# Understanding responses to environments for the Prisoner's Dilemma: A machine learning approach

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Submitted in partial fulfillment of the requirements for the degree of Doctor of Philosophy.



# Abstract

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# Acknowledgements

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### Dissemination of Work

### Publications (4 Published & 5 In preparation)

- 2018: Reinforcement learning produces dominant strategies for the Iterated Prisoner's Dilemma. Marc Harper, Vincent Knight, Martin Jones, Georgios Koutsovoulos, Nikoleta E. Glynatsi, Owen Campbell - PLOS One - Preprint arXiv:1707.06307
- 2018: An evolutionary game theoretic model of rhino horn devaluation. Nikoleta
   E. Glynatsi, Vincent Knight, Tamsin Lee. Ecological Modelling Preprint arXiv:1712.07640
- 2017: Evolution reinforces cooperation with the emergence of self-recognition mechanisms: an empirical study of the Moran process for the Iterated Prisoner's dilemma. Vincent Knight, Marc Harper, Nikoleta E. Glynatsi, Owen Campbell - PLOS ONE - Preprint arXiv:1707.06920
- 4. 2016: An open framework for the reproducible study of the Iterated prisoner's dilemma. Vincent Knight, Owen Campbell, Marc Harper et al Journal of Open Research Software

### IN PREPARATION

- 2019: A meta analysis of tournaments and an evaluation of performance in the Iterated Prisoner's Dilemma. Nikoleta E. Glynatsi and Vincent A. Knight - In preparation to be submitted - Preprint arXiv:2001.05911
- 2. 2019: A bibliometric study of research topics, collaboration and influence in the field of the Iterated Prisoner's Dilemma. Nikoleta E. Glynatsi and Vincent A. Knight - In preparation to be submitted - Preprint arXiv:1911.06128
- 3. 2019: Game Theory and Python: An educational tutorial to game theory and repeated games using Python. Nikoleta E. Glynatsi and Vincent A. Knight
   Submitted to the Journal of Open Source Education Available on GitHub Nikoleta-v3/Game-Theory-and-Python
- 4. 2019: A theory of mind: Best responses to memory-one strategies. The limitations of extortion and restricted memory. Nikoleta E. Glynatsi and Vincent A. Knight Submitted to Scientific Reports Nature Preprint arXiv:1911.12112
- 2019: Recognising and evaluating the effectiveness of extortion in the Iterated Prisoner's Dilemma. Vincent Knight, Marc Harper, Nikoleta E. Glynatsi, Jonathan Gillard - Submitted to Nature Communications - Preprint arXiv:1904.00973

### Talks & Posters

INVITED TALKS (KEYNOTES)

- How does a smile make a difference?, PyCon UK, Cardiff, 2018.
- The Fallacy of Meritocracy, PyCon Balkan, Belgrade, 2019.

#### OTHERS

- Accessing open research literature with Python PyCon Namibia, Namibia 2017.
- Writing tests for research software PyCon Namibia, Namibia 2017.
- Optimisation of short memory strategies in the Iterated Prisoners Dilemma Wales Mathematics Colloquium, Gregynog Hall 2017.
- PIP INSTALL AXELROD (poster) Euroscipy, Erlangen Germany 2017.
- Arcas: Using Python to access open research literature Euroscipy, Erlangen Germany 2017.
- A trip to earth science with python as a companion PyConUK, Cardiff 2017.
- The power of memory (poster) SIAM UKIE Annual Meeting, Southampton 2018.
- Rhinos with a bit of Python PyConNA, Namibia 2018.
- Memory size in the Prisoners Dilemma Wales Mathematics Colloquium, Gregynog Hall 2018.
- Memory size in the Prisoners Dilemma SIAM UKIE National Student Chapter, Bath University 2018.
- Stability of defection, optimisation of strategies and testing for extortion in the Prisoner's Dilemma (poster) STEM for Britain, London 2019.
- Stability of defection, optimisation of strategies and testing for extortion in the Prisoner's Dilemma - 18th International Conference on Social Behaviour, Sedona, Arizona 2019.
- An introduction to Time Series Joint workshop between CUBRIC & Mathematics Departments, Cardiff 2019.

### Software Development

- Arcas, an open source package designed to help users collect academic articles' metadata from various prominent journals and pre print serves. **Contribution:** Main developer
- Axelrod-Python library, an open source framework decided to the study of the Iterated Prisoner's Dilemma. **Contributions:** Implementation of spatial tournaments functionality, implementation/addition of strategies (from the literature) to the library, reviewing of code contributed by other contributors
- SymPy, a Python library for symbolic mathematics. **Contributions:** Implementation of Dixon's and Macaulay's resultants which were developed for my 2019 publication "Sta-

bility of defection, optimisation of strategies and the limits of memory in the Prisoner's Dilemma"

• Pandas, an open source library providing high-performance, easy-to-use data structures and data analysis tools. **Contribution:** Fix bug which converted NaN values to strings

# Contents

$\mathbf{A}$	bstra	ct		i
A	ckno	wledge	ements	iii
Di	issen	ninatio	n of Work	iv
1	Tra	ining l	ong short-term memory networks to play the Iterated Prisoner'	s
	Dile	emma		1
	1.1	Introd	luction	1
	1.2	Artific	cial, recurrent neural and long short-term memory networks	2
	1.3	Traini	ng LSTM networks to play the Iterated Prisoner's Dilemma	7
		1.3.1	Building the networks with Keras	9
		1.3.2	Training on GPU	11
		1.3.3	Training data sets	12
	1.4	Valida	ation of LSTM base strategies using a meta tournament analysis	15
		1.4.1	Fingerprinting the LSTM based strategies	16
	1.5	Chapt	er Summary	23

# List of Figures

1.1	Graphical representation of an ANN
1.2	Graphical representation of a RNN
1.3	An LSTM hidden layer at time step $t$
1.4	Graphical representation of an LSTM network
1.5	A graphical representation of a sequence to sequence LSTM model 8
1.6	A graphical representation of a sequence to probability LSTM model 8 $$
1.7	Keras sequence to sequence
1.8	Keras sequence to probability
1.9	Validation of sequence to sequence
1.10	Validation of sequence to probability

# List of Tables

1.1	The interactions between Adaptive and one of the best response sequences to	
	the strategy	7
1.2	The interactions between Adaptive and one of the best response sequences to	
	the strategy were $0 \to D$ and $1 \to C$	7
1.3	Subset of data set that the networks are trained on. The APL project performs	
	a round robin tournament with all the strategies implemented on every release.	
	The top strategies and the strategies across are based on the results of the	
	tournament with APL version 4.5.0	12
1.4	Number of epochs each both networks, sequence to sequence and sequence to	
	probability, were trained for each subset	12
1.5	Median normalised ranks of th 24 LSTM strategies over standard tournaments.	16

### Chapter 1

# Training long short-term memory networks to play the Iterated Prisoner's Dilemma

The research reported in this Chapter has been carried out with:

Axerod-Python library version: 4.2.0 Associated data set:

### 1.1 Introduction

In Chapter ?? it was mentioned that consepsualating and introducing new strategies has been an important aspect of research to the field. The aim of this Chapter is to introduce new IPD strategies based on an archetype that has not received much attention in the literature.

In Chapter ?? it was concluded that one of the properties successful strategies in a IPD competition need to have is cleverness/complexity. Complexity can confer to adaptability, and adaptability is important in performing well in diverse set of environments. This was established not only in Chapter ?? but also from the results of Chapter ??. The set of complex strategies that ranked highly across distinct tournaments in Chapter ?? included strategies based on archetypes such as finite state automata, hidden Markov models and artificial neural networks (ANNs).

ANNs have successfully been trained to play games exterior to the IPD, such as checkers [4] and chess [8]. The usage of feed forward ANNs in the IPD was firstly introduced in [11], and have been used ever since [1, 2, 7, 9]. Possibly, the most successful ANNs strategies are the one introduced in [10], as they were ranked  $7^{th}$ ,  $9^{th}$ , and  $11^{th}$  in a tournament of 223 strategies.

A type of ANNs that have not received much attention in the literature are the recurrent neural

networks (RNNs). RNNs are a type of neural networks that include a feedback connection and are designed to work with inputs in the form of sequences. RNNs were firstly considered as an archetype in 1996. In [21] a RNN which considered a single previous step as an input was trained via a Q learning algorithm to play against a single strategy. The results of [21] were limited only in the network learning to play against a single strategy.

The limitations of [21] could potentially have been due to the limitations of the RNNs themselves. As it will be discussed later the RNNs quickly became unable to learn due to the vanishing gradient problem. To improve on the standard recurrent networks a new network called the long short-term memory networks (LSTMs) were introduced in [12]. LSTMs do not suffer from the vanishing gradient, can be successfully trained and are being used in a number of innovative applications such as time series analysis [15], speech recognition [20] and prediction in medical care pathways [22].

LSTMs have not received attention in the IPD literature. This Chapter trains and introduced a number of strategies based on LSTMs. The training has been possible due to the collection of best response sequences generated in Chapter ??. The performance of the new strategies is evaluated and compared in a meta tournament analysis of standard tournaments. The results demonstrate

This Chapter is structured as follows:

- section 1.2, presents an introduction to artificial, recurrent and long short-term memory networks.
- section 1.3, cover the specifics of the LSTMs trained in the Chapter. Their architectural details are presented as well as the different training sets they have been trained on.
- section 1.4, evaluates and compares the performance of 18 LSTMs based strategies in 300 standard IPD computer tournaments.

# 1.2 Artificial, recurrent neural and long short-term memory networks

ANNs are computing systems vaguely inspired by the biological neural networks that constitute animal brains. As stated in [13] the research on ANNs has experienced three periods of extensive activity. The first peak was in the 1940s. The work of [17] opened the subject by creating a computational model for neural networks. The second peak occurred in 1960s with the introduction of the perceptron [19]. The perceptron is a linear binary classifier and in the 1960s in was shown that it could be used to learn to classify simple shapes with  $20 \times 20$  pixels input. However, it was impossible for the perceptron to be extended to classification tasks with many categories. The limitations of the model were presented in [18] which alted the enthusiasm of most researchers in ANNs, resulting in no activity in field for almost 20 years. The third peak occurred in the early 1980s with the introduction of the back propagation algorithm for multilayered networks [23]. Though the algorithm was initially proposed by [23] it has been reinvented several times and it was popularizes by [16]. Following the introduction of the algorithm and the ability to now train more complex models, ANNs have received considerable interest, and have been used in a number of innovative applications such as [6, 14].

ANNs based on the connection pattern/architecture can be grouped into two categories:

- Feed forward networks.
- Recurrent or feedback networks.

ANNs can be viewed as weighted directed graphs in which artificial neurons are nodes and directed edges, with weights, are connections between the neuron outputs and the neuron inputs [13]. In feed forward networks the graphs have no loops whereas in recurrent loops occur because of feedback connections. A graphical representation of a feed forward network with a single hidden layer is given by Figure 1.1.

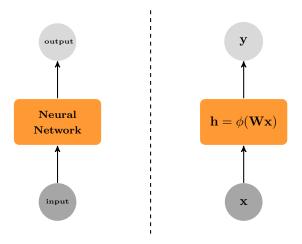


Figure 1.1: Graphical representation of an ANN.

A feed forward network is composed by an input layer, hidden layers, and an output layer. The networks' input is a vector x. The dimensionality, or number of nodes of the input layer is depended on the dimensionality of x. Each element of the input vector is connected to the hidden layer via a set of learned weights. The hidden layer also consists of nodes and the number of node/dimensionality of it is an architectural decision. The  $j^{\text{th}}$  hidden neuron outputs,

$$h_j = \phi(\sum_i w_{ij}x_i)$$
, where  $\phi$  is an activation function.

Common activation functions for  $\phi$  include:

- the sigmoid function  $\sigma(x)$  which squashes numbers into the range [0, 1]
- the hyperbolic tangent, tanh(x) which squashes numbers into the range [-1,1]
- the rectified linear unit, ReLU(x) = max(0, x)

In turn, the hidden layer is fully connected to an output layer. The  $j^{\rm th}$  output neuron outputs,

$$y_j = \sum_i v_{ij} h_i.$$

The output layer transforms the raw scores to probabilities via a softmax function.

Feed forward networks make predictions using forward propagation,

$$h = \phi(Wx) \tag{1.1}$$

$$y = \operatorname{softmax}(Vh) \tag{1.2}$$

where W is a weight matrix connecting the input and hidden layers, V is a weight matrix connecting the hidden and output layers. W and V are the learning parameters of the network. Their dimensionality depends on the dimensionality of networks layers. If x and x are 2 dimensional and the hidden layer had 500 nodes then  $W \in \mathbb{R}^{2 \times 200}$  and  $V \in \mathbb{R}^{2 \times 200}$ . Increasing the dimensionality of the hidden layer corresponds to more parameters.

Training a ANN corresponds to finding the learning parameters that minimise the error on the training data. The function that measures error is called the *loss function*. A common loss function with the choice of a softmax output is the *cross entropy* loss. Consider N training examples and K classes, the function for a prediction  $\hat{y}$  with respect to the true output y us give by,

$$L(y, \hat{y}) = -\frac{1}{N} \sum_{n \in N} \sum_{k \in K} y_{n,k} \log(y_{\hat{n},k})$$

The minimum of the loss function is calculated using the gradients decent algorithm. The gradient decent algorithm relies on the gradients. The gradients are the vector of derivatives of the loss function with respect to the learning parameters. Thus,  $\frac{\partial L}{\partial W}$  and  $\frac{\partial L}{\partial V}$ . These gradients are calculated with the back propagation algorithm [24], which is a way to efficiently calculate the gradients starting from the output.

An extension of feed forward networks are the recurrent networks. RNNs are capable of processing variable length sequences of inputs. More importantly, RNNs use their internal state/knowledge to a make a prediction on the input at time t based on the decisions made at the previous time steps. A graphical representation of a RNN is given by Figure 1.2. RNN are networks with loops. The loop represents that information can be passed down from one step of the network to the next. In fact they can be thought of as multiple copies of the same network, each passing a message to a successor, as demonstrated by Figure 1.2.

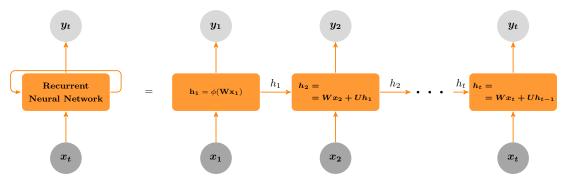


Figure 1.2: Graphical representation of a RNN.

The hidden layers pass down information by considering,

$$h_t = \phi(Wx_t + Uh_{t-1})$$

where  $h_{t-1}$  is the hidden state computed at time t-1 multiplied by some weight vector U. In matrix notation RNNs are described by,

$$h_t = \phi(Wx_t + Uh_{t-1}) \tag{1.3}$$

$$y_t = Vh (1.4)$$

Unfortunately, in practice RNNs quickly become unable to learn to connect the information. The more time steps the higher the chance that the back propagation gradient either accumulate and explode or vanish down to zero. The fundamental difficulties of RNNs were explored in depth by [3].

A network specifically designed to avoid the long-term dependency problem. was introduced by [12] called the long short-term memory network (LSTM). LSTMs work tremendously well on a large variety of problems, and are now widely used. Remembering information for long periods of time is practically their default behavior, not something they struggle to learn.

The core idea to LSTMs is the *cell state* also referred to as the long term memory, denote as  $C_t$ . The cell state is designed to pass down information with only a few carefully regulated changes being applied to it by structures called *gates*. Gates are composed out of a sigmoid neural net layer and a point wise multiplication operator. In order to explain the cell gate and subsequently how LSTMs make predictions consider a LSTM's hidden layer at time step t give by Figure 1.3.

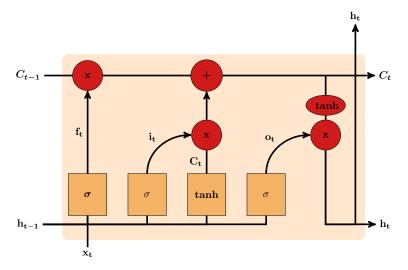


Figure 1.3: An LSTM hidden layer at time step t.

The cell state and hidden state from from the previous time step are feed back into the network. Initially, the network decides what information from the previous cell state is to be discarded. This decision is made by the *forget state*. The forget state considers the hidden state at time

t-1 and the input at time t,

$$f_t = \sigma(W_f x_t + U_f h_{t-1}) \tag{1.5}$$

Secondly, the network decides what information is going to be stored at the cell gate. There are two parts to this. Firstly, the input gate decides which values are going to be updated,

$$i_t = \sigma(W_i x_t + U_i h_{t-1} + b_i). (1.6)$$

and secondly, a tanh layer creates a vector of candidate values denoted as  $\tilde{C}_t$  that could be added to the cell state.  $\tilde{C}_t$  is given by Equation (1.9).

$$\tilde{C}_t = tanh(W_c x_t + U_c h_{t-1} + b_c).$$
 (1.7)

The cell state  $C_{t-1}$  is multiplied by  $f_t$ , forgetting the values which have been decided to discard. The new candidate values are scaled by how much information has been decided to keep from the input. Thus,

$$C_t = f_t * C_{t-1} + i_t * \tilde{C}_t. \tag{1.8}$$

The LSTM outputs a hidden state at each time step. This is either used to through an activation function to output  $y_t$ , and it is passed to the next time step cell. The output is based on the current cell state. The cell state goes through a tanh function and it is multiplied by a sigmoid gate that decides which parts to output from the cell state.

$$o_t = \sigma(W_o x_t + U_o h_{t-1} + b_o), \tag{1.9}$$

$$h_t = o_t * tanh(C_t) \tag{1.10}$$

This process is being carried out though each time step of the input sequence. At each the cell state and hidden state are feed back into the network and the hidden state can also be used to make a prediction for each time step as demonstrated by Figure 1.4.

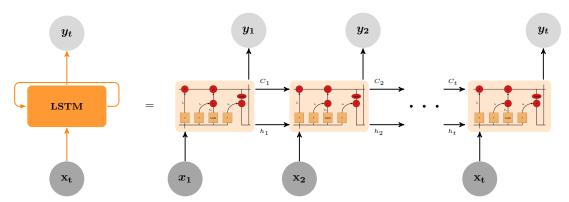


Figure 1.4: Graphical representation of an LSTM network.

The LSTM is a complex architecture and contains a large number of parameters that need to be trained. The LSTM networks have managed to achieve remarkable results. The training of an LSTM model in this Chapter is carried with the open source project Keras. The details for the training are presented in section 1.3.

# 1.3 Training LSTM networks to play the Iterated Prisoner's Dilemma

The LSTMs of this Chapter are trained based on the best response sequence data set that was generated in Chapter ??. Consider the strategy Adaptive and a best response to it as presented in section ??.

	1	2	3	4	5	6	7	8	9	10	11	<b>12</b>	 204	205
Adaptive	C	C	C	C	C	C	D	D	D	D	D	C	 C	C
$S^*$	C	C	D	D	D	D	D	D	D	D	D	D	 D	D

Table 1.1: The interactions between Adaptive and one of the best response sequences to the strategy.

For  $0 \to D$  and  $1 \to C$  Table 1.1 is given by,

	1	2	3	4	5	6	7	8	9	10	11	12	 204	205
Adaptive	1	1	1	1	1	1	1	0	0	0	0	1	 1	1
$S^*$	1	1	0	0	0	0	0	0	0	0	0	0	 0	0

Table 1.2: The interactions between Adaptive and one of the best response sequences to the strategy were  $0 \to D$  and  $1 \to C$ .

In order to train the LSTM as a player the network should predict the best response to strategy's action. Thus, the input to the network is the actions of the opponents Q, discarding the last turn.

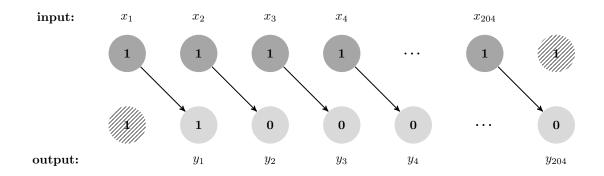


and the expected output it's the response of the  $S^*$  to those actions. Thus the genes from t=2 instead of t=1.

Two variants of LSTM network have been training in this Chapter. These are referred to:

- sequence to sequence
- sequence to probability

The sequence to sequence model as an input of actions a each time step by a strategy an it outputs a response for each at each time step. The sequence to probability has an input the action of strategy up to time t, and outputs a response to the history at time t.



A graphical representation of the two networks are given by Figures 1.5 and 1.6.

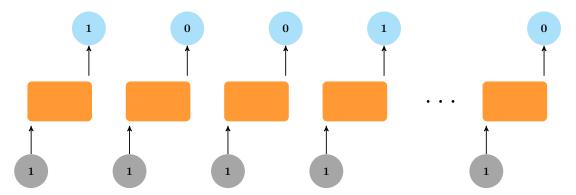


Figure 1.5: A graphical representation of a sequence to sequence LSTM model

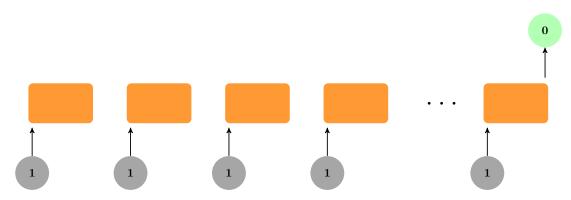


Figure 1.6: A graphical representation of a sequence to probability LSTM model

The networks are constructed in such way as the length of sequences they receive it is not fixed. The are not trained for the 5258 but each best response sequence is broken down to 204 sets

of input.

### 1.3.1 Building the networks with Keras

To construct and train the network model the open-source neural-network package Keras [5] is used. Its library written in Python, designed to enable fast experimentation with neural networks. Neural layers, cost functions, optimizers, initialization schemes, activation functions, and regularization schemes are all standalone modules that you can combine to create new models.

The code for implementing the sequence to sequence model is given by Figure 1.7. There are two main types of models available in Keras: the Sequential model, and the Model class. The Sequential model is a linear stack of layers. There is a single layer of an LSTM model that returns the hidden state at each time step. The input shape is (None, 1) because there is not a fixed length and there is a single time step between the of the input. There are 100 the outputs turn into a probability with a sigmoid . The dropout layer There are a total of 41301 parameters that need to be trained for the sequence to sequence model.

Similarly, for the sequence to probability network the implementation of the network using Keras is given by Figure 1.8. There are now two LSTM layers. The first one outputs a hidden state at each time step and the second layer outputs only at time t the output then goes

```
>>> from keras.models import Sequential
1
    >>> from keras.layers import (
     ... Dense,
4
           Dropout,
         CuDNNLSTM,
    ...)
    >>> num_hidden_cells = 100
10
    >>> drop_out_rate = 0.2
11
    >>> model = Sequential()
12
    >>> model.add(
14
15
     ... CuDNNLSTM(
             num_hidden_cells, return_sequences=True, input_shape=(None, 1))
16
    ...)
17
18
    >>> model.add(Dropout(rate=drop_out_rate))
19
20
     >>> model.add(Dense(1, activation="sigmoid"))
21
22
    >>> model.summary()
    Model: "sequential_1"
24
25
     Layer (type)
                             Output Shape
26
    cu_dnnlstm_1 (CuDNNLSTM) (None, None, 100) 41200
27
     dropout_1 (Dropout) (None, None, 100) 0
29
     dense_1 (Dense) (None, None, 1) 101
31
32
     Total params: 41,301
33
34
     Trainable params: 41,301
     Non-trainable params: 0
36
```

Figure 1.7: Keras sequence to sequence.

through Dropout then turned into probability with a sigmoid function. There are total of 122101 parameters.

```
>>> from keras.models import Sequential
     >>> from keras.layers import (
            Dense,
             Dropout,
             CuDNNLSTM.
5
     >>> num hidden cells = 100
     >>> drop out rate = 0.2
10
11
     >>> model = Sequential()
12
13
     >>> model.add(
14
             CuDNNLSTM(num_hidden_cells, return_sequences=True, input_shape=(None, 1))
15
16
     >>> model.add(CuDNNLSTM(num_hidden_cells))
17
18
     >>> model.add(Dropout(rate=drop_out_rate))
19
20
     >>> model.add((Dense(1, activation="sigmoid")))
21
     >>> model.summary()
     Model: "sequential 2"
22
23
     Layer (type)
                                Output Shape
                                                          Param #
^{24}
25
     cu_dnnlstm_2 (CuDNNLSTM) (None, None, 100)
                                                          41200
27
                                                           80800
28
     cu_dnnlstm_3 (CuDNNLSTM) (None, 100)
29
30
     dropout_2 (Dropout) (None, 100)
                                                           Ω
31
     dense 2 (Dense)
32
                               (None, 1)
                                                           101
33
     Total params: 122,101
34
35
     Trainable params: 122,101
     Non-trainable params: 0
36
37
```

Figure 1.8: Keras sequence to probability.

Keras contains two implementations on an LSTM model. That is the class LSTM and the CC-NLSTM which was used here. CCNLSTM is a fast LSTM implementation with NVIDIA CUDA Deep Neural Network library (cuDNN) which run on a graphics processing unit (GPU).

### 1.3.2 Training on GPU

Conventionally most training and running computer code happens on the central processing unit (CPU). A CPU also called a central processor or main processor is essentially the brain of any computing device, carrying out the instructions of a program by performing control, logical, and input/output (I/O) operations. CPUs, can have a number of cores, and its basic design and purpose—to process complex computations—has not really changed. A CPU core is designed to support an extremely broad variety of tasks.

A graphical processing unit (GPU), on the other hand, has smaller-sized but many more logical cores (arithmetic logic units or ALUs, control units and memory cache) whose basic design is to process a set of simpler and more identical computations in parallel.

Initially designed mainly as dedicated graphical rendering workhorses of computer games, GPUs were later enhanced to accelerate other geometric calculations (for instance, transforming polygons or rotating verticals into different coordinate systems like 3D). Nvidia created a parallel computing architecture and platform for its GPUs called CUDA, which gave developers access and the ability to express simple processing operations in parallel through code.

While a CPU core is more powerful than a GPU core, the vast majority of this power goes unused by ML applications. whereas a GPU core is optimized exclusively for data computations. Because of this singular focus, a GPU core is simpler and has a smaller die area than a CPU, allowing many more GPU cores to be crammed onto a single chip. Consequently, ML applications, which perform large numbers of computations on a vast amount of data, can see huge (i.e., 5 to 10 times) performance improvements when running on a GPU versus a CPU. As many have said GPUs are so fast because they are so efficient for matrix multiplication.

Keras CCNLSTM class makes running on GPU easy. Cardiff's super computer was used for the training of the models.

### 1.3.3 Training data sets

It was explained in the the best response data set from Chapter 6 is used as training data. Moreover, each best responses has broken down to 204 different inputs. In order to understand the difference the effect of the training data the models have been trained on different subsets of the data set. More specifically in total of four different sub sets. These with a brief explanation and number are given by Table 1.3.

Data set	number of $Q$	Explanation	number of $S^*$
all strategies	192	The data set as presented in Chapter ??	5258
top performing strategies	18	Top 18 performing strategies in a standard tournaments performed with APL $$	714
strategies across ranks	15	A set 15 that there rank is across	212
basic strategies	11	A set 11 which are classified as the basic in APL	84

Table 1.3: Subset of data set that the networks are trained on. The APL project performs a round robin tournament with all the strategies implemented on every release. The top strategies and the strategies across are based on the results of the tournament with APL version 4.5.0.

Due to the different size of the data set the networks have trained for a different number of epochs. The number of epochs is the number of complete passes through the training data set. The number of epochs each network has trained for each data set is given by Table 1.4.

	all strategies	top strategies	strategies across ranks	basic strategies
sequence to sequence	19999	83999	39999	129999
sequence to probability	7999	74999	39999	84999

Table 1.4: Number of epochs each both networks, sequence to sequence and sequence to probability, were trained for each subset.

The data are split into training and test training set. The test set is 20% of the data. The validation and loss is calculated at each epoch. The validation is the measure that. The loss is based on the.

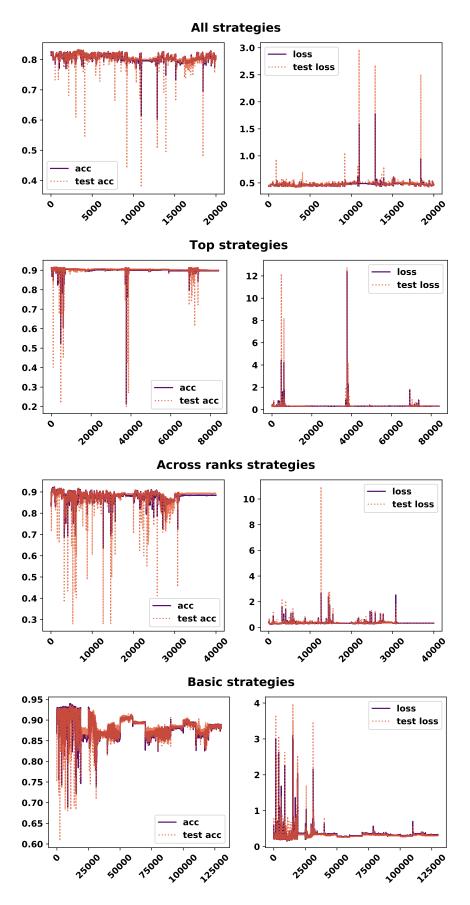


Figure 1.9: Validation of sequence to sequence

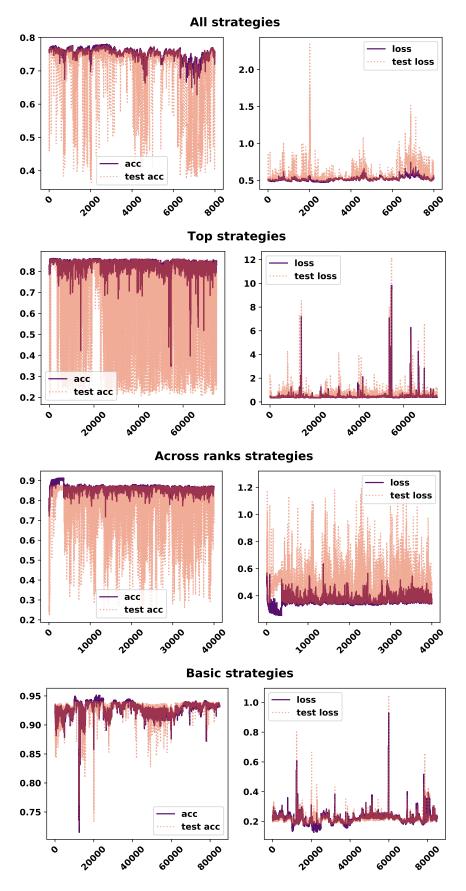


Figure 1.10: Validation of sequence to probability

Discussion on validation. In section 1.4 the LSTM networks are validated by as an IPD strategy using a meta tournament analysis.

### 1.4 Validation of LSTM base strategies using a meta tournament analysis

To evaluate the LSTM players as strategies, a strategy class was implemented that takes as an argument the a Keras model. The player then uses the history of the match translate it to an input in the form that the networks can then translate into an action. Note that the network was not trained to consider it's opening move. The opening move is also an input to the strategy class. In the meta tournament analysis different strategies that have a different opening play are considered.

```
import numpy as np
     import axelrod as axl
 3
 4
     from axelrod.random_ import random_choice
     from keras.layers import LSTM, Dense, Dropout
     from keras.models import Sequential
     C, D = axl.Action.C, axl.Action.D
10
     class LSTMPlayer(ax1.Player):
11
       name = "The LSTM player"
12
         classifier = {
13
              "memory_depth": float("inf"),
14
              "stochastic": True.
15
16
             "inspects_source": False,
17
              "manipulates_source": False,
              "manipulates_state": False,
         }
19
20
21
         def __init__(self, model, reshape_history_funct, opening_probability=0.78):
             self.model = model
22
              self.opening_probability = opening_probability
             self.reshape_history_function = reshape_history_funct
24
25
              super().__init__()
             if opening_probability in [0, 1]:
26
                  self.classifier["stochastic"] = False
27
         def strategy(self, opponent):
29
             if len(self.history) == 0:
31
                  return random_choice(self.opening_probability)
32
              history = [action.value for action in opponent.history]
              prediction = float(
34
35
                  {\tt self.model.predict(self.reshape\_history\_function(history))[0][-1]}
36
37
              return axl.Action(round(prediction))
39
          def __repr__(self):
40
41
              return self.name
```

The implemented class can interact with any opponent from APL. An example of usage of the LSTMPlayer is given by Figure.

To evaluate the performance of the LSTM strategy a meta tournament analysis, similar to Chapter ??, is carried out. The tournament type considered here are standard tournaments.

The data collections is given by Algorithm ??. At each trial a number of opponents, a list of opponent From 5 to 10. The number of turns and repetitions are fixed to 200 and 50.

The number 205 was hard coded. The evaluation does not want to take into account the last turns. 205 was selected so the last turns could have beed discarded and the match length would be 200 which is the number of turns commonly used in the literature.

### Algorithm 1: Data collection Algorithm

return result standard;

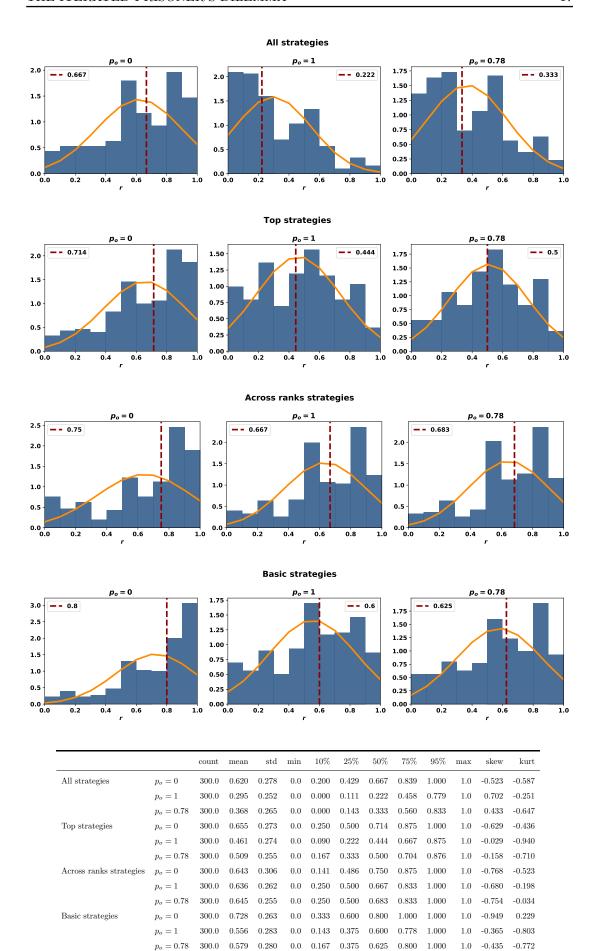
A total of 300 trials of Algorithm 1 have been performed. Similarly to Chapter ?? each trial outputs a result summary of the tournament. The performance of the strategies are evaluated on the normalised rank r, and more specifically on the median normalised rank  $\bar{r}$ . A total of different strategies competed in the tournaments, and the strategy with the maximum participation, excluding the LSTM strategy, was.

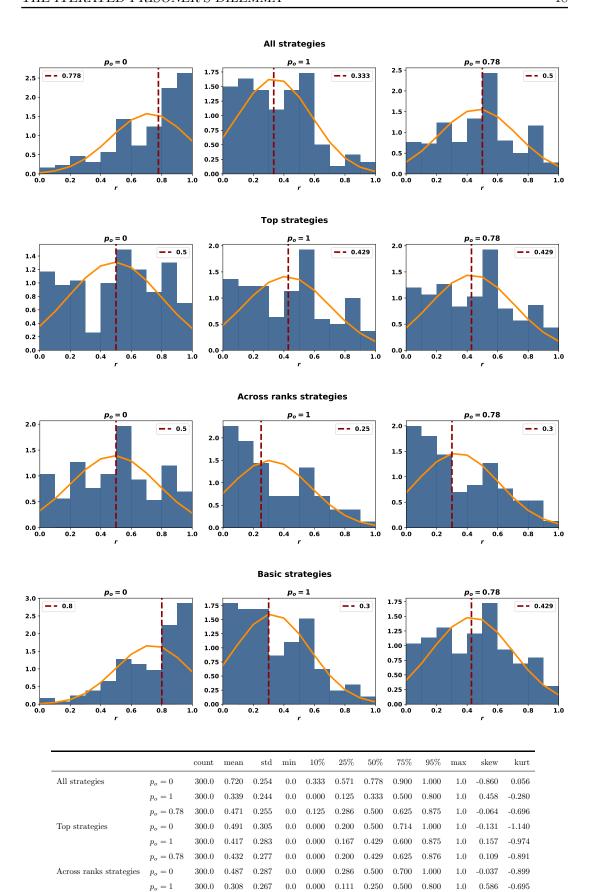
There are two distinct networks that have been trained on four different data sets. Moreover, each networks translates into a player that the opening probability can have three different values. Thus a total of  $2 \times 4 \times 3$  LSTM players performance is evaluated in this section.

	seque	ence to s	equence	sequence to probability				
	$p_o = 0$	$p_o = 1$	$p_o = 0.78$	$p_o = 0$	$p_o = 1$	$p_o = 0.78$		
All strategies	0.667	0.222	0.333	0.778	0.333	0.500		
Top strategies	0.714	0.444	0.500	0.500	0.429	0.429		
Across ranks strategies	0.750	0.667	0.683	0.500	0.250	0.300		
Basic strategies	0.800	0.600	0.625	0.800	0.300	0.429		

Table 1.5: Median normalised ranks of th 24 LSTM strategies over standard tournaments.

#### 1.4.1 Fingerprinting the LSTM based strategies





 $p_o=0.78$ 

 $p_o = 0.78$ 

 $p_o = 0$ 

Basic strategies

300.0

290.0

290.0

290.0

0.335

0.738

0.323

 $0.432 \quad 0.269$ 

0.272

0.239

0.249

0.0 - 0.000

0.0 - 0.400

0.0 - 0.000

0.0 0.000

0.125

0.600

0.125

0.200

0.300

0.800

0.300

0.429

0.556

0.975

0.500

0.625

0.800

1.000

0.778

0.875

1.0

1.0 -0.881

1.0

1.0

0.465

0.491

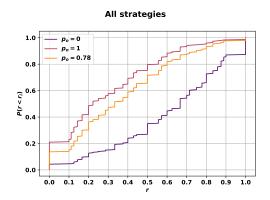
0.096

-0.867

0.315

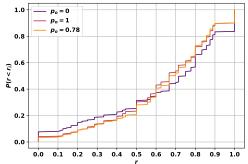
-0.535

-0.916



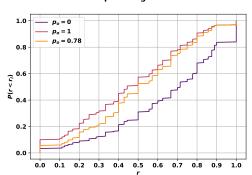
(a) P(r < 0.5) : 0.350 P(r < 0.5) : 0.797 P(r < 0.5) : 0.717

### Across ranks strategies



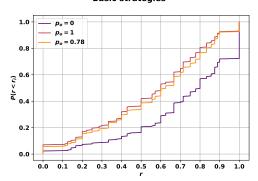
(c) ( P(r<0.5):0.313P(r<0.5):0.303P(r<0.5):0.280)

#### Top strategies

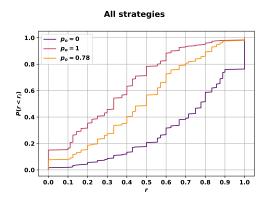


(b) P(r < 0.5) : 0.310 P(r < 0.5) : 0.573 P(r < 0.5) : 0.523

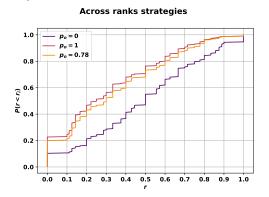
### Basic strategies



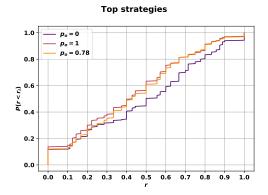
(d) P(r<0.5):0.207P(r<0.5):0.420P(r<0.5):0.387



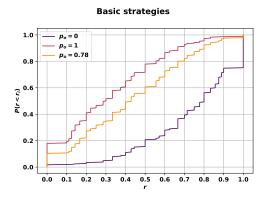
(a) P(r < 0.5) : 0.207 P(r < 0.5) : 0.783 P(r < 0.5) : 0.567



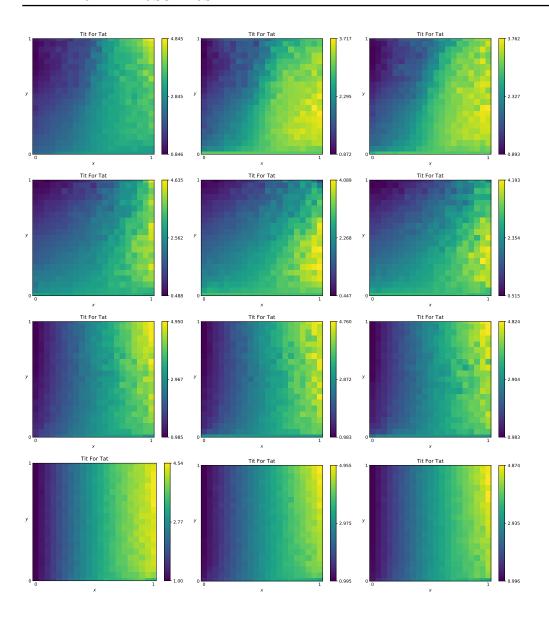
(c) ( P(r<0.5):0.550P(r<0.5):0.763P(r<0.5):0.733)

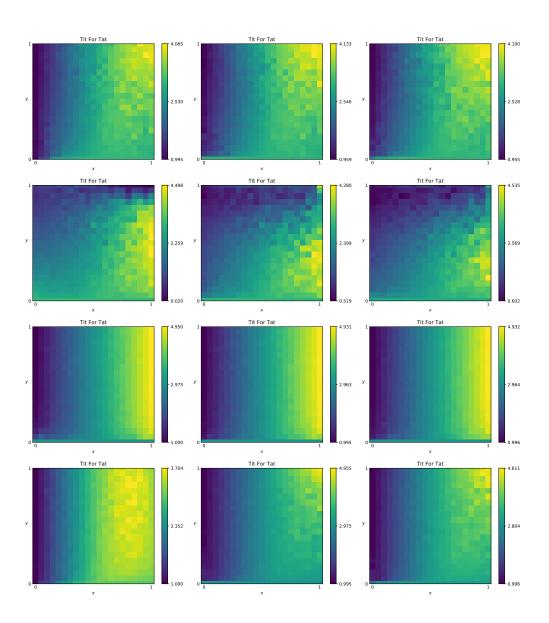


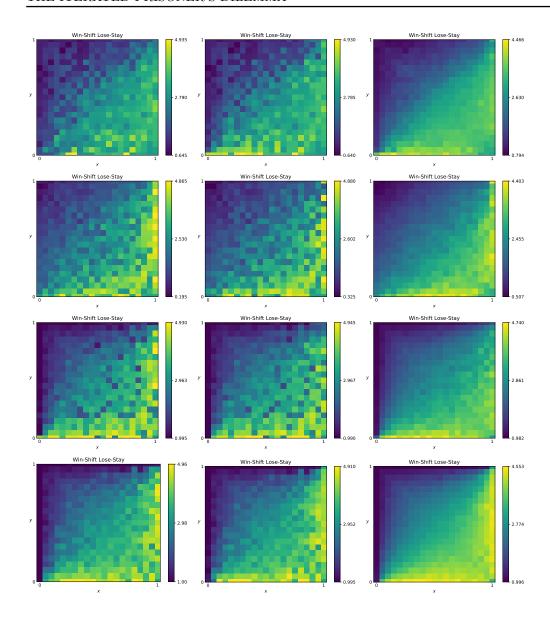
(b) P(r < 0.5) : 0.503 P(r < 0.5) : 0.633 P(r < 0.5) : 0.610



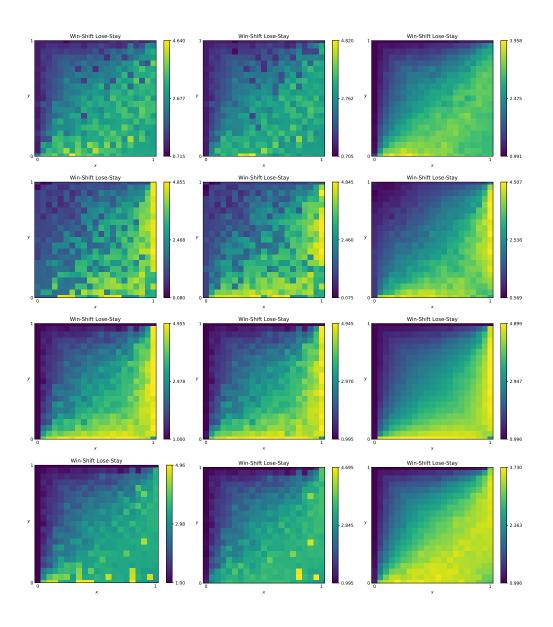
(d) P(r < 0.5) : 0.207, P(r < 0.5) : 0.779, P(r < 0.5) : 0.607



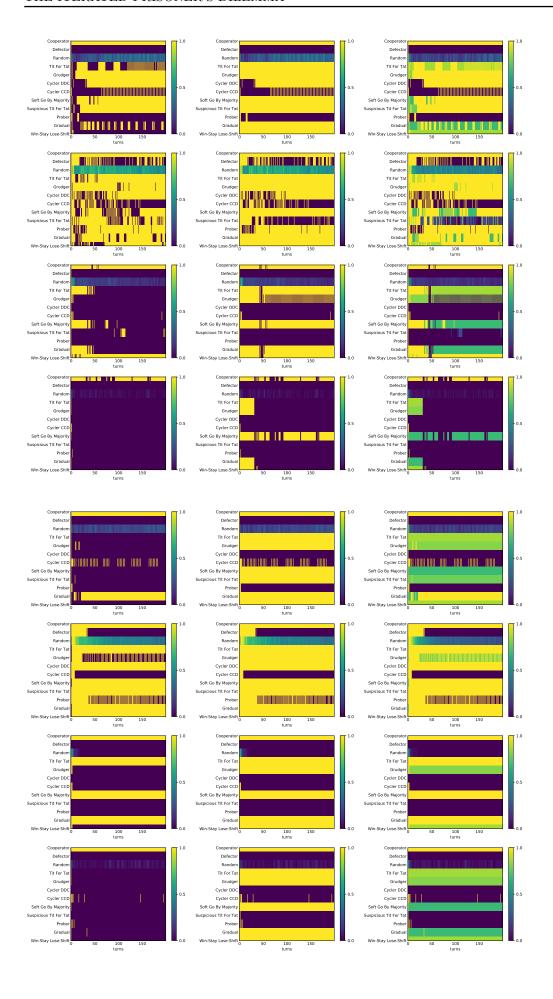




# 1.5 Chapter Summary







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