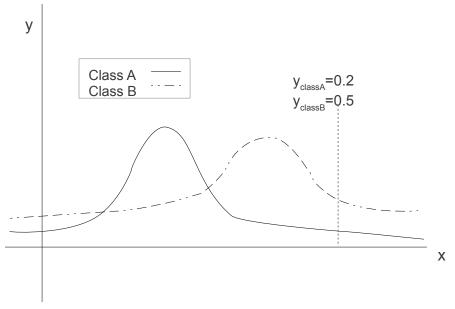
CIS501 Final exam, Fall 2012

Answer all questions. Unless stated otherwise, select a single **best** answer to each question. Questions carry one mark each unless marked otherwise.

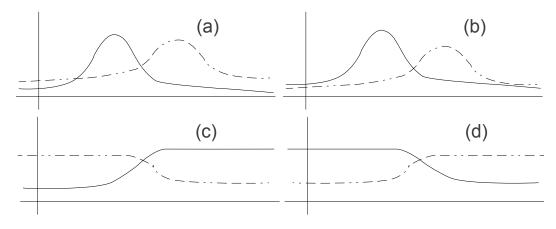
- 1. Broadly speaking there are two ways in which complexity can be controlled by constraining the data, and by constraining the model. For the second option, which amongst these is NOT a valid approach:
 - a. Reducing the number of parameters in the model
 - b. Limiting the size of the model
 - c. Limiting the magnitude of model parameter values
 - d. Using second order information when configuring the model parameters
 - e. Removing elements of the model which do not contribute to performance
- 2. In terms of "constraining the data", one approach is to reduce the dimensionality of the input space. The following are benefits of doing this EXCEPT for:
 - a. Reducing the dimensionality simply reduces the amount of data to be processed, thus saving computational resources
 - b. Removing redundant channels can help to remove noise from the data
 - c. By reducing the parameter space, reducing the dimensionality can help to avoid overfitting
 - d. Reducing the dimensionality can help to improve the validity and reliability of distance measurements
 - e. All are valid reasons



The curves depict the *likelihoods* for classes A and B respectively. Also, you are given that $p(A)=3\times p(B)$

Based on this figure, please answer the following two questions

- 3. At the point depicted by the vertical dotted line, what are the values of the *posterior distribution* for classes *A* and *B*?
 - a. 0.15,0.125
 - b. 0.20,0.50
 - c. 0.55,0.45
 - d. 0.75,0.25
 - e. 0.90,0.10
- 4. Approximately, what would the curves of the posterior distributions looks like?



- 5. Consider the following three applications: (i) Fraud detection (ii) Disease diagnosis (iii) Web search for each one, which is *more* important, Precision or Recall?
 - a. (i) Precision (ii) Recall (iii) Recall
 - b. (i) Precision (ii) Recall (iii) Precision
 - c. (i) Precision (ii) Precision (iii) Recall
 - d. (i) Recall (ii) Precision (iii) Recall
 - e. (i) Recall (ii) Recall (iii) Precision
- 6. The non-disjoint and lazy discretization methods are significant improvements over equal width discretization (i.e. the "histogram method"). Why is this?
 - a. They place the point of interest at (or near) the center of the bin used to perform density estimation
 - b. They give better density estimates in regions with low data density
 - c. They have lower computational requirements
 - d. They allow contributions from all points in the data set to be taken into account
 - e. They work well for sparse data sets
- 7. Consider the following distribution of values taken from the training set with a single continuous feature:



When designing a decision node in a tree, which of the following suggested thresholds would be a reasonable choice:

- a. -2
- b. -1
- c. +1
- d. +2
- e. +3
- 8. The following procedures are valid methods of pruning a decision tree EXCEPT for:
 - a. Stopping training when the number of training instances in a particular node drops below a threshold
 - b. Stopping training when the depth of a tree exceeds a particular threshold

- c. Deleting nodes where the information gain fails to exceed a particular threshold.
- d. Deleting nodes that do not improve (or only result in small improvements) to the test error
- e. None of the above

The following three questions address the following (feature:label) pairs:

[(0:'A'),(0:'A'),(0:'A'),(0:'B'),(1:'B'),(1:'B'),(1:'B'),(1:'B')]

- 9. calculate the information gain (IG), using base-2 logarithms:
 - a. 0.14
 - b. 0.23
 - c. 0.31
 - d. 0.45
 - e. 0.55
- 10. For the same set of pairings, please calculate the Gini impurity index:
 - a. 1/16
 - b. 2/16
 - c. 3/16
 - d. 4/16
 - e. 5/16
- 11.Let's say that this is one feature (let's call it "F1") in a Naive Bayes Classifier. What is P(F1|"B")?
 - a. 1/5
 - b. 2/5
 - c. 3/5
 - d. 4/5
 - e. 1

The following description applies to the next two questions

The following two are expressions designed for evaluating cluster quality. For simplicity, they are written for the two cluster case:

$$\frac{\sum\limits_{i \in c_{1}} \sum\limits_{j \in c_{2}} d_{ij}}{\sum\limits_{k,l \in c_{1}} d_{kl} + \sum\limits_{m,n \in c_{2}} d_{mn}} \quad , \quad \frac{\min\limits_{i \in c_{1},j \in c_{2}} d_{ij}}{\max\{\max\limits_{k,l \in c_{1}} d_{kl}, \max\limits_{m,n \in c_{2}} d_{mn}\}}$$

Given the following:

 $d_{ij} \rightarrow Distance$ between points i and j

 $c_h \rightarrow set$ of points in cluster h (for simplicity, $h \in \{1,2\}$)

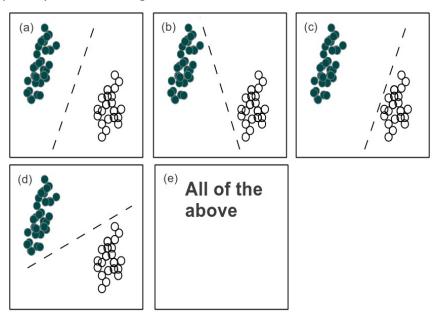
- 12. For *better* clustering performance, we would want the two measures to be:
 - a. Lower, Lower
 - b. Lower, Higher
 - c. Higher, Lower
 - d. Higher, Higher
 - e. Unable to tell
- 13. For the second expression, what is the purpose of applying the *max* operations (in the denominator)?
 - a. So as to minimize the *worst case distances* between points in the same clusters.
 - b. So as to maximize the *worst case distances* between points in the same clusters.
 - c. So as to minimize the *best case distances* between points in the same clusters.
 - d. So as to maximize the *best case distances* between points in the same clusters.
 - e. None of the above
- 14. The "C-index" for evaluating clusterings is defined as:

$$C = \frac{S - S_{min}}{S_{max} - S_{min}}$$

where a smaller value of C is "better". For the following two clusters of points, what is the value of the C-index:

- a. 0
- b. 0.2
- c. 0.4
- d. 0.6
- e. 0.8
- f. 1
- 15. The following figures depict decision boundaries induced by a single perceptron. The task is a two class classification problem, where the training data points have been clearly labeled using filled and empty circles.

Which of the figures denote a "valid" boundary – i.e. one corresponding to a converged perceptron, assuming that the perceptron learning rule was used.



- 16.If you wanted to use gradient descent to train the standard perceptron, what modification is required?
 - a. Use a "smooth" activation function
 - b. Addition of a hidden layer
 - c. Addition of hidden units
 - d. Removal of bias term
 - e. No modification is required

The passage below applies to the following three questions

The "error" term for a particular optimization problem takes the following expression:

$$E(w_1, w_2) = w_1^2 + 2w_2^2 - w_1$$

where, w_1 and w_2 are the parameters to be optimized.

- 17. What is the minimum value of the error function E, and at what values of w_1 and w_2 does this occur?
 - a. -0.22,1/3,0
 - b. -0.22,2/3,0
 - c. -0.22,0,1/3
 - d. -0.25,0,0.5

- e. -0.25,0.5,0
- 18. Iterative techniques are frequently used to solve optimization problems. You decide to use gradient descent with the initial parameter values of

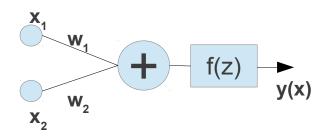
$$(w_1, w_2) = (1,1)$$

and a learning rate of 0.1. What is the initial update term?

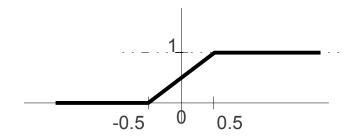
- a. (0,-0.2)
- b. (-0.1,-0.2)
- c. (-0.1, -0.4)
- d. (-0.2, -0.4)
- e. (-0.3,-0.6)
- 19. Does this point to the minimum of the error function? Why is this?
 - a. Yes, the vector gradient term indicates the direction of steepest descent, which will lead to the minimum
 - b. Yes, the error surface is convex and has no local minima, and so gradient descent leads right to the minimum of the error function
 - No, the path to the minimum is obstructed by numerous local minima, hence gradient descent does not provide an optimum path
 - d. No, the error surface has been "stretched", resulting in w_2 being over-emphasized by the gradient function.
 - e. None of the above

The description below applies to the following three questions

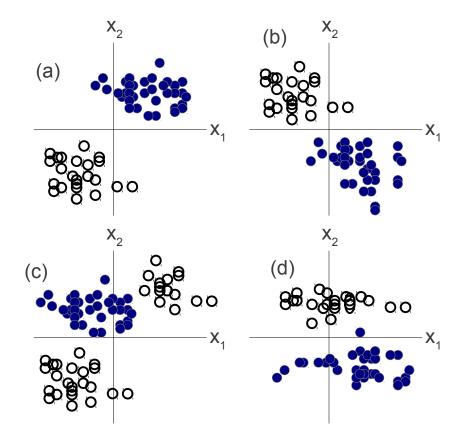
Consider the following network architecture:



Where, the activation function f(z) is depicted by:



- 20.If $x_1=1$, $x_2=2$, $w_1=w_2=0.5$, what is the network's output?
 - a. 0.5
 - b. 1
 - c. 1.5
 - d. 2
 - e. 2.5
- 21. Assuming that the target output is 0, what is the gradient of the error function w.r.t. w_2 ?
 - a. 0
 - b. 0.5
 - c. 1
 - d. 1.5
 - e. 2
- 22. Which of the following classification problems can be appropriately handled by this particularly network?



23. Consider the following three descriptions:

- (i) "Finding clusters by combining individual elements into larger collections"
- (ii) "A form of learning where only parameters related to large inputs and erroneous outputs are changed"
- (iii) "A process by which input data points are matched with network weights"

Which three terms match the descriptions above:

- a. Divisive clustering; Hebbian Learning; Projection
- b. Agglomerative clustering; Gradient descent optimization; Distance Calculation
- c. Agglomerative clustering; Hebbian Learning; Projection
- d. Divisive clustering; Gradient descent optimization; Projection
- e. Divisive clustering; Hebbian Learning; Distance Calculation

This description applies to the following two questions:

The following is the distance matrix between a set of five points:

	1	2	3	4	5
1	0	1	3.5	5	6
2	1	0	2.2	4	5
3	3.5	2.2	0	2	3.3
4	5	4	2	0	1
5	6	5	3.3	1	0

- 24. You decide to use the k-centers (also known as k-medoids) algorithm to cluster these points. Set points 1 and 2 to be the initial centroids. What will the initial clustering be?
 - a. (1),(2,3,4,5)
 - b. (1,3),(2,4,5)
 - c. (1,3,4),(2,5)
 - d. (1,3,4,5),(2)
 - e. (1,3,5),(2,4)
- 25.At some point during a run of the k-centers algorithm, we are left with the following two clusters:

$$C_1:(1,2,3,4) C_2:(5)$$

For cluster **C**₁, what should be the centroid be?

- a. 1
- b. 2
- c. 3
- d. 4
- e. 5
- 26. Which of the following is a valid reason to choose the k-centers algorithm over k-means:
 - a. Works better with high dimensional data
 - b. Lower computational complexity
 - c. Lower memory requirements
 - d. Better able to handle missing data points
 - e. Better resistance to outliers
- 27. The self organizing map is similar to the k-means clustering algorithm in some ways, but is different in others. Name one similarity and one difference:

- a. **S**: Nodes (or centroids) strive to be representative of a group of data points; **D**: The nodes of a SOM are topologically linked
- b. **S**: Nodes (or centroids) strive to be representative of a group of data points; **D**: SOM is trained using an iterative rule
- c. **S**: Both methods provide nonlinear visualization capabilities; **D**: The nodes of a SOM are topologically linked
- d. **S**: Both methods provide nonlinear visualization capabilities; **D**: SOM is trained using an iterative rule
- e. **S**: Both methods provide nonlinear visualization capabilities; **D**: SOM can work in multiple dimensions
- 28. When training the SOM, the learning rate and neighborhood size need to be gradually:
 - a. Increased; Increased
 - b. Increased; Decreased
 - c. Decreased; Increased
 - d. Decreased; Decreased
 - e. None of the above
- 29. With reference to the previous question, why do these parameters need to be changed?
 - a. Early in the training, the map may be overfitting the data, while later on this is no longer a problem
 - b. Early in the training, the map may be underfitting the data, while later on this is no longer a problem
 - c. Early in the training the parameters have very low variability and as such need a higher degree of tuning
 - d. Early in the training, the map needs to be "fine-tuned", while later on it simply needs to be "folded".
 - e. Early in the training, the map needs to be "unfolded", while later on, it needs to be "fine-tuned"
- 30. Consider the following three statements:
 - (i) "Principle Component Analysis works by finding the subspace which contains the largest proportion of the variance in the data"
 - (ii) "Provided that there are no local minima, the direction of steepest descent gives the most direct path to the global minimum"
 - (iii) "The Levenberg-Marquardt algorithm gives better performance than backpropagation by using second order information about the error surface"

Indicate whether each statement is "true" or "False":

a. True; True; True

b. False; True; True

c. True; False; True

d. True; True; False

e. False; False; False