power of kernels in SVMs
Stephen D. Scott	<ul> <li>We also know from Chapter 2 about some basic kernels and simple ways to build new kernels out of old ones</li> <li>Linear scaling, addition, multiplication, etc. of existing kernels</li> </ul>
March 23, 2006	<ul> <li>We'll look at other ways to construct new ker- nels from existing ones, plus other completely different types of kernels</li> </ul>
*Most figures ©2002 MIT Press, Bernhard Schölkopf, and Alex Smola. 1	• Some of them might look familiar 2
Outline	Tricks for Constructing Kernels
<ul> <li>Tricks for constructing kernels</li> </ul>	• If $k_1$ and $k_2$ are kernels, then so are $\alpha_1 k_1 + \alpha_2 k_2$ for $\alpha_1, \alpha_2 \ge 0$
• String kernels	⇒ If input vectors can be partitioned into subvectors of different types (e.g. strings and real values), can apply <u>direct sum</u> :
Spectrum kernels	$(k_1 \oplus k_2)(x_1, x_2, x_1', x_2') = k_1(x_1, x_1') + k_2(x_2, x_2')$
Locality-improved kernels	where $x_1, x_1' \in \mathcal{X}_1$ (e.g. $\mathbb{R}^n$ ) and $x_2, x_2' \in \mathcal{X}_2$ (e.g. strings)
<ul> <li>Kernels defined on graphs</li> </ul>	$k_{1}k_{2}$
• Sections 13.1–13.3, 13.5, assorted papers	⇒ Similar to application of direct sum, use <u>tensor product</u> : $(k_1 \otimes k_2)(x_1, x_2, x'_1, x'_2) = k_1(x_1, x'_1) k_2(x_2, x'_2)$

# Tricks for Constructing Kernels

**Conformal Transformations** 

- For a real-valued function f, k'(x, x') = f(x)f(x') is a kernel
- This leads to <u>conformal transformations</u>:

$$k_f(x, x') = f(x)k(x, x')f(x')$$

- If k is a kernel, then so is  $k_f$
- Recall that if ||x|| = ||x'|| = 1, then  $\langle x, x' \rangle = \cos(\angle(x, x'))$ ; thus

 $\begin{aligned} \cos(\angle(\Phi_f(x), \Phi_f(x'))) &= \frac{f(x)k(x, x')f(x')}{\sqrt{f(x)k(x, x)f(x)}\sqrt{f(x')k(x', x')f(x')}} \\ &= \frac{k(x, x')}{\sqrt{k(x, x)}\sqrt{k(x', x')}} \\ &= \cos(\angle(\Phi(x), \Phi(x'))) \end{aligned}$ 

I.e. angles in feature space are preserved in a conformal transformation

# Tricks for Constructing Kernels

Convolution Kernels (cont'd)

• Generally, define the set of allowed decompositions as a relation  $R(x_1,\ldots,x_D,x)$  and define the *R*-convolution

$$(k_1 \star \cdots \star k_D)(x, x') := \sum_R \prod_{d=1}^D k_d(x_d, x'_d)$$

(i.e. sum over all allowable decompositions of x into  $x_1, \ldots, x_D$ , etc.)

- Based on earlier results, we know this to be a valid kernel
- A special case: ANOVA kernel of order D

$$k_D(x, x') := \sum_{1 \le i_1 < \dots < i_D \le N} \prod_{d=1}^D k^{(i_d)}(x_{i_d}, x'_{i_d})$$

 $(D = N \Rightarrow \text{tensor prod}, D = 1 \Rightarrow \text{direct sum})$ 

Tricks for Constructing Kernels Convolution Kernels

- Notions of tensor products and direct sums lead to <u>*R*-convolution kernels</u>
- E.g. consider partitioning the string x = ATG into two distinct, contiguous, nonemtpy substrings:

$$R_1: \quad x_{1,R_1} = A \quad \underline{\text{AND}} \quad x_{2,R_1} = TG$$
$$\underline{OR}$$

 $R_2: x_{1,R_2} = AT \quad \underline{\text{AND}} \quad x_{2,R_2} = G$ (similarly, decompose x')

• Now can compute a kernel for each substring of each partitioning and combine:

$$k(x, x') = k_1(x_{1,R_1}, x'_{1,R_1})k_2(x_{2,R_1}, x'_{2,R_1}) +k_1(x_{1,R_2}, x'_{1,R_2})k_2(x_{2,R_2}, x'_{2,R_2})$$

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### String Kernels

- To apply SVMs to text classification, can map documents to <u>bag-of-words</u> representation and use kernels defined on  $\mathbb{R}^n$ 
  - Each dimension is one word, value in that dimension is word frequency
  - Ignores word ordering
- Alternatively, can use a <u>string kernel</u>, which computes similarities between two strings based on their common substrings
- Related to *R*-convolution kernel

#### String Kernels (cont'd)

- Let  $\Sigma$  be a finite alphabet,  $\Sigma^n$  be set of all length-*n* strings over  $\Sigma$ , and  $\Sigma^* = \bigcup_{n=0}^{\infty} \Sigma^n$
- Given  $s \in \Sigma^*$ , let  $\mathbf{i} := (i_1, \dots, i_{|u|})$  be an index sequence with  $1 \leq i_1 < \dots < i_{|u|} \leq |s|$  and  $u := s(\mathbf{i}) := s(i_1) \cdots s(i_{|u|})$  be a (possibly non-contiguous) subsequence of s
- $l(\mathbf{i}) := i_{|u|} i_1 + 1$  is the length of u in s
  - E.g. if s = ABBA, then l(1,2,3) = 3 (for ABB), l(1,4) = 4 (for AA)
  - $\Phi_n(s)$  defines one dimension per substring  $u \in \Sigma^n$ , and the *u*th component of  $\Phi_n(s)$  is

$$[\Phi_n(s)]_u := \sum_{\mathbf{i}:s(\mathbf{i})=u} \lambda^{l(\mathbf{i})}$$

for 0 <  $\lambda \leq$  1

String Kernels

(cont'd)

- To efficiently compute the kernel, define for  $i=1,\ldots,n-1$ 

$$k'_i(s,t) = \sum_{u \in \Sigma^i} \sum_{(\mathbf{i},\mathbf{j}):s(\mathbf{i})=t(\mathbf{j})=u} \lambda^{|s|+|t|-i_1-j_1+2}$$

• Then if  $x \in \Sigma^1$ , can recursively compute  $k_n(s,t)$ :

 $\begin{array}{l} k_0'(s,t) = 1 \ \, \text{for all } s,t \\ k_i'(s,t) = 0 \ \, \text{if } \min(|s|,|t|) < i \\ k_i(s,t) = 0 \ \, \text{if } \min(|s|,|t|) < i \end{array}$ 

$$\begin{aligned} k_i'(sx,t) &= \lambda k_i'(s,t) + \sum_{j:t_j=x} k_{i-1}'(s,t[1,\ldots,j-1])\lambda^{|t|-j+2} \\ k_n(sx,t) &= k_n(s,t) + \sum_{j:t_j=x} k_{n-1}'(s,t[1,\ldots,j-1])\lambda^2 \end{aligned}$$

String Kernels (cont'd)

- E.g. if s = ABBA, then  $[\Phi_2(s)]_{AB} = \lambda^2 + \lambda^3$
- $[\Phi_n(s)]_u$  larger if u (nearly) contiguous and common in s
- The string kernel is then

$$k_n(s,t) = \sum_{u \in \Sigma^n} [\Phi_n(s)]_u [\Phi_n(t)]_u$$
$$= \sum_{u \in \Sigma^n} \sum_{(\mathbf{i},\mathbf{j}):s(\mathbf{i})=t(\mathbf{j})=u} \lambda^{l(\mathbf{i})} \lambda^{l(\mathbf{j})}$$

- If want to vary n, use  $k := \sum_n c_n k_n$
- Since value of  $k_n$  (and therefore k) depend on lengths of s and t, normalize k in feature space

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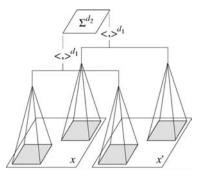
### Spectrum Kernel

- Another type of string kernel
- For a fixed integer  $\gamma \ge 1$ , define the  $\gamma$ -spectrum of a sequence to be the set of all length- $\gamma$  contiguous sequences it contains
- Feature map for spectrum kernel is indexed by all possible length- $\gamma$  subsequences from alphabet  $\Sigma$  (similar to bag of words)
- For each  $a \in \Sigma^{\gamma}$ , let  $\phi_a(x) =$  number of times a occurs in x contiguously
- Now define  $\Phi_{\gamma}(x) = (\phi_a(x))_{a \in \Sigma^{\gamma}}$ 
  - This is a weighted representation of x 's  $\gamma\text{-}$  spectrum
  - A sparse vector

### Locality-Improved Kernels

- Spectrum Kernels (cont'd)
- Can compute  $k_{\gamma}(x, x') = \langle \Phi_{\gamma}(x), \Phi_{\gamma}(x') \rangle$  in time O(|x| + |x'|)
  - 1. Collect set of length- $\gamma$  subsequences of x into array  $A_x$  and sort it (same with x')
    - $A_x$  contains non-zero entries of  $\Phi_{\gamma}(x)$
  - 2. Scan  $A_{x}$  and  $A_{x^{\prime}}\text{,}$  multiplying entries that match, and sum the products

- A variation on existing kernels to emphasize local correlations over long-range (global) ones
- E.g. in image processing, replace polynomial kernel  $\langle x,x'\rangle^d$  with a variant that focuses on subimages first
- Generally, take the dot product over all corresponding subimages of the two images, raise to the  $d_1$  power, sum these values, then raise to the  $d_2$  power



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Locality-Improved Kernels Image Processing (cont'd)

- Specifically:
  - 1. Compute (x. \* x'), the pixel-wise product of x and x'
  - 2. Sample (x.\*x') with pyramidal receptive fields:

$$z_{ij} := \sum_{i',j'} w(\max(|i-i'|,|j-j'|))(x_{\cdot} * x')_{i'j'}$$

where e.g. weighting function  $w(n) = \max(q-n, 0)$ ; i.e. only include pixels in a width-p window (p = 2q + 1) centered at (i, j)

- 3. Raise each  $z_{ij}$  to the  $d_1$  power (this gives local correlations)
- 4. Sum  $z_{ij}^{d_1}$  over entire image and raise this sum to the  $d_2$  power (long-range correlations)
- If  $d_1 = 1$ , get standard polynomial kernel  $\langle x, x' \rangle^{d_2}$

Locality-Improved Kernels Image Processing (cont'd)

Classifier	Error on MNIST (%)
$k^{1,4}$	4.0
$k_{9}^{2,2}$ $k_{9}^{4,1}$	3.1
$k_{9}^{4,1}$	3.4
Virt SV	2.8
VSV $k_9^{2,2}$	2.0

# Locality-Improved Kernels DNA Start Codon Recognition

• Problem: in a DNA sequence (from alpha-

- bet  $\{A, C, T, G\}$ ), identify subsequences that encode genes
  - Typically such a coding region begins with  $\ensuremath{\mathit{ATG}}$
  - But not all ATG occurrences imply a coding region
  - Thus the learning problem is to take a length-200 window centered at an ATG and predict if it's a coding region
- For this problem, long-range dependencies aren't very important, so use a kernel to emphasize local correlations

# Locality-Improved Kernels

DNA Start Codon Recognition (cont'd)

• We'll consider correlations inside small windows of length  $2\ell+1$ :

$$\operatorname{win}_{p}(x, x') = \left(\sum_{j=-\ell}^{+\ell} v_{j} \operatorname{match}_{p+j}(x, x')\right)^{d_{1}}$$

where  $\operatorname{match}_{p+j}(x, x') = 1$  if x and x' match at position p + j and 0 otherwise, and  $v_j$  is a weight for window position j (larger near 0)

• Now we sum the values of win<sub>p</sub>:

$$k(x, x') = \left(\sum_{p=1}^{\ell} \operatorname{win}_p(x, x')\right)^{d_2}$$

(Should summation really be only to  $\ell$ ?)

Classifier	Error (%)
ANN	15.4
Poly kernel, $d = 1$	13.8
L-I kernel, $d_1 = 4, \ell = 4$	11.9
Codon-improved kernel, $d_1 = 2, \ell = 3$	12.2
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## **Complete Graph Kernels**

- A complete graph kernel k is one whose implicit remapping  $\Phi : \mathcal{G} \to \mathcal{H}$  distinguishes all pairs of graphs  $(G, G') \in \mathcal{G} \times \mathcal{G}$ , i.e.  $\Phi$  is injective
- Example (Subgraph feature space): Let each dimension in  $\Phi(G)$  correspond to a distinct connected graph  $H \in \mathcal{G}$ . Then  $[\Phi(G)]_H =$  number of times an isomorphism of H appears in G.
- Gärtner et al. [2003] showed that for injective  $\Phi$ ,  $k(G,G) + k(G',G') - 2k(G,G') = \langle \Phi(G) - \Phi(G'), \Phi(G) - \Phi(G') \rangle = 0$  iff  $G \simeq G'$ 
  - $\Rightarrow$  Computing k is as hard as graph isomorphism, for which no efficient algorithm is currently known
- Further, the kernel for the subgraph mapping is in fact NP-hard to compute (reduce from Hamiltonian path), even to approximate and/or if *H* comes from a restricted class of graphs

# Kernels on Graphs

- Very general form of structured data
- Can represent many data types, including chemical structures
- Will consider directed graphs with labels on edges and nodes
- Let *G* be the space of all graphs, modulo isomorphism

### Kernels Based on Label Pairs

- Now consider more restrictive kernels that can be efficiently considered
- Focus on graphs with labels on nodes but not edges; labels come from \$\mathcal{L} = \{\ell\_1, \ldots, \ell\_m\}\$
- Let <u>label matrix</u> L be such that  $[L]_{ri} = 1$  if node  $v_i$ 's label is  $\ell_r$  and  $[L]_{ri} = 0$  otherwise
- Let adjacency matrix E be such that  $[E]_{ij} = 1$ if directed edge  $(v_i, v_j)$  exists in graph G and  $[E]_{ij} = 0$  otherwise;  $[E^n]_{ij}$  is number of lengthn walks from  $v_i$  to  $v_j$
- $\left[LL^{\top}\right]_{rr}$  = number of times label  $\ell_r$  is assigned to a vertex in G
- $\left[ LE^nL^\top \right]_{ij}$  = number of walks of length n between vertices labeled  $\ell_i$  and vertices labeled  $\ell_j$

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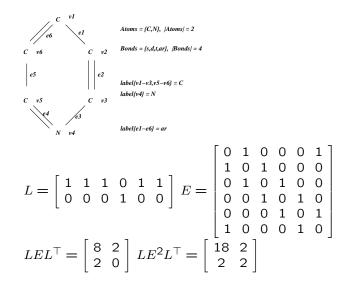
# Kernels Based on Label Pairs (cont'd)

- $\mathcal{W}_n(G) = \text{set of all } n\text{-edge walks in } G$
- For walk  $w \in W_n(G)$ ,  $l_1(w) =$  label of first vertex of w and  $l_{n+1}(w) =$  label of last vertex
- $\lambda =$  sequence of nonnegative weights  $\lambda_0, \lambda_1, \dots$
- Define mapping Φ(G) to have one feature per pair of labels (ℓ<sub>i</sub>, ℓ<sub>j</sub>): [Φ(G)]<sub>ℓ<sub>i</sub>,ℓ<sub>j</sub></sub> =

 $\sum_{n=0}^{\infty} \lambda_n \left| \left\{ w \in \mathcal{W}_n(G) : l_1(w) = \ell_i \wedge l_{n+1}(w) = \ell_j \right\} \right|$ 

i.e. the weighted sum of the number of lengthn walks from an  $\ell_i$ -labeled vertex to an  $\ell_j$ labeled vertex, weighted by  $\lambda_n$ , summed over all  $n\to\infty$ 

# Matrix Example



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### Kernels Based on Label Pairs (cont'd)

• Thus kernel is  $\langle \Phi(G), \Phi(G') \rangle =$ 

$$\left(L\left(\sum_{i=0}^{\infty}\lambda_{i}E^{i}\right)L^{\top},L'\left(\sum_{i=0}^{\infty}\lambda_{i}E'^{i}\right)L'^{\top}\right)$$

- Under certain conditions, can efficiently compute the matrix power series
- E.g. if  $\lambda_i = \beta^i/i!$  for some  $\beta > 0$  and if E can be diagonalized such that  $E = T^{-1}DT$ , then  $E^n = T^{-1}D^nT$  and  $[D^n]_{ii} = [D_{ii}]^n$  since D is diagonal
- Now we can compute

as

$$T^{-1}\left(\lim_{n\to\infty}\sum_{i=0}^n \frac{\beta^i D^i}{i!}\right)T$$
,

 $\lim_{n \to \infty} \sum_{i=0}^{n} \frac{(\beta E)^{i}}{i!}$ 

where limits are taken component-wise

## Kernels Based on Contiguous Label Sequences

- Previous kernel's mapping  $\Phi$  has a low-dimensional feature space:  $|\mathcal{L}|^2$ 
  - $\Rightarrow$  E.g. if all node labels are C or N, then feature space has dimension 4
- For a more expressive feature mapping, will use mapping with one dimension per <u>label sequence</u> rather than <u>label pair</u>
- Assume we have labels for both nodes and edges; if nodes or edges are not labeled, use generic symbol '#'

# Kernels Based on Contiguous Label Sequences

(cont'd)

- Let  $S_n$  be set of all possible label sequences of walks with n edges and let  $\lambda$ ,  $W_n(G)$ , and  $l_i(w)$  be as before
- Define mapping  $\Phi(G)$  to have one feature per possible label sequence  $s \in \bigcup_n S_n$ :

 $[\Phi(G)]_s = \sqrt{\lambda_n} |\{w \in \mathcal{W}_n(G) : \forall i \ s_i = l_i(w)\}|$ 

i.e. the number of walks in G with n edges whose (vertex and edge) label sequences match  $s = s_1, s_2, \ldots, s_{2n+1} \in S_n$ , weighted by  $\sqrt{\lambda_n}$ 

Topic summary due in 1 week!

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#### Kernels Based on Contiguous Label Sequences (cont'd)

• To compute the kernel, use the notion of a product graph: given  $G_1 = (\mathcal{V}_1, \mathcal{E}_1)$  and  $G_2 = (\mathcal{V}_2, \mathcal{E}_2)$ ,  $G_{\times} = G_1 \times G_2$  is defined as

$$\mathcal{V}_{\times} = \{(v_1, v_2) \in \mathcal{V}_1 \times \mathcal{V}_2 : label(v_1) = label(v_2)\}$$

$$\mathcal{E}_{\times} = \{((u_1, u_2), (v_1, v_2)) \in \mathcal{V}_{\times}^2 : (u_1, v_1) \in \mathcal{E}_1$$

 $\wedge (u_2, v_2) \in \mathcal{E}_2 \wedge label(u_1, v_1) = label(u_2, v_2) \}$ 

• One can show that

$$|\{w \in \mathcal{W}_n(G_1 \times G_2) : \forall i \ s_i = l_i(w)\}|$$
$$= |\{w \in \mathcal{W}_n(G_1) : \forall i \ s_i = l_i(w)\}|$$
$$\cdot |\{w \in \mathcal{W}_n(G_2) : \forall i \ s_i = l_i(w)\}|$$

• Since an *n*-edge walk in  $G_1 \times G_2$  corresponds to a walk in each of  $G_1$  and  $G_2$ , each with same label sequence, the dot product  $\langle \Phi(G_1), \Phi(G_2) \rangle$ can be computed as

$$k_{\times}(G_1, G_2) = \sum_{i,j=1}^{\mathcal{V}_{\times}} \left[ \sum_{n=0}^{\infty} \lambda_n E_{\times}^n \right]_{ij}$$