Artificial Neural Networks (Gerstner). Exercises for week 4

Regularization and Tricks of the Trade

Exercise 1. Bagging and dropout

Assume K variants of a model, with model k having test error $E_k = E_0 + \epsilon_k$, where ϵ_k has mean $\mathbb{E}[\epsilon_k] = 0$, auto-variance $\mathbb{E}\left[\epsilon_k^2\right] = v$ and co-variance $\mathbb{E}[\epsilon_k \epsilon_n] = c$.

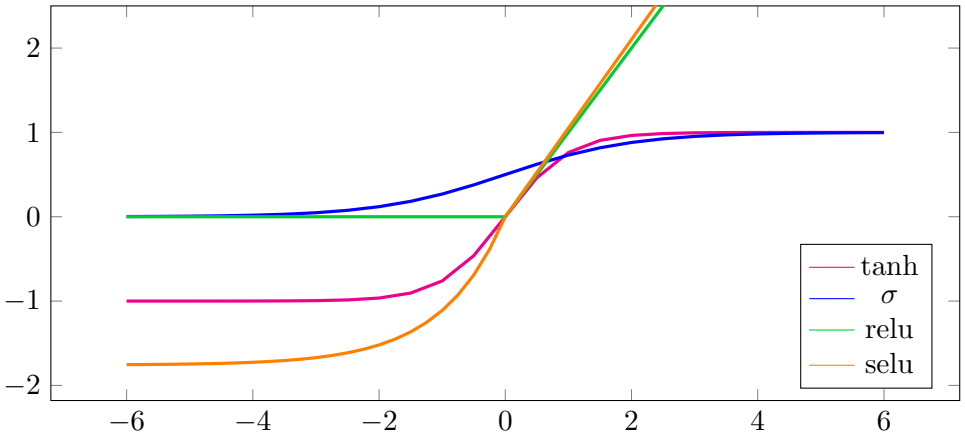
a. What is the expected value of the *bagging test error*, i.e. the expected test error of the model obtained by averaging over all variants?

b. What is the variance of the bagging test error? Does the number K of variants play a role?

c. For dropout: do you expect the co-variance c to be close to zero, between zero and v, close to v or larger than v? Justify your answer with one sentence.

Exercise 2. Different activation functions

The choice of the non-linearity function $g(x)$ can have a significant impact on learning speed and final performance. Which non-linearity is best, is still an active research question; the favorite non-linearity in the last century was probably the hyperbolic tangent $\tanh(x)$; since 2010, the rectified linear unit $\text{relu}(x) = \max(0, x)$ is highly popular and there is a fair chance that the new favorite will be the scaled exponential linear unit $\text{selu}(x) = \lambda x$ if $x > 0$ and $\text{selu}(x) = \alpha(\exp(x) - 1)$ otherwise, with $\lambda \approx 1.0507$ and $\alpha \approx 1.75814$. Currently, it seems that the key concepts to discuss the different non-linearities are, first, *linearity problem*, second the *vanishing gradient problem* and, third, the *bias shift problem*.



a. Linearity problem

(i) Show that a multi-layer neural network with linear activation function $g(x) = x$ is equivalent to a single layer linear network. Hint: the product of two matrices is again a matrix.

(ii) Assume that in each layer the inputs follow a Normal distribution with mean zero and small variance, i.e. $\sigma^2 \ll 1$. For which of the activation functions $\sigma(x) = 1/(1 + \exp(-x))$, $\tanh(x)$, $\text{relu}(x)$ and $\text{selu}(x)$ is a deep network basically equivalent to a linear network for this input distribution?

b. Vanishing gradient problem

(i) Assume now the inputs are such that they also fall into the non-linear regimes. For simplicity we assume that in each layer the activations are $a_1 = -10, a_2 = -5, a_3 = -1, a_4 = 1, a_5 = 5, a_6 = 10$. Without a calculator, determine the fraction of values close to zero of $g(a_i)$ and $g'(a_i)$ for all i and $g = \sigma, \tanh, \text{relu}, \text{selu}$. For example, for tanh none of the values $\tanh(-10), \tanh(-5), \ldots, \tanh(10)$ is close to zero but $4/6 = 2/3$ of the values of $\tanh' = 1 - \tanh^2$ are close to zero.

(ii) The update of a weight w_{ij} is proportional to $g'(a_i) \cdot g(a_j)$. Determine the fraction of $g'(a_i) \cdot g(a_j)$ that are close to zero considering all combinations of a_i and a_j and all activations $g = \sigma, \tanh, \text{relu}, \text{selu}$.

(iii) The δ's in backpropagation are in each layer multiplied with g'. Consider backpropagation through 3 layers, i.e. terms like $g'(a_i)g'(a_j)g'(a_k)$. Determine the fraction of such terms that are close to zero for $g = \sigma, \tanh, \text{relu}, \text{selu}$.

c. Bias shift problem

Consider a simple classification task. The data exist in $\mathcal{R}^N$. Data points from C_0 (with target $t = 0$) are uniformly distributed in each dimension such that $x_i \in [1, 2]$ for $i = 1 \dots N$. Data points from C_1 (with target $t = 1$) are uniformly distributed in each dimension such that $x_i \in [3, 4]$ for $i = 1 \dots N$. We want to learn to classify points using a logistic sigmoid unit trained with the cross–entropy loss; from last week, this results in the weight update rule

$$\Delta w_i = \eta \cdot (t - y) \cdot x_i$$

where $y = \sigma\left(\sum_i^N w_i x_i\right)$.

Points are presented one at a time (i.e. stochastic gradient descent).

(i) Assume we start with all weights $w_i = 0$ and present the point $\mathbf{x}^a$ from C_0, update the weights, then present $\mathbf{x}^b$. Give the drive $a = \sum_i^N w_i x_i^b$ of the output unit in response to $\mathbf{x}^b$, in terms of η, $\mathbf{x}^a$ and $\mathbf{x}^b$. Note: we do not yet need to specify which class $\mathbf{x}^b$ belongs to.

(ii) In general, we can encounter oscillations in stochastic gradient descent if a single training example strongly affects the network output – for instance, if it results in the same network output for any possible input.

We assume that if $a < -5$, $y \approx 0$, and if $a > 5$, $y \approx 1$. Under what conditions will the network output y be the same for all possible inputs $\mathbf{x}^b$ after the first training step? Can we choose a small enough η to prevent this, independent of N? What if we had chosen $\mathbf{x}^a$ from C_1 instead?

(iii) A common input normalization technique to to remove the mean from the dataset, such that $E[x_i] = 0$ across all dimensions x_i. Assume that each data point has an equal probability of coming from either C_0 and C_1. What are the new data ranges for C_0 and C_1 after removing the mean? Repeating step (ii), do we get the same result?

(iv) Consider a deep network where each hidden layer uses either the hyperbolic tangent or logistic sigmoid nonlinearity. Given what we've seen above, can you suggest one non–linearity or the other between hidden layers? Note that one layer's output is another layer's input.

d. Summarize your results by ranking the different activation functions for each of the problems discussed in this exercise.

	linearity problem	vanishing gradient problem	bias shift problem
tanh			
σ			
relu			
selu			

Exercise 3. Normalization of activations across multiple layers

In class we have seen, that by an appropriate normalization of the input patterns (for each input component: zero mean, unit standard deviation) combined with a Gaussian distribution of input weights ($\langle w_{ij}^{(1)} \rangle = 0$ and $\langle [w_{ij}^{(1)}]^2 \rangle = 1/N$] we can ensure that the activation variable of neurons in

the first layer has mean $\langle a_i^{(1)} = 0$ and variance $\langle [a_i^{(1)}]^2 = 1$. In the following we assume that the distribution of activations in layer 1 is Gaussian,i.e., a normal distribution $N(0,1)$.

The aim of the exercise is to go by induction from layer n to layer $n+1$. We start in layer 1.

Assume that neuron j in layer 1 has a rectified linear activation function, i.e., $x_j^{(1)} = [a_j^{(1)}]_+$.

a. What is the mean $\langle x_j^{(1)} \rangle$?

b. Assume that the weights in layer 2 are initialized with zero mean and variance $\langle [w_{kj}^{(2)}]^2 \rangle = [c^2]/N_1]$ where N_1 is the number of hidden neurons in the first layer.

   What is the mean activation $\langle \tilde{a}_k^{(2)} \rangle$ in layer 2? Here $\tilde{a}_k^{(2)} = \sum_{j=1}^{N_1} w_{kj}^{(2)} x_j^{(1)}$. The total activation of neuron k in layer 2 is $a_k^{(2)} = \tilde{a}_k^{(2)} - \theta_k$.

   What value should you choose for the threshold θ_k in layer 2, so that $\langle a_k^{(2)} \rangle = 0$ in layer 2?

c. Assume that you found a threshold so that $\langle a_k^{(2)} \rangle = 0$. Calculate the variance $\langle [a_k^{(2)}] \rangle$ as a function of the constant c.

d. Choose c such that the variance is one.

e. Can you now go from layer 2 to layer 3?