Bayesian Logistic Classification

3F8: Inference Coursework
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Abstract

Abstract goes here

1 Introduction

2 Model definition

2.1 Logistic classification

The logistic classifier is a binary classifier that takes a D-dimensional input $\boldsymbol{x}_n$ and outputs a class label $y_n = \{0, 1\}$, where the class labels are modelled as being independent and identically generated from a Bernoulli distribution:

$$\begin{aligned} p(y_n = 1|\tilde{\boldsymbol{x}}_n) &= \sigma(\boldsymbol{w}^T\tilde{\boldsymbol{x}}_n) \\ p(y_n = 0|\tilde{\boldsymbol{x}}_n) &= 1 - \sigma(\boldsymbol{w}^T\tilde{\boldsymbol{x}}_n) = \sigma(-\boldsymbol{w}^T\tilde{\boldsymbol{x}}_n) \end{aligned} \tag{1}$$

with $\tilde{\boldsymbol{x}}_n = \left[1, \boldsymbol{x}_n^T\right]^T$, $\boldsymbol{w}$ as a vector of D+1 model weights, and $\sigma(x) = \frac{1}{1+e^{-x}}$ (the logistic function).

2.2 Bayesian logistic classification

A posterior distribution of the model weights is required to perform fully Bayesian classification:

$$p(\boldsymbol{w}|\boldsymbol{X}, \boldsymbol{y}) = \frac{p(\boldsymbol{X}|\boldsymbol{w}, \boldsymbol{y})p(\boldsymbol{w})}{p(\boldsymbol{X}|\boldsymbol{y})} \tag{2}$$

From the definition of the logistic classifier, the log-likelihood of $\mathbf{w}$ is:

$$\begin{aligned} \mathcal{L}(\boldsymbol{w}) = \log p(y|\tilde{\boldsymbol{X}}, \boldsymbol{w}) &= \log \prod_{n=1}^{N} \sigma(\boldsymbol{w}^T\tilde{\boldsymbol{x}}_n)^{y_n}\sigma(-\boldsymbol{w}^T\tilde{\boldsymbol{x}}_n)^{1-y_n} \\ &= \sum_{n=1}^{N} y_n \log\sigma(\boldsymbol{w}^T\tilde{\boldsymbol{x}}_n) + (1 - y_n)\log\sigma(-w^T\tilde{\boldsymbol{x}}_n) \end{aligned} \tag{3}$$

The prior distribution of the model weights is chosen to be Gaussian, $p(\boldsymbol{w}) = \mathcal{N}(\boldsymbol{w}; \boldsymbol{m}_0, \boldsymbol{S}_0)$. This gives the log-prior as:

$$\mathcal{P}(\boldsymbol{w}) = \log p(\boldsymbol{w}) = -\frac{1}{2\sigma_0^2}(\boldsymbol{w} - \boldsymbol{m}_0)^T\boldsymbol{S}_0^{-1}(\boldsymbol{w} - \boldsymbol{m}_0) + \text{const} \tag{4}$$

The log-posterior is:

$$\log p(\boldsymbol{w}|\boldsymbol{X}, \boldsymbol{y}) = \mathcal{L}(\boldsymbol{w}) + \mathcal{P}(\boldsymbol{w}) + \text{const} \tag{5}$$

Calculating the model evidence $p(\boldsymbol{X}|\boldsymbol{y})$ exactly requires integrating the product of the likelihood and the prior, which is intractable. We present two methods to circumvent this problem: MAP estimation and the Laplace approximation.

2.2.1 'Semi-Bayesian' classification: MAP estimation

One method of performing Bayes-driven logistic classification is to use the MAP estimate of the model weights, $\boldsymbol{w}_{\text{MAP}}$, which can be found by applying gradient ascent to the log-posterior. This avoids needing the model evidence, but discards a lot of the information contained in the posterior and does not involve calculating the distribution of the model weights. As such, it is not 'true' Bayesian classification. The gradient of the log-posterior is calculated as follows:

$$\begin{aligned}\frac{\partial}{\partial \boldsymbol{w}} \log p(\boldsymbol{w}|\tilde{\boldsymbol{X}}, \boldsymbol{y}) &= \frac{\partial}{\partial \boldsymbol{w}}\mathcal{L}(\boldsymbol{w}) + \frac{\partial}{\partial \boldsymbol{w}}\mathcal{P}(\boldsymbol{w}) \\ &= \sum_{n=1}^{N} \left(y_n - \sigma(\boldsymbol{w}^T \tilde{\boldsymbol{x}}_n)\right) \tilde{\boldsymbol{x}}_n + \frac{1}{\sigma_0^2}\boldsymbol{w} \\ &= \tilde{\boldsymbol{X}}(\mathbf{y} - \sigma(\tilde{\boldsymbol{X}}^T \boldsymbol{w})) + \frac{1}{\sigma_0^2}\boldsymbol{w} \end{aligned} \tag{6}$$

where $\tilde{\boldsymbol{X}} = [\tilde{\boldsymbol{x}}_1 \dots \tilde{\boldsymbol{x}}_N]$. A standard gradient-based solver can be used with Equation 6 to find a value for $\boldsymbol{w}_{\text{MAP}}$, which can be used as a setting for the weights in Equation 1 to classify data points. The performance of the MAP-based classifier will be compared to a fully Bayesian classifier later in the report (**??**).

2.2.2 'True' Bayesian classification: the Laplace approximation

The Laplace approximation allows us to approximate the model evidence by finding a Gaussian distribution $q(\boldsymbol{w})$ that closely models the posterior $p(\boldsymbol{w}|\tilde{\boldsymbol{X}}, \boldsymbol{y})$ around a local maximum. The normalising constant of a Gaussian is well defined, so we can use this as the approximation of the evidence. For brevity, rewrite Equation 2 as:

$$p(\boldsymbol{w}|\boldsymbol{X}, \boldsymbol{y}) = \frac{f(\boldsymbol{w})}{K} \approx q(\boldsymbol{w}) \tag{7}$$

where $f(\boldsymbol{w}) = p(\boldsymbol{X}|\boldsymbol{w}, \boldsymbol{y})p(\boldsymbol{w})$ and $K = p(\boldsymbol{X}|\boldsymbol{y}) = \int f(\boldsymbol{w})d\boldsymbol{w}$. To find $q(\boldsymbol{w})$, we start with the truncated Taylor expansion of $\log f(\boldsymbol{w})$ around a local maximum $\boldsymbol{w}_0$:

$$\log f(\boldsymbol{w}) \approx \log f(\boldsymbol{w}_0) + \nabla \log f(\boldsymbol{w})\big|_{\boldsymbol{w}=\boldsymbol{w}_0}(\boldsymbol{w} - \boldsymbol{w}_0) + \frac{1}{2}(\boldsymbol{w} - \boldsymbol{w}_0)^T \nabla^2 \log f(\boldsymbol{w})\big|_{\boldsymbol{w}=\boldsymbol{w}_0}(\boldsymbol{w} - \boldsymbol{w}_0)$$

At a maximum of $f(\boldsymbol{w})$, $\nabla \log f(\boldsymbol{w}) = 0$ as the logarithm is a monotonic function. Hence, close to $\boldsymbol{w}_0$:

$$\begin{aligned}\log f(\boldsymbol{w}) &\approx \log f(\boldsymbol{w}_0) + \frac{1}{2}(\boldsymbol{w} - \boldsymbol{w}_0)^T \nabla^2 \log f(\boldsymbol{w})\big|_{\boldsymbol{w}=\boldsymbol{w}_0}(\boldsymbol{w} - \boldsymbol{w}_0) \\ f(\boldsymbol{w}) &\approx f(\boldsymbol{w}_0) \exp\left(\frac{1}{2}(\boldsymbol{w} - \boldsymbol{w}_0)^T \nabla^2 \log f(\boldsymbol{w})\big|_{\boldsymbol{w}=\boldsymbol{w}_0}(\boldsymbol{w} - \boldsymbol{w}_0)\right) \end{aligned} \tag{8}$$

Equation 8 is of the form of an un-normalised Gaussian centred on the maximum at $\boldsymbol{w}_0$. Set $\boldsymbol{w}_0 = \boldsymbol{w}_{\text{MAP}}$, and let $\boldsymbol{S}_N = -\nabla^2 \log f(\boldsymbol{w})\big|_{\boldsymbol{w}=\boldsymbol{w}_{\text{MAP}}}$, then normalise Equation 8 to obtain the result for $q(\boldsymbol{w})$:

$$\begin{aligned}p(\boldsymbol{w}|\tilde{\boldsymbol{X}}, \boldsymbol{y}) &\approx \frac{1}{(2\pi)^{\frac{N}{2}} \det \boldsymbol{S}_N^{-\frac{1}{2}}} \exp\left(\frac{1}{2}(\boldsymbol{w} - \ \boldsymbol{w}_{\text{MAP}})^T \boldsymbol{S}_N (\boldsymbol{w} - \boldsymbol{w}_{\text{MAP}})\right) \\ &= \mathcal{N}(\boldsymbol{w}; \boldsymbol{w}_{\text{MAP}}, \boldsymbol{S}_N^{-1}) \end{aligned} \tag{9}$$

where N is the number of data points $\boldsymbol{x}_n$. For Equation 9 to hold, $\boldsymbol{S}_N$ must be positive definite, which is equivalent to saying $\boldsymbol{w}_{\text{MAP}}$ must be a maximum (which is true by definition).

Now we can obtain an approximate value of the normalising constant K, by substituting Equation 9 into the definition of K:

$$\begin{aligned}K = \int f(\boldsymbol{w})d\boldsymbol{w} &\approx f(\boldsymbol{w}_0) \int \exp\left(\frac{1}{2}(\boldsymbol{w} - \boldsymbol{w}_0)^T \boldsymbol{S}(\boldsymbol{w} - \boldsymbol{w}_0)\right) d\boldsymbol{w} \\ &= \sqrt{\frac{(2\pi)^N}{\det \boldsymbol{S}_N}} f(\boldsymbol{w}_{\text{MAP}}) \end{aligned} \tag{10}$$

Using Equation 9, the covariance matrix of $q(\boldsymbol{w})$ can be found:

$$\begin{aligned}\boldsymbol{S}_N^{-1} &= -\nabla^2 \log p(\boldsymbol{w}|\boldsymbol{X}, \boldsymbol{y})\big|_{\boldsymbol{w}=\boldsymbol{w}_{\text{MAP}}} \\ &= \boldsymbol{S}_0^{-1} + \sum_{n=1}^{N} \sigma(\boldsymbol{w}^T \tilde{\boldsymbol{x}}_n)\sigma(-\boldsymbol{w}^T \tilde{\boldsymbol{x}}_n)\tilde{\boldsymbol{x}}_n \tilde{\boldsymbol{x}}_n^T \end{aligned} \tag{11}$$

Defining $\boldsymbol{\sigma} = \sigma(\tilde{\boldsymbol{X}}^T \boldsymbol{w})$, this can be written in vector form as:

$$\boldsymbol{S}_N^{-1} = \boldsymbol{S}_0^{-1} + \tilde{\boldsymbol{X}} \tilde{\boldsymbol{X}}^T \boldsymbol{\sigma}(1 - \boldsymbol{\sigma}) \tag{12}$$

Hence, using Equation 9, the posterior distribution is:

$$p(\boldsymbol{w}|\boldsymbol{X}, \boldsymbol{y}) \approx \mathcal{N}(\boldsymbol{w}; \boldsymbol{w}_{\text{MAP}}, \boldsymbol{S}_N) \tag{13}$$

And using Equation 10, the normalizing constant is:

$$p(\boldsymbol{X}|\boldsymbol{y}) = \frac{(2\pi)^{\frac{N}{2}}}{\det \boldsymbol{S}_N^{-\frac{1}{2}}} p(\boldsymbol{X}|\boldsymbol{w} = \boldsymbol{w}_{\text{MAP}}, \boldsymbol{y}) p(\boldsymbol{w} = \boldsymbol{w}_{\text{MAP}}) \tag{14}$$

2.3 Predictive distribution

The predictive distribution can also be approximated using the Laplace method:

$$\begin{aligned} p(y^* = 1|\boldsymbol{x}^*, \boldsymbol{y}, \boldsymbol{X}) &= \int p(y^* = 1|\boldsymbol{x}^*, \boldsymbol{w}) p(\boldsymbol{w}|\boldsymbol{y}, \boldsymbol{X}) d\boldsymbol{w} \\ &\approx \int \sigma(\boldsymbol{w}^T \boldsymbol{x}^*) q(\boldsymbol{w}) d\boldsymbol{w} \end{aligned} \tag{15}$$

Using the sifting property of the delta function:

$$\sigma(\boldsymbol{w}^T \boldsymbol{x}) = \int \delta(a - \boldsymbol{w}^T \boldsymbol{x}) \sigma(a) da \tag{16}$$

Hence:

$$\begin{aligned} p(y^* = 1|\boldsymbol{x}^*, \boldsymbol{y}, \boldsymbol{X}) &\approx \int \int \delta(a - \boldsymbol{w}^T \boldsymbol{x}^*) \sigma(a) q(\boldsymbol{w}) d\boldsymbol{w} da \\ &= \int \sigma(a) \int \delta(a - \boldsymbol{x}^{*T} \boldsymbol{w}) q(\boldsymbol{w}) d\boldsymbol{w} da \end{aligned}$$

The inner integral applies a linear constraint to $q(\boldsymbol{w})$, as the argument of the delta function is 0 unless $a = \boldsymbol{x}^{*T} \boldsymbol{w}$. Hence, the approximate predictive distribution is:

$$\begin{aligned} p(y^* = 1|\boldsymbol{x}^*, \boldsymbol{y}, \boldsymbol{X}) &\approx \int \sigma(a) \mathcal{N}(a; \boldsymbol{x}^{*T} \boldsymbol{w}_{\text{MAP}}, \boldsymbol{x}^{*T} \boldsymbol{S}_N \boldsymbol{x}^*) da \\ &= \int \sigma(a) \mathcal{N}(a; \mu_p, \sigma_p^2) da \end{aligned} \tag{17}$$

Where

$$\mu_p = \boldsymbol{x}^{*T} \boldsymbol{w}_{\text{MAP}} \tag{18}$$

$$\sigma_p^2 = \boldsymbol{x}^{*T} \boldsymbol{S}_N \boldsymbol{x}^* \tag{19}$$

This integral cannot be expressed analytically, so another approximation is required. The logistic function can be approximated well by a probit function scaled such that the gradient of the two functions at the origin are equal. It can be shown that this gives:

$$\sigma(x) \approx \Phi^{-1}\left(\sqrt{\frac{\pi}{8}} x\right) = \Phi^{-1}(\lambda x) \tag{20}$$

Substituting into Equation 17 and evaluating using properties of the probit function gives:

$$p(y^* = 1|\boldsymbol{x}^*, \boldsymbol{y}, \boldsymbol{X}) \approx \int \Phi^{-1}(\lambda x)\mathcal{N}(a; \mu_p, \sigma_p)da$$
$$= \Phi^{-1}\left(\frac{\mu_p}{\sqrt{\lambda^{-2} + \sigma_p^2}}\right) \tag{21}$$

Using Equation 20, this can be converted back into a logistic function:

$$p(y^* = 1|\boldsymbol{x}^*, \boldsymbol{y}, \boldsymbol{X}) \approx \sigma\left(\frac{\mu_p}{\sqrt{1 + \sigma_p^2\lambda^2}}\right) \tag{22}$$

Summary of the Laplace approximation:

- Posterior distribution of $\boldsymbol{w}$:

$$p(\boldsymbol{w}|\boldsymbol{X}, \boldsymbol{y}) \approx \mathcal{N}(\boldsymbol{w}; \boldsymbol{w}_{\text{MAP}}, \boldsymbol{S}_N) \tag{23}$$

- Predictive distribution for new points $\boldsymbol{x}_*$:

$$p(y_* = 1|\boldsymbol{x}_*, \boldsymbol{y}, \boldsymbol{X}) \approx \sigma\left(\frac{\mu_p}{\sqrt{1 + \sigma_p^2\lambda^2}}\right) \tag{24}$$
$$p(y_* = 0|\boldsymbol{x}_*, \boldsymbol{y}, \boldsymbol{X}) = 1 - p(y_* = 1|\boldsymbol{x}_*, \boldsymbol{y}, \boldsymbol{X})$$

- Model evidence:

$$p(\boldsymbol{X}|\boldsymbol{y}) \approx \sqrt{\frac{(2\pi)^N}{\det \boldsymbol{S}_N^{-1}}} \exp\left(\mathcal{L}(\boldsymbol{w}_{\text{MAP}}) + \mathcal{S}(\boldsymbol{w}_{\text{MAP}})\right) \tag{25}$$

$$\boldsymbol{S}_N = \left(\frac{1}{\sigma_0^2}\boldsymbol{I} + \tilde{\boldsymbol{X}}\tilde{\boldsymbol{X}}^T\boldsymbol{\sigma}(1 - \boldsymbol{\sigma})\right)^{-1} \tag{26}$$
$$\boldsymbol{\sigma} = \sigma(\tilde{\boldsymbol{X}}^T\boldsymbol{w}) \tag{27}$$
$$\mu_p = \tilde{\boldsymbol{x}}_*^T\boldsymbol{w}_{\text{MAP}} \tag{28}$$
$$\sigma_p^2 = \tilde{\boldsymbol{x}}_*^T\boldsymbol{S}_N\tilde{\boldsymbol{x}}_* \tag{29}$$
$$\lambda^2 = \frac{\pi}{8} \tag{30}$$
$$\mathcal{L}(\boldsymbol{w}_{\text{MAP}}) = \log p(\boldsymbol{X}|\boldsymbol{w}, \boldsymbol{y})\big|_{\boldsymbol{w}=\boldsymbol{w}_{\text{MAP}}} \tag{31}$$
$$\mathcal{P}(\boldsymbol{w}_{\text{MAP}}) = \log p(\boldsymbol{w})\big|_{\boldsymbol{w}=\boldsymbol{w}_{\text{MAP}}} \tag{32}$$
$$N = \text{number of data points } \boldsymbol{x}_n \tag{33}$$

2.4 Implementation in Python

Firstly, the data is loaded and split into training and test sets.

```
import numpy as np
from sklearn.model_selection import train_test_split
from sklearn.utils import shuffle

X_data, y_data = shuffle(np.loadtxt('data/X.txt'),
                         np.loadtxt('data/y.txt'))

X_train, X_test, y_train, y_test = train_test_split(X_data, y_data, train_size=800)
```

Functions are defined to expand the data through a set of radial basis functions centred on the training points, and prepend a column of ones:

```
def prepend_ones(M):
    return np.column_stack((np.ones(M.shape[0]), M))

def expand_rbf(l, X, Z=X_train):
    X2 = np.sum(X**2, 1)
    Z2 = np.sum(Z**2, 1)
    ones_Z = np.ones(Z.shape[ 0 ])
    ones_X = np.ones(X.shape[ 0 ])
    r2 = np.outer(X2, ones_Z) - 2 * np.dot(X, Z.T) + np.outer(ones_X, Z2)
    return prepend_ones(np.exp(-0.5 / l**2 * r2))
```

A numerical approximation of the MAP estimate for the weights is found:

```
from scipy.optimize import fmin_l_bfgs_b as minimise

def logistic(x):
    return 1 / (1 + np.exp(-x))

def log_prior(w, variance):
    return -1 / (2 * variance) * (w.T @ w)

def log_likelihood(w, X, y):
    sigma = logistic(X @ w)
    return np.sum(y * np.log(sigma)
                  + (1 - y) * np.log(1 - sigma))

def negative_log_posterior(w, X, y, prior_variance):
    return -(log_likelihood(w, X, y) + log_prior(w, prior_variance))

def negative_posterior_gradient(w, X, y, prior_variance):
    return -((y - logistic(X @ w)) @ X - w / prior_variance)

def find_w_map(X, y, w0=None, prior_variance=1):
    if w0 is None:
        w0 = np.random.normal(size=X.shape[1])
    w_map, posterior_at_wmap, d = minimise(negative_log_posterior,
                                           w0,
                                           negative_posterior_gradient,
                                           args=[X, y, prior_variance])
    return w_map

expanded_training_set = expand_rbf(rbf_width, X_train)
w_map = find_w_map(expanded_training_set, y_train)
```

Note that `scipy.optimize.fmin_l_bfgs_b` is a minimisation function, so to maximise the posterior we have to work with the negative posterior and negative posterior gradient. Once $\boldsymbol{w}_{\text{MAP}}$ is found, the Laplace approximation for the model evidence and the predictive distribution can be found:

```
def S_N_inv(w, rbf_width, prior_variance):
    M = w.shape[0]
    X_tilde = expand_rbf(rbf_width, X_train)
    sigma = logistic(X_tilde @ w)
    return np.identity(M) / prior_variance + X_tilde @ X_tilde.T @ sigma @ (1 - sigma)

def log_evidence(w, X, y, rbf_width, prior_variance):
    d = np.linalg.det(S_N_inv(w, rbf_width, prior_variance))
    return (M/2)*np.log(2*np.pi) - 0.5*np.log(d) + log_likelihood(w, X, y) + log_prior(w, prior_
```

```
def laplace_prediction(inputs, weights, rbf_width, prior_variance):
    X_tilde = expand_rbf(rbf_width, X_train)
    sigma = logistic(X_tilde @ weights)
    S_N = np.linalg.inv(S_N_inv(weights, rbf_width, prior_variance))
    predictive_mean = inputs @ weights
    predictive_variance = np.array([x.T @ C_N @ x for x in inputs])
    return logistic(predictive_mean / np.sqrt(1 + predictive_variance*np.pi/8))
```

3 Performance of the Laplace Approximation