

Consider the following training data points and their corresponding  $\alpha$  values

$$\mathbf{x}^1 = \begin{bmatrix} 2 \\ 0 \\ -1.5 \end{bmatrix}, y^1 = +1, \alpha^1 = 0.5 \quad \mathbf{x}^2 = \begin{bmatrix} 1 \\ -0.5 \\ 0 \end{bmatrix}, y^2 = +1, \alpha^2 = 2$$

$$\mathbf{x}^3 = \begin{bmatrix} 5 \\ 0 \\ 2 \end{bmatrix}, y^3 = -1, \alpha^3 = 2 \quad \mathbf{x}^4 = \begin{bmatrix} 13 \\ -8 \\ 7 \end{bmatrix}, y^4 = -1, \alpha^4 = 0$$

$$\mathbf{x}^5 = \begin{bmatrix} 7 \\ -6 \\ 0 \end{bmatrix}, y^5 = -1, \alpha^5 = 0.5 \quad \mathbf{x}^6 = \begin{bmatrix} -5 \\ 0 \\ -5 \end{bmatrix}, y^6 = +1, \alpha^6 = 0$$

Which of the above are considered support vectors?

Any point with  $\alpha > 0$ . In other words:  $\mathbf{x}^1$ ,  $\mathbf{x}^2$ ,  $\mathbf{x}^3$ , and  $\mathbf{x}^5$

Use the above data to compute the separating hyperplane  $\mathbf{w}$  found by the linear SVM.

$$\begin{aligned} \mathbf{w} &= \sum_i \alpha^i y^i \mathbf{x}^i = 0.5 \times 1 \times \begin{bmatrix} 2 \\ 0 \\ -1.5 \end{bmatrix} + 2 \times 1 \times \begin{bmatrix} 1 \\ -0.5 \\ 0 \end{bmatrix} + 2 \times -1 \times \begin{bmatrix} 5 \\ 0 \\ 2 \end{bmatrix} + 0.5 \times -1 \times \begin{bmatrix} 7 \\ -6 \\ 0 \end{bmatrix} \\ &= \begin{bmatrix} 10.5 \\ 2 \\ -4.75 \end{bmatrix} \end{aligned}$$

Consider a Bayesian classifier and a data set of containing 30 features. Assume there are 100 training data points. We report the highest Likelihood classifier given  $n$  features below:

#Feats	1	2	3	4	5	6	7	...	21
MaxL	$6 \times 10^{-23}$	$1 \times 10^{-21}$	$2 \times 10^{-20}$	$4 \times 10^{-20}$	$8 \times 10^{-19}$	$2 \times 10^{-19}$	$4 \times 10^{-19}$		$1 \times 10^{-9}$

#Feats	22	23	24	25	26	27	28	29	30
MaxL	$1 \times 10^{-8}$	$1 \times 10^{-7}$	$6 \times 10^{-7}$	$4 \times 10^{-6}$	$1 \times 10^{-5}$	$5 \times 10^{-5}$	$9 \times 10^{-5}$	$2 \times 10^{-4}$	$3 \times 10^{-4}$

What is the first feature removal step (how many features left?) when AIC no longer improves?

AIC:  $\log(L) - k$ :

k	30	29	28	27	26	
AIC	-38.1	-37.5	-37.3	-36.9	-37.5	

**Removing 4th feature, no longer benefits AIC.**

What is the first feature selection step (how many features chosen?) when BIC no longer improves?

BIC:  $\log(L) - 0.5 \times k \times \log(m)$ :

$\log(100) = 4.6$

k	1	2	3	4		
AIC	-53.5	-53.0	-52.3	-53.9		

**Adding 4th feature, no longer benefits BIC**

Consider the following Principal Components:

$$\mathbf{u}_1 = \begin{bmatrix} 0.41 \\ -0.41 \\ 0 \\ 0.82 \end{bmatrix} \quad \mathbf{u}_2 = \begin{bmatrix} 0 \\ 0.53 \\ 0.80 \\ 0.27 \end{bmatrix}$$

For each data point below, calculate the weights  $z$  for each principal component.

$$\mathbf{x}^1 = \begin{bmatrix} -4.4 \\ 3.8 \\ -1.0 \\ -9.2 \end{bmatrix}$$

$$z_j^i = \mathbf{u}_j^T \mathbf{x}^i \quad z_1^1 = \mathbf{u}_1^T \mathbf{x}^1 = -4.4 \times 0.41 + 3.8 \times -0.41 - 1.0 \times 0 - 9.2 \times 0.82 \approx -11$$

$$z_2^1 = \mathbf{u}_2^T \mathbf{x}^1 = -4.4 \times 0 + 3.8 \times 0.53 - 1.0 \times 0.8 - 9.2 \times .27 \approx -1.3$$

$$\mathbf{x}^2 = \begin{bmatrix} 0.1 \\ 0.7 \\ 1.2 \\ 0.7 \end{bmatrix}$$

$$z_j^i = \mathbf{u}_j^T \mathbf{x}^i \quad z_1^2 = \mathbf{u}_1^T \mathbf{x}^2 = 0.1 \times 0.41 + 0.7 \times -0.41 + 1.2 \times 0 + 0.7 \times 0.82 \approx 0.3$$

$$z_2^2 = \mathbf{u}_2^T \mathbf{x}^2 = 0.1 \times 0 + 0.7 \times 0.53 + 1.2 \times 0.8 + 0.7 \times .27 \approx 1.5$$

$$\mathbf{x}^3 = \begin{bmatrix} 2.3 \\ -2.5 \\ -0.3 \\ 4.4 \end{bmatrix}$$

$$z_j^i = \mathbf{u}_j^T \mathbf{x}^i \quad z_1^2 = \mathbf{u}_1^T \mathbf{x}^2 = 2.3 \times 0.41 - 2.5 \times -0.3 + 1.2 \times 0 + 4.4 \times 0.82 \approx \mathbf{5.3}$$

$$z_2^1 = \mathbf{u}_2^T \mathbf{x}^1 = 2.3 \times 0 - 2.5 \times 0.53 - 0.3 \times 0.8 + 4.4 \times .27 \approx \mathbf{-0.4}$$

Using both principal components, compute the reconstruction for each of the above data points.

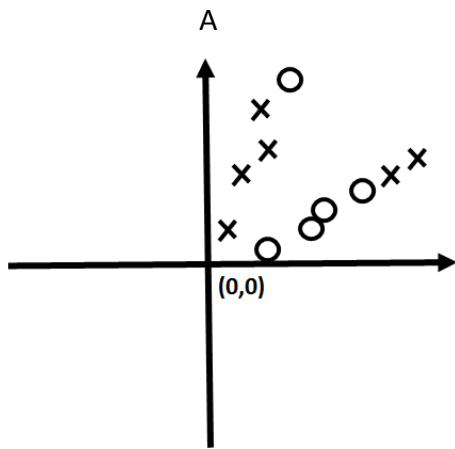
$$\tilde{\mathbf{x}}^i = z_1^i \mathbf{u}_1 + z_2^i \mathbf{u}_2$$

$$\tilde{\mathbf{x}}^1 = \begin{bmatrix} -4.5 \\ 4.5 \\ 0 \\ -9.02 \end{bmatrix} + \begin{bmatrix} 0 \\ -0.69 \\ -1.0 \\ -0.35 \end{bmatrix} = \begin{bmatrix} \mathbf{-4.5} \\ \mathbf{3.81} \\ \mathbf{-1.0} \\ \mathbf{-9.37} \end{bmatrix}$$

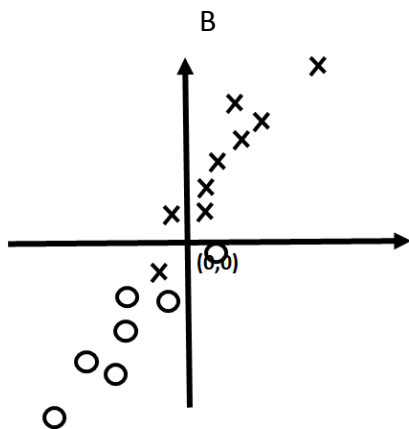
$$\tilde{x}^2 = \begin{bmatrix} 0.12 \\ -0.12 \\ 0 \\ 0.25 \end{bmatrix} + \begin{bmatrix} 0 \\ 0.80 \\ 1.2 \\ 0.41 \end{bmatrix} = \begin{bmatrix} \mathbf{0.12} \\ \mathbf{0.68} \\ \mathbf{1.2} \\ \mathbf{0.66} \end{bmatrix}$$

$$\tilde{x}^3 = \begin{bmatrix} 2.2 \\ -2.2 \\ 0 \\ 4.3 \end{bmatrix} + \begin{bmatrix} 0 \\ -0.21 \\ -0.32 \\ -0.11 \end{bmatrix} = \begin{bmatrix} 2.2 \\ -2.41 \\ -0.32 \\ -4.19 \end{bmatrix}$$

We observe the following data points in two dimensions. We wish to define new dimensions to better describe the data. Should we use PCA, ICA, or NMF?

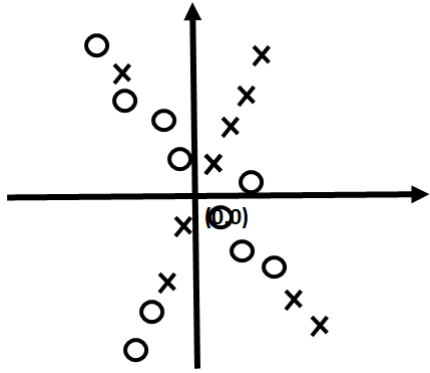


**NMF** – all features are positive



**PCA** – one main direction of variance

C



**ICA** – two non-orthogonal directions of variance, both positive and negative

Consider an HMM with 3 states:

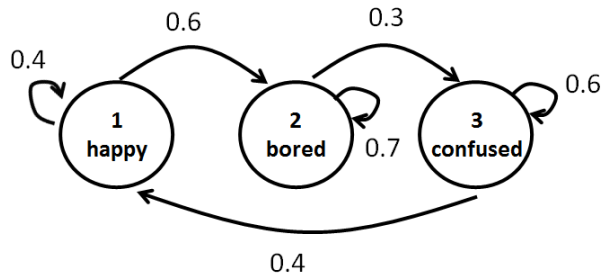
1: happy, 2: bored, 3: confused

And 4 potential outputs at each time (each output is punctuation mark)

1: ! 2: ? 3: . 4: -

The transition matrix is capture by the diagram

The initial probabilities are:



$$\begin{aligned}\pi_{happy} &= 0.2 \\ \pi_{bored} &= 0.6 \\ \pi_{confused} &= 0.2\end{aligned}$$

The emission matrix is  $\phi$ :

State\Obs	!	?	.	-
Happy	0.8	0.1	0.1	0
Bored	0	0.2	0.7	0.1
Confused	0.1	0.7	0	0.2

Which of the following are impossible state sequences:

a) happy, confused, happy, happy

**impossible**

b) bored, bored, bored, confused

**possible**



c) happy, happy, bored, confused, happy

**possible**

d) bored, bored, happy, happy

**impossible**

Presume the sequence:

?, ?, ?, ?

What is the forward value  $\alpha_1(\text{happy})$

$$\phi_{?,\text{happy}}\pi_{\text{happy}} = 0.1 \times 0.2 = \mathbf{0.02}$$

What is the forward value  $\alpha_2(\text{confused})$

$$\alpha_2(\text{confused}) = \phi_{?,\text{confused}} \sum_{\text{mood}} A_{\text{confused},\text{mood}} \alpha_1(\text{mood})$$

Compute  $\alpha_1(\text{mood})$ :

$$\alpha_1(\text{happy}) = 0.02 \text{ (see above)} \quad \alpha_1(\text{bored}) = 0.2 \times 0.6 = 0.12$$

$$\alpha_1(\text{confused}) = 0.7 \times 0.1 = 0.07$$

$$\begin{aligned} \alpha_2(\text{confused}) &= \phi_{?,\text{confused}} \sum_{\text{mood}} A_{\text{confused},\text{mood}} \alpha_1(\text{mood}) \\ &= 0.7 \times [0 \times 0.2 + 0.3 \times 0.12 + 0.6 \times 0.07] \approx \mathbf{0.06} \end{aligned}$$

Final answer: **0.06**

**Note: Viterbi was only briefly covered in class. I include this problem for review/clarification of the process but will make Viterbi worth few points (if any) on the exam.**

Using the Viterbi algorithm, what is the sequence of most probable states for each observation sequence below:

!, ?, !

First compute  $\delta_t(i)$  for all  $t$  and  $i$

$$\delta_1(i) = \pi_i \phi_{!,i} \quad \delta_1(happy) = 0.2 \times 0.8 = 0.16 \quad \delta_1(bored) = 0.6 \times 0 = 0$$

$$\delta_1(confused) = 0.2 \times 0.1 = 0.02$$

$$\delta_2(i) = \phi_{?,i} \max_j A_{i,j} \delta_1(j)$$

$$\delta_2(happy) = 0.1 \times \max_j A_{happy,j} \delta_1(j) = 0.1 \times \max \begin{cases} 0.4 \times 0.16 \\ 0 \times 0 \\ 0.4 \times 0.02 \end{cases} = 0.0064$$

$$\delta_2(bored) = 0.2 \times \max_j A_{bored,j} \delta_1(j) = 0.2 \times \max \begin{cases} 0.6 \times 0.16 \\ 0.7 \times 0 \\ 0 \times 0.02 \end{cases} = 0.0192$$

$$\delta_2(confused) = 0.7 \times \max_j A_{confused,j} \delta_1(j) = 0.7 \times \max \begin{cases} 0 \times 0.16 \\ 0.3 \times 0 \\ 0.6 \times 0.02 \end{cases} = 0.0084$$

$$\delta_3(i) = \phi_{!,i} \max_j A_{i,j} \delta_2(j)$$

$$\delta_3(happy) = 0.2 \times \max_j A_{happy,j} \delta_2(j) = 0.2 \times \max \begin{cases} 0.4 \times 0.0064 \\ 0 \times 0.0192 \\ 0.4 \times 0.0084 \end{cases} = 0.00269$$

$$\delta_3(bored) = 0.6 \times \max_j A_{bored,j} \delta_2(j) = 0 \times \max \begin{cases} 0.6 \times 0.0064 \\ 0.7 \times 0.0192 \\ 0 \times 0.0084 \end{cases} = 0$$

$$\delta_3(confused) = 0.2 \times \max_j A_{confused,j} \delta_2(j) = 0.1 \times \max \begin{cases} 0 \times 0.0064 \\ 0.3 \times 0.0192 \\ 0.6 \times 0.0084 \end{cases} = 0.00058$$

Find largest  $\delta_3(i)$ : **q<sub>3</sub>=happy**

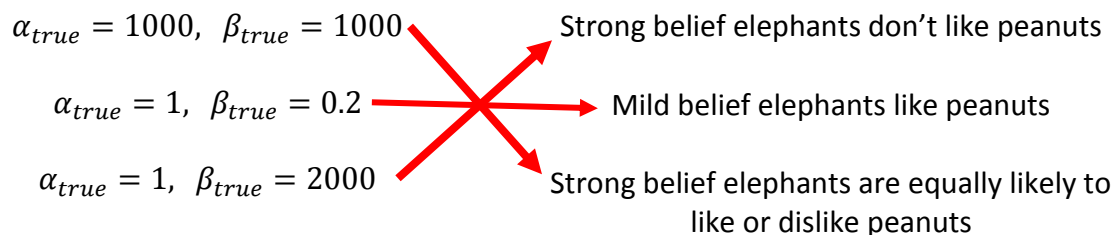
What  $\delta_2(i)$  maximizes  $\delta_3(happy)$ ?: **q<sub>2</sub>=confused**

What  $\delta_1(i)$  maximizes  $\delta_2(confused)$ ? **q<sub>1</sub>=confused**

How can we compute the emission probabilities for each state if we have a set of training sequences where both the observations and the underlying states are known? (Describe in a sentence and/or write a mathematical expression.)

For each entry  $\phi_{obs,state}$  count the number of times that *state* occurs and the number of times the *state* occurs with the *obs* observation. Then  $\phi_{obs,state} = \frac{\#D(obs \wedge state)}{\#D(state)}$

We seek to apply MAP to learn  $P(X|Y; \theta)$ , where X is a binary feature “likes peanuts?” (yes or no) and Y represents the class “is an elephant” (true or false). We use  $\alpha_{true}$  and  $\beta_{true}$  to represent the probability prior belief X=yes and X=no given Y=true. **Match each  $\alpha_{true}, \beta_{true}$  pairing on the left with its corresponding meaning on the right.**



What are the effects of changing the following parameters in gradient ascent learning for logistic regression?

1:  $\varepsilon$

**Decrease in  $\varepsilon$  decreases the amount of change to the learned classifier parameters at each update. Increasing  $\varepsilon$  will increase the amount of change.**

2:  $\lambda$

**Increase in  $\lambda$  will decrease the effect of the corresponding regularization term. Decreasing  $\lambda$  will increase the effect of the corresponding regularization term.**

We have data points with 10 features each. Each feature is a numeric value along the real number line. We wish to learn a classifier to label each data point as one of four classes. Using logistic regression (including +b term!), how many parameters must we learn?

**We have  $(4-1) \times (10+1) = 3 \times 11 = 33$  features to learn. 11 parameters per classifier, 3 classifiers total. 11 parameters correspond to 10 features plus an offset (+b).**

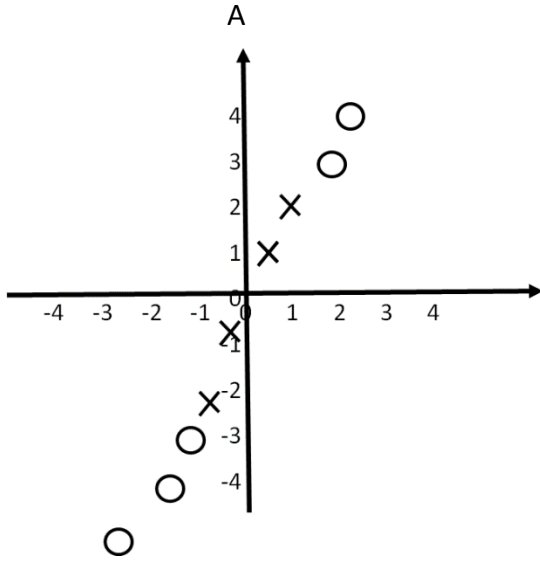
How many parameters must we learn if we wish to classify the data above into only **two** classes using a linear SVM?

**11 parameters**, including **w** (10 features) and **b**.

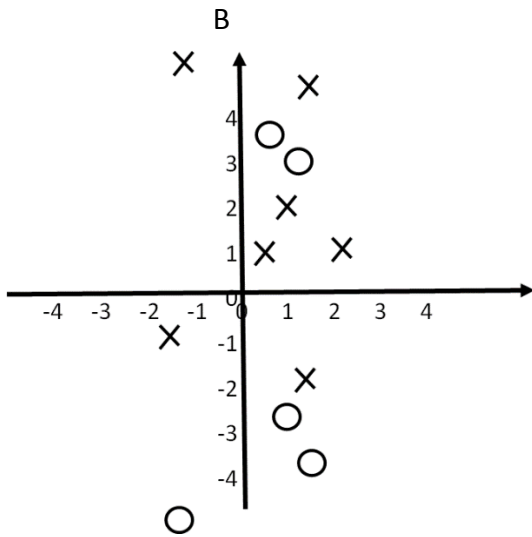
How many parameters must we learn if we wish to classify the data above into only two classes using a quadratic kernel SVM?

**There are two potential interpretations. Either we can say we learn  $10^2+1 = 101$  parameters, for all quadratic combinations of all features, or we say the number of parameters are the number of training point weights  $\alpha^i$ . Most of these weights will be drive to zero and the remaining ones will be used to identify support vectors for classification.**

Define a mapping function that will allow a linear separator to distinguish between the two classes for each set of data points. ( $x_1$  is horizontal axis,  $x_2$  is vertical axis)



$$\varphi(x_1, x_2) = (2x_1 + x_2)^2$$

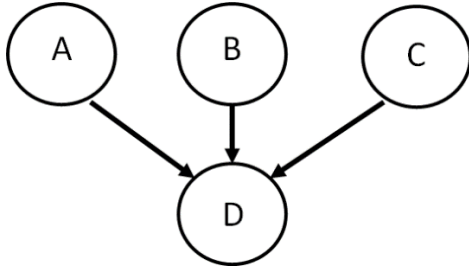


$$\varphi(x_1, x_2) = (x_2, x_2^2, x_2^3)$$

This will allow us to find a cubic function along  $x_2$  that will rise above 0 and fall below 0 four times, corresponding to the four alternating regions in the y axis.

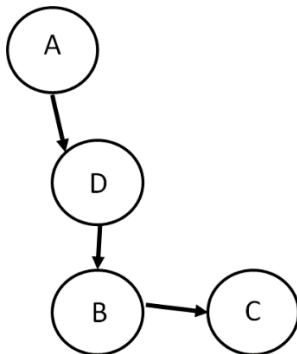
Write the formula to find the total joint probability for each of the following Bayesian Networks (to find  $P(A,B,C,D)$ ).

A:



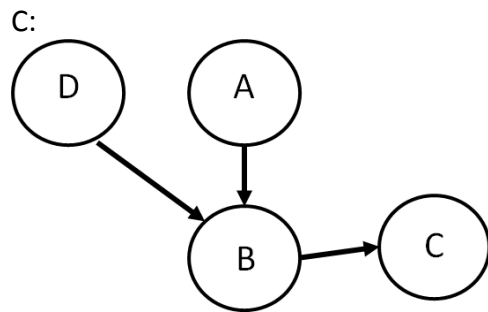
$$P(A)P(B)P(C)P(D|A,B,C)$$

B:



$$P(A)P(D|A)P(B|D)P(C|B)$$

This is basically a Markov model!



Compute  $P(\text{subset of variables})$

$\sum_{\{vars \text{ not in subset}\}} \mathbf{P(D)P(A)P(B|A, D)P(C|B)}$  – marginalize over variables not in subset

For each Bayes net above, how can we find the value of  $P(B)$

Example A:

$\mathbf{P(B)}$

Example B:

$\sum_{A,D} \mathbf{P(A)P(D|A)P(B|D)}$



Example C:

$$\sum_{A,D} P(D)P(A)P(B|A,D)$$

Let us presume we wish to classify whether an animal is a mammal or a bird, and we use 30 features  $x_1, \dots, x_{30}$  as a basis for classification. (For example,  $x_1$  can be size,  $x_2$  can be typical speed of motion,  $x_3$  can be blood temperature, etc.) For a given animal we observe the

following feature vector:  $\mathbf{x} = \begin{bmatrix} 10 \\ 1 \\ 55 \\ \vdots \\ 2 \\ 16 \\ 240 \end{bmatrix}$

and we are informed of the following probabilities:

$$P(x_1=10|Y=\text{bird})=0.01 \quad P(x_2=1|Y=\text{bird})=0.2 \quad \dots \quad P(x_{29}=16|Y=\text{bird})=0.05 \\ P(x_{30}=240|Y=\text{bird})=0.2$$

$$P(x_1=10|Y=\text{mammal})=0.06 \quad P(x_2=1|Y=\text{mammal})=0.08 \quad \dots \quad P(x_{29}=16|Y=\text{mammal})=0.1 \\ P(x_{30}=240|Y=\text{mammal})=0.07$$

If we use Naïve Bayes on a standard computer, we will find  $P(x_1, \dots, x_{30}|Y=\text{bird})=0$  and  $P(x_1, \dots, x_{30}|Y=\text{mammal})=0$ , despite the fact that **none** of the terms  $P(x_i|Y=\text{mammal})=0$  nor  $P(x_i|Y=\text{bird})=0$

a) Why does this happen on a computer?

**When fractional values become sufficiently low, the computer will round to 0.**

b) What mathematical operation can we use to prevent this problem?

**We can use logarithms to convert the product of low probabilities, like  $10^{-20}$  to the sum of 2-3 digit negative numbers like -50.**