#### Lecture 6

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#### Administrative

- Homework from 15/06/2019 due now!
- Please complete and hand in the declarations of authorship.

# Questions?

#### Solution to homework from lecture 4

### Bayesian learning and neural networks

Given data of the form

```
d = \{(x_1, c(x_1)), ..., (x_m, c(x_m))\}
```

where c is an unknown function, Bayesian learning attempts to find the most probable hypothesis  $h \in H$  that can be used to determine c.

When using neural networks, h represents an assignment of the weights and tresholds of the neurons.

### Bayesian learning and neural networks

It can be shown (see slides of Lecture 5) that the most probable hypothesis is, under common circumstances, the one that minimises the *error* (or *loss*) function

```
E(w) = \sum (x_i - f w (x_i))^2
```

where  $f: Weights \rightarrow X \rightarrow \mathbb{R}$  is the function implemented by the network.

# Hill climbing

Consider the following problem: given a function

```
f: ([-5, 5], [-5, 5]) \rightarrow \mathbb{R}
```

find x,  $y \in [-5, 5]$  such that f(x, y) is minimal.

We could, e.g., try some values:

$$f(1, 1) = 0.0$$

$$f(2, 1) = 0.0$$

$$f(3, 1) = 2.0$$

$$f(3, 2) = -2.0$$

# Hill climbing

We need a way of exploring the domain of f more systematically. We can, for example, start in (0, 0) and test the corners of the square of side 1 centered in the origin:

```
f(0, 0) = 5

f(-0.5, -0.5) = 8.25

f(-0.5, 0.5) = 4.75

f(0.5, 0.5) = 2.25

f(0.5, -0.5) = 6.75
```

We can now use (0.5, 0.5) as a starting point and continue, perhaps making the side of the square smaller (e.g., 0.9 instead of 1.

# Hill climbing and neural networks

Attempts to use hill climbing to minimise **E** are usually unsuccessful. That is because the space to be explored is exponential in the number of parameters (for two dimensions we have 4 points, for three dimensions 8, for n dimensions 2<sup>n</sup> points).

We need a "guide" to point us in the right direction.

#### Gradient

For a function  $f: X \to \mathbb{R}$  where  $X \subseteq \mathbb{R}^n$ , the gradient (if it exists) is the vector

$$\nabla$$
 f (x) = [D<sub>1</sub> f (x), ..., D<sub>n</sub> f (x)]

where we use  $D_{\text{i}}~\text{f}~(x)$  instead of the more frequent  $\delta~\text{f}~(x)~/~\delta~x_{\text{i}}$ 

 $\nabla$  f (x) points towards the direction of steepest increase in f at x.

#### Gradient descent

If we have information about the gradient of f, then we can do **much** better than hill climbing:

```
f(0, 0) = 5
gradf(0, 0) = [-2, -4]
```

So if we move in the direction [-2, -4] we should see an increase in f; if we move in the opposite direction, a decrease:

$$f(-2, -4) = 37$$
  
 $f(2, 4) = -3$ 

We now move to (2, 4) and start again, perhaps choosing a different step size (e.g., 0.9).

This procedure is a variant of *gradient descent*. In "real" gradient descent, the step size nis chosen to minimise

$$f(x - \eta * D_1 f(x, y), y - \eta * D_2 f(x, y))$$

#### Revisiting perceptrons

```
perceptron : (\mathbb{R}^n, \mathbb{R}) \to \mathbb{R}^n \to \{-1, 1\}

perceptron ([w_1, ..., w_n], \theta) [x_1, ..., x_n] = if s \ge \theta then 1

else -1

where s = w_1 * x_1 + ... + w_n * x_n

We are given

d = \{(x_1, c(x_1)), ..., (x_m, c(x_m))\}
```

Can we use gradient descent to determine  $w_i$  and  $\theta$ ?

We can compute **E**, but the problem is that it is not differentiable w.r.t.

```
Because perceptron ([w_1, \ldots, w_n], \theta) : \mathbb{R}^n \to \{-1, 1\} is discontinuous.
```

We can approximate the step function required by perceptron with a differentiable function. For example:

```
perceptron ([w_1,..., w_n], \theta) [x_1,..., x_n] = \sigma(s)

where s = w_1*x_1 + ... + w_n * x_n

\sigma(y) = 1 / 1 + \exp(-y)
```

Note that this scales the output to [0, 1].

```
\sigma'(y) = \sigma(y) * (1 - \sigma(y))
```

#### More generally

```
perceptron ([w_1, ..., w_n], \theta) [x_1, ..., x_n] = f(s)

where s = w_1*x_1 + ... + w_n * x_n

f(y) = ...
```

where f is some nicely differentiable function.

In the following, we consider a perceptron with no  $\theta$ , since

```
perceptron ([w_1, ..., w_n], \theta) [x_1, ..., x_n] = perceptron ([w_1, ..., w_n, \theta]) [x_1, ..., x_n, 1] and abbreviate
```

```
WS = [W_1, \ldots, W_n]
```

Consider m = 1 (only one element in the data), so that

```
E(ws) = (c(x) - perceptron ws (x))^{2}= (t - per ws (x))^{2}
```

We can apply gradient descent if we can compute  $D_i \in (w_i)$ . This is easy:

Notice the similarity with the learning rule from the last lecture.

### Stochastic gradient descent

In practice, m > 1.

"Proper" gradient descent demands the computation of E.

However, this would be too time-consuming, and usually an approximate **E** is computed instead, using a "batch" size between 1 and m. This approximate version is known as *stochastic gradient descent*.

# More than one layer

What if we have more than one layer? Consider a simple example:

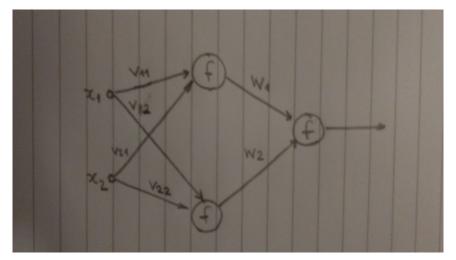


Figure 1: Two layer network

### More than one layer

We can compute the error as before

```
E(ws, vs) = (t - per ws [y<sub>1</sub>, y<sub>2</sub>])<sup>2</sup>

where y<sub>1</sub> = per vs<sub>1</sub> [x<sub>1</sub>, x<sub>2</sub>]

y<sub>2</sub> = per vs<sub>2</sub> [x<sub>1</sub>, x<sub>2</sub>]

vs<sub>1</sub> = [v<sub>11</sub>, v<sub>21</sub>]

vs<sub>2</sub> = [v<sub>12</sub>, v<sub>22</sub>]

D<sub>1</sub> E(ws, vs) = -2 * (t - f(s)) * f'(s) * y<sub>1</sub>

But how can we compute, e.g., D<sub>3</sub> E(ws, vs)?
```

### Backpropagation

#### Homework

Consider a two-layer perceptron as in Figure 1, with f(y) = y, f'(y) = 1 (such a perceptron is called *linear*). Suppose we start with

```
V_{11} = 1, V_{12} = -1, V_{21} = -1, V_{22} = 1, W_{1} = 1, W_{2} = 2
```

Assume the data is ([-1, 1], 0). How does the backpropagation algorithm adjust  $v_{11}$ ?