(1) (1.0 pt.) Assume that there is a labeling function $f : \mathcal{X} \to \mathcal{Y}$, where probability distribution over \mathcal{X} is denoted by \mathcal{D} . The Error of a classifier $h : \mathcal{X} \to \mathcal{Y}$ can be defined as:

 $L_{\mathcal{D},f}(h) \stackrel{\text{def}}{=} \mathbb{P}_{x \sim ?}[h(x) \neq f(x)]$, where in the learner, x is assumed to be a randomly chosen example from: 1- known distribution \mathcal{D}

2- unknown distribution ${\mathcal D}$

3- unknown distribution $\mathcal{D}' \neq \mathcal{D}$

(2) (1.0 pt.) What is the probability of having more than 10 noisy examples, in a sample of 35 drawn uniformly from a distribution with 20% inherent noise.

Hint 1: you need to first compute probability of having 10 or less noisy examples in a sample of size 35 Hint 2: you will need to sum up Binomial trials, where each drawn example can be noisy or not Hint 3: you can compute it by hand, but it could be faster computationally

1-0.07

2-0.02

 $3. \ 0.10$

(3) (1.0 pt.) In each of the following problems which measure should be prioritized for evaluating a classifier.A) Cancer prediction (positive: cancer, negative: healthy). The patients with positive prediction will go through more analysis, and the negative ones will be sent home. Hint: sending a cancer patient home, can be dangerous.

B) Spam email prediction (positive:spam, negative: non-spam). The spam detected emails would be automatically removed. Hint: we do not want to miss any important email.

$$precision = \frac{TP}{TP+FP}$$
$$recall = \frac{TP}{TP+FN}$$
$$1- A: precision, B: recall\\2- A: recall, B: precision\\3- A: precision, B: precision$$

(4) (1.0 pt.) Import the given files X.csv and Y.csv in the Materials section, as the inputs and targets. The provided files can also be imported from Boston dataset available in sklearn, where the features LSTAT and RM are used as inputs and target is used as output. Split the dataset to training and test set using the following Python code:

from sklearn.model_selection import train_test_split

 $X_{train}, X_{test}, Y_{train}, Y_{test} = train_{test} \\ split(X, Y, test_{size} = 0.3, random_{state} = 1)$

function description

using Linear Regression (LR) without considering bias (intercept), which option is the root-mean-square error (RMSE) over test data:

- 1-5.26
- 2-6.34
- 3-4.62

(5) (1.0 pt.) In the previous question, again using LR without bias, compute average (K-folds cross-validation error) and variance of the root-mean-square error (RMSE) on test folds, for K = 2 and K = 5. To generate the folds, use the following Python code:

from sklearn.model_selection import KFold
KFold(n_splits=k, random_state=1, shuffle=True)

function description

- 1- K=2 : average = 6.23, variance = 0.21 / k=5: average = 6.12 , variance = 0.64
- 2- K=2 : average = 5.38 , variance = 0.43 / k=5: average = 5.53, variance = 0.38
- 3- K=2 : average = 6.03 , variance = 0.37 / k=5: average = 5.72 , variance = 0.50