Final Review

Anurag Nagai

Logistic Regression

Model Evaluation

Instance Based Learning

Ensemble Methods

Unsupervised

Bayes Net

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Final Review

** This is the review of some post-midterm topics.

This review is NOT exhaustive.

You are responsible for covering the entire course.

The final will be comprehensive **

Anurag Nagar

Machine Learning Class

Topics Covered

Final Review

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List of topics covered post midterm:

- Logistic Regression
- Model Evaluation
- Instance Based Learning (k-Nearest Neighbor)
- Ensemble Learning
- Unsupervised Learning, PCA
- Bayes Net
- Hidden Markov Models (HMM)

Outline

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Logistic Regression

- 1 Logistic Regression

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Logistic Regression

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Creates a probabilistic model for classes i.e. tries to predict $P(Y = y_k)$ where y_k could be 0 or 1 for a binary dataset.

$$p = P(Y = 1) = \frac{1}{1 + e^{-z}}$$
 where $z = w_0 + \sum_{i=1}^{n} w_i x_i$ and

$$1 - p = P(Y = 0) = \frac{e^{-z}}{1 + e^{-z}}$$

$$ln(\frac{p}{1-p}) = z = w_0 + \sum_{i=1}^{n} w_i x_i$$

so log of odds ratio is a linear function of x's

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As always,

$$W = \begin{bmatrix} w_0 \\ w_1 \\ \dots \\ w_n \end{bmatrix} \quad \text{and} \quad X = \begin{bmatrix} x_0 \\ x_1 \\ \dots \\ x_n \end{bmatrix}$$

where $x_0 = 1$

In some questions, you are given W and X and are asked if the data is more likely to belong to class 1 or class 0. You may also be asked to find out the dividing boundary. You need to check if

$$P(Y=1) > P(Y=1)$$

for given values.

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Note: z is a dot product of vectors W and X

$$z = W \cdot X$$

Assign to class 1 if:

$$\frac{p}{1-p} > 1$$
 or $z > 0$ or

$$w_0 + \sum_{i=1}^n w_i x_i > 0$$

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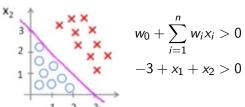
Unsupervised Learning

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$$W = \begin{bmatrix} -3\\1\\1 \end{bmatrix} \quad \text{and} \quad X = \begin{bmatrix} 1\\x_1\\x_2 \end{bmatrix}$$

Class is 1 if



Red is class 1

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Difference between generative and discriminative classifiers?

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Difference between generative and discriminative classifiers?

Generative classifiers estimate P(Y|X) from the training data using Bayes rule. In other words, they estimate joint probability of y and x from the data.

Discriminative classifiers make a model of P(Y|X)

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If there are n Booleans $X_1, X_2, \ldots X_n$, and they are all dependent on each other, how many parameters would be needed to describe their joint probability $P(X_1, X_2, \ldots X_n)$

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If there are n Booleans $X_1, X_2, \ldots X_n$, and they are all dependent on each other, how many parameters would be needed to describe their joint probability $P(X_1, X_2, \ldots X_n)$

$$2^{n} - 1$$

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If there are 2 Boolean classes Y_1, Y_2 Booleans, and each class has n Boolean features. We assume inter-class feature independence i.e. feature of a class depend on each other, but are independent of features in the other class. How many parameters are needed to describe the system.

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If there are 2 Boolean classes Y_1 , Y_2 Booleans, and each class has n Boolean features. We assume inter-class feature independence i.e. feature of a class depend on each other, but are independent of features in the other class. How many parameters are needed to describe the system.

$$2 \times (2^{n} - 1)$$

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Bayes Ne

If there are 2 Boolean classes Y_1 , Y_2 each consisting of n Boolean features, and each class is modeled using the traditional logistic regression function

$$P(Y = Y_i | (X_1, X_2, ..., X_n)) = \frac{1}{1 + exp(w_0 + \sum_{j=1}^n w_j X_j)}$$

How many total parameters are needed to describe the system?

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Bayes Ne

If there are 2 Boolean classes Y_1 , Y_2 each consisting of n Boolean features, and each class is modeled using the traditional logistic regression function

$$P(Y = Y_i | (X_1, X_2, ..., X_n)) = \frac{1}{1 + exp(w_0 + \sum_{j=1}^n w_j X_j)}$$

How many total parameters are needed to describe the system?

$$(n + 1)$$

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Suppose you are playing a game that involves hitting a target. You are given 3 chances, but the games gets tougher every time you miss. The probabilities of hitting in the first, second, and third attempts is 0.6, 0.5, 0.4 Find the odds ratio for this game.

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Bayes Net

Suppose you are playing a game that involves hitting a target. You are given 3 chances, but the games gets tougher every time you miss. The probabilities of hitting in the first, second, and third attempts is 0.6, 0.5, 0.4 Find the odds ratio for this game.

Probability of failure:

$$1 - p = (1 - 0.6) \times (1 - 0.5) \times (1 - 0.4) = 0.12$$

Probability of success: p = 0.88

Odds Rato =
$$\frac{0.88}{0.12}$$
 = 7.33

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Logistic Regression is what type of model? Linear or Non-Linear

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Logistic Regression is what type of model? Linear or Non-Linear

Linear (Proved in class)

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Logistic Regression

Suppose the logistic regression for binary class Y = 0 and Y = 1 and two feature X_1, X_2 problem is defined as below:

$$P(Y = 1 | (X_1, X_2)) = \frac{1}{1 + exp(w_0 + \sum_{j=1}^{2} w_j X_j)}$$

The values of w_0 , w_1 , w_2 are -1, 2, 0 respectively. What will be the separation rule for the two classes.

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Suppose the logistic regression for binary class Y=0 and Y=1 and two feature X_1,X_2 problem is defined as below:

$$P(Y = 1 | (X_1, X_2)) = \frac{1}{1 + exp(w_0 + \sum_{j=1}^{2} w_j X_j)}$$

The values of w_0 , w_1 , w_2 are -1, 2, 0 respectively. What will be the separation rule for the two classes.

For data to be in class 1: P(Y = 1) > P(Y = 0) or $w_0 + w_1x_1 + w_2x_2 > 0$ or $1 + 2 \times x_1 + 0 \times x_2 > 0$ or $-1 + 2x_1 > 0$ or $x_1 > \frac{1}{2}$

Final Review

Logistic Regression

For estimating best parameters of logistic regression, which technique is used:

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technique is used:

MLE + Gradient Ascent

For estimating best parameters of logistic regression, which

Logistic Regression

Logistic Regression Question

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Logistic Regression

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Consider the entropy function for a binary class problem:

$$H = p \log_2(p) + (1-p) \log_2(1-p)$$

where p is probability of class 1. A plot of H vs p would be what type of curve?

- 1 Concave
- 2 Linear
- 3 Convex
- 4 Polynomial of degree 4

Logistic Regression Question

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Consider the entropy function for a binary class problem:

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- 1 Concave
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- 5 Unsupervised Learning

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Consider the confusion matrix shown below:

| | Predicted = YES | Predicted = NO |
|--------------|-----------------|----------------|
| Actual = YES | 100 | 40 |
| Actual = NO | 10 | 50 |

Calculate:

- Accuracy
- Recall
- Precision
- F-Statistic

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Consider the confusion matrix shown below:

| | Predicted = YES | Predicted = NO |
|--------------|-----------------|----------------|
| Actual = YES | 100 | 40 |
| Actual = NO | 10 | 50 |

Calculate:

• Accuracy =
$$\frac{150}{200}$$

Precision (p) =
$$\frac{100}{110}$$

• Recall (r) =
$$\frac{100}{140}$$

■ F-Statistic =
$$\frac{2rp}{r+p}$$
 = 0.80

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Consider the following confusion matrix.

Number of cases: 100,000

| Actual Predicted State Negative | | | | dicted sitive |
|---------------------------------|----|-------|----|------------------|
| Negative | TN | 97750 | FP | 150 |
| Positive | FN | 330 | TP | 1770 |

Compute precision, recall, F-statistic

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Consider the following confusion matrix.

Number of cases: 100,000

| Actual | Predicted | | Predicted | |
|----------|-----------|-------|-----------|------|
| State | Negative | | Positive | |
| Negative | TN | 97750 | FP | 150 |
| Positive | FN | 330 | TP | 1770 |

Compute precision, recall, F-statistic

Should be easy

ROC Curve

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Remember the definitions:

- True Positive Rate (TPR) = $\frac{TP}{TP + FN}$
- False Positive Rate (FPR) = $\frac{FP}{FP + TN}$
- Receiver Operative Characteristic (ROC) curve is a plot of TPR vs FPR

Model Evaluation Question

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I am testing a model that predicts the probability of a person having disease A. I have obtained the following results:

| Person | Predicted probability | True Value |
|--------|-----------------------|------------|
| 1 | 0.95 | 1 |
| 2 | 0.80 | 1 |
| 3 | 0.60 | 1 |
| 4 | 0.45 | 1 |
| 5 | 0.15 | 0 |

Calculate area under ROC curve by setting the positive threshold i.e. probability equal to or greater than which prediction is labeled as positive :

- **0.0**
- **0.25**
- **0.50**
- **0.75**
- **1.0**

What according to you should be the best threshold for predicted probability?

Model Evaluation Question

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| Threshold | 0.0 | 0.25 | 0.5 | 0.75 | 1.0 |
|-----------|-----|------|------|------|-----|
| TP | 4 | 4 | 3 | 2 | 0 |
| FP | 1 | 0 | 0 | 0 | 0 |
| TN | 0 | 1 | 1 | 1 | 1 |
| FN | 0 | 0 | 1 | 2 | 4 |
| TPR | 1.0 | 1.0 | 0.75 | 0.50 | 0.0 |
| FPR | 1.0 | 0.0 | 0.0 | 0.0 | 0.0 |

Area under ROC = 100.0%

Best threshold = 0.25

Model Evaluation Question

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I am testing a model that predicts the probability of a person having disease A. I have obtained the following results:

| Person | Predicted probability | True Value |
|--------|-----------------------|------------|
| 1 | 0.80 | 1 |
| 2 | 0.60 | 0 |
| 3 | 0.40 | 1 |
| 4 | 0.20 | 0 |

Calculate area under ROC curve by setting the positive threshold i.e. probability equal to or greater than which prediction is labeled as positive :

- **0.0**
- **0.2**
- **0.4**
- **0.6**
- **8.0**
- **1.0**

1.0

What according to you should be the best threshold for predicted probability?



Model Evaluation Question

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Calculate area under ROC curve by setting the positive threshold i.e. probability equal to or greater than which prediction is labeled as positive :

- 0.0
- **0.2**
- **0.4**
- **0.6**
- 0.8
- **1.0**

What according to you should be the best threshold for predicted probability? Try yourself

Model Evaluation

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Logistic Regression

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Instance

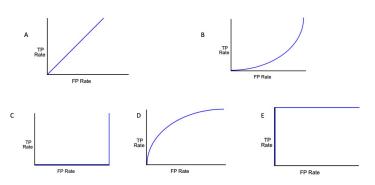
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Consider the following ROC curves.



Arrange the curves from best to worst

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Evaluation

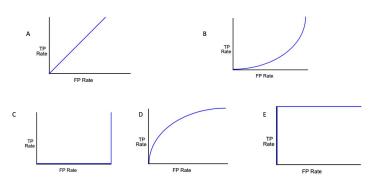
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Consider the following ROC curves.



Arrange the curves from best to worst

E, D, A, B, C



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- In k-NN, the parameter k is a hyperparameter not a model parameter. Understand the differences.
- Understand training and test phases.
- Practice questions on predictions using k-NN
- Bias and Variance versus k in k-NN

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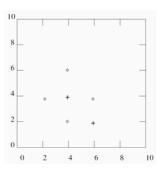
Methods

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Consider the dataset shown below:



What will be the error of 1-NN and 3-NN applied to this dataset?

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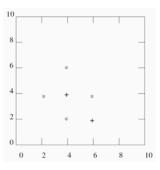
Ensemble Methods

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Consider the dataset shown below:



What will be the error of 1-NN and 3-NN applied to this dataset?

$$1-NN = 6/6$$
, $3-NN = 4/6$

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Consider the dataset shown below:

| | Solid | Temp | Force Applied | Class |
|----|-------|------|------------------|-------|
| 1 | Yes | 0 | 0.406 | No |
| 2 | No | 1 | 0.25 | No |
| 3 | No | 0 | 0.063 | No |
| 4 | Yes | 1 | 0.375 | No |
| 5 | No | 0.5 | 0.219 | Yes |
| 6 | No | 1 | 0 | No |
| 7 | Yes | 0.5 | 1 | No |
| 8 | No | 0 | 0.156 | Yes |
| 9 | No | 1 | 0.094 | No |
| 10 | No | 0 | 0.188 | Yes |

Using 1-NN, what will be the prediction for item (No, 0, 0.375)? You can assume Manhattan distance for each of the attributes. For the first attribute, you can assume Yes = 1 and No = 0

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Consider the dataset shown below:

| | Solid | Temp | Force Applied | Class |
|----|-------|------|------------------|-------|
| 1 | Yes | 0 | 0.406 | No |
| 2 | No | 1 | 0.25 | No |
| 3 | No | 0 | 0.063 | No |
| 4 | Yes | 1 | 0.375 | No |
| 5 | No | 0.5 | 0.219 | Yes |
| 6 | No | 1 | 0 | No |
| 7 | Yes | 0.5 | 1 | No |
| 8 | No | 0 | 0.156 | Yes |
| 9 | No | 1 | 0.094 | No |
| 10 | No | 0 | 0.188 | Yes |

Using 1-NN, what will be the prediction for item (No, 0, 0.375)? You can assume Manhattan distance for each of the attributes. For the first attribute, you can assume Yes = 1 and No = 0

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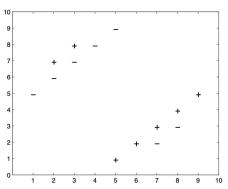
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Consider the visual dataset shown below:



What value of k in k-NN will give the best and the worst training error? Can you obtain 0 training error for any value of k?

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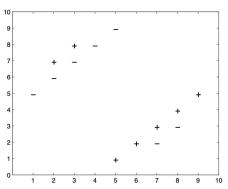
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Consider the visual dataset shown below:



What value of k in k-NN will give the best and the worst training error? Can you obtain 0 training error for any value of k?

Try various values of k, and investigate

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-MN

Suppose I create a model for regression, such as $y = w_0 + \sum_{i=1}^n w_i x^i$. I can fit the *training data* nicely by making a complex polynomial i.e. increasing the value of n. What effect will it cause to bias and variance

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Suppose I create a model for regression, such as $y = w_0 + \sum_{i=1}^n w_i x^i$. I can fit the *training data* nicely by making a complex polynomial i.e. increasing the value of n. What effect will it cause to bias and variance

Decrease bias and increase variance

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4MN

Suppose I create a model for regression, such as $y = w_0 + \sum_{i=1}^{n} w_i x^i$. I can fit the *training data* nicely by making a complex polynomial i.e. increasing the value of n. How can I prevent the model from becoming too complex?

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Suppose I create a model for regression, such as $y = w_0 + \sum_{i=1}^n w_i x^i$. I can fit the *training data* nicely by making a complex polynomial i.e. increasing the value of n. How can I prevent the model from becoming too complex?

Add a penalty term called regularization: Total Error = $\sum_{i=1}^{M} (t_i - y_i)^2 + \lambda \sum_{i=1}^{n} w_i^2$ M is the number of data points and n is the degree of polynomial used.

https://towards datascience.com/regularization-an-important-concept-in-machine-learning-5891628907 earlier (as a concept-in-machine-learning-5891628907) and the concept-in-machine-learning-5891628907 earlier (as a concept-in-machine-learning-5891628907). The concept-in-machine-learning-5891628907 earlier (as a concept-in-machine-learning-5891628907) and the concept-in-machine-learning-5891628907 earlier (as a concept-in-machine-learning-5891628907). The concept-in-machine-learning-5891628907 earlier (as a concept-in-machine-learning-5891628907) and the concept-in-machine-learning-5891628907 earlier (as a concept-in-machine-learning-section-learni

https://datanice.github.io/machine-learning-101-what-is-regularization-interactive.html

https://towardsdatascience.com/regularization-in-machine-learning-76441ddcf99a

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How do you control bias and variance in the following models:

- Linear Regression
- Neural Networks
- Decision Trees
- Logistic Regression
- k-Nearest Neighbors

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How do you control bias and variance in the following models:

- Linear Regression
- Neural Networks
- Decision Trees
- Logistic Regression
- k-Nearest Neighbors

Do your research and find answers.

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How does an ensemble (set of classifiers working together) prevent overfitting and still give a good accuracy?

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How does an ensemble (set of classifiers working together) prevent overfitting and still give a good accuracy?

Averaging of models gives a more stable model overall i.e. a model with low variance.

It's like taking the opinion of multiple people before taking a decision.

https://towards datascience.com/simple-guide-for-ensemble-learning-methods-d87cc68705a2

https://machinelearning mastery.com/ensemble-methods-for-deep-learning-neural-networks/

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Practice examples of Adaboost. Like the ones solved in class, and

http://ais.informatik.uni-freiburg.de/teaching/ws11/robotics2/pdfs/rob2-18-adaboost.pdf
http://www.authorstream.com/Presentation/aSGuest79199-727683-animated-adaboost-example/

Conceptual knowledge of Adaboost, such as:

Can Adaboost using stumps (i.e. decision tree of depth 1) always give 0 *training* error ?

Will Adaboost always converge?

Random Forest

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In Random forest you can generate large number of trees (say T_1, T_2, \ldots, T_n) and then aggregate the results of these trees. Which of the following is true about individual(T_k) tree in Random Forest?

- 1 Individual tree is built on a subset of the features
- 2 Individual tree is built on all the features
- 3 Individual tree is built on a subset of observations
- 4 Individual tree is built on full set of observations

Random Forest

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We would like to perform K-means clustering on the four point shown below:

$$x1 = (1, 1)$$

$$x2 = (2, 2)$$

$$x3 = (6, 6)$$

$$x4 = (7, 7)$$

The initial centers are x1 and x2 for clusters C1 and C2 respectively. We will use Manhattan distance.

For the first iteration, which points will be assigned to C1 and C2?

New centers after first iteration:

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We would like to perform K-means clustering on the four point shown below:

$$x1 = (1, 1)$$

$$x2 = (2, 2)$$

$$x3 = (6, 6)$$

$$\times 4 = (7, 7)$$

The initial centers are x1 and x2 for clusters C1 and C2 respectively. We will use Manhattan distance.

For the first iteration, which points will be assigned to C1 and C2?

New centers after first iteration:

$$C1 = (1, 1), C2 = (5, 5)$$

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We would like to perform K-means clustering on the four point shown below:

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$$x4 = (7, 7)$$

The initial centers are x1 and x2 for clusters C1 and C2 respectively. We will use Manhattan distance.

For the second iteration, which points will be assigned to C1 and C2?

New centers after second iteration:

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We would like to perform K-means clustering on the four point shown below:

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$$x3 = (6, 6)$$

$$\times 4 = (7, 7)$$

The initial centers are x1 and x2 for clusters C1 and C2 respectively. We will use Manhattan distance.

For the second iteration, which points will be assigned to C1 and C2?

 $\times 1$, $\times 2$ to C1, and $\times 3$, $\times 4$ to C2

New centers after second iteration:

$$C1 = (1.5, 1.5), C2 = (6.5, 6.5)$$

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We would like to perform K-means clustering on the four point shown below:

$$x1 = (1, 1)$$

$$x2 = (2, 2)$$

$$\times 3 = (6, 6)$$

$$\times 4 = (7, 7)$$

The initial centers are x1 and x2 for clusters C1 and C2 respectively. We will use Manhattan distance.

For the third iteration, which points will be assigned to C1 and C2?

New centers after third iteration:

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We would like to perform K-means clustering on the four point shown below:

$$x1 = (1, 1)$$

$$x2 = (2, 2)$$

$$x3 = (6, 6)$$

$$\times 4 = (7, 7)$$

The initial centers are x1 and x2 for clusters C1 and C2 respectively. We will use Manhattan distance.

For the third iteration, which points will be assigned to C1 and C2?

x1, x2 to C1, and x3, x4 to C2

New centers after third iteration:

C1 = (1.5, 1.5), C2 = (6.5, 6.5) No change, so can stop

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We would like to perform K-means clustering on the four point shown below:

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The initial centers are x1 and x2 for clusters C1 and C2 respectively. We will use Manhattan distance.

What will be the value of Within Cluster Sum of Squares (WCSS) after the end of clustering?

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We would like to perform K-means clustering on the four point shown below:

$$x1 = (1, 1)$$

$$x2 = (2, 2)$$

$$x3 = (6, 6)$$

$$\times 4 = (7, 7)$$

The initial centers are x1 and x2 for clusters C1 and C2 respectively. We will use Manhattan distance.

What will be the value of Within Cluster Sum of Squares (WCSS) after the end of clustering?

Cluster 1:
$$1^2 + 1^2$$
, Cluster 2: $1^2 + 1^2$

$$Total = 4.0$$

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Suppose I have the following dataset:

| Point | <i>X</i> ₁ | <i>X</i> ₂ |
|-------|-----------------------|-----------------------|
| Α | 1 | 1 |
| В | 1 | 0 |
| C | 3 | 3 |
| D | 2 | 4 5 |
| Е | 3 | 5 |

Perform k-means clustering with k=2, with A and C as initial centers. Use Manhattan distance.

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First Iteration

Starting distance:

from A:
$$A = 0$$
, $B = 1$, $C = 4$, $D = 4$, $E = 6$

from C:
$$A=4$$
, $B=5$, $C=0$, $D=2$, $E=2$

Update Centers:

$$C1 = (1, 0.5)$$

$$C2 = (2.67, 4)$$

Second Iteration

Starting distance:

from C1:
$$A = 0.5$$
, $B = 0.5$, $C = 4.5$, $D = 4.5$, $E = 6.5$

from C2:
$$A = 4.67$$
, $B = 5.67$, $C = 0.67$, $D = 0.67$, $E = 1.33$

Update Centers:

$$C1 = (1, 0.5)$$

$$C2 = (2.67, 4)$$

No change in assignment or centers, so can stop

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Suppose I have the following dataset:

| Point | <i>x</i> ₁ | <i>X</i> ₂ |
|-------|-----------------------|-----------------------|
| Α | 1 | 1 |
| В | 1 | 0 |
| C | 1 3 | 3 |
| D | 2 | 4 5 |
| E | 3 | 5 |

Perform single link hierarchical clustering and draw the final dendrogram. Use Manhattan distance.

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Suppose I have the following dataset:

| Point | <i>x</i> ₁ | <i>X</i> ₂ |
|-------|-----------------------|-----------------------|
| Α | 1 | 1 |
| В | 1 | 0 |
| C | 1 3 | 3 |
| D | 2 | 4 5 |
| E | 3 | 5 |

Perform single link hierarchical clustering and draw the final dendrogram. Use Manhattan distance.

Try yourself

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Which of the following are aims of PCA analysis?

- **1** To convert a set of correlated attributes of a dataset into features that are orthonormal to each other.
- 2 To discover a set of principal components that explain variance in data.
- 3 To rename columns.
- To remove duplicate rows.

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Which of the following are aims of PCA analysis?

- 1 To convert a set of correlated attributes of a dataset into features that are orthonormal to each other.
- 2 To discover a set of principal components that explain variance in data.
- 3 To rename columns.
- To remove duplicate rows.

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Which of the following is true about the first principal component?

- 1 It is a linear combination of other principal components and explains the largest percentage of variance in the dataset.
- 2 It explains the least variance in the dataset.
- 3 It is orthogonal to all other principal components and explains the largest percentage of variance in the dataset.
- Principal components are randomly arranged.

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Which of the following is true about the first principal component?

- 1 It is a linear combination of other principal components and explains the largest percentage of variance in the dataset.
- 2 It explains the least variance in the dataset.
- It is orthogonal to all other principal components and explains the largest percentage of variance in the dataset.
- 4 Principal components are randomly arranged.

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ИΜЬ

Understand the concept of PCA, explained variance, etc. No numerical calculations will be asked.

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AMA

Suppose I have a Bayes net of 5 nodes, where each node can take 4 possible values i.e. the nodes are not binary. How many parameters would be needed if they were all dependent, and when they were all fully independent.

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4ML

Practice more questions involving joint probability, summing out etc, dependency separation (d-separation), etc

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Practice Forward and Viterbi algorithm questions.