Single Option Correct Questions

- 1. Which of the following is NOT true about multicollinearity?
 - a. It is represented by a high correlation coefficient between pairs of variables
 - b. It causes the parameter estimates to be biased
 - c. It leads to the feature matrix not being a full-rank matrix
 - d. It causes one or more features to be estimable from other features
- 2. Which of the following is NOT a feature of PCA?
 - a. It creates uncorrelated features
 - b. It creates mutually orthogonal features
 - c. The features created are linear combinations of the older features
 - d. None of the above
- 3. What is meant by "preserving maximum information" in context of PCA?
 - a. The sum of squares of residuals is maximized
 - b. The variance captured is maximized
 - c. The mean squared error is minimized
 - d. The variance captured is minimized
- 4. Which one of the following is NOT a pair of basis vectors of the 2D space?
 - a. \hat{i}, \hat{j}

 - b. $-\hat{i}, -\hat{j}$ c. $\frac{1}{\sqrt{2}}(\hat{i}+\hat{j}), \frac{1}{\sqrt{2}}(\hat{i}-\hat{j})$
- 5. Which is NOT the fundamental property of basis vectors in a vector space?
 - a. Linear combinations of basis vectors are basis vectors
 - b. They have unit magnitude
 - c. They are orthogonal
 - d. They can uniquely describe each point in the vector space
- 6. How do you determine the covariance matrix for a dataset?
 - a. By calculating pairwise correlation coefficients and using eigendecomposition
 - b. By calculating pairwise correlation coefficients and normalizing the matrix
 - c. By calculating pairwise covariances and normalizing the matrix
 - d. By calculating pairwise covariances
- 7. How are the directions of the principal components determined?
 - a. By eigen-decomposing the feature matrix
 - b. By eigen-decomposing the covariance matrix
 - c. By factorizing the covariance matrix
 - d. By factorizing the feature matrix
- 8. Which of the following statements regarding PCA is NOT right?

- a. Eigenvalue corresponding to a particular eigenvector represent the percentage variance explained by that eigenvector
- b. Eigenvalues when performing PCA can be sometimes negative
- c. Eigenvectors represent the directions of principal components
- d. Eigenvectors represent linear transformation of the original basis vectors
- 9. Arrange the following steps for the PCA algorithm in the correct order
 - i. Represent all data in a covariance matrix after normalizing them
 - ii. Eigen-decompose the covariance matrix to get set of eigenvectors and eigenvalues
 - iii. New basis vectors are created from eigenvectors and assigned principal components based on decreasing magnitude of its corresponding eigenvalues iv. Points are assigned coordinates based on the new basis vectors
 - a. I-> II-> III-> IV
 - b. II -> I -> III -> IV
 - c. III -> II -> I -> IV
 - d. IV -> I -> II -> III
- 10. How do you choose the value of k in KMeans?
 - a. Kink in the elbow plot
 - b. Minimizing silhouette score
 - c. Maximizing area under the elbow curve
 - d. All of the above
- 11. What pre-processing step do you need to perform before fitting the data to KMeans?
 - a. Standard Scaling
 - b. Label encoding of categorical variables
 - c. One hot encoding of categorical variables
 - d. All of the above
- 12. What is NOT one of the drawbacks of the KMeans method?
 - a. The value of k has to be set beforehand
 - b. The clusters are spherical in shape
 - c. The loss function is the within cluster sum squared value
 - d. Categorical variables need to be dropped before
- 13. To solve which hindrance of the KMeans method was the KMeans++ method devised?
 - a. KMeans suffers from the problem of the loss function being composed of Euclidean distances
 - b. KMeans clustering is susceptible to cluster centroid initialization
 - c. KMeans requires input of k, the number of clusters
 - d. KMeans cannot return non-spherical clusters
- 14. What is/are the advantage(s) of hierarchical clustering over KMeans?
 - a. The number of clusters do not need to be passed

- b. The distance measure can be non-euclidean
- c. We can select arbitrary number of clusters based on threshold values
- d. All of the above

15. What does DBSCAN stand for?

- a. Density Based Statistical Clustering with Amplified Noise
- b. Density Based Spatial Clustering of Applications with Noise
- c. Density Based Statistical Clustering of Applications with Noise
- d. Density Based Spatial Clustering with Amplified Noise
- 16. What is the complexity vs interpretability tradeoff?
- 17. What is the bias vs variance tradeoff?
- 18. What is NOT true in the context of overfitting?
 - a. It occurs when the model has learnt the test data really well, but cannot generalize to train data
 - b. It is undesirable
 - c. It is usually associated with more complex models
 - d. We can reduce overfitting by creating bootstrapped training datasets
- 19. Why do we need cross-validation?
- 20. What are the types of regularization present in Ridge and Lasso and ElasticNet?
 - a. L1 for Ridge, L2 for Lasso, L1 & L2 for ElasticNet
 - b. L1 for Lasso, L2 for Ridge, L1 & L2 for ElasticNet
 - c. L1 for ElasticNet, L2 for Ridge, L1 & L2 for Lasso
 - d. L1 for ElasticNet, L2 for Lasso, L1 & L2 for Ridge