

ASSIGNMENT 4

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Q1.

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Q1. To extend LDA to non-linear mappings, the data can be mapped to a new feature space, F , via some function ϕ . In this new feature space, the function that needs to be maximised is:

$$J(w) = \frac{w^T S_B \phi w}{w^T S_w \phi w}$$

where

$$S_B \phi = (m_2 \phi - m_1 \phi)(m_2 \phi - m_1 \phi)^T$$

$$S_w \phi = \sum_{i=1,2} \sum_{n=1}^{L_i} (\phi(x_n^i) - m_i \phi)(\phi(x_n^i) - m_i \phi)^T$$

$$m_i \phi = \frac{1}{L_i} \sum_{j=1}^{L_i} \phi(x_j^i)$$

Note $w \in F$. We use kernel trick in which the dot product in the new feature space is replaced by a kernel function.

$$K(x, y) = \phi(x) \cdot \phi(y).$$

LDA now can be reformulated in terms of dot products by noting that

$$w = \sum_{i=1}^L \alpha_i \phi(x_i).$$

$$w^T m_i \phi = \frac{1}{L_i} \sum_{j=1}^{L_i} \sum_{k=1}^{L_i} \alpha_j K(x_j, x_k^i) = \alpha^T M_i.$$

where $(M_i)_j = \frac{1}{L_i} \sum_{k=1}^{L_i} K(x_j, x_k^i).$

numerator of $T(w)$ can be written as:

$$w^T S_B \phi w = w^T (m_2 \phi - m_1 \phi) (m_2 \phi - m_1 \phi)^T w \\ = \alpha^T M \alpha$$

$$\text{where } M = (M_2 - M_1) \cdot (M_2 - M_1)^T$$

$$\text{denominator } \rightarrow w^T S_w \phi w = \alpha^T N \alpha$$

$$\text{where } N = \sum_{j=1,2} K_j (I - 1_{L_j}) K_j^T$$

with n^{th} , m^{th} component of K_j defined as $k(x_n, x_m^j)$.

I is the identity matrix, and 1_{L_j} the matrix with all entries equal to $1/L_j$.

$$\begin{aligned} w^T S_w \phi w &= \left(\sum_{i=1}^L \alpha_i \phi^T(x_i) \right) \left(\sum_{j=1,2}^{L_j} \sum_{n=1}^L (\phi(x_n^j) - m_j \phi) (\phi(x_n^j) - m_j \phi)^T \right) \\ &= \sum_{j=1,2} \sum_{n=1}^L \sum_{p=1}^{L_j} \sum_{k=1}^L \alpha_i \phi^T(x_i) (\phi(x_n^j) - m_j \phi) (\phi(x_n^j) - m_j \phi)^T \alpha_k \phi(x_k) \\ &= \sum_{j=1,2} \sum_{n=1}^L \sum_{p=1}^{L_j} \sum_{k=1}^L \left(\alpha_i k(x_i, x_n^j) - \frac{1}{L_j} \sum_{p=1}^{L_j} \alpha_i k(x_i, x_p^j) \right) \\ &\quad \left(\alpha_k k(x_k, x_n^j) - \frac{1}{L_j} \sum_{p=1}^{L_j} \alpha_k k(x_k, x_p^j) \right) \\ &= \sum_{j=1,2} \left(\sum_{i=1}^L \sum_{n=1}^{L_j} \sum_{k=1}^L \left(\alpha_i \alpha_k k(x_i, x_n^j) - \frac{\alpha_i \alpha_k}{L_j} \sum_{p=1}^{L_j} k(x_i, x_p^j) \right) \right. \\ &\quad \left. k(x_k, x_p^j) \right) \end{aligned}$$

$$= \sum_{j=1,2} \alpha^T x_j k_j^T \alpha - \alpha^T x_j \frac{1}{l_j} k_j^T \alpha$$

$$= \alpha^T N \alpha.$$

With these equations for the numerator & denominator of $J(\omega)$, the equation for J can be rewritten as

$$J(\alpha) = \frac{\alpha^T M \alpha}{\alpha^T N \alpha}.$$

Differentiating w.r.t α and setting equal to 0 gives.

$$(\alpha^T M \alpha) N \alpha = (\alpha^T N \alpha) \cdot M \alpha.$$

Since only the direction of ω , and hence direction of α , matters, the above can be solved for α as

$$\alpha = N^{-1} (M_2 - M_1)$$

N is usually singular, so multiple of I is added to it

$$N_\epsilon = N + \epsilon \cdot I.$$

Given the solution for α , the projection of a new data point is given by.

$$y(x) = (\omega \cdot \phi(x)) = \sum_{i=1}^L \alpha_i \cdot k(x_i, x).$$

Q2.

Arcene- Kernel PCA

```
a = importdata('arcene_train.data');
b = importdata('arcene_train.labels');
c = importdata('arcene_valid.data');
d = importdata('arcene_valid.labels');
sigma = 10000;
[x1 y1] = size(a);
[x2 y2] = size(c);
kernel = zeros(x1,x1);

for i=1:x1
    for j=1:x1
        kernel(i,j) = exp(-norm(a(i,:)-a(j,:))^2/sigma^2);
    end
end
temp = ones(x1,x1)/x1;
kernel_new = kernel - temp*kernel - kernel*temp + temp*kernel*temp;
[eigenvectors1 eigenvalues1] = eig(kernel_new);
eigenvalues1 = diag(eigenvalues1);

%eigenvalues1 = eigenvalues1(end:-1:1);
%eigenvectors1 = fliplr(eigenvectors1);

kernel_t = zeros(x2,x1);
for i=1:x2
    for j=1:x1
        kernel_t(i,j) = exp(-norm(c(i,:)-a(j,:))^2/sigma^2);
    end
end

for i=1:x1
    eigenvectors1(:,i) = eigenvectors1(:,i)/eigenvalues1(i);
```

end

for t=1:2

v1 = eigenvectors1(:,1:10^t);

train1 = kernel_new*v1;

test1 = kernel_t*v1;

trainmodel1 = svmtrain(train1,b);

acc_linear=100*(size(find(svmclassify(trainmodel1,test1)==d),1)/size(train1,1));

disp(acc_linear);

trainmodel1 = svmtrain(train1,b,'kernel_function','rbf','rbf_sigma',5);

acc_rbf=100*(size(find(svmclassify(trainmodel1,test1)==d),1)/size(train1,1));

disp(acc_rbf);

end

Arcene- Kernel LDA

a = importdata('arcene_train.data');

b = importdata('arcene_train.labels');

c = importdata('arcene_valid.data');

d = importdata('arcene_valid.labels');

sigma = 10000;

[x1 y1] = size(a);

[x2 y2] = size(c);

kernel = zeros(x1,x1);

for i=1:x1

for j=1:x1

kernel(i,j) = exp(-norm(a(i,:)-a(j,:))^2/sigma^2);

end

end

```
temp = ones(x1,x1)/x1;  
kernel_n = kernel - temp*kernel - kernel*temp + temp*kernel*temp;
```

```
kernel_t = zeros(x2,x1);  
for i=1:x2  
    for j=1:x1  
        kernel_t(i,j) = exp(-norm(c(i,:)-a(j,:))^2/sigma^2);  
    end  
end
```

```
m1ind = find(b==1);  
m2ind = find(b==-1);  
M1 = mean(kernel_n(m1ind,:));  
M2 = mean(kernel_n(m2ind,:));  
N =  
kernel_n(m1ind,:)*(eye(size(m1ind,1))-(1/size(m1ind,1)))*kernel_n(m1ind,:) +  
kernel_n(m2ind,:)*(eye(size(m2ind,1))-(1/size(m2ind,1)))*kernel_n(m2ind,:);  
N1 = N + 644*eye(size(kernel_n,1));  
N = N + 8000*eye(size(kernel_n,1));  
train = kernel_n*inv(N)*(M1-M2)';  
test = kernel_t*inv(N)*(M1-M2)';  
train1 = kernel_n*inv(N1)*(M1-M2)';  
test1 = kernel_t*inv(N1)*(M1-M2)';  
trainmodel = svmtrain(train,b);  
accuracy = size(find(svmclassify(trainmodel, test)==d),1);  
trainmodel = svmtrain(train1,b,'kernel_function','rbf');  
accuracy1 = size(find(svmclassify(trainmodel, test1)==d),1);  
accuracy  
accuracy1
```

	PCA	LDA																
Arcene	<table><tr><td>K=10</td><td>Linear</td><td>56.0</td></tr><tr><td>K=10</td><td>RBF</td><td>70.0</td></tr><tr><td>K=100</td><td>Linear</td><td>56.0</td></tr><tr><td>K=100</td><td>RBF</td><td>68.0</td></tr></table>	K=10	Linear	56.0	K=10	RBF	70.0	K=100	Linear	56.0	K=100	RBF	68.0	<table><tr><td>Linear</td><td>68.0</td></tr><tr><td>RBF</td><td>69.0</td></tr></table>	Linear	68.0	RBF	69.0
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Observation:

- More dimension PCA increases the predictability of data.
- Gaussian Kernel give poor accuracies in case of PCA for K =10 , better results for K =100.
- Gaussian kernel gives poor results in LDA compared to Linear Kernel.