

MACHINE LEARNING

- Which of the following in sk-learn library is used for hyper parameter tuning?
A) GridSearchCV()
B) RandomizedCV()
C) K-fold Cross Validation
D) All of the above
- In which of the below ensemble techniques trees are trained in parallel?
A) Random forest
B) Adaboost
C) Gradient Boosting
D) All of the above
- In machine learning, if in the below line of code:
`sklearn.svm.SVC (C=1.0, kernel='rbf', degree=3)`
we increasing the C hyper parameter, what will happen?
A) The regularization will increase
B) The regularization will decrease
C) No effect on regularization
D) kernel will be changed to linear
- Check the below line of code and answer the following questions:
`sklearn.tree.DecisionTreeClassifier(*criterion='gini',splitter='best',max_depth=None, min_samples_split=2)`
Which of the following is true regarding max_depth hyper parameter?
A) It regularizes the decision tree by limiting the maximum depth up to which a tree can be grown.
B) It denotes the number of children a node can have.
C) both A & B
D) None of the above
- Which of the following is true regarding Random Forests?
A) It's an ensemble of weak learners.
B) The component trees are trained in series
C) In case of classification problem, the prediction is made by taking mode of the class labels predicted by the component trees.
D)None of the above
- What can be the disadvantage if the learning rate is very high in gradient descent?
A) Gradient Descent algorithm can diverge from the optimal solution.
B) Gradient Descent algorithm can keep oscillating around the optimal solution and may not settle.
C) Both of them
D) None of them
- As the model complexity increases, what will happen?
A) Bias will increase, Variance decrease
B) Bias will decrease, Variance increase
C)both bias and variance increase
D) Both bias and variance decrease.
- Suppose I have a linear regression model which is performing as follows:
Train accuracy=0.95 and Test accuracy=0.75
A) model is underfitting
B) model is overfitting
C) model is performing good
D) None of the above

Q9 to Q15 are subjective answer type questions, Answer them briefly.

- 9. Suppose we have a dataset which have two classes A and B. The percentage of class A is 40% and percentage of class B is 60%. Calculate the Gini index and entropy of the dataset.**

ANS: The Gini index and entropy of the dataset with 40% A and 60% B classes can be calculated as follows:

$$\text{Gini Index} = 1 - (0.4)^2 - (0.6)^2 = 0.48$$

$$\text{Entropy} = -(0.4\log_2(0.4) + 0.6\log_2(0.6)) = 0.97$$

- 10. What are the advantages of Random Forests over Decision Tree?**

Random Forests have the following advantages over Decision Tree:

Random Forests are less prone to overfitting.

They are faster to train.

They can handle missing values and imbalanced datasets better.

They can give feature importance.

- 11. What is the need of scaling all numerical features in a dataset? Name any two techniques used for scaling.**

Ans: Since the features have different scales, there is a chance that higher weightage is given to features with higher magnitude. This will impact the performance of the machine learning algorithm and obviously, We do not want our algorithm to be biased towards one feature. We scale our data before employing a distance based algorithm so that all the features contribute equally to the result. o Normalization is a scaling technique in which values are shifted and rescaled so that they end up ranging between 0 and 1. It is also known as Min-Max scaling. o Standardization is another scaling technique where the values are centred around the mean with a unit standard deviation. This means that the mean of the attribute becomes zero and the resultant distribution has a unit standard deviation.

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- 12. Write down some advantages which scaling provides in optimization using gradient descent algorithm.**

Ans: Machine learning algorithms like linear regression, logistic regression, neural network, etc. that use gradient descent as an optimization technique require data to be scaled. The presence of feature value X in the formula will affect the step size of the gradient descent. The difference in ranges of features will cause different step sizes for each feature. To ensure that the gradient descent moves smoothly towards the minima and that the steps for gradient descent are updated at the same rate for all the features, we scale the data before feeding it to the model.

- 13. In case of a highly imbalanced dataset for a classification problem, is accuracy a good metric to measure the performance of the model. If not, why?**

ANS: When the class distribution is slightly skewed, accuracy can still be a useful metric. When the skew in the class distributions are severe, accuracy can become an unreliable measure of model performance. The reason for this unreliability is centered around the average machine learning practitioner and the intuitions for classification accuracy. Typically, classification predictive modeling is practiced with small datasets where the class distribution is equal or very close to equal. Therefore, most practitioners develop an intuition that large accuracy score (or conversely small error rate scores) are good, and values above 90 percent are great. Achieving 90 percent classification accuracy, or even 99 percent classification accuracy, may be trivial on an imbalanced classification problem. This means that intuitions for classification accuracy developed on balanced class distributions will be applied and will be wrong, misleading the practitioner into thinking that a model has good or even excellent performance when it, in fact, does not.

- 14. What is “f-score” metric? Write its mathematical formula.**

ANS: F-score is a metric that is used to evaluate the performance of a model in binary classification problems. It is the harmonic mean of precision and recall. Its mathematical formula is as follows:

$$\text{F-score} = 2 * (\text{Precision} * \text{Recall}) / (\text{Precision} + \text{Recall})$$

- 15. What is the difference between fit(), transform() and fit_transform()?**

Ans: Fit() - In the fit() method, where we use the required formula and perform the calculation on the feature values of input data and fit this calculation to the transformer. For applying the fit() method we have to use .fit() in front of the transformer object. Transform() - For changing the data we probably do transform, in the transform() method, where we apply the calculations that we have calculated in fit() to every data point in feature F. We have to use .transform() in front of a fit object because we transform the fit calculations. We use the example that is used above section when we create an object of the fit method then we just put it in front of the .transform and transform method uses those calculations to transform the scale of the data points, and the output will we get is always in the form of sparse matrix or array. Fit_transform() - This fit_transform() method is basically the combination of fit method and transform method, it is equivalent to fit().transform(). This method performs fit and transform on the input data at a single time and converts the data points. If we use fit and transform separate when we need both then it will decrease the efficiency of the model so we use fit_transform() which will do both the work.