

# Problem Sheet 1

#### 1 Nearest Neighbour Classification

In the lectures, we studied the perceptron, a linear classifier of the form  $y = \operatorname{sign}(\mathbf{w} \cdot \mathbf{x} + w_0)$ , where  $\operatorname{sign}(z) = 1$  if  $z \geq 0$  and  $\operatorname{sign}(z) = 0$  otherwise. The parameters to be learnt are  $\mathbf{w}$  and  $w_0$ . The "Nearest neighbour classifier" (NN) is a different approach to learning from data. Suppose we are given N points  $(\mathbf{x}_1, y_1), \ldots, (\mathbf{x}_N, y_N)$  where  $y_i \in \{0, 1\}$ ; for a parameter k and given a new point  $\mathbf{x}^*$ , the k-NN approach does the following: find  $\mathbf{x}_{j_1}, \ldots \mathbf{x}_{j_k}$  the k-closest points to  $\mathbf{x}^*$ , then output  $\widehat{y}^*$  as the majority label from the set  $\{y_{j_1}, \ldots, y_{j_k}\}$ , i.e., the most commonly occurring label among the k-nearest neighbours.

- 1. What advantage does the k-NN approach offer over a linear classifier like the perceptron?
- 2. How many parameters does the nearest neighbour model have? How much memory do you need to store the model? What is the computational cost of predicting the label  $\hat{y}^*$ ?
- 3. In this part, we'll look at the setting where the vectors  $\mathbf{x}$  are points on the boolean hypercube, *i.e.*,  $\mathbf{x} \in \{0,1\}^D$ . Fix  $\mathbf{x}^* = (0,0,\ldots,0)$  to be the origin and imagine that data consists of points drawn uniformly at random from the boolean hypercube. What is the distribution of the Hamming distance of data points from  $\mathbf{x}^*$ ? What happens as  $D \to \infty$ ? (*Hint*: Use the central limit theorem.)
- 4. Let us now fix some numbers. Suppose the dimension of the data D=10,000; let  $\mathbf{x}^*=(0,0,\ldots,0)$  and suppose we generated N=10,000 data points. What do you expect the distance of  $\mathbf{x}^*$  from the nearest data-point to be? the furthest? How large does N need to be to get points that are reasonably close to  $\mathbf{x}^*$ , say within Hamming distance 50?

**Remark**: You do not have to write precise numbers or even mathematical expressions for the answers to part 4 above. Make sure you understand the behaviour qualitatively. The phenomenon explored in the last two parts of the question is referred to as the *curse of dimensionality*.

#### 2 Normalization constant for a 1D Gaussian

The normalization constant for a zero-mean Gaussian is given by

$$Z = \int_{-\infty}^{\infty} \exp\left(-\frac{x^2}{2\sigma^2}\right) dx. \tag{2.1}$$

To compute this, consider its square

$$Z^{2} = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} \exp\left(-\frac{x^{2} + y^{2}}{2\sigma^{2}}\right) dx dy.$$
 (2.2)





Let us change variables from cartesian (x, y) to polar  $(r, \theta)$  using  $x = r \cos \theta$  and  $y = r \sin \theta$ . Since  $dx dy = r dr d\theta$  (recall that r is the determinant of the Jacobian matrix in 2D) and  $\cos^2 \theta + \sin^2 \theta = 1$ , we have

$$Z^{2} = \int_{0}^{2\pi} \int_{0}^{\infty} r \exp\left(-\frac{r^{2}}{2\sigma^{2}}\right) dr d\theta \tag{2.3}$$

Evaluate this integral and thus show that  $Z = \sqrt{2\pi\sigma^2}$ .

**Hint 1:** Separate the integral into a product of two integrands, the first of which (involving  $d\theta$ ) is constant, so is easy.

**Hint 2:** If  $u = \exp\left(-\frac{r^2}{2\sigma^2}\right)$  then  $\frac{du}{dr} = -\frac{r}{\sigma^2} \cdot \exp\left(-\frac{r^2}{2\sigma^2}\right)$ , so the second integral is also easy (since  $\int u'(r) dr = u(r)$ ).

## 3 Reducing the cost of linear regression for large D, small N

The ridge method is a regularized version of least squares with objective function:

$$\min_{\mathbf{w} \in \mathbb{R}^D} \|\mathbf{y} - \mathbf{X}\mathbf{w}\|_2^2 + \lambda \|\mathbf{w}\|_2^2$$
 (3.1)

Here  $\lambda$  is a scalar, the input matrix  $\mathbf{X} \in \mathbb{R}^{N \times D}$  and the output vector  $\mathbf{y} \in \mathbb{R}^{N}$ . The parameter vector  $\mathbf{w} \in \mathbb{R}^{D}$  is obtained by differentiating the cost function, yielding the normal equations

$$(\mathbf{X}^T \mathbf{X} + \lambda \mathbf{I}_D) \mathbf{w} = \mathbf{X}^T \mathbf{y}, \tag{3.2}$$

where  $\mathbf{I}_D$  is the  $D \times D$  identity matrix. The predictions  $\hat{\mathbf{y}} = \hat{\mathbf{y}}(\mathbf{X}_*)$  for new test points  $\mathbf{X}_* \in \mathbb{R}^{N^* \times D}$  are obtained by evaluating the hyperplane

$$\widehat{\mathbf{y}} = \mathbf{X}_* \mathbf{w} = \mathbf{X}_* (\mathbf{X}^T \mathbf{X} + \lambda \mathbf{I}_D)^{-1} \mathbf{X}^T \mathbf{y} = \mathbf{H} \mathbf{y}.$$
 (3.3)

The matrix  $\mathbf{H}$  is known as the *hat matrix* because it puts a "hat" on y.

- 1. Show that the solution can be written as  $\mathbf{w} = \mathbf{X}^T \widetilde{\mathbf{w}}$ , where  $\widetilde{\mathbf{w}} = \lambda^{-1} (\mathbf{y} \mathbf{X} \mathbf{w})$ .
- 2. Show that  $\widetilde{\mathbf{w}}$  can also be written as follows:  $\widetilde{\mathbf{w}} = (\mathbf{X}\mathbf{X}^T + \lambda \mathbf{I}_N)^{-1}\mathbf{y}$  and, hence the predictions can be written as follows:

$$\widehat{\mathbf{y}} = \mathbf{X}_* \mathbf{w} = \mathbf{X}_* \mathbf{X}^T \widetilde{\mathbf{w}} = [\mathbf{X}_* \mathbf{X}^\mathsf{T}] ([\mathbf{X} \mathbf{X}^T] + \lambda \mathbf{I}_N)^{-1} \mathbf{y}.$$
(3.4)

(This an awesome trick because if N=20 patients with D=10,000 gene measurements, the computation of  $\widetilde{\mathbf{w}}$  only requires inverting the  $N\times N$  matrix, while the direct computation of  $\mathbf{w}$  would have required the inversion of a  $D\times D$  matrix.)



### Machine Learning

Michaelmas Term 2016

Week 3

## 4 Logical Gates Using Perceptrons

Recall that a perceptron with input features  $x_1, \ldots, x_D$ , weights  $w_1, \ldots, w_D$  and bias  $w_0$  outputs the value:

$$y = \begin{cases} 1 & \text{if } w_0 + \sum_{i=1}^D w_i x_i \ge 0\\ 0 & \text{otherwise} \end{cases}$$
 (4.1)

- 1. Suppose there are at most two inputs and the inputs always take binary values, *i.e.*,  $x_i \in \{0,1\}$ . Show how to construct AND, OR and NOT gates by suitably adjusting weights.
- 2. The constructions for AND and OR gates required only the bias term  $w_0$  to be negative, all other weights were positive. Can you achieve a similar construction for the NOT gate? Why?
- 3. Can you construct an XOR (exclusive or) gate? If not, give reasons.
- 4. Often, instead of using a hard threshold we would like to use a continuous approximation. Recall the hyperbolic tangent function  $\tanh(z) = \frac{e^z e^{-z}}{e^z + e^{-z}}$ . We consider another type of artificial neuron whose output is defined as

$$y = \tanh\left(w_0 + \sum_{i=1}^{D} w_i x_i\right). \tag{4.2}$$

Suppose you treat outputs above 0.99 as true and those below -0.99 as false. Show that similar constructions to the ones you had earlier can still be used to construct logic gates.