

# Assignment - 1

## SECTION A – True / False

- 1. Bagging reduces the chance of overfitting, making the model more adaptable to unseen data. **True**
- 2. Averaging predictions reduces fluctuations in data. **True**
- 3. In boosting, the weights of data points change at every step to make the next gradient descent more accurate. **True**
- 4. Sampling without randomness can introduce sampling bias. **True**
- 5. If a model is overfit, training error can be 0% while testing error can be 100%. **True**
- 6. Regularization simplifies the model, which decreases bias. **False**
- 7. Increasing the depth of a decision tree always prevents overfitting. **False**
- 8. Every classifier makes assumptions about the data. **True**
- 9. Random Forests are more accurate than single decision trees because they combine bagged trees. **True**
- 10. Irreducible error is the lower bound on error due to inherent noise in the data. **True**

## SECTION B – ERM and SVM

1. Fill the following table:

Model	Loss Function	Regularizer
SVM	<u>Hinge Loss</u> : $\max(0, 1 - y(w \cdot x + b))$	<u><math>L_2   w^2  </math></u> : loss func. + $\lambda \sum_{i=1}^m w_i^2$
LASSO	<u>Mean Square Error (MSE)</u> : $(y - \hat{y})^2$	<u><math>L_1   w,   </math></u> : loss func. + $\lambda \sum_{i=1}^m  w_i $
RIDGE	<u>Mean Square Error (MSE)</u> : $(y - \hat{y})^2$	<u><math>L_2   w^2  </math></u> : loss func. + $\lambda \sum_{i=1}^m w_i^2$

Here;  $\lambda$  = regularization const.

3. Short Answer:

- (a) Which loss functions can be optimized using gradient descent? Why?
- (b) Which loss functions can be optimized using Newton's method? Why?

→ (a) Gradient descent works with first derivative of the loss function. So every differentiable (at least one) loss function can be optimized by gradient descent method. like - MSE (Convex Upward), logistic loss can be optimized by gradient descent.

b) While Newton's method deals with gradient (first order derivative) as well as Hessian



(second order derivative). So loss like that are twice differentiable can be optimized by Newton's method.  
like: MSE; as it is twice differentiable.

## SECTION C – Bias and Variance

1. Explain one major reason why underfitting occurs.

→ Underfitting occurs when the model is **too simple** to capture the underlying pattern in the data, leading to **high bias** and poor performance on both training and test data.

2. If both training and test error is high, what does this imply about the data?

→ If both the training & test error is high, it simply implies that the model suffers from **high bias**. Hence it is a **underfitted** model and unable to learn the structure of the data.

3. Explain how bagging reduces variance.

→ Bagging first divide the total dataset into several random samples or **bootstrap**; then it make prediction on the samples and finally take **average** of these predictions. By this method, it can cancel out individual model fluctuations and stabilize the final prediction.

4. Explain the effect of boosting on bias and variance.

→ Boosting also divide the dataset into some

random samples. Then it starts from assumed initial parameters, go through each sample, calculate loss function, update parameters accordingly. By this sequential learning method; it can **reduces bias**. It may **slightly increase variance**, but overall improves model accuracy.

## SECTION E – Decision Trees

1. Derive that the optimal prediction at a leaf (with square loss) is the mean.

→ Let a leaf contain targets  $y_i, \forall i, i = \{1, \dots, n\}$   
 Let the optimal prediction of the leaf be  $C$ ;  
 where  $C$  is a constant.  
 Square loss:

$$L(C) = \sum_{i=1}^n (y_i - C)^2$$

$$\Rightarrow \frac{\partial L(C)}{\partial C} = -2 \sum_{i=1}^n (y_i - C) \quad \left( \begin{array}{l} \text{differentiate} \\ \text{wrt } C \end{array} \right)$$

As a  $C$  is the optimal prediction;

then;  $L(C) \Big|_{C=C} = \text{minimum}$

$$\Rightarrow \frac{\partial L(C)}{\partial C} \Big|_{C=C} = 0 \Rightarrow \sum_{i=1}^n y_i - nC = 0$$

$$\Rightarrow C = \frac{1}{n} \sum_{i=1}^n y_i$$

∴ Hence; the optimal prediction is the mean of the targets



in the leaf.

2. What are the max/min Gini impurity values for 3 classes?

→ Gini impurity -

$$G = 1 - \sum_{k=1}^3 p_k^2$$

$p_k$  = Probability for each  $k$

Now we know that -

A lower gini index indicates more homogeneity on pure distribution.

While a higher gini index indicates a more heterogeneous on impure distribution.

Hence; minimum gini possible for a pure node:

→ pure node:  $p = (1, 0, 0)$

$$G_{\min} = 1 - (1^2 + 0^2 + 0^2) = 0$$

Similarly maximum gini possible for a perfect impure node. A perfect impure or heterogeneous node is a node with uniform distribution.

→ Uniformly distributed node:  $p = (1/3, 1/3, 1/3)$

$$\begin{aligned} G_{\max} &= 1 - \left( \left(\frac{1}{3}\right)^2 + \left(\frac{1}{3}\right)^2 + \left(\frac{1}{3}\right)^2 \right) \\ &= 1 - \left( 3 \times \frac{1}{9} \right) = \frac{2}{3} \end{aligned}$$

∴ Hence  $G_{\min} = 0$  &  $G_{\max} = \frac{2}{3}$  for 3 classes.

3. Why are decision trees myopic learners?

→ Myopic means short-sighted or greedy.

Decision tree acts myopic or greedy alike because at each node it chooses the best split to get immediate reduction in impurity independently **without considering any previous or future split.**

It means as other algorithms try to reach global minima in loss function; decision tree does not follow that.

4. Explain two methods to avoid overfitting in decision trees.

→ 1. Limiting maximum depth; we can prevent over-complexity of model.

2. Remove features, branches; we simplify the model and improve generalization.

## SECTION F – Boosting & Bagging

1. Can Random Forest use the same data for training and testing? Justify.

→ YES – Random Forest can use the same dataset for training and testing.

Because, Random Forest first does a sampling (random) over the dataset for training. So a chunk amount of data remains untouched after the training. So we can use the rest of the dataset as a test set. Hence Random Forests obtain an **unbiased estimate of test error** without requiring a separate test dataset.



## 2. Explain the key differences between bagging and boosting.

- 1. Bagging learn parallelly over all the random sampled dataset. While, boosting learn sequentially.
- 2. Bagging aim to reduce variance productively. While Boosting aim to reduce bias.
- 3. Hence, bagging deal with overfitting & boosting deal with underfitting.
- 4. Bagging; each model lead to diff. optima; finally take average of them. While Boosting update parameter sequentially.

