MACHINE LEARNING

in Q1	to Q5, only one option is correct, Choose the (correct option:
1.	In which of the following you can say that the mode. A) High R-squared value for train-set and High R B) Low R-squared value for train-set and High R C) High R-squared value for train-set and Low R D) None of the above Answer-C	R-squared value for test-set. R-squared value for test-set.
2.	Which among the following is a disadvantage of A) Decision trees are prone to outliers. B) Decision trees are highly prone to overfitting. C) Decision trees are not easy to interpret D) None of the above. Answer-B	decision trees?
3.	Which of the following is an ensemble technique A) SVM C) Random Forest Answer-C	B) Logistic Regression D) Decision tree
4.	disease is most important. In this case which of A) Accuracy	for detection of a fatal disease where detection of the the following metrics you would focus on? B) Sensitivity None of the above.
5.	The value of AUC (Area under Curve) value for 0.85. Which of these two models is doing better A) Model A C) both are performing equal Answer-B	
In Q6	to Q9, more than one options are correct, Cho	ose all the correct options:
6.	Which of the following are the regularization tech A) Ridge C) MSE Answer-A&D	nnique in Linear Regression?? B) R-squared D) Lasso

B) Decision Tree

D) Xgboost.

C) Random Forest

A) Adaboost

Answer-A&D

8. Which of the techniques are used for regularization of Decision Trees?

7. Which of the following is not an example of boosting technique?

- A) Pruning
- C) Restricting the max depth of the tree
- B) L2 regularization
- D) All of the above

Answer-A&C

- 9. Which of the following statements is true regarding the Adaboost technique?
 - A) We initialize the probabilities of the distribution as 1/n, where n is the number of data-points
 - B) A tree in the ensemble focuses more on the data points on which the previous tree was not performing well
 - C) It is example of bagging technique
 - D) None of the above

Answer- A&B

Q10 to Q15 are subjective answer type questions, Answer them briefly.

10. Explain how does the adjusted R-squared penalize the presence of unnecessary predictors in the model?

Answer-

The adjusted R-squared is a modified version of R-squared that accounts for predictors that are not significant in a regression model. In other words, the adjusted R-squared shows whether adding additional predictors improve a regression model or not. To understand adjusted R-squared, an understanding of R-squared is required. R-squared comes with an inherent problem – additional input variables will make the R-squared stay the same or increase (this is due to how the R-squared is calculated mathematically). Therefore, even if the additional input variables show no relationship with the output variables, the R-squared will increase. An example that explains such an occurrence is provided below.

Essentially, the adjusted R-squared looks at whether additional input variables are contributing to the model. Consider an example using data collected The idea is that you can change the value of one independent variable and not the others. However, when independent variables are correlated, it indicates that changes in one variable are associated with shifts in another variable. The stronger the correlation, the more difficult it is to change one variable without changing another. It becomes difficult for the model to estimate the relationship between each independent variable and the dependent variable *independently* because the independent variables tend to change in unison.

11. Differentiate between Ridge and Lasso Regression.

Answer-

Ridge:- Similar to the lasso regression, ridge regression puts a similar constraint on the coefficients by introducing a penalty factor. However, while lasso regression takes the magnitude of the coefficients, ridge regression takes the square.

Lasso :- In this shrinkage technique, the coefficients determined in the linear model from equation above are shrunk towards the central point as the mean by introducing a penalization factor called the alpha α (or sometimes lamda) values. Alpha (α) is the penalty term that denotes the amount of shrinkage (or constraint) that will be implemented in the equation. With alpha set to zero, you will find that this is the equivalent of the linear regression model from equation 1.2, and a larger value penalizes the optimization function. Therefore, lasso regression shrinks the coefficients and helps to reduce the model complexity and multi-collinearity.

Considering the geometry of both the lasso (left) and ridge (right) models, the elliptical contours (red circles) are the cost functions for each. Relaxing the constraints introduced by the penalty factor leads to an increase in the constrained region (diamond, circle). Doing this continually, we will hit the center of the ellipse, where the results of both lasso and ridge models are similar to a linear regression model.

However, both methods determine coefficients by finding the first point where the elliptical contours hit the region of constraints. Since lasso regression takes a diamond shape in the plot for the constrained region, each time the elliptical regions intersect with these corners, at least one of the coefficients becomes zero. This is impossible in the ridge regression model as it forms a circular shape and therefore values can be shrunk close to zero, but never equal to zero.

12. What is VIF? What is the suitable value of a VIF for a feature to be included in a regression modelling?

Answer-

A variance inflation factor (VIF) is a measure of the amount of multicollinearity in regression analysis. Multicollinearity exists when there is a correlation between multiple independent variables in a multiple regression model. This can adversely affect the regression results. Thus, the variance inflation factor can estimate how much the variance of a regression coefficient is inflated due to multicollinearity. A variance inflation factor (VIF) provides a measure of multicollinearity among the independent variables in a multiple regression model. Detecting multicollinearity is important because while multicollinearity does not reduce the explanatory power of the model, it does reduce the statistical significance of the independent variables.

Multicollinearity creates a problem in the multiple regression model because the inputs are all influencing each other. Therefore, they are not actually independent, and it is difficult to test how much the combination of the independent variables affects the dependent variable, or outcome, within the regression model.

13. Why do we need to scale the data before feeding it to the train the model?

Answer-

scaling of the data comes under the set of steps of data pre-processing when we are performing machine learning algorithms in the data set. As we know most of the supervised and unsupervised learning methods make decisions according to the data sets applied to them and often the algorithms calculate the distance between the data points to make better inferences out of the data. In real life, if we take an example of purchasing apples from a bunch of apples, we go close to the shop, examine various apples and pick various apples of the same attributes. Because we have learned about the attributes of apples and we know which are better and which are not good also we know which attributes can be

compromised and which cannot. So if most of the apples consist of pretty similar attributes we will take less time in the selection of the apples which directly affect the time of purchasing taken by us. The moral of the example is if the apples every apple in the shop is good, we will take less time to purchase or if the apples are not good enough we will take more time in the selection process which means that if the values of attributes are closer we will work faster and the chances of selecting good apples also strong.

14. What are the different metrics which are used to check the goodness of fit in linear regression?

Answer-

Mean absolute error (MAE): Represent average error(You are going to take error from every single data points, and take average)

Mean Squared error (MSE): Similarly to MAE but noise is exaggerated and larger error are "punished", it is harder to interpret than MAE as it not in base units, however it is generally more popular. (There are some outliers in data, it can ignore all outliers and it will go with majority) Root Mean Squared Error: Most popular metric, similar to MSE, however, the result is square rooted to make it more, it interpretable as it's in base units, it is recommended that RMSE be used as primary metric to interpret your model.

15. From the following confusion matrix calculate sensitivity, specificity, precision, recall and accuracy.

Actual/Predicted	True	False
True	1000	50
False	250	1200

Answer-

1000 = True positive 50 False Positive 250 = False Negative 1200 True Negative 1) Accuracy

TP + TN / (TP + TN + FP + FN) = 1000 + 1200 / (1000 + 1200 + 50 + 250) = 2200 / 2500 = 0.88

- 2) Recall
- = TP / (TP + FN)
- = 1000 / (1000 + 250)
- = 1000/1250
- = 0.8
- 3) Precision
- = TP / (TP + FP)
- = 1000 / (1000 + 50)
- = 1000 / 1050
- = 0.95
- 4) Specificity
- = TN / (TN + FP) = 1200 / (1200 + 50)
- = 1200 / 1250
- = 0.96