Please show all your work! Answers without supporting work will not be given credit. Write your answers in spaces provided.

There are total 13 pages including two blank pages at the end for scratch work.

You have 1 hour and 15 minutes to complete this exam.

You may use any result from the lectures or the homeworks without proof.

Name & UNI:
 (15 points) State True or False. No justification needed! (+1 for correct, 0 for blank, -1 for incorrect answer)
(a) The second of the project X_1, \ldots, X_n , the joint distribution $P(X_1, \ldots, X_n) = Assignment Project Exam Help$
(b) If two random variables are conditionally independent, then they are not necessarily independent; but if the two random variables are independent then they are necessarily conditionally herendent as yellowcoder.com
(c) In a Hidden Markov Model, the observed variable X_{t-1} (at time $t-1$) is conditionally independent of the viscosity of the X_t p at time X_t .
(d) $\frac{1}{B_1 \subseteq B_2 \subseteq B_\infty}$ Consider the unit L_p -ball in \mathbb{R}^d , that is the set $\{x \in \mathbb{R}^d : x _p \le 1\} =: B_p$. Then,
(e) A non-linear kernel transform in sufficiently high dimensions can always achieve zero test error.
(f) The dual variables (α_i) in Support Vector Machines (SVMs), take non-zero values only if the corresponding datapoints x_i are in fact "support" vectors that dictate the margin.
(g) VC dimension of decision trees in \mathbb{R}^2 is infinite.
(h) The decision boundary induced by a Logistic regression classifier on a two-class problem is always linear.

(i)	The Lloyd's algorithm for k -means clustering for $k=2$ can give solutions that are arbitrarily bad in terms of the clustering cost compared to the optimal $(k=2)$ k -means solution
(j)	The maximum likelihood setting of the parameters for a mixture model often yields undesirable results.
(k)	A directed graphical model on n variables that has a structure of a fully connected directed acyclic graph (DAG) admits no independencies (conditionally amongst the n variables.
(1)	The notation " $A \perp B \mid C$ " means " A and B are independent given C ". Then $X \perp Y \mid W, Z \text{ and } X \perp W \mid Y, Z \implies X \perp W, Y \mid Z.$
(m)	$\underline{}$ L_2 (Euclidean) distances can always be computed efficiently in Kernel space via the kernel trick.

- $\text{\tiny (n)} \ \, \underline{Assignment}_{\mathrm{is}} \underbrace{Paroject}_{\mathrm{e}} \underbrace{Exam_{\mathrm{C}}}_{\mathrm{Hel}} \underline{p}_{|\mathcal{F}|)}.$
- (o) ______When compared to batch tearning, exting learning causignificantly reduce the number of labelled samples needed to learn a concept to a desired lever of accuracy.

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2. [Maximum Likelihood Estimation of fully observed HMMs (15 points)] Let the distribution of $(X_t, Y_t)_{t \in \{1, 2, ...\}}$ be from a discrete space HMM, where each Y_i (the hidden state at time t) takes in values from $[K] := \{1, 2, ..., K\}$ and each X_t (the observation at time t) takes values in $[D] := \{1, 2, ..., D\}$. Recall that the HMM parameters $\theta = (\pi, A, B)$ have the following semantics: $P[Y_1 = i] = \pi_i, P[Y_{t+1} = j|Y_t = i] = A_{i,j}$ and $P[X_t = j|Y_t = i] = B_{i,j}$.

Suppose you have as training data a labeled sequence $((x_1, y_1), (x_2, y_2), \dots, (x_T, y_T)) \in ([D] \times [K])^T$. What is the maximum likelihood estimation of the HMM parameters θ given this data? (Hint: if needed, use the convention 0/0 = 0 in deriving your estimates.)

- 3. [Optimization of sparse regression] Let $\mathbf{X} \in \mathbb{R}^{n \times d}$ be a data matrix with the i^{th} row of \mathbf{X} is the i^{th} data observation $x_i \in \mathbb{R}^d$. Let $y \in \mathbb{R}^n$ be the real-valued labels. Let $\mathcal{W} := \{w \in \mathbb{R}^d : w \text{ has at most one non-zero entry}\}$. Suppose you want to find the weight vector $\hat{w} \in \mathcal{W}$ that minimizes the objective function $f(w) := \|y \mathbf{X}w\|^2$ over all $w \in \mathcal{W}$.
 - (a) **(5 points)** Formulate this as an optimization problem. (Make sure to clearly indicate the dimension and the variables).

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(b) (3 points) Is this a convex optimization problem? (why or why not)?

(c) (10 points) Describe an algorithm for computing such a vector \hat{w} . Make sure that your pseudocode is clear and precise, and specify what is exactly returned by your algorithm. (Hint: all but one entry of \hat{w} are zero).

⁽d) (2 points) What is the time complexity of your algorithm (give it in terms of the parameters n and d)?

4. [Principal Components Analysis (PCA) and beyond]

Recall from lecture that given a data matrix $X = \begin{bmatrix} & & & | \\ x_1 & \cdots & x_n \\ | & & | \end{bmatrix}$, where each observation $x_i \in \mathbb{R}^d$,

the best k-dimensional linear mapping that minimizes the squared reconstruction error is given by the eigenvectors corresponding to the top k eigenvalues of the outer product matrix XX^{T} . This is also called the k-dimensional PCA subspace.

(a) Suppose you are informed that the function to compute the eigenvectors and eigenvalues in your favorite language is buggy (so you cannot use this function). As an alternative you explore the language documentation and find a function that has the the ability to decompose any $d \times n$ matrix X into a summation over three sets variables: σ_i (a scalar), u_i (a vector in \mathbb{R}^d), and v_i (a vector in \mathbb{R}^n). That is, the data matrix X can be written as:

$$X = \sum_{i=1}^{d} \sigma_i u_i v_i^\mathsf{T},$$

These variables have special properties:

- The scalars σ_i 's are all non-negative, such that: $0 \le \sigma_1 \le \ldots \le \sigma_d$.
- The vectors u_1, \ldots, u_d are orthonormal. That is, $||u_i|| = 1$ and $u_i \cdot u_j = 0$ (for $i \neq j$).

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i. (5 points) Compute XX^{T} in terms of σ_i , u_i and v_i . Simplify your answer as much as possible.

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ii. (5 points) Using the definition of eigenvector of a matrix A: " $Ax = \lambda x$ ", show that each u_i is an eigenvector of the outerproduct XX^{T} . What are the corresponding eigenvalues?

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iii. (2 points) Which k eigenvectors (from u_1, \ldots, u_d) of XX^{T} should be used to form a k-dimensional PCA subspace?

(b) (3 points) PCA yields a subspace that gives the best reconstruction error, but this subspace can be arbitrarily bad for some predictive tasks such as classification. Suppose we want to do linear classification on binary labelled data in \mathbb{R}^2 . To reduce representational complexity, we wish to project this data onto a 1D subspace via PCA and perform linear classification in 1D.

Depict an example binary labelled dataset in \mathbb{R}^2 such that even though there exists a 1D projection of the dataset that can perfectly linearly separate the two classes, the best linear separator on the 1D PCA projection will give poor classification accuracy. (You must justify your answer why PCA projection would not pick a good label separating subspace by outlining the kinds of properties your depicted dataset must have.)

5. [Bayes Nets]

- (a) (5 points) Draw a Bayes net over the random variables $\{A, B, C, D\}$ where the following conditional independence assumptions hold. Here, $X \perp Y \mid Z$ means X is conditionally independent of Y given Z, and $X \not\perp Y | Z$ means X and Y are not conditionally independent given Z, and \emptyset stands for the empty set.
 - A⊥B|∅
 - $A \not\perp D|B$
 - A⊥D|C
 - $A \not\perp C | \emptyset$
 - B ∠ C | ∅
 - A ⊥ B | D
 - $B \perp D \mid A, C$

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(b) (2 points) Write a simple expression for the joint distribution over the variables $\{A, B, C, D\}$ that captures all the conditional (in)dependencies asserted by the Bayes net you depicted in part (a).

Pr[A, B, C, D] https://powcoder.com

(c) (3 points) List all the conditional independence asserted by the Bayes net you depicted in part (a) but with all the arrows reversed.

- 6. [A better output Perceptron algorithm guarantee]
 - In class, we saw that when the training sample S is linearly separable with a maximum margin $\gamma > 0$, then the Perceptron algorithm run cyclically over S is guaranteed to converge after $T \leq (R/\gamma)^2$ updates, where R is the radius of the sphere containing the sample points. This does not guarantee however that the hyperplane solution returned by Perceptron, i.e. w_T achieves a margin close to γ .
 - (a) (5 points) Show an example training dataset S in \mathbb{R}^2 that has margin γ , and an order of updates made by the Perceptron algorithm where the hyperplane solution returned has arbitrarily bad margin on S.

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(b) Consider the foliation and that the that cap and the consider the foliation of the consideration of the consideratio

Modified Perceptron Algorithm

Input: training dataset $S = (x_i, y_i)_{i=1,...,n}$ Output: learned vector w- Initialize $w_0 := 0, t := 0$

- while there exists an example $(x,y) \in S$, such that y (w_t · x) ≤ 0 y (w_t · x) ≤ 0 y (w_t · x) $\leq \gamma ||w_t||$

- set $w_{t+1} := w_t + yx$
- set t := t + 1
- return w_t .
 - i. (2 points) If the Modified Perceptron Algorithm (MPA) terminates after T rounds, what margin guarantee is achieved by the hyperplane w_T returned by MPA? Justify your answer.

ii. (3 points) It turns out that for a training sample S of margin γ one can also show that at any iteration t where the Modified Perceptron Algorithm (MPA) makes a mistake, the following is true.

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A.
$$w_t \cdot w^* \ge (w_{t-1} \cdot w^*) + \gamma$$

B. $||w_t|| \le \frac{4R^2}{\gamma} + \frac{3}{4}t\gamma$

From properties A and B, what mistake bound can you derive for MPA? That is, bound the maximum iterations T, or equivalently, bound the number of mistakes made by MPA.

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