PSY9511: Seminar 7

Deep learning for computer vision tasks

Esten H. Leonardsen 07.11.24



Outline

- 1. Exercise 4
- 2. Deep learning
 - Motivation
 - · (Deep) neural networks
 - · Training procedure
- 3. Convolutional neural networks for computer vision

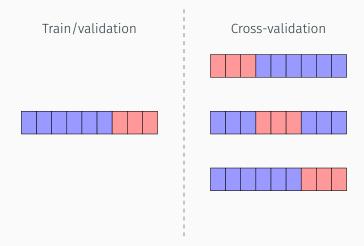


Weekly exercises

- The weekly exercises are mandatory
- · The deadlines are strict

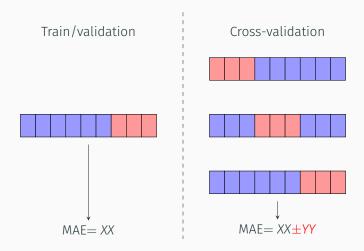


Validation procedures





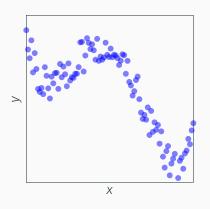
Validation procedures



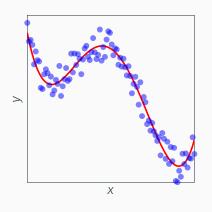


Deep learning



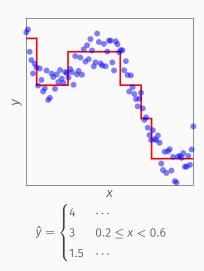


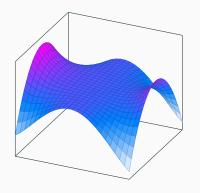




$$\hat{y} = s(x)$$





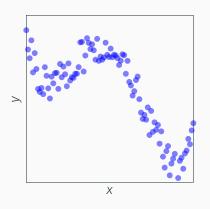




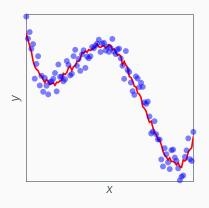


The cat wagged its tail

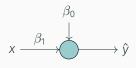




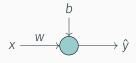




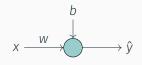




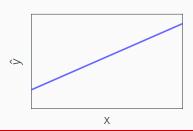
$$\hat{y} = \beta_0 + \beta_1 x$$

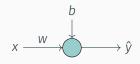


$$\hat{y} = wx + b$$

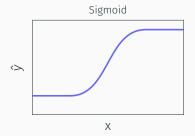


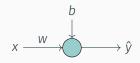
$$\hat{y} = wx + b$$





$$\hat{y} = \frac{e^{wx+b}}{1 + e^{wx+b}}$$



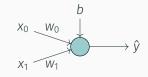


$$\hat{y} = max(0, wx + b)$$

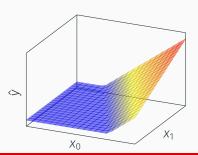
Rectified Linear Unit (ReLU)

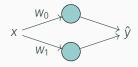




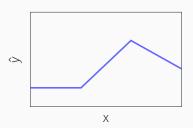


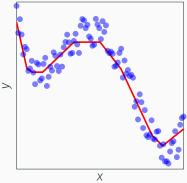
$$\hat{y} = max(0, w_0x_0 + w_1x_1 + b)$$





$$\hat{y} = max(0, w_0x + b_0) + max(0, w_1x + b_1)$$





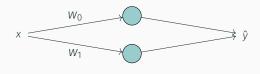
Piecewise linear function



Universal approximation theorem:

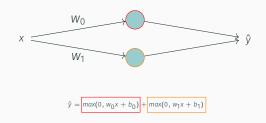
"Any relationship that can be described with a polynomial function can be approximated by a neural network with a single hidden layer."



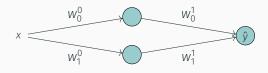


 $\hat{y} = max(0, w_0x + b_0) + max(0, w_1x + b_1)$



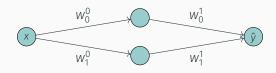






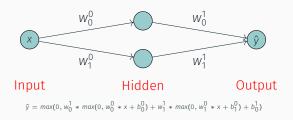
$$\hat{y} = \max(0, w_0^1 * \max(0, w_0^0 * x + b_0^0) + w_1^1 * \max(0, w_1^0 * x + b_1^0) + b_0^1)$$



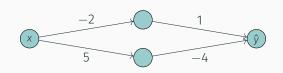


$$\hat{y} = \max(0, w_0^1 * \max(0, w_0^0 * x + b_0^0) + w_1^1 * \max(0, w_1^0 * x + b_1^0) + b_0^1)$$

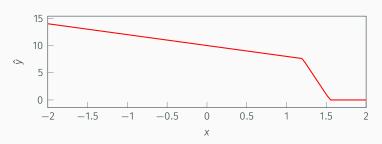




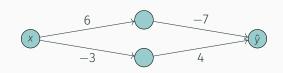




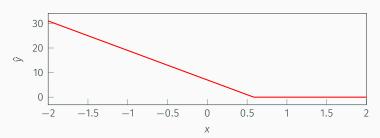
$$\hat{y} = max(0, 1 * max(0, (-2) * x + 3) + (-4) * max(0, 5 * x + (-6)) + 7)$$

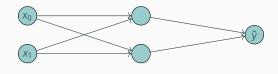






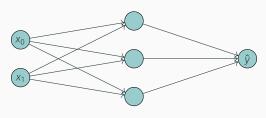
$$\hat{y} = max(0, -7 * max(0, 6 * x + (-5)) + 4 * max(0, (-3) * x + 2) - 1)$$



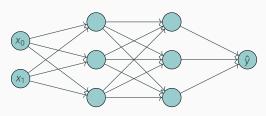


$$\begin{split} \hat{y} &= max(0, w_{0,0}^0 * max(0, w_{0,0}^0 * x_0 + w_{1,0}^0 * x_1 + b_{0,0}) + \\ & w_{1,0}^1 * max(0, w_{0,1}^0 * x_0 + w_{1,1}^0 * x_1 + b_{0,1}) + \\ & b_1) \end{split}$$





$$\begin{split} \hat{y} &= max(0, w_{0,0}^1 * max(0, w_{0,0}^0 * x_0 + w_{1,0}^0 * x_1 + b_{0,0}) + \\ w_{1,0}^1 * max(0, w_{0,1}^0 * x_0 + w_{1,1}^0 * x_1 + b_{0,1}) + \\ w_{2,0}^1 * max(0, w_{0,2}^0 * x_0 + w_{1,2}^0 * x_1 + b_{0,2}) + \\ b_1) \end{split}$$



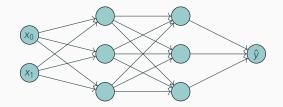
$$\begin{split} \hat{y} &= \max(0, w_{0,0}^2 * \max(0, w_{0,0}^1 * \max(0, w_{0,0}^0 * x_0 + w_{1,0}^0 * x_1 + b_{0,0}) + \\ & w_{1,0}^1 * \max(0, w_{0,1}^0 * x_0 + w_{1,1}^0 * x_1 + b_{0,2}) + \\ & w_{2,0}^2 * \max(0, w_{0,2}^0 * x_0 + w_{1,2}^0 * x_1 + b_{0,2}) + \\ & b_{1,0}) + \\ & w_{1,0}^2 * \max(0, w_{0,1}^1 * \max(0, w_{0,0}^0 * x_0 + w_{1,0}^0 * x_1 + b_{0,0}) + \\ & w_{1,1}^1 * \max(0, w_{0,1}^0 * x_0 + w_{1,1}^0 * x_1 + b_{0,1}) + \\ & w_{2,1}^2 * \max(0, w_{0,2}^0 * x_0 + w_{1,2}^0 * x_1 + b_{0,2}) + \\ & b_{1,1}) + \\ & w_{2,0}^2 * \max(0, w_{0,2}^1 * x_0 + w_{1,0}^0 * x_1 + b_{0,0}) + \\ & w_{1,2}^1 * \max(0, w_{0,1}^0 * x_0 + w_{1,1}^0 * x_1 + b_{0,1}) + \\ & w_{1,2}^2 * \max(0, w_{0,1}^0 * x_0 + w_{1,1}^0 * x_1 + b_{0,1}) + \\ & w_{2,2}^1 * \max(0, w_{0,2}^0 * x_0 + w_{1,2}^0 * x_1 + b_{0,2}) + \\ & b_{1,2}) + \end{split}$$

<u>Artificial neural networks</u>: Combines artificial neurons, simple computational units that compute a non-linear function of their inputs, in a computational graph

- Can approximate arbitrarily complex polynomial functions (given enough neurons)
- Organized in layers. We can expand a model in width (e.g. more neurons per layer) or depth (e.g. more layers)

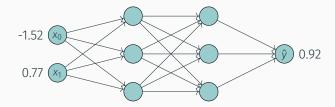


Deep learning: Loss functions



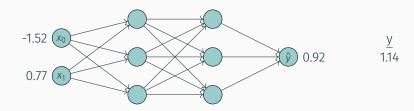


Deep learning: Loss functions

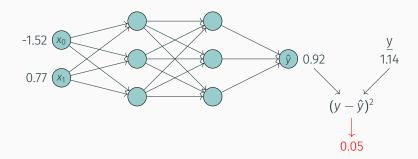




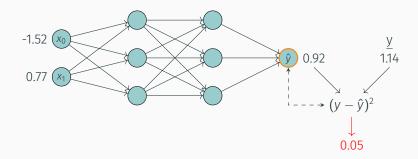
Deep learning: Loss functions













```
In[1]: from tensorflow.keras.layers import Input, Dense
    from tensorflow.keras import Model

inputs = Input(shape=(2,))
    hidden1 = Dense(units=3, activation='relu')(inputs)
    hidden2 = Dense(units=3, activation='relu')(hidden1)
    outputs = Dense(units=1, activation=?)(hidden2)

model = Model(inputs, outputs)
    model.compile(loss=?)
```

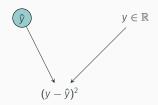


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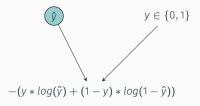
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```

Regression



Binary classification



Multiclass classification



 $y \in \{cat, dog, bat\}$



Multiclass classification







Multiclass classification



x ₀	cat	dog	bat
	1	0	1
	0	1	0
	0	0	1
	0	1	0

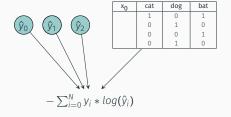
Multiclass classification







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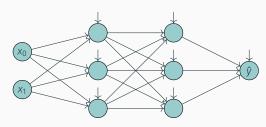




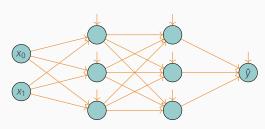
<u>Loss functions</u>: Behaves for neural networks as any other statistical learning model. However, important to **configure the final layer** correctly

- · Regression: Mean squared error
 - · No activation
- Binary classification: Binary cross-entropy
 - · Sigmoid activation
- Multiclass classification: Categorical cross-entropy
 - · Softmax activation

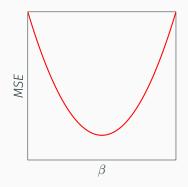




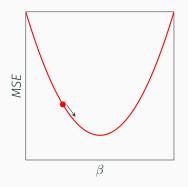
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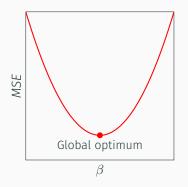
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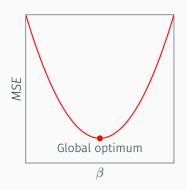












$$|\beta| \to 10^6 - 10^{12}$$

$$\beta_x \implies \beta_y$$



Learning representations by back-propagating errors

David E. Rumelhart*, Geoffrey E. Hinton† & Ronald J. Williams*

We describe a new learning procedure, back-propagation, for networks of neurone-like units. The procedure repeatedly adjusts the weights of the connections in the network so as to minimize a measure of the difference between the actual output vector of the net and the desired output vector. As a result of the weight adjustments, internal 'hidden' units which are not part of the input or output come to represent important features of the task domain, and the regularities in the task are captured by the interactions of these units. The ability to create useful new features distinguishes back-propagation from earlier, simpler methods such as the perceptron-convergence procedure'.



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[†] Department of Computer Science, Carnegie-Mellon University, Pittsburgh, Philadelphia 15213, USA

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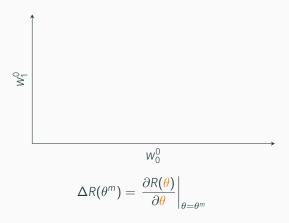


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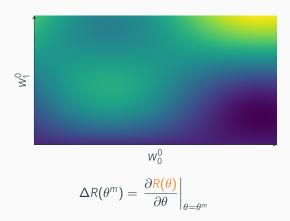
$$\Delta R(\theta^m) = \left. \frac{\partial R(\theta)}{\partial \theta} \right|_{\theta = \theta^m}$$





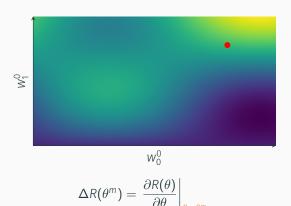
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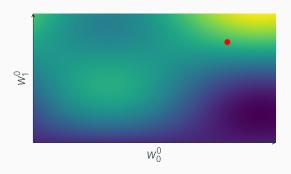
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- · $R(\theta)$ is the loss as a function of the parameters





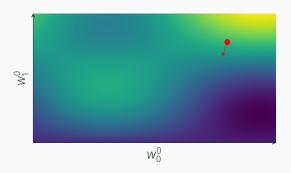
- \cdot θ are the parameters of the model
- · $R(\theta)$ is the loss as a function of the parameters
- θ^m is a specific configuration of parameters





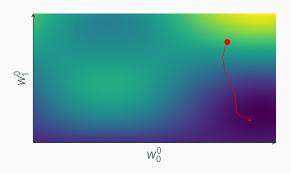
$$\frac{\partial R_i(\theta)}{\partial w_{jk}} = \frac{\partial R(\theta)}{\partial f_{\theta}(x_i)} \cdot \frac{\partial f_{\theta}(x_i)}{\partial g(z_{ik})} \cdot \frac{\partial g(z_{ik})}{\partial z_{ik}} \cdot \frac{z_{ik}}{w_{kj}}$$





$$\frac{\partial R_i(\theta)}{\partial w_{jk}} = \frac{\partial R(\theta)}{\partial f_{\theta}(x_i)} \cdot \frac{\partial f_{\theta}(x_i)}{\partial g(z_{ik})} \cdot \frac{\partial g(z_{ik})}{\partial z_{ik}} \cdot \frac{z_{ik}}{w_{kj}}$$





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<u>Backpropagation</u>: Uses gradient descent to iteratively determine how the model weights should be updated to minimize the loss function

 Gradient descent: Calculate gradient based on all data points



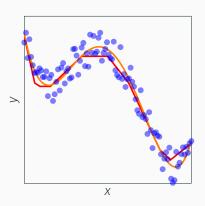
<u>Backpropagation</u>: Uses gradient descent to iteratively determine how the model weights should be updated to minimize the loss function

- Gradient descent: Calculate gradient based on all data points
- Stochastic gradient descent: Calculate gradient based on a batch of data points



<u>Splines</u>: A smooth curve implemented via piecewise polynomial functions

<u>Neural networks</u>: A piecewise linear function implemented as a hierarchy of artificial neurons





<u>Splines</u>: A smooth curve implemented via piecewise polynomial functions

• Requires us to carefully balance the complexity of the function

<u>Neural networks</u>: A piecewise linear function implemented as a hierarchy of artificial neurons



Splines: A smooth curve implemented via piecewise polynomial functions

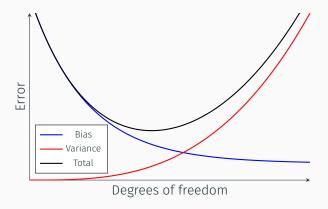
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Neural networks: A piecewise linear function implemented as a hierarchy of artificial neurons

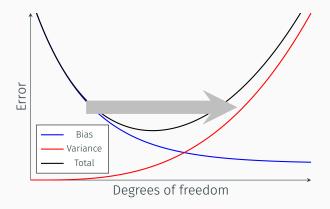
Overparameterization



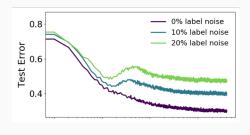


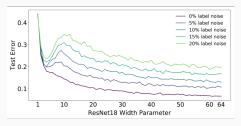




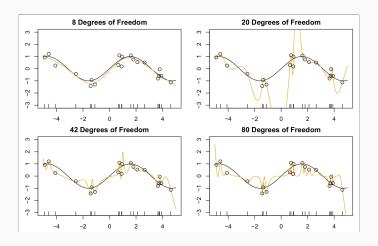














Overparameterization: Deep artificial neural networks generally have far more parameters than necessary (and often more than the number of data points)

- At face value, it is surprising that this does not yield severe overfitting
- However, it can be shown that neural networks, after perfectly fitting their training data, generally become more well-behaved and less wild



• Weight decay: Applies an ℓ_2 -penalty to the weights, similarly to ridge regression

$$R(\theta; \lambda) = R(\theta) + \lambda \sum \theta^2$$

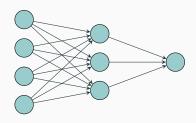


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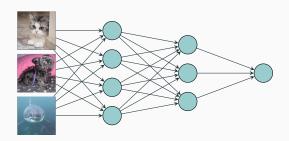
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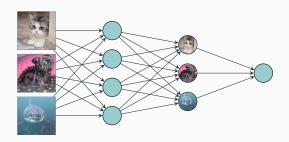
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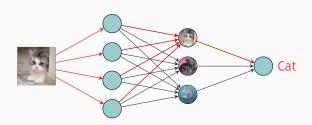
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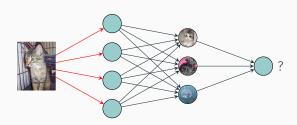
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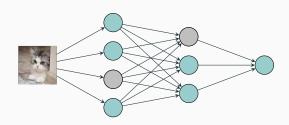
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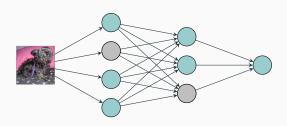
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- <u>Dropout:</u> Randomly kills a fraction of the neurons during training
- · Data augmentation: Attempts to generate more data

