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THESIS TOPIC DECLARATION

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Implementation of a platform game

Topic of the Thesis work:

Developing an implementation of the classic game "Prince of Persia" in Java using libgdx library to create executables for Android, Windows and other operating systems.

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1. Introduction

1.1. Abstract

The goal of this thesis is to train a neural network for one task and then apply the trained neural network on the other task. This will be achieved by expressing the second task in terms of the first.

In the first part of the work, several artificial neural networks will be created and trained for the Snake game and the networks' properties, such as weights and biases, will be compared to their results. Some of the networks will be trained only briefly, while others will have an extensive training.

In the second part, the best performing network will play the Maze game and the results will be compared to two networks, one of which was exclusively trained for the Maze game and the other was trained for the Snake game first and for the Maze game second.

Some networks will have different structures, such as different number of layers and different number of neurons in them, but there will also be similarly structured networks, which will be trained with a varying number of test games to observe if more data and training bring the best results.

1.2. Snake game history

The Snake game is the classic video game concept where the player plays for a line that represents the snake which grows in length after consuming other points. The game's main difficulty is staying alive without touching the walls or itself. This concept was originally seen in the arcade maze game Blockade developed by Gremlin and published by Sega in October 1976. [1]

There are many variations of The Snake Game, both single player and multiplayer, with the most known being Nibbler, Tron, Slither.io and Nokia Snake.

Soon after its creation in 1978 Snake, which was programmed by Peter Trefonas for TRS-80 microcomputer and Apple II, was instantly widespread on arcade machines, personal computers and other platforms due to its ease of implementation and replayability and in 1996 Next Generation magazine ranked it number 41 on their "Top Hundred Games of All Time." [2]

In 1998 Nokia resurged the Snake popularity after preloading it as an app on their monochrome mobile phones. This helped the game tremendously as it was open to a much larger audience than before and people could play the game to spend their time in queues and bus stops. Snake was associated with a Nokia app ever since.

1.3. Artificial neural network history

The original idea of a neural network was described in 1943 by neuropsychologist Warren McCulloch and mathematician Walter Pitts in their “A logical calculus of the ideas immanent in nervous activity”, which was on how neurons in the brain might work. The writers have modeled a simple neural network using electrical circuits. [3]

Later in 1949, Donald Hebb introduced his theory, also known as “Hebb’s Postulate”, about the neural bases of learning in his book “Organization of Behavior”, which has indicated that neural connections are enhanced every time they are being used and if two nerves fire at the same time, the connection between them becomes stronger. [4] Thus, learning is not done by brain passively, which would be an incredible feat, but rather it is a process during which the cellular structure of the brain is altered permanently. This theory has a classic status within science and is backed by recent research. [5]

In 1959, Bernard Widrow and Marcian Hoff invented the Widrow-Hoff least mean squares filter (LMS) adaptive algorithm that was the basis for the ADALINE and MADALINE (Multiple ADaptive LINear Elements) networks. MADALINE was the first neural network to ever be applied to a real-world problem. It is an adaptive filter that eliminates the echoes on phone lines, which is still in commercial use. [6]

The success of ADALINE and MADALINE networks unfortunately did more harm than good to neural network research, particularly due to high expectations and outrageous promises on the network capabilities and severe limitations of hardware back then. Unfulfilled promises and increasing popularity of von Neumann architecture led to neural network research funding being heavily reduced.

There were a few advances in the field, however it was not until 1982 when John Hopfield presented “Neural networks and physical systems with emergent collective computational abilities” paper to the National Academy of Sciences. In this work he proposed using bidirectional connections between neurons, since previously the neural networks have only used single “forward” direction. [7] This has led to an invention of an associative neural network, also known as Hopfield network, which consists of long-term and short-term memories and the new training is done using the short-term memory.

In the same year, Reilly, Cooper and Elbaum have published “A neural model for category learning” that used the “hybrid” multiple layer neural network and in 1986 Rumelhart and McClelland described the use of parallel distributed processing (connectionism) in the neural network. [8] This is now known as a backpropagation network and, unlike hybrid networks, it uses more than a couple of layers, however that makes its training much slower, as it needs many more iterations over the input to produce results.

Nowadays, networks with multiple hidden layers, also known as deep neural networks (DNN), have achieved state-of-the-art performance on computer vision problems and the current goal is

to reduce both the model size and computational cost of the networks to allow a more widespread deployment. [9]

1.4. Neural networks in games

Neural networks are quite a rare occurrence in games as the latter are usually fully scripted and use 30-year old artificial intelligence (AI) technology, such as A* and finite state machines instead. [10] There are exceptions to this, of course, such as the 1996 game Creatures, where the player assumes control over the small furry animals and teaches them how to behave. These creatures are using the neural networks to learn. Another example would be the 2013 Forza Motorsport 5 by Microsoft where the neural networks learn how the human players control the in-game vehicles and react to events, such as crashes or overtakes by other players and then the trained networks, also known as drivatars, play versus human players for more realistic competition. [11]

On the other hand, the neural networks are quite commonly used for playing the games instead. One of the most famous examples of this is AlphaGo, a Google DeepMind's DNN which in the span of 2016-2017 won official matches against two world champions of Go game (ranked 1st and 4th at the time) to become the first Go program, not only to surpass the amateur level, but also to reach the professional nine dan level in the game. Even though nine dan is the highest possible level in Go, many top Go players believe that AlphaGo is above that. [12]

1.5. Personal motivation

One of the limitations of neural networks is the fact that the networks are problem-specific, and a trained network cannot be applied to another task after it learned to solve one. [13,14]

In this work, however, I aim to create and train a neural network that can play a game of one genre and then apply it onto a game of another genre to show that it is possible to reuse the network, instead of developing a new one altogether.

My personal goal is to familiarize myself with the neural networks playing the games and later apply that knowledge in game development field to create an alternative to the current AI technology and make the techniques of scripting and AI cheating obsolete. This will also make the AI controlled game characters' actions be more human-like and give AI the ability to learn from its mistakes and adjust accordingly.

I hope that this will bring us one step closer to more realistic artificial intelligence, which in turn will provide a better gaming experience to the players.

2. Preparation

2.1. Abstract concepts

2.1.1. Neural network

So far there was a lot of history of neural networks in this work, but there was no explanation on what the neural network is.

A neural network is a massively parallel distributed processor made up of simple processing units, which has a natural propensity for storing experiential knowledge and making it available for us. It resembles the brain in two respects: the knowledge is acquired by the network from its environment through a learning process and then stored using interneuron connection strengths, known as synaptic weights. [15]

There are two types of neural networks: biological neural networks and artificial neural networks (ANN). The biological neural networks could be any group of connected biological nerve cells, known as neurons. An example of such a network is our brain.

An artificial neural network, on the other hand, is a mathematical structure designed to simulate biological networks. ANN is formed of three different types of layers: input layer, hidden layer and output layer. Each layer consists of one or more neurons, where each neuron represents a relatively simple function that takes an input and produces an output. The neurons of an input layer are passive, meaning they do not modify the data used for training, while the neurons of hidden and output layers are active instead. A visual representation of an arbitrary ANN with one hidden layer can be seen in Figure 1.

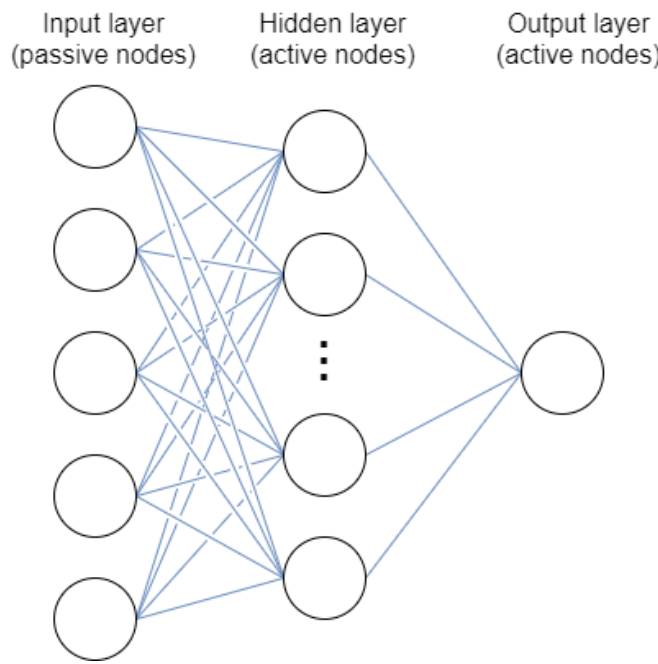


Figure 1. An arbitrary artificial neural network with one hidden layer

When the data is provided to an ANN for training, each value in an input layer is duplicated and each neuron in the hidden layers receives a copy. This is called a fully interconnected structure. After that, the values are multiplied by a set of predetermined numbers, also known as the weights, and then weighted inputs are added to produce a single number. This can be expressed as

$$y = f(x) = \sum_{i=1}^n x_i w_i + b$$

where x is the input, w is the weight and b is the bias, which is used to adjust the output. The number is passed through a non-linear function, also known as an activation function or a transfer function, which could be a sigmoid, rectifier, \tanh or any other activation function. This function takes a value between $-\infty$ and $+\infty$ and produces an output within the limits of the used function, which specifies if the neuron should fire or not. [16]

This work uses the state-of-the-art rectified linear unit (ReLU) as the activation function for a couple of reasons. First is the ease of calculation compared to other activation functions, since ReLU is a simple $f(x) = \max(0, x)$, while, for example, sigmoid uses expensive operations, such as exponentials: $f(x) = \frac{1}{1+e^{-x}}$. The other reason to use ReLU is because it returns the value of 0 more often than sigmoid or \tanh , which makes the neural network sparser, since fewer neurons fire, which more realistically simulates a biological neural network and allows for the smaller number of calculations needed.

The training of an ANN can be supervised, unsupervised or a mixture of both. Unsupervised training needs a lot more data, but the data can be “unlabeled”, which means it can lack categorization. This eliminates the need to label the data by hand, however, that also means that the network should have an effective algorithm to use the unlabeled data efficiently. The problem with this approach is that the accuracy of the said algorithm’s output cannot be evaluated. However, this way of training is closer to the true artificial intelligence and can produce the results which humans do not take into consideration due to pre-existing biases.

Supervised learning is done using “labeled” data. The input is presented to the network as the pairs of data and desired outputs. The goal of this approach is to analyze the training data and produce an accurate function, which can be used both for expected and unexpected inputs to get reasonable outputs. In this work, the neural network training will be supervised.

2.1.2. Backpropagation

During the ANN training, there will be a discrepancy between the expected and the actual outputs. This discrepancy is an error, which will be used to update the neuron weights for a better result.

It is not possible to get around this issue and choose “perfect” neuron weights during the initialization step of the neural network due to how the weights are calculated. A common

practice for choosing the initial weights since 1986, when Rumelhart and McClelland published their classic “Parallel Distributed Processing”, is to generate random non-zero numbers, which are very small. [17] As the weights are randomly generated, they will not produce the best immediate results and will need to be adjusted. Here is when backpropagation comes in.

Backpropagation, also known as backward propagation of errors or a feedback step, is a method of calculating an error made by each neuron in a network after the data is processed, to then send the values to an optimization method, such as gradient descent, which will adjust the neuron’s weight accordingly. This process requires multiple iterations, also known as epochs, to reach its highest potential to reduce the difference between the expected and the actual output of the network. It is most commonly used in deep neural networks.

2.1.3. Optimization algorithms

As it was mentioned above in section 2.1.2, the optimization algorithm adjusts the ANN’s neuron weights to produce more accurate results. There are two types of such algorithms: first order optimization algorithms and second order optimization algorithms, where the order specifies the derivative order.

First order algorithms minimize the error using the gradient values with respect to the parameters, with an example being the gradient descent algorithm. The regular (batch) gradient descent will perform an update after calculating the gradient of the whole data, which makes it very slow and inefficient if the data is too large. There is another variant of gradient descent called stochastic gradient descent (SGD), which is more popular, as it calculates the gradient for each training input data. The SGD algorithms which are most commonly used are momentum, Nesterov accelerated gradient, AdaGrad, AdaDelta and Adam, with the last being used in this work.

The second order optimization algorithms depend on the second order derivatives and, thus, are not that popular now, due to high cost of computation in terms of time and memory. However, they have a potential to outperform the first order algorithms if their derivatives are calculated in advance. Second order algorithms include Newton’s method, Quasi-Newton and Broyden–Fletcher–Goldfarb–Shanno (BFGS) algorithm.

2.1.4. Adam optimization algorithm

Adam, short for adaptive moment estimation, is a method which computes the adaptive learning rates for different parameters. The learning rate is a measure of how fast the network arrives to the final solution, so if the learning rate is too high, the network might fluctuate around the solution without reaching it or miss it completely (diverge). If the learning rate is too low, however, the network will take a very long time to reach the final solution (converge), so the best practice is to find a good middle ground.

Adam was introduced in 2015 as a SGD method that only requires first-order gradients with little memory requirement and it was designed to combine the advantages of two other popular

optimization algorithms: AdaGrad, which works well with the sparse gradients and Root mean square propagation (RMSProp), which does well in realistic settings. [18]

At this moment Adam is a state-of-the-art algorithm that compares favorably to other SGD methods, as it can be seen on Figure 2.

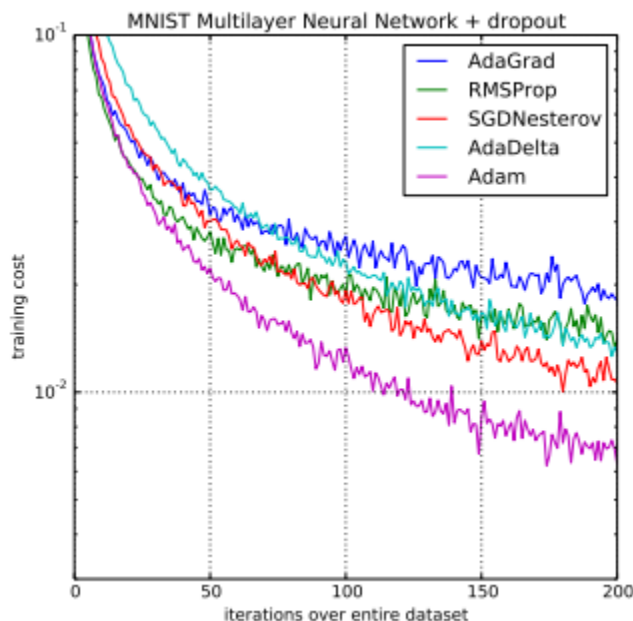


Figure 2. Comparison of Adam to other optimization algorithms training a multilayer neural network. Taken from Adam: A Method for Stochastic Optimization, 2015 [18]

2.2. Software

This work uses Python 3 programming language [19] and the publicly available libraries which can be found on “PyPI – the Python Package Index.” [20]

2.2.1. TensorFlow

TensorFlow [21] is an open-source library for numerical computation using data flow graphs. The graph nodes represent mathematical operations, while the graph edges represent the multidimensional data arrays (tensors) that flow between them. TensorFlow is available on all major operating systems, including Android and iOS. It was originally developed by Google Brain team for the purposes of conducting machine learning and deep neural networks research internally, but was released to public in 2015.

2.2.2. TFLearn

TFLearn [22] is a modular and transparent deep learning library for distributed machine learning built on top of TensorFlow. It features a higher-level API for implementing deep neural networks, as well as helper functions to train the TensorFlow graphs and offers easy graph visualization. It was created by Aymeric Damien in 2016.

2.2.3. NumPy

NumPy [23] is the fundamental Python package for scientific computations which is used massively. It contains N-dimensional array objects, tools for integrating C/C++ and Fortran code, as well as the implementations of high-level mathematical functions. The ancestor of NumPy, Numeric, was created by Jim Hugunin, which was later competing with another package, Numarray. Numarray was faster while doing operations with large arrays and Numeric was faster with small ones. In 2005, Travis Oliphant created NumPy to unify the community around a single package which had the best of both worlds.

3. Research

3.1. Neural network structure

The structure of the neural network that will be used the most throughout this work is as follows: an input layer, one hidden layer and an output layer. There will be other types of networks with more hidden layers, but those networks will only be used for comparison with the main one.

3.1.1. Input layer

The input layer receives the data, which it later passes on to the following layers without any changes. This data will be a list of five values, with first three values specifying if there are obstacles immediately to the left, front and right of the snake. The values can be either one (1) or zero (0), where zero means that there is no obstacle in specified direction and one means that the obstacle is there.

The fourth value of an input is an angle between snake's current direction vector and direction to an apple. This angle will be normalized and divided by 180 degrees, so that the resulting value ranges from -1 to 1.

The last value will be the direction the snake is suggested to take, which can be one of the three predetermined values: -1 specifies a turn to the left, 0 is to continue moving in the current direction and 1 is a turn to the right. This is what makes the neural network training supervised, as the network is told what outputs are expected for certain inputs.

3.1.2. Hidden layer

The hidden layer is the first layer where the values are expected to be modified, however the results of this layer will be sent to another hidden layer or to an output layer for additional calculations and, thus, will remain hidden outside the network, hence the name.

Since the neural network created for playing the Snake game does not have to be overly complicated, there was a decision to only have one single hidden layer with 25 neurons in it, with ReLU being the activation function.

The other network, however, will consist of two hidden layers with 100 neurons each and it will be seen if a deeper neural network with more neurons produces better results.

3.1.3. Output layer

The output layer produces the final results on the neural network and in our case the return value will be one of the three choices: -1 if the snake did not survive a turn, 0 if the snake has survived, but the direction of the snake is wrong in relation to the apple and 1 if the snake has survived and chose the correct direction in its last move.

This layer contains only one neuron and uses the linear activation function (the default one).

3.1.4. Code implementation

Described neural network is trivial to implement in Python using TFLearn package. The code is as follows:

```
network = input_data(shape=[None, 5, 1], name='input')
network = fully_connected(network, 25, activation='relu')
network = fully_connected(network, 1, activation='linear')
network = regression(network, optimizer='adam', learning_rate=0.01,
loss='mean_square', name='target')
model = tflearn.DNN(network)
```

3.2. Neural network training

The training of the neural network is a time and memory consuming process depending on the amount of games to train with. In our case, we will use different networks, which will be trained with a varying number of games and then the results of these networks will be compared after 5,000 test games are played by each.

The games are generated automatically, which means that each turn of the game will be chosen randomly out of three possible directions (left, front, and right), as going backwards is impossible, and if the snake is alive after the previous turn, new turn is generated until the game is over.

Since the computers used for generation of these initial sets of games were designed for personal use and were not designed to contain a big amount of data in the memory at the same time (there are around 6,000,000 moves for 100,000 games only), the training part of the network was designed to have an ability to load an existing network and continue training it with a new set of data. At first, however, this feature was absent, and the network was trained from scratch every time the training process was activated. This quickly became a problem due to the limitations of computer memory, as the network could not be trained with more than 100,000 games, because there was a memory error otherwise. Therefore, the network with 1,000,000 initial games was trained during 10 different training sessions of 100,000 games each.

Time spent on training the networks, as well as their performance comparison over the course of 5,000 test games, is in the table below. The networks are single layered and have 25 hidden layer neurons, unless stated otherwise and were trained with a different number of initial games. For the network with 1,000,000 training data the time was recorded over the course of 10 training sessions and calculated later. Time spent on generation of training data set is omitted, however it was the bottleneck part of the training and slowed the process immensely.

Table 1. Trained neural network performance comparison

Number of training games	1,000	10,000	100,000	1,000,000	100,000 (two-layered, 100 neurons per layer)
1st epoch (s)	4.401	35.271	362.135	3636.172	602.469
2nd epoch (s)	3.402	36.311	378.288	3843.051	614.335
3rd epoch (s)	4.144	36.273	376.790	3861.501	617.867
Total time (s)	11.947	107.855	1117.213	11340.724	1834.671
Average steps in 5,000 games	283.2438	263.3626	426.6786	400.9064	547.1236
Average score in 5,000 games	19.6092	18.1684	28.7722	27.3626	35.485
Worst score	1	2	4	2	6
Best score	53	60	75	64	88

As it can be seen from the table, the number games to train from is an important distinction between the networks and their performances. The less games there are, the worse the network performs, however, there is also a performance drop going from 100,000 games to 1,000,000. This could be attributed to the fact that the set of 5,000 games is a very small sample, however, the tests were run many times and the results were consistent across all runs.

The reason of this performance drop is most probably caused by a phenomenon known as the overtraining. Overtraining happens when there are too many iterations done on the training data and the network has increased performance over the training set, but loses an ability to generalize on the new data.

The best results were produced by the neural network of two layers. This is not surprising at all, because the state-of-the-art approach in neural network creation is to have multiple layers, with some well-performing networks having over 150! [24] Since operations which are performed by the network's neurons are very simple, combining them allows for much more complex calculations to be made possible. This, in turn, allows for a higher level of abstraction in deeper layers, where every next layer will recognize more and more features of the data set. An example is the image recognition. First layer of the neural network will only recognize the very simple things like edges, second layer will be trained to recognize more difficult shapes, such as ellipses

and circles, third layer will be able to differentiate between ears and eyes, while the fourth layer will already learn very complex features, like human faces or other objects.

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