

ETC3555

Statistical Machine Learning

Deep Neural Networks

4 September 2018

Deep learning in R with Keras

```
library(tidyverse)
 2
     library(keras)
 3
 4
     keras_model_seauential() %>%
 5
       layer_dense(units = 256, activation = "relu", input_shape = c(784)) %>%
 6
       layer_dropout(rate = 0.4) %>%
       layer_dense(units = 128, activation = "relu") %>%
 8
       layer_dropout(rate = 0.3) %>%
 9
       laver_dense(units = 10, activation = "softmax") %>%
10
       compile(loss = "categorical_crossentropy".
11
               optimizer = optimizer_rmsprop().
12
               metrics = c("accuracy")) %>%
13
       fit(x_train, y_train, epochs = n_epochs, batch_size = 128, validation_split = 0.2)
```

Introduction 2/21

Some design choices

- Network architecture
- Activation functions for hidden and output layers
- Weight initialization
- Optimization algorithm
- Regularization method
- Hyperparameter selection procedure

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Introduction 3/21

Deep neural networks

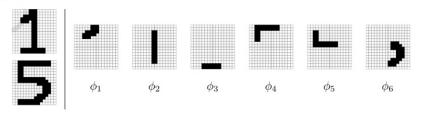
- The universal approximation theorem states that a single hidden layer with enough hidden units can approximate any target function with any desired accuracy.
- In other words, regardless of what the target function is, a large network will be able to represent this function.

Introduction 4/21

Deep neural networks

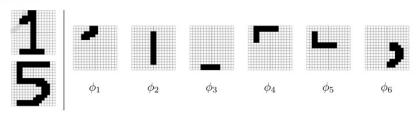
- We are not guaranteed that the training algorithm will be able to *learn* that function. The algorithm can fail due to overfitting and the hard optimization problem.
- That may not be a natural way to represent the target function. Many layers (i.e. a composition of several simpler functions) more closely mimics human learning.
- 3 Empirically, a deeper network does seem to result in better generalization for a wide variety of tasks. This is an area of active research.

Introduction 5/21



We consider the digit recognition problem to classify 1 versus 5. How to classify '1' and '5' from these basic shapes/features?

Introduction 6/21



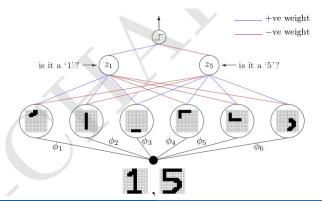
We consider the digit recognition problem to classify 1 versus 5. How to classify '1' and '5' from these basic shapes/features?

- lacksquare '1' should contain a ϕ_1 , ϕ_2 and ϕ_3
- lacksquare '5' should contain a ϕ_3 , ϕ_4 , ϕ_5 , ϕ_6 and perhaps a little ϕ_1

We want the value of these features to be large (close to 1) if its corresponding feature is in the input image and small (close to -1) if not.

Introduction 6/21

• We build complex Boolean functions from the 'basic' functions AND and OR. We want to mimic that process here. We assume we have feature functions ϕ_i which compute the presence (+1) or absence (-1) of the corresponding feature.



Introduction 7/21

Exercise

Exercise 7.18

Since the input x is an image it is convenient to represent it as a matrix $[x_{ij}]$ of its pixels which are black $(x_{ij} = 1)$ or white $(x_{ij} = 0)$. The basic shape ϕ_k identifies a set of these pixels which are black.

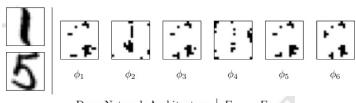
(a) Show that feature ϕ_k can be computed by the neural network node

$$\phi_k(\mathbf{x}) = \tanh\left(w_0 + \sum_{ij} w_{ij} x_{ij}\right).$$

- (b) What are the inputs to the neural network node?
- (c) What do you choose as values for the weights? [Hint: consider separately the weights of the pixels for those $x_{ij} \in \phi_k$ and those $x_{ij} \notin \phi_k$.]
- (d) How would you choose w_0 ? (Not all digits are written identically, and so a basic shape may not always be exactly represented in the image.)
- (e) Draw the final network, filling in as many details as you can.

Introduction 8/21

- **A** deep network architecture with $[d^{(0)}, d^{(1)}, d^{(2)}, d^{(3)}] = [256, 6, 2, 1]$
- The following figure shows what the 6 hidden units in the first hidden layer **learned**. The pixels corresponding to the top 20 incoming weights are shown.



Deep Network Architecture	$E_{\rm in}$	E_{test}
[256, 3, 2, 1]	0	0.170%
[256, 6, 2, 1]	0	0.187%
[256, 12, 2, 1]	0	0.187%
[256, 24, 2, 1]	0	0.183%

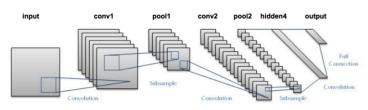
Introduction 9/21

- Our 'deep network' has just 2 hidden layers. In a more complex problem, there would be a hierarchy of "basic" features.
- "The first layer constructs a low-level representation of basic shapes"
- "The next layer builds a higher level representation from these basic shapes"
- We obtain more complex representations in terms of simple parts from previous layers. An 'intelligent' decomposition of the problem.

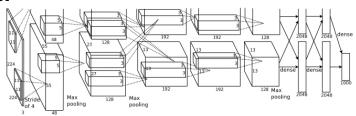
Introduction 10/21

Deep architectures

LeNet



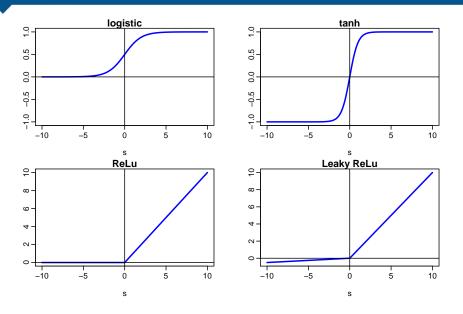
AlexNet



Many others. Source: https://bit.ly/2mWCUw1

Introduction 11/21

Activation functions



Introduction 12/21

Rectified linear unit (ReLu)

- $\theta(s) = max(0,s)$
- The most popular activation function in modern deep learning
- Better gradient propagation compared to sigmoidal activation functions that saturate in both directions (less vanishing gradient problems)
- Sigmoidal activation functions have an almost zero gradient at certain regions. This is an undesirable property since it results in slow learning.
- Efficient computation
- In a randomly initialized network, only about half of the hidden units are activated (having a non-zero output)

Introduction 13/21

Rectified linear unit (ReLu)

- Potential problems: Non-differentiable at zero, unbounded, etc.
- Dying ReLU problem: if a ReLu unit become inactive for all inputs, no gradients flow backward through that unit, and it 'dies'
- Leaky ReLU: $\theta(s) = \begin{cases} s, & \text{if } s \ge 0 \\ as & \text{otherwise} \end{cases}$ where a > 0 is a small positive number
 - Other variants

Introduction 14/21

Softmax (output) activation function

In logistic regression, we compute

$$P(y|\mathbf{x}) = egin{cases} h(\mathbf{x}) = heta(w^T \mathbf{x}) = heta(s) & ext{for } y = +1; \ 1 - h(\mathbf{x}) & ext{for } y = -1, \end{cases}$$

where $\theta(\cdot)$ is the logistic function.

The *softmax* function is a generalized logistic activation function used for multiclass classification. It has both multivariate input and output.

For a classification problem with K classes, we compute

$$P(y=j|\mathbf{x}) = \theta(\mathbf{s})_j = \theta(\mathbf{s})_j = \frac{e^{\mathbf{s}_j}}{\sum_{k=1}^K e^{\mathbf{s}_k}},$$

where $\mathbf{s}=(s_1,s_2,\ldots,s_K)^T$, $s_j=\mathbf{w}_i^T\mathbf{x}$ and $j=1,2,\ldots,K$.

Introduction 15/21

Softmax (output) activation function

- Softmax vs "hardmax "
 - Softmax: $\mathbf{s} = (5, 2, -1, 3)^T \xrightarrow{\theta} (0.84, 0.04, 0, 0.12)^T$
 - Hardmax: $\mathbf{s} = (5, 2, -1, 3)^T \stackrel{\theta}{\to} (1, 0, 0, 0)^T$

$$\theta(\mathbf{s})_j = \begin{cases} 1 & \text{if } j = \operatorname{argmax}(s_1, s_2, \dots, s_K); \\ 0 & \text{otherwise.} \end{cases}$$

- The softmax behaves as a smooth function (which can be differentiated) and approximates the indicator function.
- The outputs are represented using binary vectors and the multi-class cross-entropy (or logloss) is used as loss function

Introduction 16/21

Regularization

- L_1 and L_2 norm regularization (weight decay, weight elimination, etc)
- Early stopping
- Dropout
- Data augmentation

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Introduction 17/21

L₂ regularization: weight decay

$$E_{\text{aug}}(\mathbf{w}) = E_{\text{in}}(\mathbf{w}) + \frac{\lambda}{N} \sum_{\ell,i,j} (w_{ij}^{(\ell)})^2$$

Introduction 18/21

L₂ regularization: weight decay

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What is $\frac{\partial E_{\mathrm{aug}}}{\partial w_{ii}^{(l)}}$?

Introduction 18/21

L₂ regularization: weight decay

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What is $\frac{\partial E_{\mathrm{aug}}}{\partial w_{ii}^{(l)}}$?

$$\frac{\partial E_{\mathrm{aug}}(\mathbf{w})}{\partial \mathbf{W}^{(\ell)}} = \frac{\partial E_{\mathrm{in}}(\mathbf{w})}{\partial \mathbf{W}^{(\ell)}} + \frac{2\lambda}{N} \mathbf{W}^{(\ell)}$$

- A component in the negative direction of **w** is added to the weight update.
- We can compute the first term using backpropagation

Introduction 18/21

L₂ regularization: weight elimination

$$E_{\text{aug}}(\mathbf{w}, \lambda) = E_{\text{in}}(\mathbf{w}) + \frac{\lambda}{N} \sum_{\ell, i, j} \frac{(w_{ij}^{(\ell)})^2}{1 + (w_{ij}^{(\ell)})^2}$$

Introduction 19/21

L₂ regularization: weight elimination

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What is
$$\frac{\partial E_{\text{aug}}}{\partial w_{ii}^{(l)}}$$
 ?

Introduction 19/21

L₂ regularization: weight elimination

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 ?

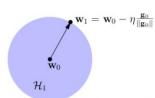
$$\frac{\partial E_{\mathrm{aug}}}{\partial w_{ij}^{(\ell)}} = \frac{\partial E_{\mathrm{in}}}{\partial w_{ij}^{(\ell)}} + \frac{2\lambda}{N} \cdot \frac{w_{ij}^{(\ell)}}{(1 + (w_{ij}^{(\ell)})^2)^2}$$

Many other weight based complexity penalties ...

Introduction 19/21

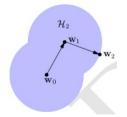
Early stopping

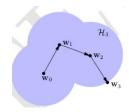
- With more gradient descent iterations, more of our hypothesis set is explored. With fewer iterations, we explore a smaller hypothesis set + better generalization
- With fixed-step gradient descent, we do:
 - **1** We start at weights \mathbf{w}_0
 - **2** We take a step size η to $\mathbf{w}_1 = \mathbf{w}_0 \eta \frac{\mathbf{g}_0}{\|\mathbf{g}_0\|}$
- We implicitly searched the following hypothesis set: $\mathcal{H}_1 = \{ \boldsymbol{w} : \| \boldsymbol{w} \boldsymbol{w}_0 \| \le \eta \}$ and picked the hypothesis \boldsymbol{w}_1 with minimum E_{in} .



Introduction 20/21

Early stopping





- $\begin{aligned} & \quad \mathcal{H}_2 = \mathcal{H}_1 \cup \{\mathbf{w} : \|\mathbf{w} \mathbf{w}_1\| \leq \eta\}, \\ & \quad \mathcal{H}_3 = \mathcal{H}_2 \cup \{\mathbf{w} : \|\mathbf{w} \mathbf{w}_2\| \leq \eta\}, \text{ and } \mathcal{H}_1 \subset \mathcal{H}_2 \subset \mathcal{H}_3 \subset \dots \end{aligned}$
- As t increases, $E_{in}(\mathbf{w}_t)$ decreases but the complexity of \mathcal{H}_t increases
- It may be better to stop early at some t^* , well before reaching a minimum of E_{in} . A validation set can be used to monitor validation error and determine when to stop training.

Introduction 21/21