# CS 461 Homework 1

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Due: Sept. 27, 11:59 pm

# 1 Problem 1: Right Choice for Empirical Metric

### 1.1 1.1 Precision Rate

Precision Rate measures the rate of positive values that are in the classification circle. The formula is

$$\frac{\text{\# of observed positive examples}}{\text{\# of samples in the circle}}$$

The classification circle trained on this data would be a small circle surrounding 1 positive example. This would ensure a precision rate of 100%.

### 1.2 Recall Rate

Recall Rate measures the rate of observed positive values by the total number of positive values in the dataset. The formula is

$$\frac{\# \text{ of observed positive examples}}{\# \text{ of total positive examples}}$$

The classification circle trained on this data would be huge. In fact, it would classify the entire dataset to make sure that no positive values are left behind.

### 1.3 Issues and Suggested Metric

The main issue with using precision rate as the only metric for classification is that it's very easy to miss a lot of data since the classification circle would be so small to ensure a 100% rate. The opposite is true of recall rate. Since it would be focused on maximizing the total amount of positive examples, the circle would be too big and it would take in 5 negative examples to take 1 positive one. The precision rate for such a classification would be horrible.

The solution is to strike a balance between the two. You could use the F1 score, which is just the harmonic mean between precision rate and recall rate.

$$F1 = 2 \cdot \frac{\text{precision} \cdot \text{recall}}{\text{precision} + \text{recall}}$$

# 1.4 Question 2.1

Since the ball only has the choice between moving left and right at each level, we can define  $\Omega_M = \{L, R\}$ . The choice is the same from  $\Omega_2 \dots \Omega_M$ .

#### 1.5 Question 2.2

 $\Omega$  is the cartesian product of all of the subspaces  $\Omega_2 \dots \Omega_M$ . We can use the following set builder notation to define:

$$\Omega = \{(x_1, x_2, \dots, x_M) \mid x_i \in \{L, R\}, i \in \mathbb{Z}, 1 \le i \le M\}$$

### 1.6 Question 2.3

The meaning of  $L_g$ , where the ball arrives, is simply the number of right turns the ball took.  $L_g$  defines the place or bin where the ball ends up, which is dictated by the number of right turns.

### 1.7 Question 2.4

We can define  $L_g$  as the random variable X. Since there are only two possibilities with an equal probability p = 0.5, we can use the following binomial distribution to describe the distribution:

$$P(X=k) = \binom{M}{k} \left(\frac{1}{2}\right)^{M}$$

### 1.8 Question 2.5

Define the PMF of your random variable for depths  $M=5,\,M=10,\,$  and M=100. Plot them and explain the phenomenon in relation to the Central Limit Theorem.

### 1.9 Central Limit Theorem (CLT) Explanation

The Central Limit Theorem (CLT) states that the distribution of the sum (or average) of independent random variables, each with finite mean and variance, approaches a normal distribution as the number of variables increases. In the context of the Galton board, as the number of levels M increases, the binomial distribution for the number of successes k begins to approximate a normal distribution. We can see that visually in the following figure:

### 1.10 Question 3.1

What is the probability that your plant will survive the week? We want to calculate P(D). Given:

$$P(D|F) = 0.8$$
,  $P(D|F') = 0.2$ ,  $P(F) = 0.3$ 

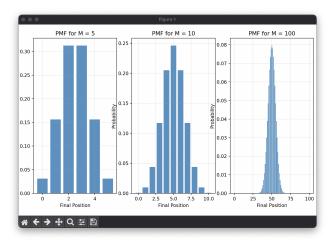


Figure 1: Illustration of the Central Limit Theorem (CLT).

We can find P(F') = 1 - P(F) = 0.7.

Now, using marginalization:

$$P(D) = P(D, F) + P(D, F') = (P(D|F) \cdot P(F)) + (P(D|F') \cdot P(F'))$$
  
$$P(D) = (0.8 \cdot 0.3) + (0.2 \cdot 0.7) = 0.38$$

Thus, the probability that the plant survives is:

$$P(D') = 1 - P(D) = 0.62$$

### 1.11 Question 3.2

If your friend forgot to water it, what is the chance that it will be dead when you return?

This is simply:

$$P(D|F) = 0.8$$

# 1.12 Question 3.3

If it is dead when you return, what is the chance that your friend forgot to water it?

Using Bayes' theorem:

$$P(F|D) = \frac{P(D|F) \cdot P(F)}{P(D)} = \frac{0.8 \cdot 0.3}{0.38} = 0.6316$$

# 2 Problem 4: Naive Bayes

# 2.1 4.1 Formula Rewrite for Naive Bayes

$$P(D=+|G=g,B=b) = \frac{P(G=g|D=+)\cdot P(B=b|D=+)\cdot P(D=+)}{(P(G=g|D=+)\cdot P(B=b|D=+)\cdot P(D=+)) + (P(G=g|D=-)\cdot P(B=b|D=-)\cdot P(D=-))}$$
 
$$P(D=-|G=g,B=b) = \frac{P(G=g|D=-)\cdot P(B=b|D=-)\cdot P(D=-)}{(P(G=g|D=+)\cdot P(B=b|D=+)\cdot P(D=+)) + (P(G=g|D=-)\cdot P(B=b|D=-)\cdot P(D=-))}$$

### 2.2 4.2 Estimation Based on Train Data

Look at mannan-nb-train.py

### 2.3 4.3 Classifier Code

Look at mannan-nb-cls.py

## 2.4 4.4 Evaluation of Classifier

Look at mannan-nb-test.py

### 2.5 4.5 Standardization Necessity

Standardization is not really helpful for Naive Bayes classification. For Gaussian Naive Bayes, we are just looking at the mean and variance. These metrics aren't affected by scaling or standardization, so it isn't needed for this classifier.

### 2.6 4.6 Data Reflection and Improvements

The dataset may not fully capture real-world diabetes diagnoses since it only includes glucose and blood pressure, while other factors like age or BMI are also important. There could also be biases if the data comes from a limited population. Without new data, feature engineering, such as creating a glucose-to-blood-pressure ratio, and using cross-validation could improve the model's accuracy and generalizability.

# Problem 5: Data Whitening

# **5.1** Expressing E[Y] and COV[Y, Y]

Let  $Y = A \cdot X + b$ , where A is a transformation matrix and b is a bias vector. The expected value of Y is:

$$E[Y] = A \cdot E[X] + b$$

Given 
$$E[X] = \begin{bmatrix} 0.2 \\ 0.3 \\ 0.1 \end{bmatrix}$$
, we have:

$$E[Y] = A \cdot \begin{bmatrix} 0.2 \\ 0.3 \\ 0.1 \end{bmatrix} + b$$

The covariance matrix of Y is:

$$COV[Y, Y] = A \cdot COV[X, X] \cdot A^T$$

Given the covariance matrix  $COV[X,X] = \begin{bmatrix} 2.75 & 0.43 & 0 \\ 0.43 & 2.25 & 0 \\ 0 & 0 & 1 \end{bmatrix}$ , the covariance

of Y becomes:

$$COV[Y,Y] = A \cdot \begin{bmatrix} 2.75 & 0.43 & 0 \\ 0.43 & 2.25 & 0 \\ 0 & 0 & 1 \end{bmatrix} \cdot A^{T}$$

# 5.2 Design A and b for Whitening

To whiten the data, we need to set E[Y] = 0 and COV[Y, Y] = I.

1. Setting E[Y] = 0:

The expected value of Y is:

$$E[Y] = A \cdot E[X] + b$$

To achieve E[Y] = 0, we set:

$$b = -A \cdot E[X]$$

Thus, the calculated b is:

$$b = \begin{bmatrix} -0.1052 \\ -0.1912 \\ -0.1 \end{bmatrix}$$

2. Setting COV[Y, Y] = I:

The covariance of Y is:

$$COV[Y,Y] = A \cdot COV[X,X] \cdot A^T$$

We want COV[Y, Y] = I, which leads to the equation:

$$A \cdot COV[X, X] \cdot A^T = I$$

Therefore, the transformation matrix A is the inverse square root of the covariance matrix COV[X,X]. The calculated A is:

$$A = \begin{bmatrix} 0.6097 & -0.0558 & 0 \\ -0.0558 & 0.6746 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

#### 3. Verification:

After applying the whitening transformation, we can verify that the transformed data has a mean close to zero and a covariance matrix close to the identity matrix. The results from the Python code are:

Mean of 
$$Y = \begin{bmatrix} 0.0023 \\ -0.0053 \\ 0.0225 \end{bmatrix}$$

$$\begin{bmatrix} 1.0053 & -0.0061 & -0.0061 \\ -0.0061 & -0.0061 \end{bmatrix}$$

Covariance of 
$$Y = \begin{bmatrix} 1.0053 & -0.0061 & -0.0088 \\ -0.0061 & 0.9814 & 0.0315 \\ -0.0088 & 0.0315 & 0.9940 \end{bmatrix}$$

As you can see, the covariance matrix is very close to the identity matrix, indicating that the whitening process was successful.

# 6.1 Maximum Likelihood Estimate (MLE)

Given that the observations  $x_1, x_2, \ldots, x_n$  are i.i.d. and follow a Gaussian distribution, the probability density function for each observation  $x_i$  is:

$$f(x_i|\mu,\sigma^2) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(x_i-\mu)^2}{2\sigma^2}\right)$$

The likelihood function for all observations is the product of the individual probabilities, and the log-likelihood is:

$$\log L(\mu, \sigma^2) = -\frac{n}{2}\log(2\pi\sigma^2) - \frac{1}{2\sigma^2}\sum_{i=1}^{n}(x_i - \mu)^2$$

To find the MLE for  $\mu$ , we differentiate the log-likelihood with respect to  $\mu$  and set it to zero:

$$\frac{d}{d\mu} \log L(\mu, \sigma^2) = \frac{1}{\sigma^2} \sum_{i=1}^{n} (x_i - \mu) = 0$$

Solving for  $\mu$ , we obtain the MLE:

$$\mu^* = \frac{1}{n} \sum_{i=1}^n x_i$$

### 6.2 Maximum A Posteriori (MAP) Estimate

Let  $X_1, \ldots, X_N$  be i.i.d. random variables with a PDF:

$$f_{X_n|\mu}(x_n|\mu) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(x_n - \mu)^2}{2\sigma^2}\right)$$

and let  $\mu$  have a prior distribution:

$$f_{\mu}(\mu) = \frac{1}{\sqrt{2\pi\sigma_0^2}} \exp\left(-\frac{(\mu - \mu_0)^2}{2\sigma_0^2}\right)$$

The MAP estimate is:

$$\hat{\mu}_{\text{MAP}} = \arg \max_{\mu} p(\mu|\mathbf{x}) \propto p(\mathbf{x}|\mu) \cdot p(\mu)$$

Taking the logarithm of the likelihood and prior:

$$\hat{\mu}_{\text{MAP}} = \arg\max_{\mu} \left\{ -\frac{1}{2\sigma^2} \sum_{n=1}^{N} (x_n - \mu)^2 - \frac{(\mu - \mu_0)^2}{2\sigma_0^2} \right\}$$

Differentiating with respect to  $\mu$ :

$$\sum_{n=1}^{N} \frac{(x_n - \mu)}{\sigma^2} - \frac{(\mu - \mu_0)}{\sigma_0^2} = 0$$

Solving for  $\mu$ , we get:

$$\mu\left(\frac{N}{\sigma^2} + \frac{1}{\sigma_0^2}\right) = \frac{\sum_{n=1}^{N} x_n}{\sigma^2} + \frac{\mu_0}{\sigma_0^2}$$

The final MAP estimate is:

$$\hat{\mu}_{\text{MAP}} = \frac{\frac{1}{N} \sum_{n=1}^{N} x_n \cdot \sigma_0^2 + \mu_0 \cdot \frac{\sigma^2}{N}}{\sigma_0^2 + \frac{\sigma^2}{N}}$$

# **6.3** Behavior of MAP for $N \to \infty$ and $N \to 0$

The MAP estimate is:

$$\mu_{\mathrm{MAP}} = \frac{\sigma^2}{N\sigma_0^2 + \sigma^2} \mu_0 + \frac{N\sigma_0^2}{N\sigma_0^2 + \sigma^2} \mu_{\mathrm{ML}}$$

- As  $N \to \infty$ : The prior  $\mu_0$  loses influence, so  $\mu_{\text{MAP}} \to \mu_{\text{ML}}$ .

$$\lim_{N \to \infty} \mu_{\text{MAP}} = \mu_{\text{ML}}$$

- As  $N \to 0$ : The data loses influence, so  $\mu_{\text{MAP}} \to \mu_0$ .

$$\lim_{N\to 0} \mu_{\text{MAP}} = \mu_0$$

In short, use MLE when N is large and MAP when N is small.