

# Homework #3

- 1.)  $X_n$  = input data sample, it's a vector  
 $y_n$  = ground-truth class label of  $X_n$   
 $\hat{y}_n$  = output "soft-label"/confidence of a logistic regression classifier given the input  $X_n$   
 $n = 1$  to  $N$   
 $K$  = # of classes

1a.)

$$L_{\text{Binary}} = -\frac{1}{N} \sum_{n=1}^K (y_n \log(\hat{y}_n) + (1-y_n) \log(1-\hat{y}_n))$$

$K=2$   
 $y_n = \text{integer}$

$$- y_n \log(\hat{y}_n) + (1-y_n) \log(1-\hat{y}_n)$$

1b.)

$$L_{\text{Binary}} = -\frac{1}{N} \sum_{n=1}^N \sum_{k=1}^K y_{(n,k)} \log(\hat{y}_{(n,k)})$$

$K \geq 2$   
 $y_n = \text{One-hot vector}$

- 1c.) 2 classes : class-0, class-1  
 $X_n$  is in class-1 ( $y_n=1$ )  
 $\hat{y}_n = 0.8$

$$L(y_n, \hat{y}_n) = -(y_n \log(\hat{y}_n) + (1-y_n) \log(1-\hat{y}_n))$$

$$L(1, 0.8) = -[(1) \log(0.8) + (1-1) \log(1-0.8)] = -\log(0.8)$$

- 1d.) 3 classes : class-0, class-1, class-2

$X_n$  is in class-2 ( $y_n=2$ )

$$\hat{y}_n = [0.1, 0.2, 0.7]^T$$

$$y_n = [0, 0, 1]$$

$$L([0, 0, 1], [0.1, 0.2, 0.7]^T) = 0 + 0 + [-(1) \log(0.7) + (1-1) \log(1-0.7)] = -\log(0.7)$$

$$1e.) \quad f(x) = -\ln\left(\frac{1}{1+e^{-x}}\right) = -\left[\ln(1) - \ln(1+e^{-x})\right] \quad \begin{aligned} u &= 1+e^{-x} \\ \frac{d}{du} &= -e^{-x} \end{aligned}$$

$$f'(u) = -\left[\ln(1) - \ln(u)\right] = -\left[0 - \frac{1}{u} du\right]$$

$$f'(x) = -\left[\frac{1}{1+e^{-x}} (-e^{-x})\right] = \frac{e^{-x}}{1+e^{-x}}$$

$$\left(\frac{w}{v}\right)' = \frac{w'v - v'w}{v^2}$$

$$w = e^{-x}$$

$$w' = -e^{-x}$$

$$v = 1+e^{-x}$$

$$v' = 0 - e^{-x}$$

$$f''(x) = \frac{-e^{-x}(1+e^{-x}) - (-e^{-x})(e^{-x})}{(1+e^{-x})^2} = -\frac{e^{-x}}{(1+e^{-x})^2}$$

$$2a.) \quad \text{MSE loss} = \frac{1}{N} \sum_{n=1}^N (y_n - \hat{y}_n)^2$$

$$2b.) \quad \text{MAE loss} = \frac{1}{N} \sum_{n=1}^N |y_n - \hat{y}_n|$$

$$2c.) \quad \text{MAPE loss} = \frac{1}{N} \sum_{n=1}^N \left| \frac{y_n - \hat{y}_n}{y_n} \right| * 100\%$$

3a.) The output of a decision tree for regression looks like a staircase because the tree partitions the input space into regions based on the feature values. Each split creates a new threshold for the feature, and the output remains constant within each region.

3b.) Building a deep tree such that every leaf-node of the tree is a pure node might not always be a good strategy as it can lead to overfitting to the outliers affecting the rest of the data.

3c.) total # of training samples: 60

3d.) max-depth: 2

3e.) # of 'pure' nodes: 2 (Node-2 and Node-4)

entropy of Node-0:  $H(p) = -\sum_{n=1}^N P_n \log_2(P_n)$

Samples = 60

[10, 20, 30]

$$P_1 = \frac{10}{60} = \frac{1}{6}$$

$$P_2 = \frac{20}{60} = \frac{1}{3}$$

$$P_3 = \frac{30}{60} = \frac{1}{2}$$

$$H = -\left[\frac{1}{6}(\log_2(\frac{1}{6})) + \frac{1}{3}(\log_2(\frac{1}{3})) + \frac{1}{2}(\log_2(\frac{1}{2}))\right] \approx 1.46$$

3f.) leaf / terminal node: 3

4a.) if the outputs are i.i.d random variables or weakly correlated. High variance low bias

4b.) Bagging

5.) The goal of Boosting is to reduce Bias whereas the goal of bagging is to reduce variance. Bagging reduces the variance by creating more training data sets; Boosting decreases bias because sample data sets are weighted on their performance.

6a.) Bagging trains multiple models on bootstrap replicates of the training set and then averages the outputs from the simple models. Stacking trains various models using the outputs from many other models to make prediction

6b.) Stacking many polynomial models of the same degree might lead to overfitting of data since they'd likely capture similar patterns and make similar predictions.

6C.) Yes, stacking models of different types/structures can be useful as they're more likely to capture different aspects of the data. This is important that the base models are diverse enough and don't correlate strongly with each other as it can lead to overfitting.

7a-Overfitting.) classification task #1 where the training accuracy is 100% and the testing accuracy is 50% is likely to be overfitting

7b-underfitting.) classification task #2 where both the training and testing accuracy are 50% is likely to be underfitting

8a.) hyper-parameters determine the structure and complexity of the model. It's value is chosen before a learning algorithm is used.

Examples: for a polynomial model, the degree.

# of epochs

learning rate

# of branches in decision tree

# of clusters

8b.) A validation set is needed to evaluate the performance of a model and tune the hyper-parameters.

8C.) This can lead to overfitting as pre-selecting the optimal hyper-parameters on the training set the model can get use to the training set and perform poorly on new data



8d.) Optimizing hyper-parameters using the testing set can also lead to overfitting. The test set should only be used to evaluate the final performance of the model, after hyperparameters have been selected using a separate validation set to avoid errors in new data.

9a.) Maximizing the margin in the input space will improve Classifier robustness against noises because it helps to separate the data points of different classes with a larger distance, reducing any errors caused by noise.

9b.) Yes, the margin in the input space will be maximized by a nonlinear SVM since it uses Kernel functions to map the input space to a higher-dimensional feature space, which can better separate data points of different classes.

9c.) SVM can potentially cause an "out-of-memory" error because it needs to compute the Kernel matrix which is memory inefficient for a large data set.

9d.) The purpose of using a kernel function is to map the original non-linear observations into a higher-dimensional space where they become separable.

10a.) We use weighted accuracy as compared to standard accuracy because the # of instances in the minority class can result in an error using standard accuracy. Weighted accuracy takes into account the class distribution giving more weight to that smaller class. This makes sure that the classifier is performing well on both classes instead of only the bigger class.

standard accuracy can mislead results by predicting the majority class for all instances

10b.) Re-sampling,  
- up-sampling  
- down sampling  
sampling with replacement  
sampling without replacement  
generate synthetic data samples

11.) PMF for a discrete variable  $X = [P_1, P_2, P_3, \dots, P_K]$   
and  $\sum_K P_K = 1$

$$H(X) = -\sum_K P_K \log_2(P_K)$$

a.) non-negative

Prove  $H(X) \geq 0$

when  $P_K = 0$  then  $H(X) = 0$

since  $\sum_K P_K = 1$  we can use Jensen's inequality

$$f(\sum_K t_K x_K) \leq \sum_K t_K f(x_K)$$

$t_K =$  non negative weights  
such that  $\sum_K t_K = 1$

$$H(X) = -\sum_K P_K \log(P_K)$$

$$\leq -\log(\sum_K P_K^2)$$

$$\leq -\log(\sum_K P_K)$$

$$= -\log(1) = 0$$

b.) reaches maximum when the PMF is a uniform distribution, i.e.  $P_K = 1/K$

$$P_K = 1/K \text{ for all } K$$

$$H(X) = \sum_K (1/K) \log(1/K)$$

$$= -(1/K) \sum_K \log(1/K)$$

$$= -(1/K) \sum_K (-\log(K))$$

$$= \log(K)$$

since  $\log(K)$  is a constant value that doesn't depend on the specific values of  $P_K$  it's maximized when  $P_K = 1/K$  for all  $K$ . This means entropy reaches the maximum when the PMF is a uniform distribution