

Minimum levels of high energy particle  
bombardment on fusion reactor vessels:  
Towards a computational multi-parameter scan for automated  
data reduction ("big data") applications

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# Contents

<b>1</b>	<b>Model</b>	<b>1</b>
1.1	EIRENE 1D . . . . .	1
1.1.1	Kinetic Boltzmann Equations . . . . .	1
1.2	Plasma Profiles . . . . .	1
1.3	Choice of sampling set . . . . .	1
1.3.1	Sobol Sequence . . . . .	2
<b>2</b>	<b>Methods</b>	<b>3</b>
2.1	Neural Networks . . . . .	3
2.1.1	General Introduction To Neural Networks . . . . .	4
2.1.2	Fully Connected . . . . .	5
2.1.3	Training . . . . .	6
2.1.4	Hyper parameters . . . . .	7
2.2	Gaussian Processes . . . . .	9
2.2.1	General Introduction . . . . .	9
2.2.2	GOGAP algorithm . . . . .	9
<b>3</b>	<b>Results</b>	<b>11</b>
3.1	Physical Results . . . . .	11
3.2	Neural Networks . . . . .	11
3.2.1	Hyper Parameters . . . . .	11
3.2.2	Derivatives . . . . .	12
3.3	Gaussian Processes . . . . .	12
3.3.1	Subdivision of parameter space . . . . .	12
3.3.2	Derivatives . . . . .	12
3.4	Comparison . . . . .	12
3.4.1	Accuracy . . . . .	12
3.5	NNGP - Maybe . . . . .	12
<b>A</b>	<b>More Data probably</b>	<b>13</b>

<b>B</b>	<b>Background Neural Network</b>	<b>15</b>
B.1	Introduction to Neural Networks . . . . .	15
B.2	Examples of neural networks in different contexts . . . . .	16
<b>C</b>	<b>Background Gaussian Processes</b>	<b>17</b>
C.1	Bayesian Statistics . . . . .	17

# Chapter 1

## Model

### 1.1 EIRENE 1D

#### 1.1.1 Kinetic Boltzmann Equations

### 1.2 Plasma Profiles

For the inputs of Eirene plasma profiles are needed, that can be dynamically provided by other algorithms like SOLPS<sup>1</sup> or EMC3<sup>2</sup>. Central piece of this work is to investigate if a substitute function can be found for the full range of possible plasma profiles by using big data methods. One can ascertain the physical limits of the parameters constituting the plasma profiles from table ???. These limits are based on different phenomenons in plasma physics, which can be

### 1.3 Choice of sampling set

Since the parameter space is high dimensional, the training points have not been selected randomly. According to <sup>3</sup> randomly sampled points might form clusters, which could skew the training towards a subsection of the parameter space. To avoid this a low discrepancy sequence, namely the Sobol sequence, has been chosen from which the training points are sampled.

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<sup>1</sup>Citation needed

<sup>2</sup>Citation needed

<sup>3</sup>Add reference

### 1.3.1 Sobol Sequence

The sobol sequence was first invented by the mathematician Ilya M. Sobol<sup>4</sup>

#### Low-discrepancy sequences

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<sup>4</sup>Add citation

# Chapter 2

## Methods

This chapter is concerned with introducing the methods used to investigate the data reduction of the previously introduced model. The focus of this work will be on artificial neural networks (ANN in short) and Gaussian processes. The following list will provide a brief overview of other machine learning techniques that could be employed to

- Support Vector Machines (SVM):  
Support Vector Machines (in short SVM) are used to classify data similar to ANNs. In contrast to ANNs SVMs are build up from theory and contain little Hyperparameters<sup>1</sup> making them easier to analyse and less prone to overfitting. Generally speaking a SVM tries to seperate data by calculating a hyperplane using given training data. In more complex situations transformations of the parameter space into a higher dimensional space wherein the p
- Random Forests:
- Boosting:

Short description of Random Forests

Short description of Boosting methods

### 2.1 Neural Networks

The following section is concerned with discussing neural networks as a means of investigating functional dependencies.<sup>2</sup>

<sup>1</sup>Please refer to section 2.1.4 for further information.

<sup>2</sup>To aid with understanding the terminology used there is a glossary in the appendix section B.1.

### 2.1.1 General Introduction To Neural Networks

An artificial neural network (ANN) in the following called neural network, abbreviated to NN, is "a computing system made up of a number of simple, highly interconnected processing elements, which process information by their dynamic state response to external inputs." [1] First concepts of learning algorithms based on neural plasticity have been introduced in the

citation needed

late 1940s by D. O. Hebb . In 1975 backpropagation became possible via an algorithm by Werbo, this led to an increase in machine learning popularity.

citation needed

During the 1980s other methods like support vector machines and linear classifiers became the preferred and/or dominating machine learning approach. With the rise in computational power in recent years neural networks have gained back a lot of popularity.

The concept idea of neural networks is to replicate the ability of the human brain to learn and to adapt to new information. The structure and naming convention reflect this origin.

A neural network is made up of small processing units called neurons. These are grouped together into so called layers. Every network needs at least two layers, the input layer and the output layer. If a network has intermediary layers between input and output, they are called hidden layers. A network with at least two hidden layers is called a deep neural network (DNN). The amount of layers in a network is called the depth of the network. While the amount of neurons in a layer is called the layers width. In a typical NN information stored in neurons is transferred into the next layer by a weighted sum. The connected neuron of the following layer then applies a non-linear function, called activation function, to calculate it's final value. This process is repeated until the output layer is reached. The activation function

reword "order"

as well as the amount and order of connections can vary in between layers. The system according to which a network is designed is called a network architecture. The most important architectures in the following work will be *dense deep feed forward* and *autoencoder*. To give insight into the basic working principle an example neural network is depicted and described in the following section 2.1.2.

Neural networks are usually used in two ways, optimization or classification. Well known examples are handwriting recognition as classification and least mean squares (LMS) optimization.

think of better example



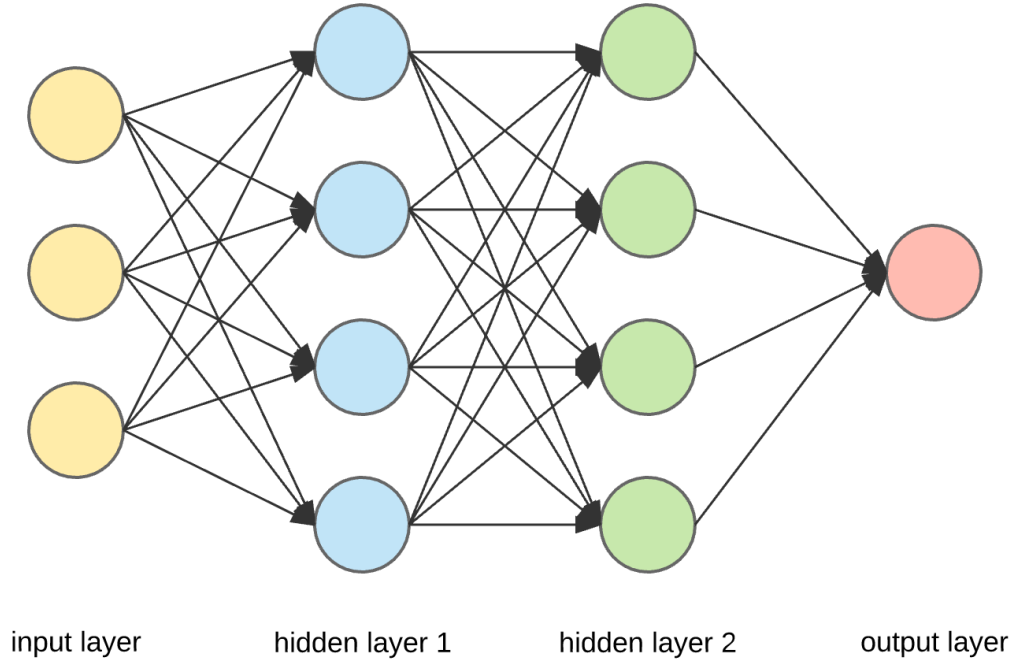


Figure 2.1: Schematic structure of the most basic fully connected deep neural network. Indicated are the input (yellow), output (red) and hidden layers (blue and green). Each neuron outputs to all neurons in the following layer, but there are no interconnection between neurons of the same layer. Note that while the network has the minimum depth (2 hidden layers) to qualify for a deep neural network, the width could be smaller.

### 2.1.2 Fully Connected

A fully connected or dense neural network like depicted in figure 2.1 is characterized by connecting every neuron from the previous to the following layer with a weight. In contrast to other networks this allows for a very high flexibility but also lacks the spatial context of data. This kind of network is especially well suited for data that is given in form of vectors or drawn from an arbitrary parameter space.

The working principle is to form a weighted sum  $\sum_{k=1}^N w_{j,k} \cdot x_k$  over the values from neurons of the previous layer  $x_k$  weighted by the connecting weight  $w_{j,k}$ . The weighted sum is then evaluated by the activation function  $\sigma()$  such that the new value  $x_j = \sigma(\sum_{k=1}^N w_{j,k} \cdot x_k)$ . Here  $k$  refers to the index of the neuron in the previous layer and  $j$  to the index of the neuron in the current layer.<sup>3</sup>

<sup>3</sup>The order of indices becomes more intuitive when talking about backpropagation and

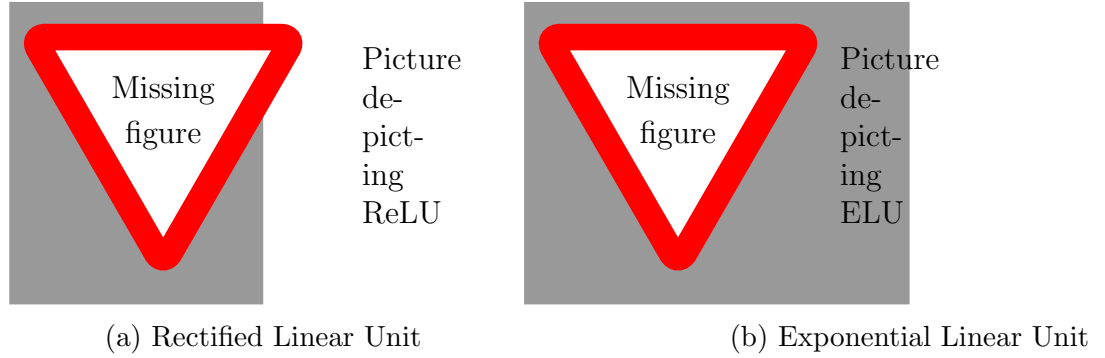


Figure 2.2: Example activation functions rectified linear unit (a) and exponential linear unit (b) used to introduce non-linearity into neural networks.

ref needed

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Since forming a weighted sum is a linear operation the activation function must be non-linear to enable the network to learn non-linear behaviour. The most common activation functions are the rectifier also called rectified linear unit (ReLU) and exponential linear unit (ELU) shown in figure 2.2. Both are inspired by the asymmetrical behaviour of biological neural connections.

### 2.1.3 Training

Before a neural network can be put to work it needs to be trained. To train a NN a set of training and test data has to be generated. This work uses the afore mentioned monte carlo simulations from the EIRENE code. The training data consists of input e.g. temperature and density of the plasma and output e.g. the sputtering rate of the first wall. The EIRENE input data is used as input of the network and the sputtering rate is compared to the output of the network via a cost function. Afterwards the weights of the network are adjusted by using backpropagation, which is a method that calculates partial weight derivatives of the output. A more detailed explanation

add citation

can be found in section 2.1.3 .

### Backpropagation

$$\frac{\partial F}{\partial w} \quad (2.1)$$

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it's matrix notation in section 2.1.3

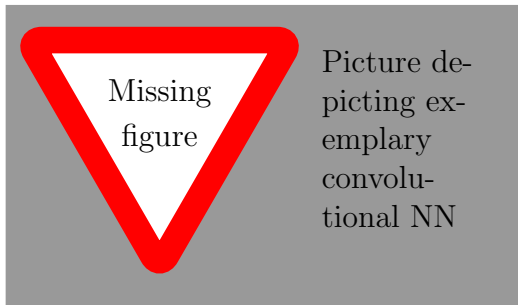


Figure 2.3: Exemplary Conv NN

In contrast to fully connected networks convolutional networks apply a so called filter to the input. These filters consider the inputs of nearby neurons as well. They are usually used in image recognition and require data that has information stored in patterns, most commonly special patterns as in images. Fig. 2.3 depicts an exemplary convolutional network for image data and gives indication to its working procedure.

### 2.1.4 Hyper parameters

A hyper parameter refers to a parameter of the network that is not changed during training. Since these can have substantial influence on the performance of the network they will be explained in the following

#### Depth

The depth of the network correlates directly to the amount of chained non-linear functions. Therefore it strongly influences the ability of a NN to learn abstract patterns. The more complicated a pattern is the more depth is "required" to learn the pattern. For example a simple classification between left and right only requires a shallow network, whereas recognizing different brands of car requires a deeper network.

#### Width

The width of a layer refers to the amount of neurons in that layer. Often networks are build of layers with the same width in each hidden layer. If that is the case one sometimes speaks of the width of the network which is then equivalent to width of a/each hidden layer.

## Architectures

### Activation or Activation Function

As previously discussed the activation function introduces non-linearity to the network. Some activation function will be better suited to model a certain problem than others. If information about the model or pattern to be predicted is known, an activation function close to this will have better performance. For example predicting outcome of a sine function will perform better with exponential activations or uneven polynomial activations than even polynomials.

### Loss Function

Choosing a loss function determines what criteria the network optimizes for which directly corresponds to which patterns it learns. For prediction one typically chooses a root mean square function. For classification cross entropy loss functions are most common.

### Batch size

The Dataset is divided into subsets called batches which are fed to the current network during training. After each batch the weights are adjusted. Splitting the dataset in this way is advantageous to the computational performance during training. Less memory is used during training and the number of epochs trained is reduced. The flip side of using a batch size smaller than the number of training data is that the gradient for optimization will be worse in comparison to the gradient calculated with the full data set.

### Epochs

An epoch describes a full training cycle of training, validating and adjusting weights for the entire training data set. If the batch size is smaller than the number of training points then multiple<sup>4</sup> adjustments are made.

### Metrics

Metrics are additional information gained from the network during training and evaluation. Metrics are not hyperparameters since they do not influence the resulting network but are an important source of information for further improving the network structure. For example a secondary loss function can

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<sup>4</sup>Number of batches in one epoch = rounded up  $\left( \frac{\text{Amount of Training Data}}{\text{Batch Size}} \right)$

be implemented as a metric to evaluate general optimization of the network in contrast to only the chosen loss quantity.

### **Regularization**

Regularization describes methods used to reduce the generalization error but not the training error. Commonly used regularization methods include L1, L2, Dropout and Early Stopping regularization. L1 and L2 regularization is applied by adding a penalty term to the loss function. This requires initial knowledge of input influences. For example an image with bad resolution might have a larger penalty term applied than an image with high resolution. Dropout regularization and early stopping are used to prevent overfitting. Since the amount of parameters in the network is often on the same order of magnitude as the amount of training data, neural networks are prone to overfitting. Early stopping interrupts the training process as soon as the validation loss stops improving by a user set minimum delta.

## **2.2 Gaussian Processes**

### **2.2.1 General Introduction**

### **2.2.2 GOGAP algorithm**



# Chapter 3

## Results

### 3.1 Physical Results

### 3.2 Neural Networks

#### 3.2.1 Hyper Parameters

The following subsets of hyper parameters have been investigated: Beginning with investigating the effect of the data amount. Using same network size per

Number	Width	Depth (hidden)	Amount of Data	Activation	Droprate
1	100	10	$2^{17}$	Relu	0.25
2	X	X	$2^{16}$	X	X
3	X	X	$2^{15}$	X	X
4	X	X	$2^{14}$	X	X
5	80	X	$2^{17}$	X	X
6	X	X	$2^{16}$	X	X
7	X	X	$2^{15}$	X	X
8	X	X	$2^{14}$	X	X

#### Depth & Width

Depth and Width together determine the total number parameters and complexity of the network. The question is how complex does the network need to be to accurately learn the model. Ideally we'd like to trim the network as much as possible without losing too much accuracy.

**Activation**

Elu, Relu, Sigmoid

**Loss Function**

SGD

**3.2.2 Derivatives****3.3 Gaussian Processes****3.3.1 Subdivision of parameter space****3.3.2 Derivatives****3.4 Comparison****3.4.1 Accuracy****3.5 NNGP - Maybe**



# Appendix A

## More Data probably



# Appendix B

## Background Neural Network

### B.1 Introduction to Neural Networks

Glossary:

- Network: A series of layers. The first layer of a network is called the input layer, the last layer is called the output layer. Any layers in between are called hidden layers. The amount of hidden layers is called the **depth** of the network.
- Layer: A collection of neurons. The amount of neurons in a layer is called the **width** of a layer.
- Neuron: A single node in a layer. It contains a single number formed by a weighted sum of it's inputs evaluated by the activation function.
- Activation function: A non-function applied to the weighted sum of a neuron. Used to introduce non linearity into the network in order to enable non linear model "fitting".
- Weight: Each connection between layers has it's own weight factor. These are adjusted during training to fit the data. Weights are often referred to as parameters.
- Regularization: Methods used to suppress overfitting.

- Metric: Additional information gathered during training/testing.
- Loss function: Function that dictates the optimization, e.g. Root Mean Squared.
- Training Data: Set of Data used during training phase. Weights are adjusted to these data.
- Validation Data: Set of Data used during training phase to evaluate training results intermediary.
- Test Data: Data set not seen during training to evaluate trained network performance.
- Input: Data fed to the network for training or evaluation.
- Output: Prediction of the network.
- Label: True output value for input data.
- Hyperparameter: A parameter not changed during training e.g. width and depth of the network.

## B.2 Examples of neural networks in different contexts

# Appendix C

## Background Gaussian Processes

### C.1 Bayesian Statistics



# Bibliography

- [1] Maureen Caudill. Neural networks primer, part i. *AI Expert*, 2(12):46–52, December 1987.