CAPTURING THE PREDICTIVE POWER OF CORTICAL LEARNING ALGORITHMS

By

ALEXANDER C MICHELS

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Westminster College, New Wilmington, PA

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advised by

Carolyn Cuff, Ph.D.

C. David Shaffer, Ph.D.

William Procasky, Ph.D.

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ABSTRACT iii

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Hierarchical Temporal Memory (HTM) is model of intelligence based on the the interactions of pyramidal neurons in the mammalian neocortex which is currently being developed by Numenta. It has stood out due to high noise tolerance and learning capacity which makes it well-suited for anomaly detection, prediction, signal processing, and sensorimotor processes. We seek to find a mathematical analogy to the predictive power of this biologically constrained theory using models from time series analysis. To do this, we statistically analyzed the predictions of HTM networks which were asked to predict outputs of autoregressive moving average (ARMA) models, varying the parameters of both the networks and the ARMA models. We hope to to find a relation between sets of HTM network parameters and ARMA model parameters to better explain cortical learning algorithms and the neocortex.

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Chapter 1

Introduction

Although we are continually bombarded with sensationalist news stories proclaiming the dawn and dangers of "artificial intelligence," it is important to first define what it means to say a machine is intelligent. This is much the same way Turing began his preponderance of artificial intelligence in his 1950 Computing Machinery and Intelligence [33]. Turing came up with an eloquent boundary for determining the point at which we call a machine 'intelligent.' His proposal, which he called "The Imitation Game," but is now referred to as "The Turing Test," is to have an interrogator question a human and an artificial intelligence, randomly labeled X and Y, in the hopes of distinguishing between the two [33]. We reach "intelligence" when an interrogator cannot reliably distinguish between the two [33].

As mathematicians, we could also take our favorite route of defining something: we look at a collection of things that we would agree to be "intelligent" and abstract the shared properties we consider to be desirable until we have a set of properties that must be met for a system to be considered intelligent. From this approach, we could say that a system **S** is intelligent if and only if it is able to "use language,"

form abstractions and concepts, solves kinds of problems now reserved for humans, and improve themselves" [24]. This is the definition from the Dartmouth AI Summer Research Project, but such a definition is obviously hard to evaluate and it would be much more difficult to form a consensus on what set of properties define intelligence than it is to get a set of properties to define other abstract mathematical concepts such as an integral domain.

It quickly becomes apparent that we also need more states than just the two Boolean "intelligent"/"not intelligent" ones to talk about intelligence in a constructive manner. Without this everything on the spectrum from a for loop with an if-else to the robots in Isaac Asimov's *I*, *Robot* are under the same label of "not intelligent" yet we know that the 'intelligence' exhibited in those cases are not at all comparable. We need a spectrum with many states of intelligence to constructively talk about the intelligence of a system.

John Searle's 1980 paper "Minds, Brains, and Programs" introduced the world to The Chinese room argument [32]. It supposes that artificial intelligence research is successful and produces an artificial intelligence that is capable of behaving as if it understands Chinese, then asks does the machine literally understand Chinese or it simulating the ability to understand Chinese [32]? Although this may seem like a pedantic distinction at first glance, the difference is truly important. The hypothesis that we can only ever simulate the ability to think is known as the weak AI hypothesis whereas the hypothesis that we can produce a machine capable of thought is the strong AI hypothesis [24]. These both stand in contrast to Artificial General Intelligence, which is an intelligence capable of performing any intellectual task that a human can [7].

1.1 The History of Artificial Intelligence

One would think that artificial intelligence would have its roots in the last century or two, but mankind has dreamed of and proposed machines with human-like intelligence for thousands years, dating back to at least Homer [7]. From our imagination and literature, artificial intelligence was brought into the realm of the academic by philosophers and mathematicians such as René Descartes' "mechanical man" and Gottfried Wilhelm Leibniz's mechanical reasoning devices [7]. Pascal and Leibniz both designed calculating machines capable of automated arithmetic, but proposing a calculator is far from what we think of as artificial intelligence today [7].

It was the rise of electronics from Turing, IBM, Bell Laboratories, and countless others in the mid-twentieth century started to change the question from a philosophical one to a practical one [7]. Artificial intelligence is a testament to the kind of interdisciplinary problem solving encouraged by the liberal arts, with contributions coming from the fields such as engineering, biology, psychology, game theory, communication theory, logic, philosophy, and linguistics [7]. Eventually, with advancements in computational power, operating systems, and language design, researchers were able to demonstrate impressive computational problem solving such as Arthur Samuel's 1952 checker-playing program written in assembly language and one of the first examples of evolutionary computation [7].

Newell, Shaw, and Simon's "Logic Theorist" program became the first artificial intelligence written for a computer in 1956 [18]. Through the use of heuristic algorithms, Logic Theorist was able to prove theorems and solve complex problems astonishingly well [18]. These early attempts at artificial intelligence were largely doing two things: searching and finding ways to represent and manipulate knowledge. Claude Shannon pointed this out in his 1950 "Programming a Computer for Playing Chess" in which

he produced what is now called the Shannon number, 10^{120} , which is a lower bound of the game tree complexity of Chess [24].

Artificial intelligence became formally recognized as a field of study and got its name from the 1956 Dartmouth Artificial Intelligence Conference [7]. Another product of the conference was a step forward in artificial intelligence's ability to represent and manipulate knowledge with John McCarthy's development of the first AI programming language, LISP [24]. The strides towards strong AI came crashing down with the publication of the 1969 paper "Perceptrons" which showed that single layer perceptrons were not able to properly handle linearly inseparable problems, leading to a steep decline in neural network research and "AI Winter" [24].

Research into artificial intelligence reemerged in the mid to late eighties, but this time with a more practical focus rather than searching for Searle's strong AI [24]. Algorithms developed and used for artificial intelligence found their way into camera auto-focus, anti-lock brakes, search engines, and medical diagnoses [24]. Another marked difference is the plethora of approaches such as agent systems and biologically inspired systems. Today research into artificial intelligence has largely remained in this practical realm, using neural networks, data mining, fuzzy logic and other tools to solve real-world problems while slowly marching towards an Artificial General Intelligence.

1.2 Approaches to Intelligence

Artificial intelligence, because of how broadly the word can be defined and how many fields contribute the its progress, can be hard to wrap one's head around. However, Connell and Livingston have proposed four categories for artificial intelligence approaches which are useful for understanding the state of artificial intelligence research and the varied potential paths to Artificial General Intelligence [9].

Their first category is labeled "Silver Bullets" and describes approaches in which much of what is needed is already believed to be present, but we are missing a crucial piece that will supposedly resolve our problems and deliver us a system with intelligence [9]. Examples include 'Fancy Logic' (second-order, non-monotonic, epistemic, deontic, modal, etc), Fuzzy Logic, Deep Language, Embodiment, and Quantum Computing [9]. Disciples of this school of thought are chasing their particular "Silver Bullet," working to formalize and perfect what they believe to be the missing link.

They next describe the "Core Values" section which puts emphasis on the central organizational scheme over other computational details, believing that this macrolevel structure has greater influence than the exact algorithms used [9]. Situatedness, Emotionality, Self-Awareness, and Hierarchy & Recursion are a few of these ideologies. There are strong arguments for this category, especially Hierarchy & Recursions argument that an intelligence needs to be able to abstract recursively [9].

Connell and Livingston's third category, "Emergence," looks at artificial intelligence approaches which believe they already have the essentials, but we haven't implemented the essentials on a large enough scale to get our intelligence yet [9]. For example, one might hold the position that intelligence is simply the ability to generate and search decision trees and we haven't realized an Artificial General Intelligence yet because our hardware doesn't allow us to do this effectively enough yet. Approaches in this category include Axiomatization, Commonsense, Learning, Evolution, and Integration [9].

Lastly, we visit "Emulation" which is the school of thought that says we are

better off copying intelligence than designing our own [9]. Neural simulation, neural networks, animal models, human development, sociality, and cortical learning algorithms all fit in this category [9]. The danger with this approach is abandoning theory in its sprint towards a functional copy, because if one does not understand the thing they have made, it is hard to see what it can do and where it can be improved. Another excellent point is that it can be very hard to correctly identify what needs to be copied, as Connell and Livingston note, "artificial feathers and flapping turn out not to be needed to create airplanes" [9].

1.3 Biologically Inspired Computing

Emulating natural processes—and even the human brain—is not a new idea to computer scientists. In fact, artificial neural networks that we still use today were originally proposed in a 1943 paper that introduced the first mathematical model of a neuron [12]. The 1950's and 60's saw the rise of evolutionary computing which attempts to simulate evolution usually for finding optimal solutions to problems. Computational neuroscience is the newest field to emerge from this interdisciplinary intersection of biologists and computer scientists.

Computational neuroscience can be described as "simulating human cognition using biologically based networks of neuronlike units" [14]. In other words, the field has tasked itself with the incredibly difficult task of trying to understand cognition by breaking the brain down into algorithms and data structures which we can then implement in a computational model. So rather than trying to create a computational model that is as intelligent as the brain, they are taking the more direct approach in attempting to model the brain itself.

One such set of computational neuroscientists are the researchers at Numenta who are attempting to replicate the mammalian neocortex. Their founder, Jeff Hawkins, realized in the early 1980's that we needed to truly understand what made the brain intelligent before we created our own intelligence [23]. Unfortunately, this view was not shared by researchers at Intel and MIT where he unsuccessfully attempted to start a research lab and study computer science respectively. To support his inquiry, Jeff Hawkins decided to supplement his computer science and engineering background with some training in biophysics and set out to develop a theoretical understanding of the neocortex. In 2002, he founded the Redwood Neuroscience Institute and Hawkins is currently developing cortical learning algorithms and Hierarchical Temporal Memory at Numenta which he founded in 2005 [23].

Chapter 2

Time Series

A time series is a "sequence of observations taken sequentially in time" and are everywhere in today's world [6]. Figure 2.1 shows two such examples, stock prices (Figure 2.1a) and levels of precipitation in Seatle, WA from 1948-1979 (Figure 2.1b). Their ubiquity drives a desire to understand, characterize, and predict time series, but it can be an extremely complex task because there is generally a dependence on prior terms due to the sequential nature and some uncertainty to their movements.



Figure 2.1: Examples of Time Series ¹

2.1 Autoregressive Models

One model for representing time series is the autoregressive (or AR) model. In an autoregressive model, the series at some time t is expressed as a finite weighted sum of previous values of the series plus a random component a_t [6]. Let $z_t, z_{t-1}, z_{t-2}, ...$ denote the series at times t, t-1, t-2, ... Further, let μ be the mean or level of the process and let $\tilde{z}_{t-n} = z_{t-n} - \mu$. We can then express an autoregressive process of order p by

$$\tilde{z}_t = \phi_1 \tilde{z}_{t-1} + \phi_2 \tilde{z}_{t-2} + \dots + \phi_p \tilde{z}_{t-p} + a_t$$
 (2.1)

where μ , ϕ_1 , ϕ_2 , ..., ϕ_p , σ_a^2 are the p+2 unknown parameters of the model with σ_a^2 being the variance of the white noise process a_t [6]. Using a backshift operator B such that $Bz_t = z_{t-1}$ we can define an autoregressive operator of order p by

$$\phi(B) = 1 - \phi_1 B - \phi_2 B^2 - \dots - \phi_p^p \tag{2.2}$$

which allows us to write an autoregressive model succinctly as $\phi(B)\tilde{z}_t = a_t$ [6].

2.2 Moving-Average Models

Suppose instead that the true process is regressed on the random shocks rather than the deviations from the level. Starting from our autoregressive model of order p we can recursively substitute the model's definition of \tilde{z}_{t-n} until we end up with an infinite summation of innovation terms [6]. However, the moving average model

¹Image (b) courtesy of Seattle Weather Blog: http://www.seattleweatherblog.com/rain-stats/

provides us a much more direct approach by representing the time series as a finite number of shocks to the process. Formally we express a moving average (MA) model of order q by

$$\tilde{z}_t = a_t - \theta_1 a_{t-1} - \theta_2 a_{t-2} - \dots - \theta_q a_{t-q} \tag{2.3}$$

just as with the autoregressive model, we have q + 2 unknown parameters and using the backshift operator we can define a moving average operator that allows us to express the model more succinctly yielding $\tilde{z}_t = \theta(B)a_t$ [6].

2.3 Autoregressive Moving-Average Models

Real world processes may require a combination of these two approaches for a succinct and accurate model, resulting in a mixed Autoregressive Moving-Average (ARMA) model [6]. Such a model includes terms from both the autoregressive and moving average model resulting in:

$$\tilde{z}_t = \phi_1 \tilde{z}_{t-1} + \dots + \phi_p \tilde{z}_{t-p} + a_t - \theta_1 a_{t-1} - \dots - \theta_q a_{t-q}$$
(2.4)

This model has p+q+2 unknown parameters and can also be written more succinctly using our autoregressive operator and moving average operator resulting in $\phi(B)\tilde{z}_t = \theta(B)a_t$ [6].

2.4 Classification of Time Series

A key question in time series analysis is how to select the correct model for a given time series. There are a few approaches, but one is to use information criterion

such as the Akaike Information Criterion or the Bayes Information Criterion [6]. For both AIC and BIC, optimal models minimize the information criterion and there is a penalty for the number of parameters used resulting in a bias towards a smaller set of parameters. The formula for BIC is

$$BIC_{p,q} = \ln(\hat{\sigma}_a^2) + r \frac{\ln(n)}{n}$$
(2.5)

where n is the sample size, $\hat{\sigma}_a^2$ is the maximum likelihood estimate of σ_a^2 , and r = p + q + 1 is the number of parameters estimated in the model.

The ubiquity of time series makes forecasting them an important and many times profitable task such as in the case of financial time series (as seen in Figure 2.1a), making them a common prediction task for machine learning and artificial intelligence systems. It allows us to test an "intelligent" systems ability to discover temporal patterns which may sometimes be fuzzy or complex: a hallmark of human intelligence.

Chapter 3

Cortical Learning Algorithms

Hierarchical Temporal Memory is a framework attempting to produce cortical learning algorithms—algorithms that mimic the processes of the human neocortex. In the human brain, the neocortex is approximately seventy-percent of the brain's volume [28]. It can be described as a "wrinkly sheet of sheet of cells, about two millimeters thick" and is responsible for the characteristics of intelligence that set humanity apart from other species [28].

Vernon Mountcastle

Overview of the neocortex

Cortical Columns and Pyramidal Neurons

3.1 Sparse Distributed Representations

Consider the binary representation of the integer 16 versus that of the integer 15. The two numbers are quite similar: they are only 1 apart, so they are as close

¹Image courtesy of Wikipedia: https://en.wikipedia.org/wiki/Neocortex

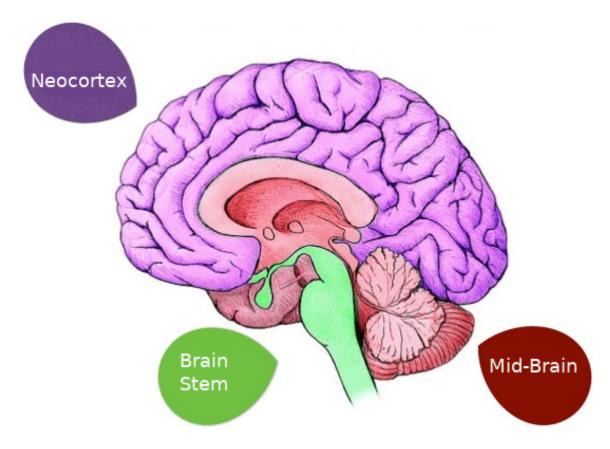


Figure 3.1: Overview of the Human Brain

as two integers can be. Yet their binary representations are [100000] and [011111] respectively. They have no shared "on" bits so despite their similarity, their binary representations reflect none at all. In fact, despite being as close as two integers can be (a Euclidean distance of 1), their binary representations have a Hamming distance, the number places in which two codewords differ, of 5, the maximum Hamming distance of two codewords of length 5 [1].

This means that our encoding does not preserve semantic similarity or a concept of distance between elements which is highly undesirable for an code because if there is some kind of error in the code we could end up decoding something meaning the

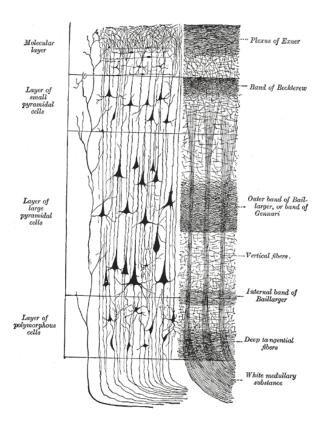


Figure 3.2: Column of the Neocortex¹

opposite of what we were trying to convey. As an example, consider \mathbb{Z} from 0 to 31 which is mapped to $GF(2)^6$ by their binary representation. The mapping of 31 is [111111] but a single error in transmission can easily lead to [011111] which would be decoded as 15. So a code Hamming distance (d_H) of one away $(\frac{1}{5}$ of the total metric) lead to an element 16 integers away $(\frac{1}{2}$ of the total metric). We would obviously like to avoid this so that errors in the transmission of our codes are either (1) correctable or (2) lead to a decoding that is as close as possible to our original element.

This is achievable by simply conveying more information in our code. In our binary representation any single error led to another valid codeword (a codeword which decoded to an element of the input/output set) which meant that no errors could be detected or corrected. By expanding our code length, we increase the number of codewords (multiplying by the cardinality of the code alphabet for each character added) meaning that fewer errors will result in other valid codewords and can possibly be detected or corrected.

A key strategy with Sparse Distributed Representations is to encode semantic similarity, such as with our idea of distance in our motivating example. This helps us achieve our second goal because even if we increase the error-tolerance of our code, there is still some probability of an uncorrectable error and we would like that error to result in a codeword as close to the original codeword as possible. To give you a real world example imagine I am sending instructions to a aircraft and I need to tell it to turn down 7° to start its descent. Obviously, an error resulting in 8° or 6° being interpretted by the pilot are both preferable to 90°.

To achieve our goal, we employ sparse distributed representations or SDRs. Just as with traditional dense binary representations, we will represent sparse distributed representations as vectors over their code alphabet, in this case GF(2). We call them sparse distributed representations because these vectors typically only have a small proportion of the components as 1. We will use Numenta's notation of letting $w_{\overrightarrow{x}}$ denote the number of components in an SDR \overrightarrow{x} that are 1, so $w_{\overrightarrow{x}} = ||x||_1$ [2].

Given our definition of distance, we could say that two decodings of sparse distributed representations, a and b, are equal if and only if the $d_H(a,b) = w_a = w_b$. This would mean that both vectors would have to have the same dimensionality, same number of on bits, and all on and off bits would have to match. This definition is good for "equals," but suppose we have a single error in transmission or a single component of our distributed system fails, equality would thus fail. In order to be able preserve

the ability to subsample and thus to preserve fault tolerance, we therefore need a less stringent definition for decoding SDRs. Numenta refers to this as the *overlap*, which is

$$overlap(\overrightarrow{a}, \overrightarrow{b}) \equiv \overrightarrow{a} \cdot \overrightarrow{b} \equiv \sum_{i=0}^{n-1} a_i b_i$$
 (3.1)

Thus, we say two SDRs \overrightarrow{d} and \overrightarrow{b} decode to the same element of the input space if and only if $overlap(a,b) \ge \theta$ where $\theta \le w_a$ and $\theta \le w_b$ [2]. Denote the function that determines if two sparse distributed representations decode to the same element of the input space using some θ and $overlap(\overrightarrow{a}, \overrightarrow{b})$, $match_{\theta}(\overrightarrow{a}, \overrightarrow{b})$ and it is a function from SDR×SDR \longrightarrow {true, false}.

Given a set of sparse distributed representations with the same dimension, $X = \{\overrightarrow{x_1}, \overrightarrow{x_2}, ..., \overrightarrow{x_n}\}$, we can union the vectors using the bitwise OR operation over the i^{th} position of the vectors in the set to produce the i^{th} position of union(X) [2]. For example, given [0100] and [0010] the union would be [0110]. We say an SDR \overrightarrow{y} is an element of the union of a set of SDRs, \overrightarrow{X} , if and only if $match_{\theta}(\overrightarrow{X}, \overrightarrow{y})$ [2].

3.2 Hierarchical Temporal Memory

Hierarchical Temporal Memory aims to implement the functions of the neocortex as faithfully as possible, and unsurprisingly uses the same atomic unit of the cell. Modeled after the pyramidal neurons of the neocortex, these cells each take their own inputs and make their own predictions which the layer combines into a cohesive prediction [21]. They have three states: active from feed-forward input, active from lateral input (predictive state), and inactive. These first two states correspond to

a short burst of action potentials in a neuron, and a slower steady burst of action potentials on a neuron respectively in our own brain.

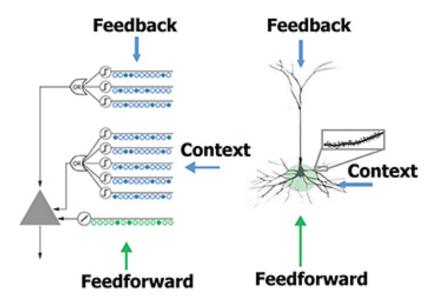


Figure 3.3: Pyramidal Neuron vs. Hierarchical Temporal Memory Cell [11]

Just like actual pyramidal neurons, HTM cells also dendrite segments. Each cell has one proximal dendrite segment and a dozen or more distal dendrite segments [21]. Proximal dendrites accept feed-forward input and all cells in the same column respond to a similar feed-forward input due to a class of inhibitory cells. On the other hand, distal dendrites receive lateral input from nearby cells and their response is unique on a cell-by-cell basis dependent on the connected synapses on each cells distal dendrites. Figure 3.3 illustrates the similarities between a pyramidal neuron and an HTM cell.

talk about synapses

3.2.1 Encoder

3.2.2 Spatial Pooler

The Spatial Pooler's job is to recieve the encoded representation of the input space and convert the input to a sparse pattern-a requirement for learning and prediction in Hierarchical Temporal Memory theory [21]. Upon intitialization, all of the potential synapses on the dentrite segments of all of the mini-columns are randomly assigned permenance values and based on these permance values a potential synapse may be valid or not [21].

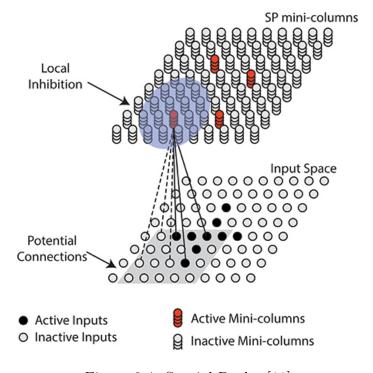


Figure 3.4: Spatial Pooler [11]

The pooler begins with an input from the encoder that has a fixed number of bits and for each column we determine how many valid synapses are connected to active input bits [21]. This number is then multipled by a "boosting" factor determined dynamically by how often a column is active relative to its neighbors. The most highly activated mini-columns after boosting then disable a percentage of the mini-columns within a dynamically determined "inhibition radius" to create a sparse set of activated mini-columns [21]. Lastly, each active column updates the permanence values of its potential synapses. The permanence values associated with active input bits are increased while those associated with inactive bits are decremented [21].

3.2.3 Temporal Pooler

The temporal pooler takes over where the spatial pooler left off, with a sparse set of activated mini-columns

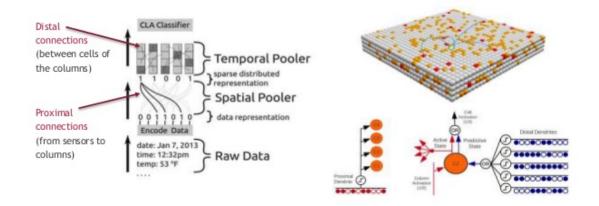


Figure 3.5: Spatial and Temporal Poolers

3.2.4 Classifier

Once the system completes its processing of a time step's input, the system outputs a sparse distributed representation which may have gone through multiple layers and it is important to be able to decode the system's prediction in order for the system to be useful. Hierarchical Temporal Memory currently uses something called an SDR Classifier to decode the predictions of an HTM [13]. At its essence, an SDR Classifier is a single layer, feed forward, neural network [13].

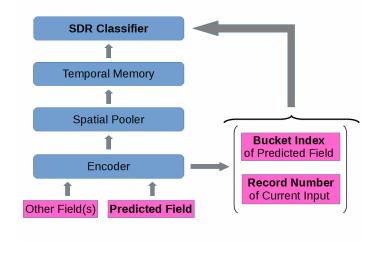


Figure 3.6: SDR Classifier

SDR Classifiers are able to decode the information HTMs give about predictions n times steps in the future by maintaining a weight matrix [13]. The matrix's weights are adjusted to "learn" after each time step to reflect the correct weighting between the output vector at time t and the probability distribution of the input/output space at time t-n [13]. This enables the matrix to reflect relationships between inputs and outputs n time steps apart. To determine the SDR Classifier's interpretation of an output at time t+n, the SDR Classifer takes in the HTM's output vector and uses Softmax to determine the most likely decoding [13]. So for each of the k classes in the output space, the certainty that the output vector is referring to it is

$$y_j = \frac{e^{a_j}}{\sum_{i=1}^k e^{a_i}} \tag{3.2}$$

where a_j is the activation level of the j^{th} element of the output space, calculated by multiplying the output vector by the j^{th} column of the weight matrix element wise [13].

Chapter 4

Literature Review

4.1 Sequence Memory for Anomaly Detection and Prediction

Numenta's work

partners such as http://numenta.com/htm-for-stocks/ grokstream.com cortical.io Masters theses on prediction

4.2 Spatial Data, Sensors, and Motor Response

Chapter 5

Experiment Design

5.1 Hardware and Software Specifications

I utilized the official Python (Python 2.7) implementation of Numenta's cortical learning algorithms, NuPIC¹ and it's Network API². At the time of my work NuPIC was at version 1.05 and I chose to use this version. It should be noted that although much of NuPIC is in Python, many of its computationally intensive algorithms are implemented in C++ using their NuPIC.core repository³ and it also relies on a number of external libraries such as numpy and pandas.

Due to the computationally intensive nature of my research and the limited time available to me, I also heavily relied on parallel processing to run my code in a reasonable time frame. In particular, the multiprocessing package in Python was indispensable to my research, allowing me to speed my code up a factor of up to 28x when running independent tests or swarming.

¹NuPIC repository: https://github.com/numenta/nupic

²Network API docs: http://nupic.docs.numenta.org/stable/api/network/

³NuPIC.core repository: https://github.com/numenta/nupic.core

Dijkstra (Asus Vivobook)			
CPU	Intel i5-8250U, 4-Core, 1.60GHz		
RAM	24 GB		
OS	Linux Mint 19 Cinnamon		
Galois (HP DL380 G7)			
CPU	2x Intel Xeon X5650, 6-Core, 2.70 GHz		
RAM	192 GB		
OS	Linux Mint 19 Cinnamon		
Aslan (Type)			
CPU	2x Intel Xeon E5-2690, 10-Core, 3.00 GHz		
RAM	252 GB		
OS	Gentoo		

Table 5.1: Hardware Specifications

On the hardware side of things, the majority of development and testing was conducted on my personal laptop, an Asus Vivobook with some small upgrades (see Table 5.1 for details). The code was then deployed to my HP DL380 G7 or [ASLAN's type] server which had more cores, more memory, and a higher clock speed allowing me to get faster results. For a synopsis of the hardware specifications, see Table 5.1.

5.2 Particle Swarm Optimization

Introduced in by Russell Eberhart and James Kennedy in their 1995 paper, A New Optimizer Using Particle Swarm Theory, Particle Swarm Optimization has changed the way the world optimizes continuous nonlinear functions [5]. Swarming is fast

and effective for optimization over complex multidimensional search spaces and lends itself well to parallelization due to its multi-agent approach, making it an excellent choice for optimizing parameters in neural networks and is implemented by Numenta for finding optimal parameters.

Usually abbreviated to PSO or *swarming*, it draws inspiration from bird flocks and schools of fish [14]. The concept is quite simple: it simulates particles on parameter space and at each time-step every particle evaluates fitness as a function of its location in parameter space, then moves towards some linear combination of its personal best and the overall best score among the particles using weights representing a cognitive constant or individuality (denoted c) and a social constant (denoted s).

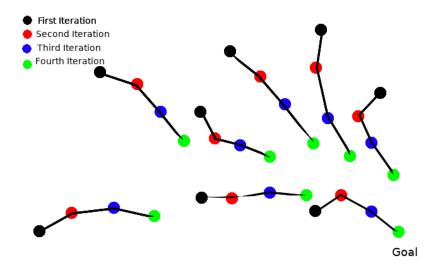


Figure 5.1: Visualization of the PSO Algorithm

The algorithm begins by initializing all particles with a random coordinate in the parameter space and a random velocity. Then at each time step, it calculates its fitness based on the function being optimized and the parameters associated with its position in parameter space. If this fitness is the best the particle has achieved, it is

recorded as its personal best, denoted **pbest** and the particle with the best fitness is chosen as global best, denoted **gbest**. We then update the velocity and particle position of each particle in our swarm using the equations described in Equation (5.1) where v_t is the velocity of a particle at time step t, p_t is the position of a particle at time step t, and β_1, β_2 are uniformly distributed random variables with $\max(\beta)$ being a parameter of the algorithm [5].

$$v_{t+1} = v_t + c \times \beta_1 \times (\text{pbest} - p_t) + s \times \beta_2 \times (\text{gbest} - p_t),$$

$$p_{t+1} = p_t + v_{t+1}.$$
(5.1)

I utilized swarming to find optimal or near-optimal sets of parameters for learning ARMA time series with a Hierarchical Temporal Memory network. Numenta themselves utilizes swarming for this very same purpose, and NuPiC comes with a swarming utility, however I wrote my own swarming utility in Python which I throughly tested on a number of complex surfaces.

5.3 Time Series Classification

A key component of this work was the classification of ARMA time series. To accomplish this, I used the statsmodels⁴ Python module which comes with a variety of tools for conducting statistical tests including time series statistics. Among them is the $arma_order_select_ic$ function which finds information criterion for ARMA models. The module also contains ARMA.fit method which takes a p and q alongside the time series and returns the (p,q)-ARMA model using exact maximum likehood via Kalman filter [31].

⁴statsmodels webpage: https://www.statsmodels.org/dev/index.html

The likelihood model used by ARMA.fit has the option to use a variety of methods [31]. I wrote a function which intially attempted to use the default solving method of limited memory Broyden-Fletcher-Goldfarb-Shanno and upon throwing an exception would then sequentially fallback upon the other methods available: Broyden-Fletcher-Goldfarb-Shanno, Newton-Rhapson, Nelder-Mead, conjugate gradient, non-conjuage gradient, and Powell, in that order [31].

Chapter 6

Results and Discussion

The first step of this project was learning how to interact with the NuPIC Network API in Python: this included writing a stream object to feed an HTM network inmemory arrays, writing training routines for HTMs, and conducting some sensitivity analysis with the model parameters.

Then we began by training HTMs on very simple deterministic sequences which did not require any temporal context past the last term-that is sequences such that you can know for certain the next term in the sequence by only knowing the last term. Examples include 1, 2, 1, 2, ..., 1, 3, 1, 3, ..., and 78, 79, 78, 79, This allowed us to ensure that we could train and predict using only an HTM's spatial pooler.

Next, we trained Hierarchical Temporal Memory networks using slighly more complex, but still deterministic sequences, that relied on some more long-range temporal dependencies. For these patterns, it is not always sufficient to know just the last term, but you may need to also know some term other previous term which would require the HTM's temporal pooler. Examples include 1, 2, 3, 2, ... and 1, 2, 2, 3, 5, 2, 2, 4, 1, In the first example, 1 and 3 are always followed by 2, but to know what is followed

by 2 you also need to know what proceeded 2.

When we transitioned to ARMA models we started with a low standard deviation for the noise terms and grew from there. We also decided to normalize the samples of the ARMA processes such that the values were bounded between [-100,100].

Autoregressive Model			Moving-Average Model		
	μ	σ	μ	σ	
AR Lag 1	0.00349	0.07462	0.00276	0.05784	MA Lag 1
AR Lag 2	-0.00185	0.03404	-0.00147	0.04248	MA Lag 2
AR Lag 3	0.00134	0.02895	-0.00049	0.02883	MA Lag 3
AR Lag 4	-0.00039	0.03909	0.00192	0.03478	MA Lag 4
AR Lag 5	0.00093	0.02925	-0.00128	0.03096	MA Lag 5
AR Lag 6	0.00387	0.03895	-0.00175	0.03532	MA Lag 6
AR Lag 7	0.00042	0.01602	-0.00016	0.01331	MA Lag 7
AR Lag 8	0.00008	0.00314	-0.00010	0.00788	MA Lag 8
MA Lag 1	-0.00461	0.06797	-0.00193	0.04801	AR Lag 1
MA Lag 2	0.00085	0.01665	0.00142	0.03004	AR Lag 2

Table 6.1: Errors on fitting coefficients of AR and MA models to instances of ARMA models

Autoregressive Model			Moving-Average Model		
	HTM	Error in	HTM	Error in	
	Output	Terms of Sigma	Output	Terms of Sigma	
AR Lag 1	0.063	0.79751	0.897	15.55602	MA Lag 1
AR Lag 2	0.000	0.05435	0.000	0.03460	MA Lag 2
AR Lag 3	0.000	13.77064	0.000	13.89143	MA Lag 3
AR Lag 4	0.000	0.00998	0.000	0.05520	MA Lag 4
AR Lag 5	0.000	10.22462	0.000	9.73127	MA Lag 5
AR Lag 6	0.000	7.60282	0.000	8.54332	MA Lag 6
AR Lag 7	0.000	0.02622	0.000	0.01202	MA Lag 7
AR Lag 8	0.000	0.02548	0.000	0.01269	MA Lag 8
MA Lag 1	0.000	0.06782	-0.817	17.05749	AR Lag 1
MA Lag 2	0.000	0.05105	0.000	0.04727	AR Lag 2

Table 6.2: Errors on fitting coefficients of AR and MA models to HTM outputs

Chapter 7

Further Work

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Appendices

Appendix A

Errors on Fitting AR and MA Models

Autoregressive Model			Moving-Average Model		
	μ	σ	μ	σ	
AR Lag 1	0.00349	0.07462	0.00276	0.05784	MA Lag 1
AR Lag 2	-0.00185	0.03404	-0.00147	0.04248	MA Lag 2
AR Lag 3	0.00134	0.02895	-0.00049	0.02883	MA Lag 3
AR Lag 4	-0.00039	0.03909	0.00192	0.03478	MA Lag 4
AR Lag 5	0.00093	0.02925	-0.00128	0.03096	MA Lag 5
AR Lag 6	0.00387	0.03895	-0.00175	0.03532	MA Lag 6
AR Lag 7	0.00042	0.01602	-0.00016	0.01331	MA Lag 7
AR Lag 8	0.00008	0.00314	-0.00010	0.00788	MA Lag 8
MA Lag 1	-0.00461	0.06797	-0.00193	0.04801	AR Lag 1
MA Lag 2	0.00085	0.01665	0.00142	0.03004	AR Lag 2

