

Machine Learning

Unit III - Supervised Learning : Regression

Data
Page

Q. 1 Q

Explain in brief techniques to reduce under fitting & over fitting.

→ - The data model is cannot capture the relationship between the i/p & o/p variables accurately called underfitting.

- It generate a high error rate on both the training set & unseen data.

- Underfitting destroy the accuracy of our ML model.

- It occurs when a model is too simple, which can be a result of a model needing more training time, more i/p features.

- Underfitting can be avoided by using more data & also reducing the features by feature selection.

- e.g - building a linear regression model over non-linear data.

• Techniques to reduce underfitting.

- ① Increase model complexity
- ② Increase number of features, performing feature engineering
- ③ Remove noise from the data
- ④ Increase the duration of training to get better result
- ⑤ High bias & low variance

2) Overfitting

- Overfitting is a opposite of underfitting.

- It occurring when the model has been overtrained or when it contains too much complexity, resulting high error rates on test data.

- Overfitting is the main problem that occur in supervised learning.

- When a model perform very well for training data but has poor performance for test data is known as overfitting.

- Reasons for overfitting - ① High variance & low bias

② The model is too complex

③ The size of training data is always in overfits.

- Techniques to reduce overfitting
- ① Increase training data
 - ② Reduce model complexity
 - ③ Early stopping during the training phase
 - ④ Ridge Regularization & Lasso Regularization
 - ⑤ Use dropout for neural networks to tackle overfitting

Find the equation of linear regression line using following data:

X	Y
1	3
2	4
3	5
4	7

$$\text{cm} = \frac{\sum xy - \bar{x}\bar{y}}{\sum x^2 - (\bar{x})^2}$$



X	Y	XY	X ²
1	3	3	1
2	4	8	4
3	5	15	9
4	7	28	16
$\Sigma x = 10$	$\Sigma y = 19$	$\Sigma xy = 56$	$\Sigma x^2 = 30$

The equation of regression line is

$$Y' = ax + b$$

$$a = n \sum xy - \bar{x} \bar{y}$$

$$n \sum x^2 - (\sum x)^2$$

$$= 4 \times 54 - 10 \times 19$$

$$4 \times 90 - 100$$

$$= 216 - 190$$

$$120 - 100$$

$$= \frac{20}{20}$$

$$a = 1.0$$

$$\begin{aligned}
 b &= \frac{1}{n} (\sum Y - \alpha \sum x) \\
 &= \frac{1}{4} (19 - 1.5 \times 10) \\
 &= \frac{5}{4} \\
 b &= 1.25
 \end{aligned}$$

$$Y = \alpha X + b$$

$$Y = 1.25X + 1.5$$

$$\begin{array}{ll}
 \text{For } X = 1 & Y = 1.25 \times 1 + 1.5 = 2.8 \\
 \text{For } X = 2 & Y = 1.25 \times 2 + 1.5 = 4.1 \\
 \text{For } X = 3 & Y = 1.25 \times 3 + 1.5 = 5.4 \\
 \text{For } X = 4 & Y = 1.25 \times 4 + 1.5 = 6.7 \\
 \hline
 & X
 \end{array}$$

(c) Write short note on :

① MAE

② RMSE

③ R²

→ ④ MAE

- MAE stand for Mean Absolute Error.
- MAE is a very simple metric.
- MAE calculates the absolute difference between actual & predicted values.

- Example :- We have I/P data & O/P data of use linear regression which draw a best fit line.

$$MAE = \frac{1}{N} \sum |Y - \hat{Y}|$$

④ sum of all the errors & divide them by a total number of observation to get a MAE.

• Advantage

- ① The MAE get same unit as the o/p variable.
- ② It is most robust to outliers

• Disadvantage:

- ① The graph of MAE is not differentiable

2) RMSE

- RMSE stands for Root Mean Squared Error
- It measures the average difference b/w a statistical model's predicted values & the actual values.
- RMSE is a simple square root of mean squared error (mse).

$$RMSE = \sqrt{MSE} = \sqrt{\frac{1}{n} \sum_{j=1}^n (y_j - \hat{y}_j)^2}$$

• Advantages:-

- ① It gives a relatively high weight to large errors
- Disadvantage:-
- ① Not robust to outliers

$$\Rightarrow R^2$$

- R^2 is also known as coefficient of determination or sometimes also known as goodness of fit
- R^2 score is used to evaluate the performance of a linear regression model

$$R^2 \text{ squared} = 1 - \frac{SS_r}{SS_m}$$

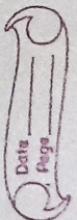
$$SS_r = \text{Sum error of regression line}$$
$$SS_m = \text{sum error of mean line}$$

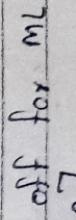
Regularization — Lasso & ridge regression

Q.2.Q

Explain in brief Lasso & Ridge Regression
→ Lasso Regression

- ① In machine learning, Lasso stands for least Absolute Shrinkage & Selection Operator

-  Data Flows
- ⑥ It is a regression analysis method.
 - ⑦ Lasso is used for the overcome of overfitting.
 - ⑧ Lasso regression is a regularization technique.
 - ⑨ It uses shrinkage [shrinkage \rightarrow Reduction].
 - ⑩ If a regression model uses the L1 regularization technique, it is called Lasso Regression.
 - ⑪ Lasso regularization has the ability to set some coefficients to zero.
 - ⑫ Lasso is applied in economics & finance.
 - ⑬ It is found to improve prediction.
- 2) Ridge Regression
- ⑭ Ridge regression is a model tuning method.
 - ⑮ It is used to analyse any data that suffer from multicollinearity.
 - ⑯ multicollinearity :- It is a statistical concept where several independent variables in a model are correlated.
 - ⑰ This method performs L2 regularization.
 - ⑱ Ridge regression model can be used on dataset that have many correlated features.
 - ⑲ Ridge regression is mostly used to reduce the overfitting in the model.
 - ⑳ It reduces the complexity of the model.
 - ㉑ Ridge regression is a type part of linear regression.

-  Data Flows
- Q) What is bias & variance trade off for ml model?
- S) First explain ridge regression.
- ① In machine learning, if we minimize one component of the error [i.e. bias], & the other component [i.e. variance] tends to increase vice versa.
 - ② Finding the right balance of bias & variance is key to creating an effective & accurate model called as bias & variance tradeoff.

③ The bias-variance tradeoff describes the relationship b/w a model's complexity.

④ Applications:-

⑤ Regression ⑥ Classification ⑦ Reinforcement learning ⑧ Human learning.

⑨ Diagram

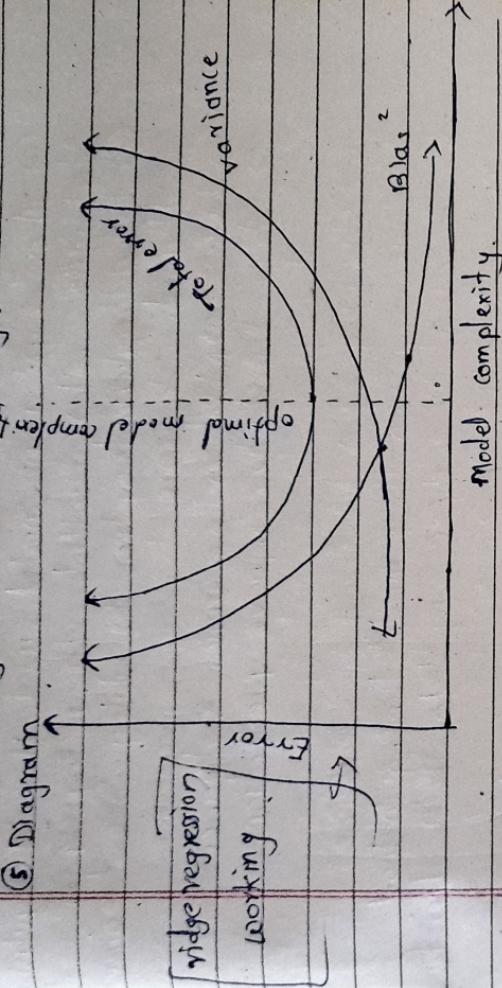


Fig - Bias-variance tradeoff.

- ⑩ Write short note on Evaluation metrics.
- ① Evaluation metrics are used to measure the quality of the statistical or machine learning models.
 - ② There are many different types of evaluation metrics available to test a model.
 - ③ The evaluation metrics includes classification accuracy, logarithmic loss, confusion matrix & so on.
 - ④ Classification accuracy is the ratio of the number of correct predictions to the total number of input samples.

⑤ Logarithmic loss also called log loss.

⑥ A confusion matrix gives us a matrix as a/p of describe the complete performance of the model.

• Applications of Evaluation metric.

① statistical analysis

② machine learning

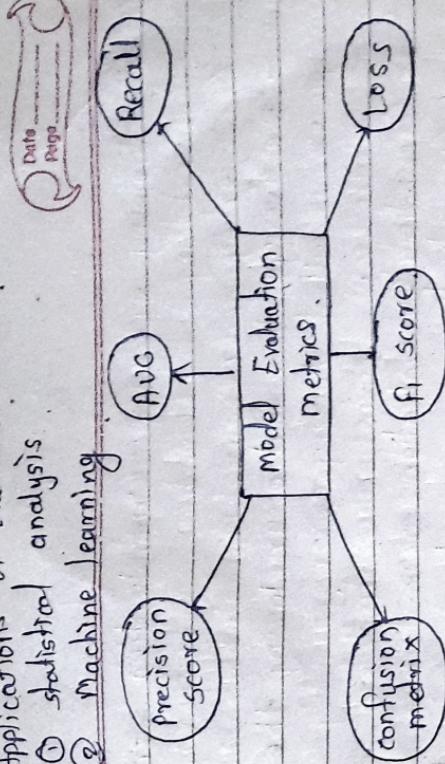


Fig :- model Evaluation metrics

* Evaluation metrics :- MAE, RMSE, R^2



Linear Regression

Regression models

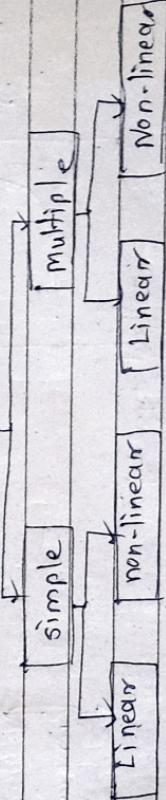


Fig . Types of regression model.

- ① It is used to predict the value of variable based on the value of another variable
- ② It is a data analysis technique , that predicts the value of unknown data by using another related & known data value

③ It provide continuous output

- ④ It is used to solve regression problem
- ⑤ It is used to draw best fit line
- ⑥ It is most popular machine learning algorithm
- ⑦ α & β are parameters of regression
- ⑧ $y = mx + c$ where y = dependent variable

x = independent variable m = slope of line

c = estimate intercept.

if identifying g predicting future outcome of

if

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#. ② Training error & Generalization error

- ↳ Training error.
 - The average loss that occur during the training process is called training error.
- ↳ Generalization error
 - Generalization error can be minimized by avoiding overfitting in the learning algorithm.
 - Types of generalization errors
 - ① Bias . ② variance, ③ irreducible error.

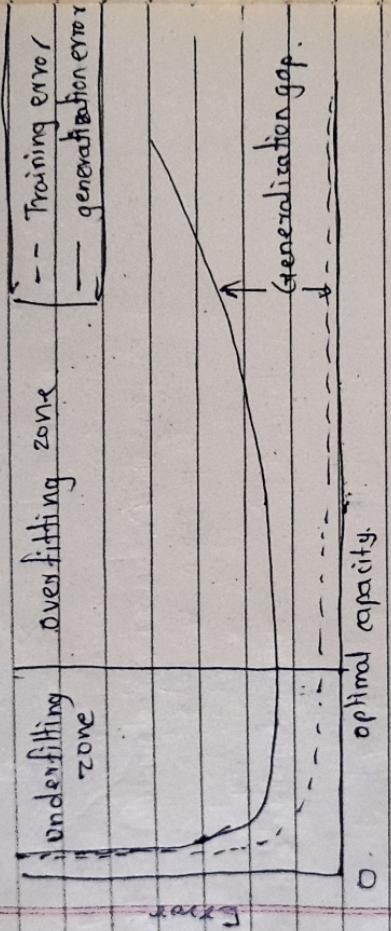


fig. Training error & generalization error.

- ## #. The Gradient Descent algorithm
- ① It is an optimisation algorithm which is commonly used to train machine learning models of neural networks.
 - ② The GD procedure is an algorithm for finding the minimum of a function.
 - ③ This method is commonly used in ML for:
 - ① to minimize a cost function.
 - ④ GD algorithm does not work for all functions. There are 2 specific requirement. a function has
 - ① Differentiable
 - ② convex

• GD method's steps

- ① choose a starting point (initialization)
- ② calculate gradient at this point
- ③ make a scaled step in the opposite direction to the gradient
- ④ repeat point 2 & 3 until one of the criteria is met:
 - ⑤ maximum no. of iterations reached
 - ⑥ step size is smaller than the tolerance

parameter tuning
parametric error

① Bias

② Variance

③ Underfitting

④ Overfitting

underfitting

- 1) Overfitting - high bias & low variance
- 2) Underfitting - low bias & high variance.

- ① Regression - performance matrix
- ② Classification - confusion matrix

regression

Loss bias - error in testing data
Variance - error in training dataset

Generalize - equal bias of variance

- Always overfit in decision tree

loss

$$\text{Ridge regression} = \text{Loss} + \alpha \|w\|_2^2$$

penalty

some feature has same value
multicollinearity

problem solve

regression

use on fit

- It minimize the complexity of problem
- penalty is reduce the loss. [penalty should be change]
- solve overfitting problem

$\|w\|$ - vector of coefficient, α = constant

- if α increases, the magnitude of coefficient will be reduce. is almost equal to 0

$$L_1 = \text{Lasso regression} = \text{Loss} + \alpha \|w\|_1$$

- If α increase, coefficient is almost zero but not 0

reduce the loss

solve overfitting problem

= Loss calculate by using cost function

Unit 4 - Supervised Learning : Classification

Q.1
SIN

Explain in brief methods used for evaluating classification models.

- ① Accuracy
- ② Precision
- ③ Error rate
- ④ Recall (sensitivity)
- ⑤ Specificity
- ⑥ F1 score
- ⑦ AUC - ROC
- ⑧ Log Loss.

-
- Accuracy is the ratio of the number of correct predictions and the total number of predictions
 - Accuracy is useful for balanced classes but not good choice for unbalanced classes
 - Formula =

$$\text{Accuracy} = \frac{\text{TP} + \text{TN}}{\text{TP} + \text{FP} + \text{TN} + \text{FN}}$$

2) Precision

- Precision is a measure of exactness
- The ability of a classification model to identify only the relevant data points
- Formula =

$$\text{precision} = \frac{\text{TP}}{\text{TP} + \text{FP}}$$

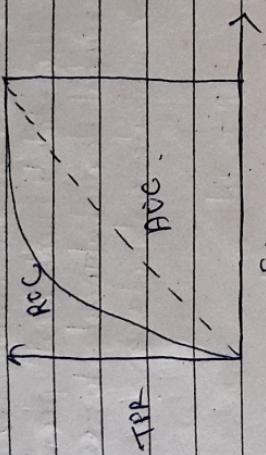
3) Recall

- Recall is a measure of completeness
- It has ability to detect positive samples
- Recall & sensitivity are similar
- Formula

$$\text{Recall} = \frac{\text{TP}}{\text{TP} + \text{FN}}$$

4) AUC - ROC

- AUC stands for Area under the Curve of ROC curve. It represents the probability curve of ROC.
- ROC is the probability curve of ROC represents the degree or measure of separability.
- It tells how much the model is capable for distinguishing between classes.



① TPR (True Positive Rate) (Recall).

$$TPR = \frac{TP}{TP + FN}$$

② FPR (False positive Rate)

$$\begin{aligned} FPR &= 1 - Recall \\ &= \frac{FP}{TN + FP} \end{aligned}$$

5) Log loss

- It is also called as logarithmic loss or logistic regression.
- Log loss measures the quality of probabilistic predictions, it score ignores the probabilistic measure of nature of classification.
- It takes input as a probability value between 0 & 1.
- Log loss is closer to '0'.
- If lower log loss value means better predictions.

Q.2

Consider the following data to predict the student pass or fail using the K-Nearest Neighbour (KNN) for the values physics = 6 marks, chemistry = 8 marks with number of neighbours $K = 9$.

Physics (marks)	Chemistry (marks)	Results
4	9	Fail
6	7	Pass
7	8	Pass
5	5	Fail
8	8	Pass

\rightarrow Physics = 6 marks, Chemistry = 8 marks
 $K=3$

- According to Euclidean distance

$$d = \sqrt{\sum_{i=1}^2 (X_{\text{observed}} - X_{\text{actual}})^2}$$

Hence, we have 2 samples,
so,

$$d = \sqrt{(X_{01} - X_{A1})^2 + (X_{02} - X_{A2})^2}$$

$$\sqrt{(6-4)^2 + (8-8)^2} = \sqrt{2^2} = 2$$

$$\begin{aligned} \textcircled{1} & \quad \sqrt{(6-7)^2 + (8-8)^2} = \textcircled{1} \\ \textcircled{2} & \quad \sqrt{(6-6)^2 + (8-8)^2} = \textcircled{1} \end{aligned}$$

out of this
select '3'
nearest value
because the value
of $X = 3$

$$\textcircled{3} \quad \sqrt{(6-5)^2 + (8-8)^2} = \sqrt{10} = 3.16$$

$$\textcircled{4} \quad \sqrt{(6-5)^2 + (8-5)^2} = \sqrt{10} = 3.16$$

$$\textcircled{5} \quad \sqrt{(6-8)^2 + (8-8)^2} = \textcircled{2}$$

out of these select '2' nearest value because the value of $k=2$

- i.e. 1, 1, 2
- Nearest selected values have "Pass result" and zero fail i.e. $g > 0$
so the student is Pass

Q.3 Write short note on Ensemble learning methods

S. i) Simple & mean
ii) Advanced OR
What do you mean by ensemble learning
of difference b/w bagging & boosting
- Types of Ensemble learning.

→ Ensemble learning - It is a technique that combines multiple ml algorithms to produce one optimal solution

- It improves helps to improve ml results & decision

- In learning models noise, variance, bias are major sources of error these methods helps to minimize these errors

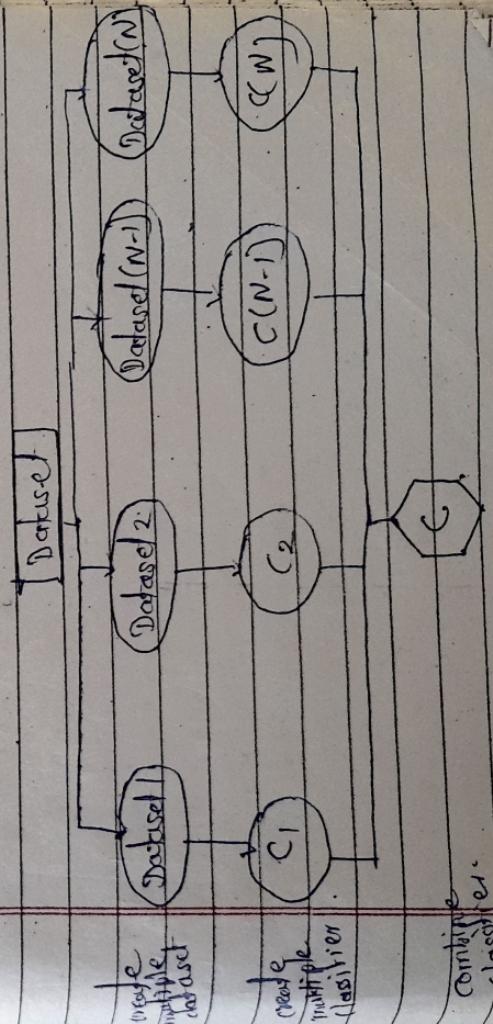


Fig. An overview of Ensemble learning.

contd.

↳ Simple :-

- ① mean / average
- ② mode
- ③ the weighted average.

* mean / average.

- In the mean / average ensemble techniques, data analysts takes average prediction made by all models when making ultimate predictions.

→ Model

- In statistical terminology, "mode" is the no. or value that most appears in a dataset.
- In ensemble learning techniques, ml professionals use a no. of models for making predictions about each data point.

→ The weighted average.

- In the weighted average ensemble method, data scientists assign different weights to all the models in order to make a prediction.

• Types of ensemble

Advanced :- ① Bagging ② Boosting ③ AdaBoost

Bagging

① Reduce variance

② High

③ Each model receives equal weight

④ Parallel

⑤ High

⑥ Random Sampling

⑦ Highly computing intensive

Example

Robust

Boosting

① Reduce bias

② Low

③ models are weighted according to their performance.

④ sequential

⑤ low.

⑥ Systematic sampling.

⑦ less computing intensive.

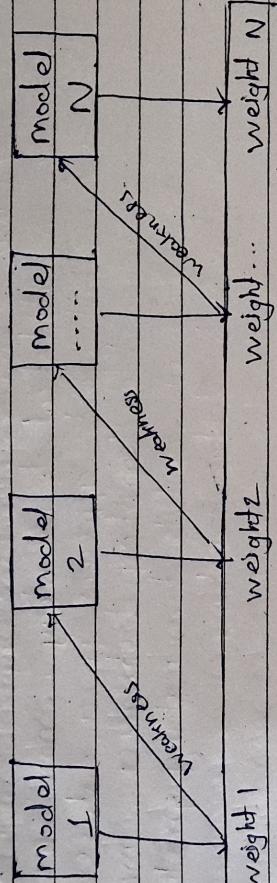
⑧ Adaboost

⑨ Not robust

5/11

③ AdaBoost

- AdaBoost stands for Adaptive boosting
- It is a popular boosting algorithm
- AdaBoost algorithms can be used for both classification & regression problem.
- AdaBoost helps you combine multiple "weak classifiers" into a "single" strong classifier.
- It is indirectly proportional to the weak learners.



Ensemble. (with all it's predecessors)

fig. AdaBoost Algorithm

④ Random Forest

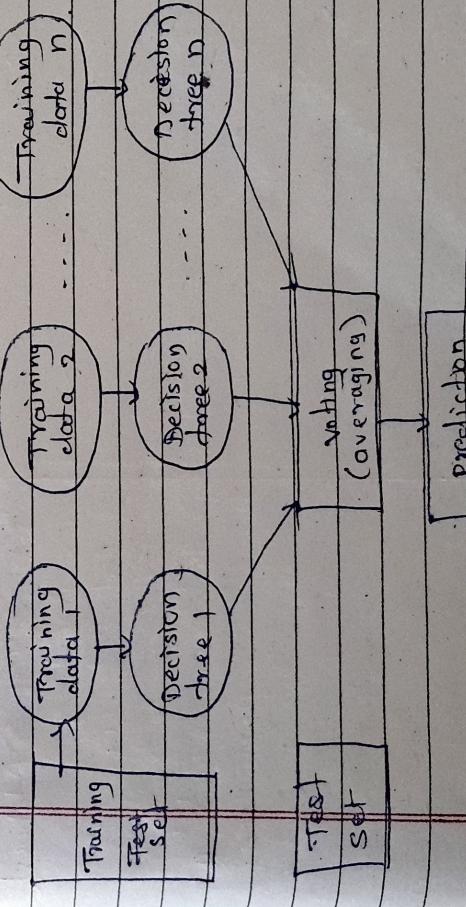


fig. Working of Random forest algorithm

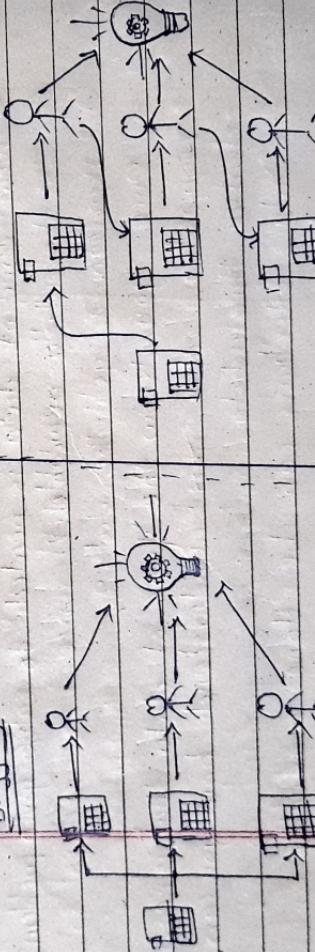
- Random forest is a popular ML algorithm that belongs to the supervised learning techniques.
- It can be used for both classification & regression problem in ml.
- It good predicted produced good prediction than can be understand easily.
- Random forest is a classifier that contains a number of decision trees of the given dataset.
- It is used as black box model in business.

- Working of Random Forest algo.

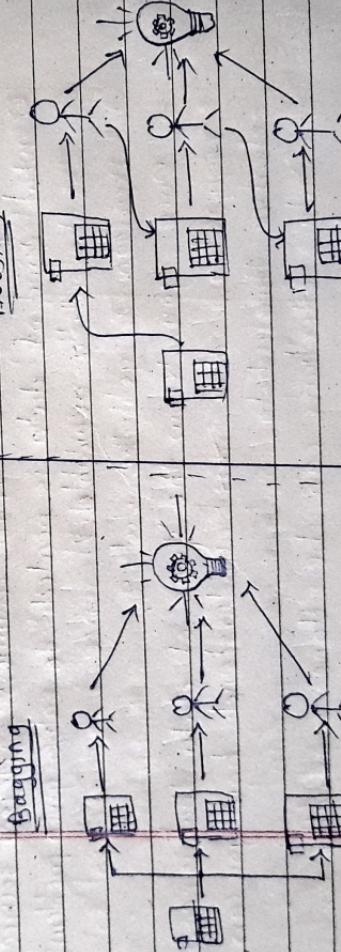
- Ensemble use of two types of methods

- ① Bagging
- ② Boosting

Bagging



Boosting



Advantages

- It takes less training time as compare to other.
- It is used in classification & regression problems.
- It solve overfitting problem.
- It predicts the o/p with high accuracy.
- It can handle large dataset efficiently.
- It automatically handle missing values.

Disadvantages

- ① Random Forest is highly complex.
- ② It can take lot of memory.
- ③ More trees slow down the model.
- ④ Random forest is much noisy.



Q. Write short note on importance of confusion matrix.

Ans. 6 M

or steps in ml modeling

- ⇒ • Steps in ml modeling
 - 1) Data collection
 - 2) Data preprocessing
 - 3) Model selection
 - 4) Model training - Training & testing K-fold cross validation
 - 5) Model Evaluation - ① Accuracy ② precision ③ Recall ④ Confusion matrix
 - 6) Hyper Parameter Tuning
 - 7) Prediction

1) Data collection

- Data collection is the process of gathering & measuring information from countless different sources.
 - Q what kind of data is needed to solve given problem?
 - Q Is the data available?
 - Q How can I get the data?
 - Once you know the type of data that is needed required, you must understand how you can derive this data
 - Data collection can be done manually or by web scraping.

2)

- Data preprocessing
 - It is a technique that is used to convert the raw data into clean dataset
 - Data preprocessing steps
 - ① Removing duplicates - It is an important part of data analysis if analyzing duplicate values of remove them
 - pandas drop_duplicates() method helps in removing duplicate values from data frame

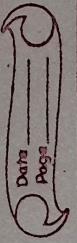
② Transformation of data using function or mapping
- Data transformation changes the format, structure or value of the data & converts them into clean, usable data.

③ Replacing values
- pandas `dataframe.replace()` function is used to replace a string, list, dictionary, series, numbers, etc from a `dataframe`.

④ Handling missing ~~to~~ data
- Missing values can be handled by deleting the rows or columns having null values
- `isna()` & `isnull()` function used to check missing data.

⑤ Model selection
- model selection is a procedure that select the most suitable model for dataset
- Common techniques for model selection includes cross validation, hyperparameter tuning & using evaluation metrics.
- choose appropriate ml algorithm based on the nature of your problem (e.g. classification, regression, clustering).
- Select a set of candidate models if necessary & consider their strength & weaknesses.

A) Model training
In ml data preprocessing, we divide our dataset into training set & test set
- This is an important step because it improve the performance of our ml model.
① Training set - we already know the op.
② Testing set - by using the test set predict the op.



• K-fold cross validation

- Cross-validation is a resampling procedure
- It is a technique used to determine / check the performance of ML model
- It involves splitting the dataset into "k" subsets or "folds"

5) Model Evaluation

Q) Precision, ② recall ③ Accuracy

A) Confusion matrix

- A confusion matrix is a table that shows how well a ML model is performing
- It compares the models prediction with the actual labels in the datasets
- The matrix has four matrix values:
 - ① True positives
 - ② True negatives
 - ③ False positives
 - ④ False negatives
- It calculate the difference both actual value & predicted value.

• Importance of confusion matrix

- The confusion matrix is important in ml if helps us evaluate how well our model is performing by showing us the no. of true positives, true negatives, false positives & false negatives
- With this information, we can calculate metrics like accuracy, precision, recall & F1 score
- It's like a road map to understand the strength & weaknesses of our models predictions

6)

Hyper Parameter tuning - It is an important step in model development process.

- Hyperparameter tuning in ML model is all about finding the best settings for the model
- By adjusting things like learning rate, regularization strength or no. of hidden units that is performing our model best

7) Predictions :

- Once the model is evaluated & improved, it is finally used to make prediction.
- The final output can be a categorical variable or it can be a continuous quantity.

Q

Explain with example the variant of SVM, the support vector regression SVM,

or

Define following terms with reference to SVM.
 ① Separating hyperplane ii Margin. 6m

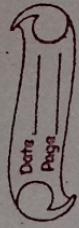
→

- SVM - Support vector machine
- SVM is a type of supervised learning that can be used for classification & regression
- Support vector regression (SVR) is a variant of SVM, that is used for regression task
- SVR is design to predict continuous numeric value.

- The goal of SVR is to find a function that best fits the training data.
- SVR is minimizing the error.

Example :- consider we have a dataset of houses with the following features & prices.

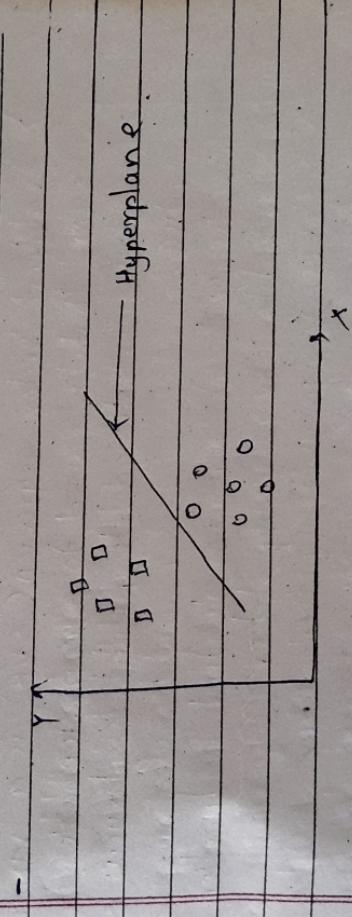
Square footage	No. of bedrooms	Location	Price
1500	3	1	400,000
2000	4	2	900,000
1800	3	1	2,50,000
2200	4	3	3,50,000
2500	5	2	4,00,000



Using SVD, we can train a model on this data to predict house prices based on the feature.

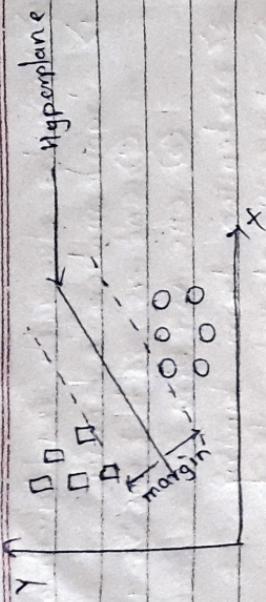
1) Separating Hyperplane

- Separating hyperplane is the line which is used to separate the dataset
- If data is n dimensional we required n-1 dimensional hyperplane
- SVM algorithm helps to find the best line or decision boundary, this best boundary is called as hyperplane
- It helps to determine which datapoint belongs to which class
- It helps to improve the generalization & robustness of the model



2) Margin

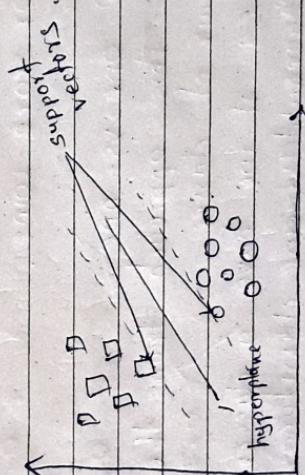
- The distance between the vectors of the hyperplane is called as margin.
- The hyperplane with maximum margin called optimal hyperplane
- we always choose the larger margin because if any mistake made in model then it can be easily solve it.



- It allows better classification & generalization.

3) Support vector

- Support vector is a data point that is closest to the decision boundary (hyperplane).
- SVM uses optimal hyperplane to create an optimal hyperplane that maximizes margin to improve performance of model.



~~Q~~ ~~SVM~~

- What are different variants of multi-class classifications
Explain them with suitable example.
- In ml multi-class classification is the problem of classifying data points into one of three or more classes
 - Many classification algorithm can be used to turn binary algorithm into multinomial classifier such as multinomial logistic regression

- Variants of multi-class classification
 - 1) Transformation to binary
 - (a) one-vs-rest [one-vs-all]
 - (b) one-vs-one
 - 2) extension from binary
 - a) hierarchical classification
 - 3) transformation to binary
 - (a) one-vs-rest [one-vs-all]
 - In multi-class problem, we create a separate binary classifier for each class
 - For example, if you have classes A, B, & C you create three binary classifiers: A vs {B, C}, B vs {A, C} and C vs {A, B}
 - The classifier with the highest probability or score wins
 - (b) one-vs-one
 - This approach trains the classifier on each pair of classes
 - e.g. If we have 3 classes A, B, & C it would train classifiers for A vs B, A vs C & B vs C.
 - The classifier with the highest score wins.

Q2 Calculate macro average precision, macro average recall & macro average F-score for the following given confusion matrix of multi-class classification

	A	B	C	D
A	100	80	10	10
B	0	9	0	1
C	0	1	8	1
D	0	1	0	9

→ Macro average precision
- First find out the precision for A, B, C & D

$$\textcircled{1} \quad P(A)$$

$$\text{Precision} = \frac{\text{TP}}{\text{TP} + \text{FP}}$$

$$\textcircled{1} \quad P(A) = \frac{100}{100+0+0+0} = \frac{100}{100} = 1$$

$$\textcircled{2} \quad P(B) = \frac{9}{9+8+1+1} = \frac{9}{19} = 0.099$$

$$\textcircled{3} \quad P(C) = \frac{8}{8+10+0+0} = \frac{8}{18} = 0.444$$

$$\textcircled{4} \quad P(D) = \frac{9}{9+10+1+1} = \frac{9}{21} = 0.429$$

Macro average precision = $P_1 + P_2 + \dots + P_n$

$$= 1 + 0.099 + 0.444 + 0.429$$

$$= \frac{1.972}{4} = \underline{\underline{0.493}}$$

2) Macro average recall

$$\text{recall} = \frac{\text{TP}}{\text{TP} + \text{FN}}$$

$$R(A) = \frac{100}{100+80+10+0} = \frac{100}{200} = \frac{1}{2} = 0.5$$

$$R(B) = \frac{9}{9+0+0+1} = \frac{9}{10} = 0.9$$

$$R(C) = \frac{8}{8+1+1+0} = \frac{8}{10} = 0.8$$

$$R(D) = \frac{9}{9+0+1+0} = \frac{9}{10} = 0.9$$

$$\text{Macro average recall} = \frac{R_1 + R_2 + \dots + R_n}{n}$$

$$= 0.5 + 0.9 + 0.8 + 0.9$$

$$= 0.75 \quad \underline{3.1}$$

$$= 0.775$$

3) Macro average F1-score

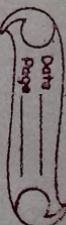
$$F_1\text{-score} = \frac{2 \times \text{precision} \times \text{recall}}{\text{precision} + \text{recall}} = \frac{2 \times \text{TP}}{2\text{TP} + \text{FP} + \text{FN}}$$

$$F_1(A) = \frac{2 \times 1 \times 0.5}{1+0.5} = \frac{1}{1.5} = 0.666$$

$$F_1(B) = \frac{2 \times 0.099 \times 0.9}{0.099+0.9} = \frac{0.1982}{0.999} = 0.1982$$

$$F_1(C) = \frac{2 \times 0.444 \times 0.8}{0.444+0.8} = \frac{0.7104}{1.244} = 0.571$$

$$F_1(D) = \frac{2 \times 0.429 \times 0.9}{0.429+0.9} = \frac{0.772}{1.329} = 0.581$$



Macro average F₁-Score = $\frac{P_{avg} + R_{avg}}{2}$

H

$$= \frac{0.666 + 0.198 + 0.571 + 0.581}{4}$$

$$= 1.9963$$

H

$$= 0.499025$$

Micro average precision =

$$= \frac{TP(A) + TP(B) + TP(C) + TP(D)}{TP(A) + TP(B) + TP(C) + TP(D) + FP(A) + FP(B) + FP(C) + FP(D)}$$

$$= \frac{100 + 9 + 8 + 9}{100 + 9 + 8 + 9 + 0 + 82 + 10 + 12}$$

$$= \frac{126}{280}$$

$$= \underline{0.4548}$$

2) Micro average recall =

$$= \frac{TP(A) + TP(B) + TP(C) + TP(D)}{TP(A) + TP(B) + TP(C) + TP(D) + FN(A) + FN(B) + FN(C) + FN(D)}$$

$$= \frac{100 + 9 + 8 + 9}{100 + 9 + 8 + 9 + 100 + 1 + 2 + 1}$$

$$= \frac{126}{280} = \underline{0.4548}$$

$$= \underline{0.548}$$

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Page _____

3) micro average F_1 -score

$$= \frac{TP(A) + TP(B) + TP(C) + TP(D)}{TP(A) + TN(A) + TP(C) + TP(D) + \frac{1}{2} (FP(A) + FP(B) + FP(C) + FP(D) + FN(A) + FN(B) + FN(C) + FN(D))}$$

$$= 100 + 9 + 8 + 9$$

$$100 + 9 + 8 + 9 + \frac{1}{2} (0 + 82 + 10 + 12 + 100 + 1 + 2 + 1)$$

$$= 126$$

$$126 + \frac{1}{2} (208)$$

$$= 126 = \frac{126}{230} = 0.548$$

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