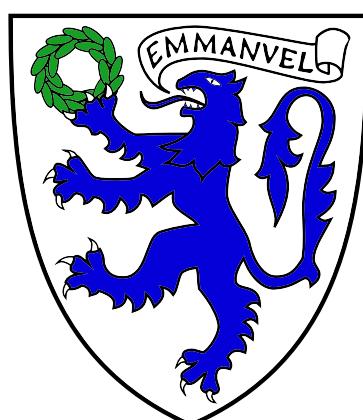


Unifying Transformers and Convolutional Neural Processes

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Chapter 1

Introduction

1.1 Motivation

Machine learning models have been immensely successful in variety of applications to generate predictions in data-driven domains such as computer vision, robotics, weather forecasting. While the success of these models is undeniable, they tend to lack the ability to understand the uncertainty in the predictions. This is a major drawback in the deployment of these models in real-world applications, for example, in weather forecasting, the uncertainty of the prediction of the weather is arguably as valuable as the prediction itself. In this work, we aim to implement a model that is **uncertainty aware** whilst also possessing further desirable properties.

1.2 Desirable Properties

On top of being uncertainty aware, we would like to insert some desirable inductive biases that help the model to generalize better and be more interpretable. These properties are:

Flexible: *The model should be able to work on a variety of data types.* As long as a data point can be represented as a vector, the model should be able to operate on it. This allows the model to be used in a variety of applications and domains.

Scalable: *The model should be able to work on large datasets.* There should be no limit on the size of the dataset that the model can be trained on. Which is not the case with many traditional models such as Large Language Models. Another aspect of scalability is the ability learn high-dimensional data with good computational efficiency.

Permutation Invariant: *The prediction of the model should not change if the order of the input data is changed.* When each data point contains the information about input and output pairs, the model should not care about the order in which they are fed into the model. For example, in the case of a weather forecasting model, which uses data from multiple weather stations, the model should not care about the order in which the data from the weather stations is fed into the model, thus making the model permutation invariant.

Translation Equivariant: *Shifting the input data by a constant amount should result in a constant shift in the predictions.*

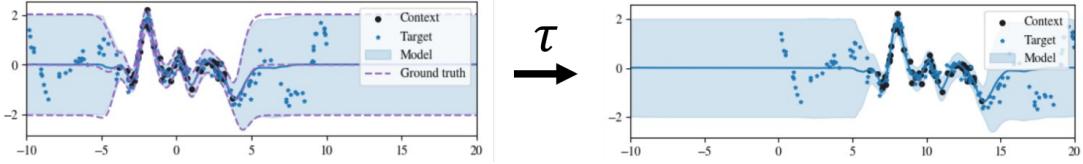


Figure 1.2.1: The Translation Equivariant property on a 1D dataset.

Figure 1.2.1 illustrates this property, when the input data on the left plot is shifted by a constant amount, the prediction should also shift by the same amount (right).

$$f : \mathbf{x} \rightarrow (\mathbf{x}, \hat{\mathbf{y}}) \quad (1.2.1)$$

$$f : \mathbf{x} + \mathbf{c} \rightarrow (\mathbf{x} + \mathbf{c}, \hat{\mathbf{y}}) \quad (1.2.2)$$

where \mathbf{x} is the input and $\hat{\mathbf{y}}$ is the output and \mathbf{c} is a constant shift in the input. This property allows the model to be more robust and generalize better to unseen data, particularly in the case of stationary data.

Off-the-Grid Generalization: *The model should be able to work on off-the-grid data points.* Off-the-grid data points are the data points that are not in a regular gridded structure, for example, images that have missing pixel values are off-the-grid. Traditional models like Convolutional Neural Networks (CNNs) are not able to operate on off-the-grid data points since they require a regular structure to apply the convolution operation. By making the model off-the-grid generalizable, we can create models that can work on many types of data and easily handle missing data points. It also allows the model to generalize to regions of the input space that it has not seen during training. Tasks like image inpainting, where the model is required to fill in missing pixels in an image, can benefit from this property.

TODO

Add a figure to illustrate off-the-grid generalization

Neural Processes (NPs) [Garnelo, Schwarz, et al. 2018] are a class of models that satisfy the above properties. The framework underlying NPs is general purpose, and thus can be modified with a variety of neural network architectures.

1.3 Aims and Objectives

In this work, we aim to implement and compare two different neural network architectures for Neural Processes, the first being based on a Convolutional Neural Network (CNN) called Convolutional Neural Processes (ConvNP) and the second being based on a Transformer architecture called Transformer Neural Processes (TNP). We aim to compare the two models on a variety of tasks and datasets to see how they perform in terms of generalization, scalability, and uncertainty estimation.

Furthermore, we investigate how the TNP can be optimized to achieve better performance. We then investigate new Transformer architectures that have better computational efficiency compared to the original Transformer architecture.

Our end goal is to develop a better understand of the properties of these models. We aim to figure out the best practices for using these models in different scenarios, where they perform well and where they do not. We also aim to understand the limitations of these models and how they can be improved.

Chapter 2

Neural Processes

2.1 Introduction

Neural Process (NP) is a meta-learning framework introduced in [Garnelo, Rosenbaum, et al. 2018; Garnelo, Schwarz, et al. 2018] that can be used for few-shot uncertainty aware meta learning. There exists two variants of the Neural Process, the Conditional Neural Process (CNP) and the Latent Neural Process (LNP), whilst we will discuss the differences between the two in this chapter, we will focus on the CNP for the majority of the project and hence we will implicitly refer to CNP as NP.

The main idea behind Neural Processes is to learn a distribution over the input locations we want to predict conditioned on the training data. In the NP literature we refer the training data as the context set and the input locations we want to predict the output for as the target set. The model is trained on a set of context-target pairs and then tested on a new set of context-target pairs to see how well it can generalize to new target sets.

2.2 Architecture

2.2.1 Conditional Neural Processes

Conditional Neural Processes (CNPs) [Garnelo, Rosenbaum, et al. 2018] was one of the two original Neural Processes introduced by Garnelo et al. in 2018. The general framework for a CNP requires us to take a context set $\mathcal{C} = \{\mathcal{C}_i\}_{i=1}^{N_c}$ where each context point \mathcal{C}_i is a input-output pair $\mathcal{C}_i = (\mathbf{x}_i^{(c)}, \mathbf{y}_i^{(c)})$ and a target set $\mathcal{T} = \{\mathbf{x}_i^{(t)}\}_{i=1}^{N_t}$ where each target point $\mathbf{x}_i^{(t)}$ is an input point we want to predict the output for.

We firstly encode each data point in the context set \mathcal{C}_i into an embedding using an encoder network.

$$\mathbf{r}(\mathcal{C}_i) = \text{Enc}_\theta(\mathcal{C}_i) = \text{Enc}_\theta([\mathbf{x}_i^{(c)}, \mathbf{y}_i^{(c)}]) \quad (2.2.1)$$

Where \mathbf{r} is the embedding of the context set \mathcal{C}_i and θ are the parameters of the encoder. Then we process the embeddings of the context sets to get a global representation of the dataset.

$$\mathbf{R}(\mathcal{C}) = \text{Process}(\{\mathbf{r}(\mathcal{C}_i)\}_{i=1}^D) \quad (2.2.2)$$

We require this ‘processing‘ to be **permutation invariant**, so typically it is a simple summation of the embeddings. The global representation \mathbf{R} is then used to condition the decoder to predict the outputs of the target set to give us a posterior distribution over the outputs $\mathbf{y}_i^{(t)}$.

$$p(\mathbf{y}_i^{(t)} | \mathbf{x}_i^{(t)}, \mathcal{C}) = \text{Dec}_\theta(\mathbf{x}_i^{(t)}, \mathbf{R}(\mathcal{C})) \quad (2.2.3)$$

The overall architecture is shown in Figure 2.2.1.

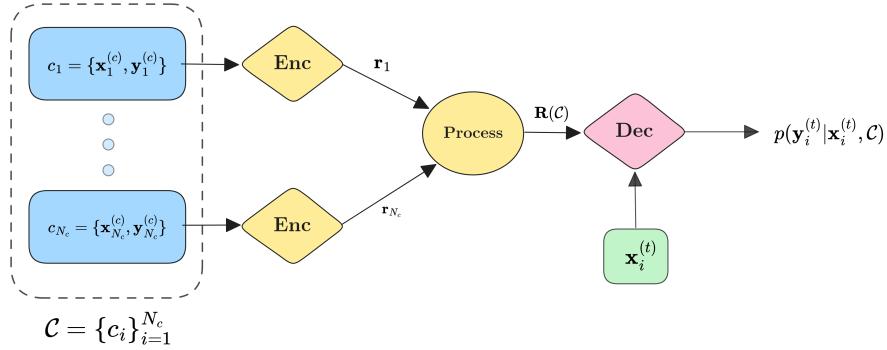


Figure 2.2.1: CNP Architecture: The model vectorizes each individual data point \mathcal{C}_i in the context set \mathcal{C} and then processes/aggregates them to obtain a global representation $\mathbf{R}(\mathcal{C})$ which is then used to condition the decoder to predict a distribution over the target points $\mathbf{y}^{(t)}$.

In the original CNP paper, the encoder and decoder are implemented as simple Multi-Layer Perceptrons (MLPs) and the processing is implemented as a mean operation, this happens to be an implementation off the ‘DeepSet’ architecture [Zaheer et al. 2018].

Importantly, CNPs make the strong assumption that the posterior distribution *factorizes* over the target points. This means that the posterior distribution over the target points is independent of each other.

$$p(\mathbf{y}^{(t)} | \mathbf{x}^{(t)}, \mathcal{C}) \stackrel{(a)}{=} \prod_{i=1}^{N_t} p(y_i^{(t)} | x_i^{(t)}, \mathbf{R}(\mathcal{C})) \stackrel{(b)}{=} \prod_{i=1}^{N_t} \mathcal{N}(y_i^{(t)} | \mu_i, \sigma_i^2) \quad (2.2.4)$$

$$(2.2.5)$$

The benefit of this factorization assumption (a) is that the model can scale linearly with the number of target points with a tractable likelihood. However, this assumption means **CNPs are unable to generate coherent sample paths, they are only able to produce distributions over the target points**. Furthermore, we need to select a marginal likelihood for the distribution (b) which is usually a Heteroscedastic Gaussian Likelihood (Gaussian with a variance that varies with the input) [Garnelo, Rosenbaum,

et al. 2018]. This adds an assumption as we have to select a likelihood for the distribution which may not be appropriate for the data we are modeling.

As the likelihood is a Gaussian, the model can be trained using simple maximum likelihood estimation (MLE) by maximizing the negative log-likelihood (\mathcal{L}) of the target points.

$$\mathcal{L} = \mathbb{E}_{(\mathcal{C}, \mathcal{T})} \left[- \sum_{i=1}^{|\mathcal{T}|} \log p(y_i^{(t)} | x_i^{(t)}, \mathcal{C}) \right] \quad (2.2.6)$$

2.3 Performance of Vanilla NP

Whilst the Vanilla CNP using DeepSets is flexible and scalable, in reality it is unable to perform well on more complicated and higher dimensional data since the model is unable to learn a good representation of the data using a simple MLP and summation operation.

Could we replace the encoder and decoder with more powerful networks? And if so, what would be the best architecture to use? We aim to answer in this project by exploring the use of a Convolutional Neural Network (CNN) and a Transformer as encoders of our NP. CNNs and Transformers have been shown to perform well on a variety of tasks and at scale, thus we hypothesize that they will be able to learn a better representation of the context set and thus improve the performance of the NP. Both are bound to have their unique advantages and disadvantages which we will explore in the following chapters by firstly exploring the Convolutional Neural Process (ConvCNP) and then the Transformer Neural Process (TNP).

Chapter 3

Convolutional Neural Processes

3.1 Motivation

Convolutional Neural Networks (CNNs) are a class of neural networks that are renowned for being translation equivariant, they have been widely used in image processing tasks and have shown to be able to learn spatial patterns in the data via convolutional filters. CNNs are very powerful and have been the state of the art in many image processing tasks making them a desirable backbone for a Neural Process. Hence, the motivation for Convolutional Neural Processes (ConvCNPs) [Gordon et al. 2020]. We will briefly discuss the ConvCNP model and its architecture however for a more detailed explanation, we refer the reader to the original paper.

3.2 Model

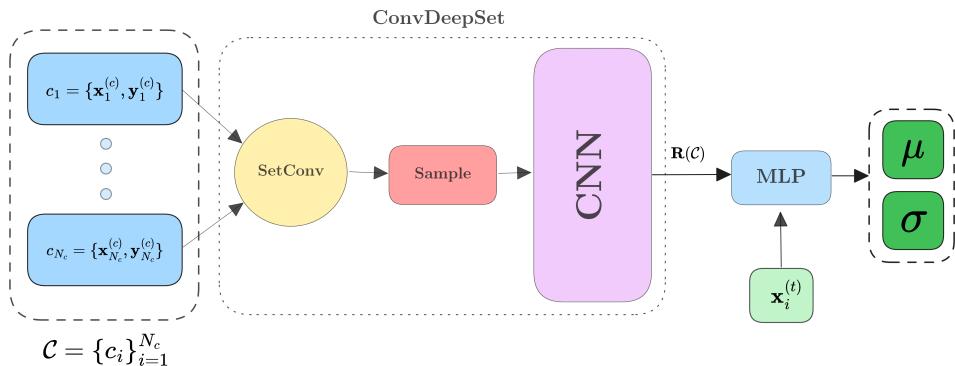


Figure 3.2.1: ConvCNP Architecture

ConvCNPs [Gordon et al. 2020] Figure 3.2.1 utilize CNNs, more specifically the UNet architecture. However, a few hurdles are preventing us from directly plugging data into a CNN.

CNNs operate on **on-grid data**, meaning that the data is on a regular grid, for example, an image. However, the data we are working with is sometimes **off-grid data**, for example time series dataset with irregular timestamps. Furthermore, to ‘bake-in’ the

TE property, we need to be able to shift the data in the input space and the output space in the same way. This is not trivial when using standard vector representations of the data. [Gordon et al. 2020] propose a solution to this problem by using **functional embeddings** to model the data. Functional embeddings are a way to represent data as a function that has a trivial translation equivariance property.

These functional embeddings are created using the **SetConv** operation. The SetConv operation is a generalization of the convolution operation that operates on sets of data, it takes a set of input-output pairs and outputs a continuous function. The SetConv operation is defined as follows:

$$\text{SetConv}(\{\mathbf{x}_i, \mathbf{y}_i\}_{i=1}^N)(x) = \sum_{i=1}^N [1, \mathbf{y}_i]^T \mathcal{K}(\mathbf{x} - \mathbf{x}_i) \quad (3.2.1)$$

Where \mathcal{K} is a kernel function that measures the distance between the query x and the data point \mathbf{x}_i .

This operation has some key properties:

- We append 1 to output \mathbf{y}_i when computing the SetConv, this acts as a flag to the model so that it knows which data points are observed and which are not. Say we have a data point of $\mathbf{y}_i = 0$, then if we did not append the 1, the model would not be able to distinguish between an observed data point and an unobserved data point (as both would be 0).
- The ‘weight’ of the Kernel depends only on the relative distance between points on the input space which means that the model is translation equivariant.
- The summation over the data points naturally introduces **Permutation Invariance** to the model, meaning that the order of the data points does not matter.

In [Gordon et al. 2020], they refer to the addition of the 1 as the **density channel**, this channel is used to distinguish between observed and unobserved data points.

Now that we have a function representation of the context set, we need to sample it/discretize it at **evenly** spaced points. Thus converting the data into a **on-grid** format. Now it can be fed into a CNN. Now that we have a CNN that operates on the data, we need to convert it back to a continuous function. This is done again by using the **SetConv** operation. The final output of the encoder is a continuous function that represents the context set which can be queried at any point in the input space using the target set.

$$R(\mathbf{x}_t) = \text{SetConv}(\text{CNN}(\{\text{SetConv}(\mathcal{C})(\mathbf{x}_d)\}_{d=1}^D))(\mathbf{x}_t) \quad (3.2.2)$$

The decoder is a simple MLP that takes the output of the encoder and the target set as input and outputs the mean and variance of the predictive distribution $\mu(\mathbf{x}_t), \sigma(\mathbf{x}_t) = \text{MLP}(R(\mathbf{x}_t))$

Chapter 4

Transformer Neural Processes

4.1 Transformers

4.1.1 Introduction

The Transformer is a deep learning architecture introduced in [Vaswani et al. 2017]. Originally devised as a sequence-to-sequence model that uses attention to learn the relationships between the input and output sequences. Transformers have been revolutionary in the field of natural language processing (NLP) leading to the development of models like BERT [Devlin et al. 2019] and GPT [Brown et al. 2020]. They are also starting to gain traction in fields like computer vision [Dosovitskiy et al. 2021] reaching becoming the new state of the art. The attractiveness of transformers is their general purpose nature and ability to learn relationships with an infinite context window whilst being massively parallelizable. A brief overview of the transformer model is given below.

4.1.2 Embedding

Data points it into a vector representation called an *embedding* or *token* ensuring postional information is added to these tokens via a positional encoding. The embedding is a simple linear transformation of the input data $\mathbf{X} \in \mathbb{R}^{N \times D}$ where N is the number of data points and D is the dimensionality of the data.

4.1.3 (Self-)Attention

The attention mechanism is a way to learn the relationships between the input and output sequences. ‘Normal’ attention infers what the most important word/phrase in an input sentence is which is not very powerful. This is where the idea of *self-attention* comes in. Self-attention is a way to learn the relationships between the data in the input sequence itself (Note when we say attention from now on, we mean self-attention). Consider a language modelling task, the sentence `The quick brown fox jumps over the lazy dog` has strong attention between the data like `fox` and `jumps` representing an action, `brown` and `fox` representing the color of the fox and so on, then there are very weak attentions between data `quick` and `dog` representing the lack of relationship between the two data. Using self-attention we can learn these relationships between the data in the input sequence, this gives a powerful mechanism to learn the relationships between the

input and output sequences and thus allows the model to learn the translation of the input sequence.

In the transformer models we will use the embeddings $\mathbf{X} \in \mathbb{R}^{N \times D}$ as the input to generate a query $\mathbf{Q} \in \mathbb{R}^{N \times d_k}$, a key $\mathbf{K} \in \mathbb{R}^{N \times d_k}$ and a value $\mathbf{V} \in \mathbb{R}^{N \times d_v}$ matrices via a simple linear transformation matrix $\mathbf{W}_q \in \mathbb{R}^{D \times d_k}$, $\mathbf{W}_k \in \mathbb{R}^{D \times d_k}$ and $\mathbf{W}_v \in \mathbb{R}^{D \times d_v}$ respectively.

Where each row of the matrices is the query, key and value vectors for each data point in the input sequence.

The query, key and value matrices are then used to compute the attention matrix $\mathbf{A} \in \mathbb{R}^{N \times N}$ as follows:

$$\mathbf{A} = \text{softmax} \left(\frac{\mathbf{Q}\mathbf{K}^T}{\sqrt{d_k}} \right) \quad (4.1.1)$$

The intuition behind this is that we want to compute the similarity between the query and the key vectors as such we use the dot product between the query and key vectors. The softmax is used to normalize the attention matrix so that the rows sum to 1. The softmax is also scaled by $\sqrt{d_k}$ to prevent the softmax from saturating. The attention matrix is then used to compute the output matrix $\mathbf{H} \in \mathbb{R}^{N \times d_v}$ as $\mathbf{H} = \mathbf{AV}$

The overall attention function for a layer is given by:

$$\mathbf{H} = \text{Attention}(\mathbf{Q}, \mathbf{K}, \mathbf{V}) = \text{softmax} \left(\frac{\mathbf{Q}\mathbf{K}^T}{\sqrt{d_k}} \right) \mathbf{V} \quad (4.1.2)$$

4.1.4 Multi-Head Self-Attention

So far we have only computed the attention matrix \mathbf{A} once so the model only learns one attention relationship, however, we can take advantage of using multiple attention ‘heads’ in parallel to learn many attention relationships, this scheme is called the *Multi-Head Attention* (MHSA).

Each attention head is computed using simple dot product attention of a transformed query, key and value matrix. They are transformed by a simple linear layer (a matrix) which is unique for each head of the MHSA, $\mathbf{W}_q^{(i)} \in \mathbb{R}^{d_k \times d_k}$, $\mathbf{W}_k^{(i)} \in \mathbb{R}^{d_k \times d_k}$ and $\mathbf{W}_v^{(i)} \in \mathbb{R}^{d_v \times d_v}$ where $i \in [1, h]$ for a head count of h . Then the attention for the particular head is computed as follows:

$$\mathbf{H}^{(i)} = \text{Attention}(\mathbf{Q}\mathbf{W}_q^{(i)}, \mathbf{K}\mathbf{W}_k^{(i)}, \mathbf{V}\mathbf{W}_v^{(i)}) \in \mathbb{R}^{N \times d_v} \quad (4.1.3)$$

Then the output of the MHSA is the concatenation of the outputs of each head $\mathbf{H}^{(i)}$ (stacked on top of each other) multiplied by a learnable matrix $\mathbf{W}_O \in \mathbb{R}^{hd_v \times D}$ which transforms the concatenated output to the original dimensionality of the input sequence.

$$\text{MHSA}(\mathbf{Q}, \mathbf{K}, \mathbf{V}) = \text{concat}(\mathbf{H}^{(1)}; \mathbf{H}^{(2)}; \dots; \mathbf{H}^{(h)}) \mathbf{W}_O = \begin{bmatrix} \mathbf{H}^{(1)} \\ \mathbf{H}^{(2)} \\ \vdots \\ \mathbf{H}^{(h)} \end{bmatrix} \mathbf{W}_O \in \mathbb{R}^{N \times D}$$

4.1.5 Encoder

Now that we have covered the MHSA block, we can move on to the encoder of the transformer.

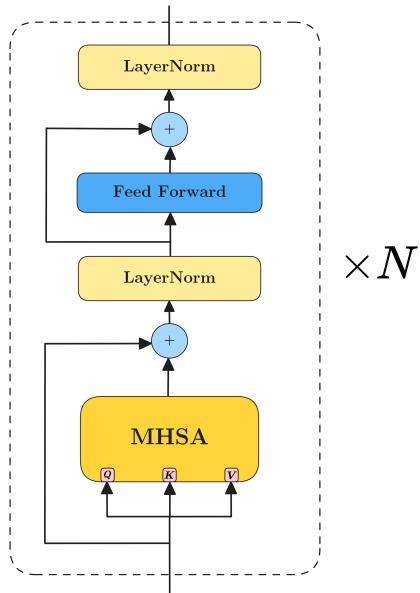


Figure 4.1.1: Transformer Encoder [Vaswani et al. 2017]

Figure 4.1.1 shows the encoder of the transformer model. The encoder is composed of a stack of N identical layers. Each layer is composed of two sub-layers, the MHSA and a simple feed-forward network. The output of each sub-layer is $\text{LayerNorm}(x + \text{Sublayer}(x))$ where x is the input to the sub-layer. The final output of the encoder is the output of the last layer which shall be denoted as $\mathbf{Y} \in \mathbb{R}^{N \times D}$.

Key Points

The Transformer Encoder Layer takes an input set of embeddings $\mathbf{X} \in \mathbb{R}^{N \times D}$ and outputs a set of embeddings $\mathbf{Y} \in \mathbb{R}^{N \times D}$ of the **same dimensionality** but with the **patterns of the input sequence learned**. It can be viewed as a function that takes a set and outputs a set of the same dimensionality.

4.2 Vanilla Transformer Neural Process

An Attention based encoder for the Neural Process was investigated in [Kim et al. 2019], though the results were impressive, the model fails to perform at larger scale and ‘tends

to make overconfident predictions and have poor performance on sequential decision-making problems’ [Nguyen and Grover 2023]. It is natural to think that the Transformer [Vaswani et al. 2017] could be used to improve the performance of the Neural Process, as Transformers have been proven to effectively model large scale data using attention mechanisms. [Nguyen and Grover 2023] introduced the Transformer Neural Process (TNP) which uses an *encoder-only* Transformer to learn the relationships between the context and target points via self-attention with appropriate masking.

4.2.1 Model Architecture

Similar to the standard Neural Process architecture we are required to encode data points within the context set into a vector representation, in language modelling literature we refer to this as tokenization. The tokenization is done by using a simple Multi-Layer Perceptron (MLP) to encode the data points into vector tokens with a configurable token dimension, D_{em} .

Where the TNP differs is that it also encodes the target points which is padded with zeros to represent the lack of value for the target data points, then both context and target tokens are passed into the transformer at the same time instead of computing a context representation and then passing the target points as in the standard NP.

A flag bit is introduced into the tokens to indicate whether the token is a context or target token. Consider the context dataset $\mathcal{C} = \{(\mathbf{x}_i, \mathbf{y}_i)\}_{i=1}^{N_c}$ and target set $\mathcal{T} = \{(\mathbf{x}_i)\}_{i=N_c+1}^N$, the model will encode a data point into a token \mathbf{t}_i as follows:

$$\mathbf{t}_i = \begin{cases} \text{MLP}(\text{cat}[\mathbf{x}_i, 0, \mathbf{y}_i]) & \text{if } i \leq N_c \\ \text{MLP}(\text{cat}[\mathbf{x}_i, 1, \mathbf{0}]) & \text{if } i > N_c \end{cases}$$

Where we use a flag bit of 0 to indicate a context token and 1 to indicate a target token, **cat** is the concatenation operation and **0** is a vector of zeros of the same dimension as \mathbf{y}_i .

Now that the data points are tokenized we can pass them through the Transformer encoder to learn the relationships between the context and target points. Importantly the Transformer encoder is masked such that the target tokens can only attend to the context tokens and previous target tokens. Alternatively one perform self attention on the context tokens then cross attention between the context and target tokens giving a more efficient implementation [Feng et al. 2022], we use this implementation in our experiments.

We can now pass the tokens through the Transformer encoder layer several times to learn the relationships between the context and target points and generate a vector output. The output of the Transformer encoder is then passed through a Multi-Layer Perceptron (MLP) to generate the mean and variance of the predictive distribution of the target points.

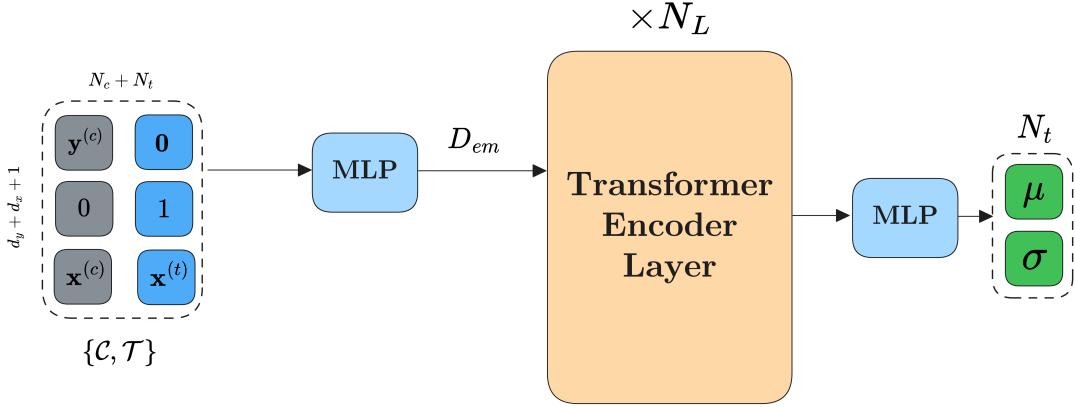


Figure 4.2.1: Vanilla Transformer Neural Process Architecture. All the context and target sets are tokenized and passed through the Transformer Layer. The output of the Transformer Layer is passed through an MLP to generate the mean and variance of the predictive distribution.

4.2.2 Performance

One of the key benefits of Transformers (and attention) is the ability to have a global view of the data, analogous to an ‘infinite receptive field’ in Convolutional Neural Networks. This allows the model to learn very complicated relationships across many length scales in the data. However, this can be a double-edged sword as the model can overfit to the data within its training region and fail to generalize to unseen data.

Translation Equivariance is a key property behind CNNs which allow them to generalize excellently to unseen data by learning features irrespective of their position in the data. Such feature learning is critical to modelling real world data where the positions of the data do not matter as much as the relative relationships between the data points, e.g. in image classification the position of the object in the image does not matter as much as the features of the object itself. Transformers lack this property, causing them to generate random predictions when the data is shifted out of context as shown in Figure 4.2.2.

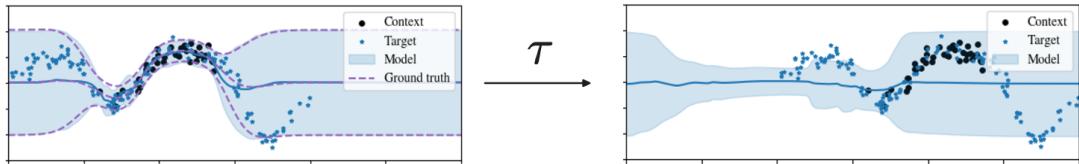


Figure 4.2.2: Vanilla Transformer Neural Process failing to generalize in out of context data. The TNP successfully models the data in the context region (left) but fails to make predictions when the data is shifted (right).

What if we could combine the best of both worlds? Could we create a Translation Equivariant Transformer Neural Process (TETNP) which possesses the global view of the Transformer and the generalization properties of the CNN? This is the question we will investigate in the next section and is the main contribution of this work and [Ashman et al. 2024].

4.3 Translation Equivariant TNP

Recall, Translation Equivariance requires the model to yield the same output when the input is shifted. Such property must imply that the model only learns the relative distances between the data points ($\mathbf{x}_i - \mathbf{x}_j$) and *not* the absolute positions of the data points. How could one enforce such a property in the Transformer Neural Process? The solution used is very simple, we add a term to the attention mechanisms in the Transformer to enforce the Translation Equivariance property!

Consider the standard attention mechanism in the Transformer, the attention weights are computed as:

$$\text{Attention}(\mathbf{Q}, \mathbf{K}, \mathbf{V}) = \text{softmax}(\mathbf{E}) \mathbf{V} \quad (4.3.1)$$

$$E_{ij} = \frac{\mathbf{q}_i \cdot \mathbf{k}_j}{\sqrt{d_k}} \quad (4.3.2)$$

To create a Translation Equivariant Attention mechanism we add a term to the attention weights which enforces the Translation Equivariance property. The new attention weights are computed with the following equation:

$$\text{Attention}(\mathbf{Q}, \mathbf{K}, \mathbf{V}, \mathbf{X}) = \text{softmax}(\mathbf{E} + F(\Delta)) \mathbf{V} \quad (4.3.3)$$

$$E_{ij} = \frac{\mathbf{q}_i \cdot \mathbf{k}_j}{\sqrt{d_k}} \quad (4.3.4)$$

$$\Delta_{ij} = \mathbf{x}_i - \mathbf{x}_j \quad (4.3.5)$$

Where F is some function that introduces non-linearity into the attention weights and is applied to each entry of the matrix independently ($F(\Delta)_{ij} = F(\Delta_{ij})$), we will investigate the effect of different functions in the experiments later on. It is clear to see that if all the input locations are shifted by a constant τ , the relative term will remain the same $F(\Delta_{ij}) = F(\mathbf{x}_i + \tau - \mathbf{x}_j - \tau) = F(\mathbf{x}_i - \mathbf{x}_j)$ and the attention weights will remain the same.

Importantly since we are enforcing the Transformer to learn the relative distances between the data points, we must remove the \mathbf{x} values from the tokenization of the data points. The tokenization of the data points is now:

$$\mathbf{t}_i = \begin{cases} \text{MLP}(\text{cat}[0, \mathbf{y}_i]) & \text{if } i \leq N_c \\ \text{MLP}(\text{cat}[1, \mathbf{0}]) & \text{if } i > N_c \end{cases}$$

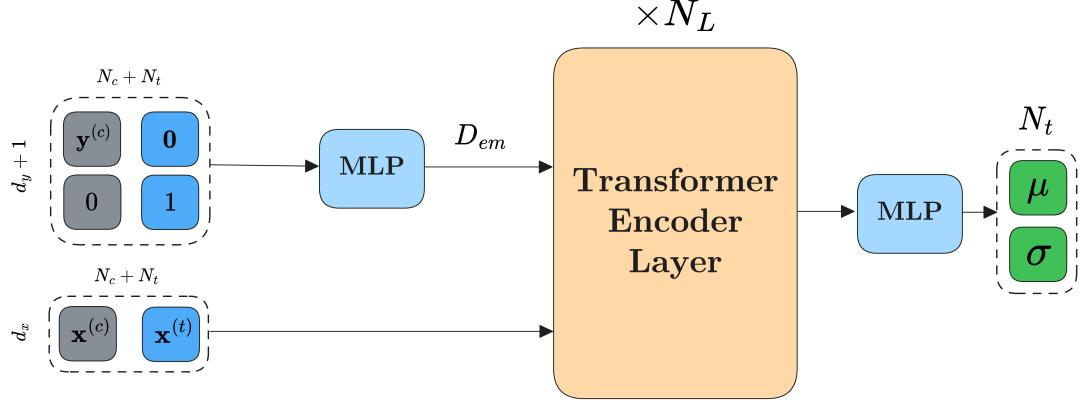


Figure 4.3.1: Translation Equivariant Transformer Neural Process Architecture. All the outputs of the context and target sets are tokenized and passed through the Transformer Layer, whilst the input locations are passed in raw to the attention mechanism to enforce the Translation Equivariance property. The output of the Transformer Layer is passed through an MLP to generate the mean and variance of the predictive distribution.

A benefit of removing the \mathbf{x} values from the tokenization is that the tokens only need to encode the \mathbf{y} values, reducing the dimensionality of the tokens whilst also separating the encoding for \mathbf{x} and \mathbf{y} values which could be beneficial for the model as in the vanilla TNP the model loses distinction between the \mathbf{x} and \mathbf{y} values since they are embedded together.

Chapter 5

Experimentation on 1D Datasets

We will use 1D datasets to firstly optimize the hyperparameters of the TNP models and use this to compare the performance of the TNP and ConvNP models. The metric used to compare the models is the validation loss of the models on unseen data. Furthermore, we will look at the plots of the predictions to observe the behavior of the models especially in the regions outside the training data.

5.1 Datasets

5.1.1 Gaussian Process

We will use samples from a Gaussian Process to generate the data. The aim will be for our model to learn the underlying function of the Gaussian Process and fit the data just like the original Gaussian Process. The Gaussian Process is defined as:

$$f(x) \sim \mathcal{GP}(0, k(x, x')) \quad (5.1.1)$$

Where we use the squared exponential kernel:

$$k(x, x') = \exp\left(-\frac{(x - x')^2}{2l^2}\right) \quad (5.1.2)$$

With length scales being sampled from a uniform distribution $l \sim \mathcal{U}(\log(-0.101), \log(0.101))$. We choose to sample N_c context points from the Gaussian Process and use these as the training data for our models. We then sample N_t target points from the Gaussian Process and use these as the validation data for our models. Depending on the specific task we will either fix or vary the number of context points N_c and target points N_t .

5.1.2 Sawtooth

Whilst the GP is useful for testing the models on a smooth function, we also want to test the models on a more complex function, particularly one with discontinuities. We will use the sawtooth function for this purpose. The sawtooth function with period T is defined as:

$$f(x) = x - T \left\lfloor \frac{x}{T} \right\rfloor + n \quad (5.1.3)$$

Where n is some random noise which is sampled from a normal distribution $n \sim \mathcal{N}(0, 0.1)$. We will sample N_c context points from the sawtooth function and use these as the training data for our models. We then sample N_t target points from the sawtooth function and use these as the validation data for our models. Depending on the specific task we will either fix or vary the number of context points N_c and target points N_t .

5.2 Relative Attention Function

As mentioned in section 4.3 we need to pass the matrix of differences (Δ) between x values through a function F to apply non-linearities to the differences acting as hyperparameter of our mode. We will investigate using a simple linear function with no bias and gradient 1 ('identity') as a baseline. For non-linear functions, we will consider using a Gaussian Radial Basis Function (RBF) and a Multi-Layer Perceptron (MLP). The functions are defined below:

$$F_{\text{identity}}(\Delta) = \Delta \quad F_{\text{RBF}}(\Delta) = \exp\left(-\frac{\Delta^2}{2\sigma^2}\right) \quad F_{\text{MLP}}(\Delta) = \text{MLP}(\Delta) \quad (5.2.1)$$

Where σ is a hyperparameter of the RBF function and $\text{MLP}(\Delta)$ is a 2-layer MLP with ReLU activation functions.

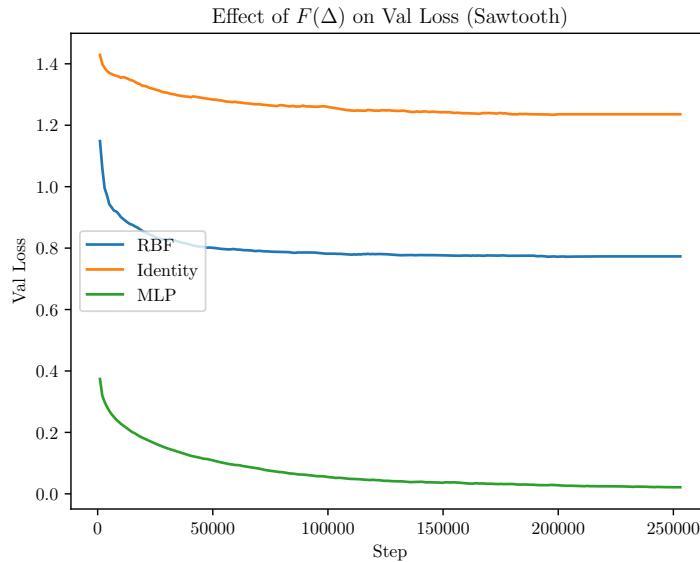


Figure 5.2.1: Relative Attention Functions on Validation Loss for the TETNP on the 1D Sawtooth Dataset. Lower validation loss is better.

Figure 5.2.1 shows the validation loss curves for the TNP with different relative attention functions. We can see that the MLP function performs significantly better than the other two functions. This is likely due to the MLP being able to **learn** whilst the other two

functions are fixed. We can also see that the RBF function performs better than the identify function since the effect of adding the raw difference affects the dot product of the attention mechanism. So we can conclude that the MLP function is the best function to use for the TNP. The computational cost of using the MLP function is also not significantly higher than the other two functions, since the MLP is very small.

5.3 Optimizing Hyperparameters

The multi-headed attention mechanism in the Transformer Encoder has three hyperparameters that we will investigate. These are the token embedding dimension of the data (D_{em}), the number of attention heads (N_h) and the embedding dimension of the attention heads (D_h). We will investigate how changing these hyperparameters affects the performance of the model. To gauge the effect of these hyperparameters, we use a 1 million parameter model and reduce the value of one of the hyperparameters and see the effect on the validation loss. This process is repeated for the other hyperparameters to see which hyperparameters have the most effect on the validation loss.

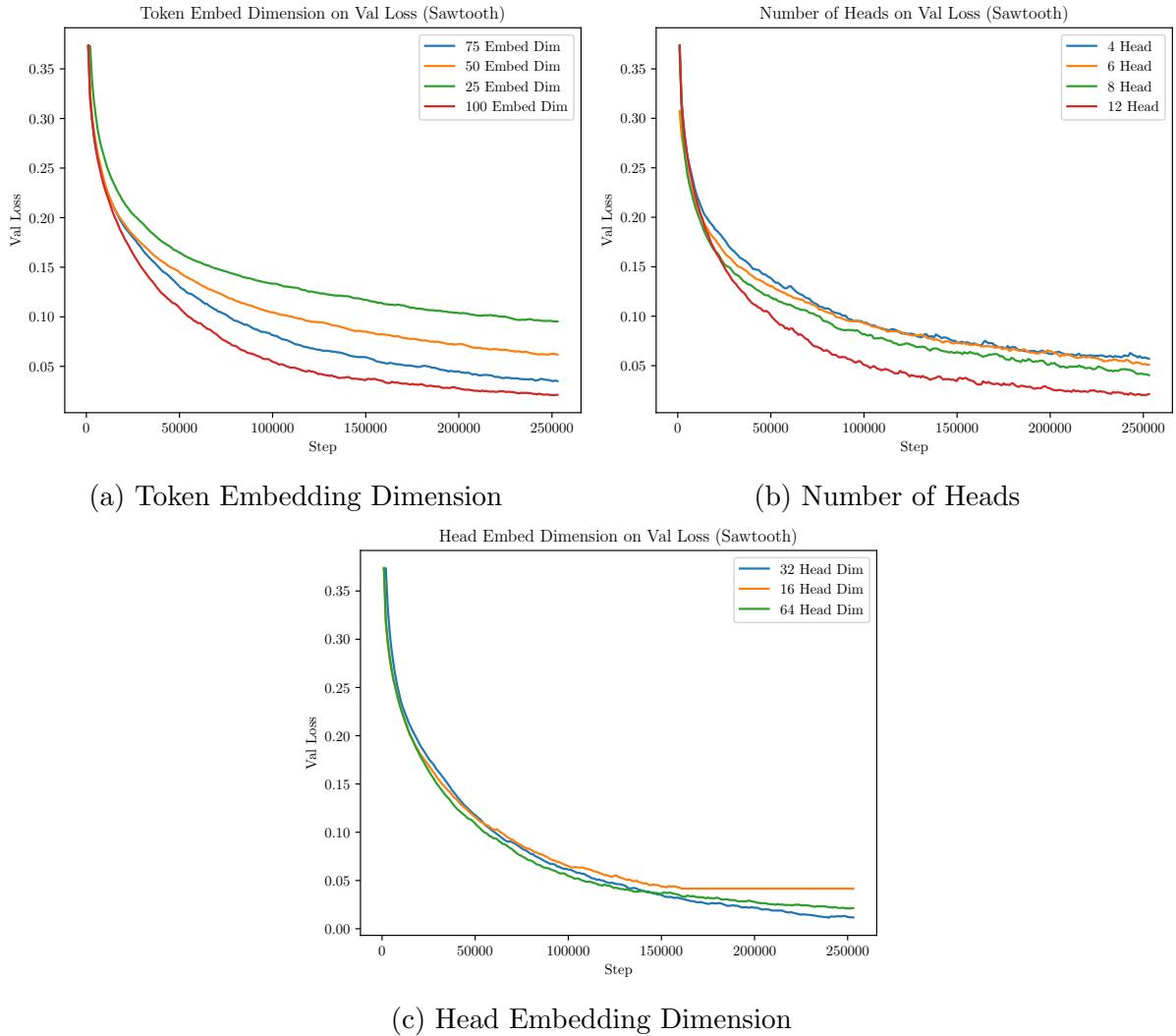


Figure 5.3.1: Hyperparameter Selection

From Figure 5.3.1 it is clear that reducing the token embedding dimension has a significant effect on the validation loss, with the number of heads making a smaller difference and the head embedding dimension making very little difference. This highlights the importance of the token embedding dimension even when using low one-dimensional data as the input data (in this case sawtooth function). Furthermore, the head embedding dimension has very little effect on the validation loss, so we can set this to a small value to reduce the number of parameters in the model and distribute the parameters to the other hyperparameters. Using this knowledge, we will now investigate how to select the hyperparameters using a *constant parameter budget* of 1 million parameters.

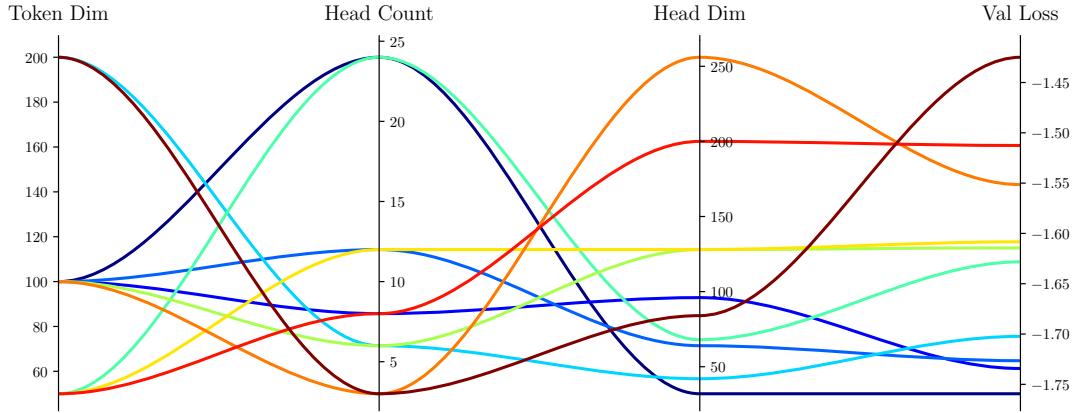


Figure 5.3.2: Constant Parameter Budget Hyperparameter Selection

Figure 5.3.2 shows a parallel coordinates plot of the validation loss for different hyperparameter configurations, where dark blue is the best and dark red is the worst. We can see that the model that performs the best has a very high token embedding dimension, high headcount and low head embedding dimension, which is consistent with the previous results. We can also see that if we go too high in the token embedding dimension (200), the model performs worse as we have to sacrifice the number of heads in the transformer. These results give us a set of best practices for selecting hyperparameters for the TNP: high token embedding dimension, high number of heads and low head embedding dimension.

5.4 TNP vs ConvNP

Finally using our ‘best’ TNP model, we will compare the performance of the ConvNP and TNP on fitting sawtooth and GP (using EQ Kernel) data using 1 million parameters.

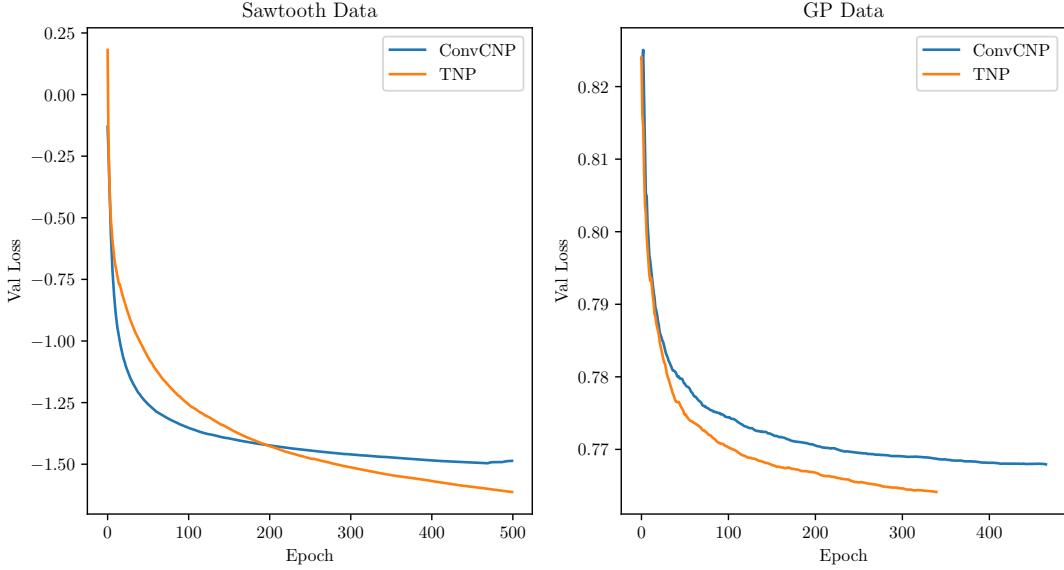


Figure 5.4.1: ConvNP vs TNP on Sawtooth and GP Data

Figure 5.4.1 shows the validation loss curves for the ConvNP and TNP on sawtooth and GP data. We can see that the validation loss for the TNP is lower than the ConvNP for both datasets which is very promising and indicates the TNP is a better model than the ConvNP.

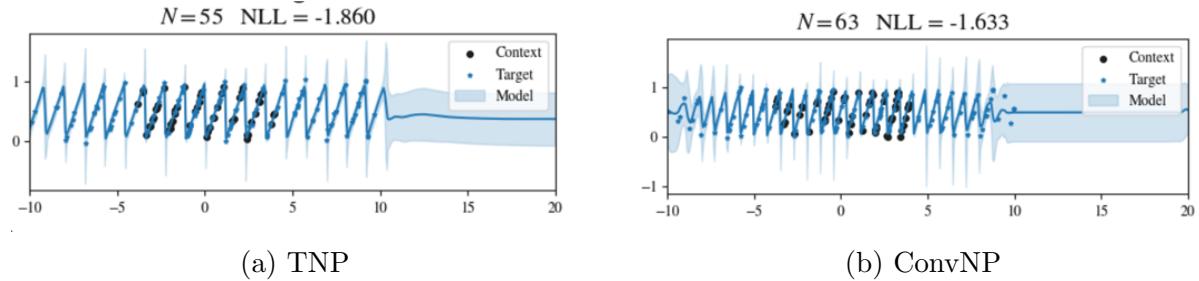


Figure 5.4.2: ConvNP vs TNP on Sawtooth Data

Inspecting the model fits on sawtooth data Figure 5.4.2, we observe that the TNP can extrapolate the structure of the sawtooth function beyond the range of the context set (black points) whilst the ConvNP performs decently but fails to retain the structure as well as the TNP, since the amplitude of the sawtooth reduces the further away from the context set. Hence we can conclude that the TNP can better understand the structure of the data than the ConvNP.

5.4.1 Computational Complexity

Table 5.4.1 shows that the TETNP uses drastically more memory than the ConvNP for the same number of context and target points. This is due to the quadratic complexity of the TETNP which scales with $\mathcal{O}(N_c^2 + N_c N_t)$ compared to the linear complexity of the ConvNP which scales with $\mathcal{O}(N_c D_x^3 + N_t D_x)$. Whilst on the 1D dataset the ConvNP is more memory efficient, the TETNPs complexity does not scale with the number of

N_c	N_t	ConvNP Memory (MB)	TETNP Memory (MB)
10	10	18	13
100	10	16	18
1000	10	16	339
5000	10	124	9357
<hr/>			
10	1000	36	399
100	1000	36	469
1000	1000	36	1510
5000	1000	124	13480

Table 5.4.1: Memory Usage in inference of the ConvNP and TETNP on 1D data using N_c context points and N_t target points.

dimensions of the data, whilst the ConvNP *scales cubically with the number of dimensions*. This may suggest that the TETNP is more suitable for high-dimensional data than the ConvNP since the TETNP has a fixed memory cost per dimension. We will investigate this in the next chapter on 2D datasets to see if the discrepancy in memory usage is still present.

Chapter 6

Experimentation on 2D Datasets

6.1 Datasets

6.1.1 Gaussian Process

The 2D Gaussian Process is the natural extension of the 1D Gaussian Process described in subsection 5.1.1 where we use the squared exponential kernel. We continue to use the same range of lengthscale across both input dimensions as the 1D Gaussian Process.

The following plots show some samples from the GP dataset.

TODO

Add plots of GP dataset

6.1.2 Sawtooth

The 2D Sawtooth dataset is the natural extension of the 1D Sawtooth dataset described in subsection 5.1.2. We continue to use the same period T and noise n across both input dimensions as the 1D Sawtooth dataset.

The following plots show some samples from the Sawtooth dataset.

TODO

Add plots of Sawtooth dataset

6.1.3 Restricted Sawtooth

By accident when generating the 2D Sawtooth dataset, we ended up restricted the ‘direction of travel’ of the sawtooth function to the line of $x_1 = x_2$ or $x_1 = -x_2$. This was not intentional but when training both models on this dataset, we found very interesting results. As the models only learn a subset of the ‘full sawtooth’ function, we can see how well the models can generalize to samples from the full sawtooth function.

The following plots show some samples from the Restricted Sawtooth dataset.

6.2 Post or Pre Relative Attention Function

TODO

Highlight MLP

In our original formulation of the TETNP (section 4.3) we pass the matrix of differences (Δ) between x values through a function F to apply non-linearities then add it to the dot product attention Equation 4.3.3, whilst this performs well we can also consider applying this non-linearity after combining the dot product attention and the relative attention, this method is called the ‘Post Relative Attention Function’.

$$\text{Attention}(\mathbf{Q}, \mathbf{K}, \mathbf{V}, \mathbf{X}) = \text{softmax}(\mathbf{E}) \mathbf{V} \quad (6.2.1)$$

$$\text{Pre: } \mathbf{E}_{ij} = \mathbf{D}_{ij} + \text{MLP}(\Delta_{ij}) \quad (6.2.2)$$

$$\text{Post: } \mathbf{E}_{ij} = \text{MLP}(\text{cat}[\mathbf{D}_{ij}, \Delta_{ij}]) \quad (6.2.3)$$

Where

$$\mathbf{D}_{ij} = \mathbf{q}_i \cdot \mathbf{k}_j / \sqrt{d_k} \quad \Delta_{ij} = \mathbf{x}_i - \mathbf{x}_j \quad (6.2.4)$$

We will investigate the performance of the TETNP with the ‘Post Relative Attention Function’ compared to the original ‘Pre Relative Attention Function’. We choose to use the Sawtooth dataset as it is more difficult to learn than the Gaussian Process dataset and will show the differences between the two models more clearly.

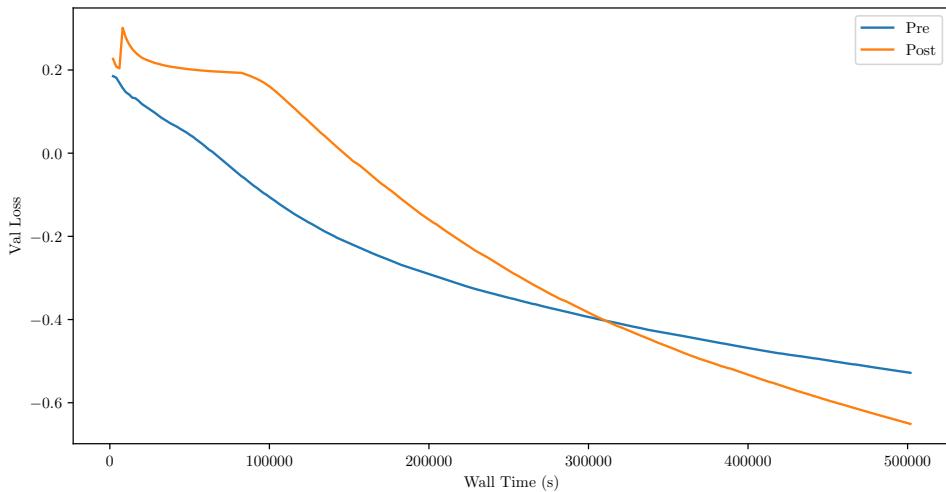


Figure 6.2.1: Validation Loss of TETNP with the ‘Post Relative Attention Function’ and ‘Pre Relative Attention Function’ on the 2D Sawtooth Dataset. Lower validation loss is better.

The results show that the TETNP with the ‘Post Relative Attention Function’ outperforms the TETNP with the ‘Pre Relative Attention Function’ by quite a large margin. Trivially this makes a lot of sense as the Post function can further refine the dot product attention through the MLP whilst in the ‘Pre’ function the MLP is *only* applied to the Δ matrix. The computational complexity of these two functions are not too different as the MLP are small and applied to the same size matrices.

6.3 ConvNP vs TETNP

We have discovered in the 1D section that the TETNP outperforms the vanilla TNP in all cases, hence for the 2D experiments we will only compare the ConvNP to the TETNP. When performing our experiments we will use models which are both 1 million parameters in size, to ensure a fair comparison.

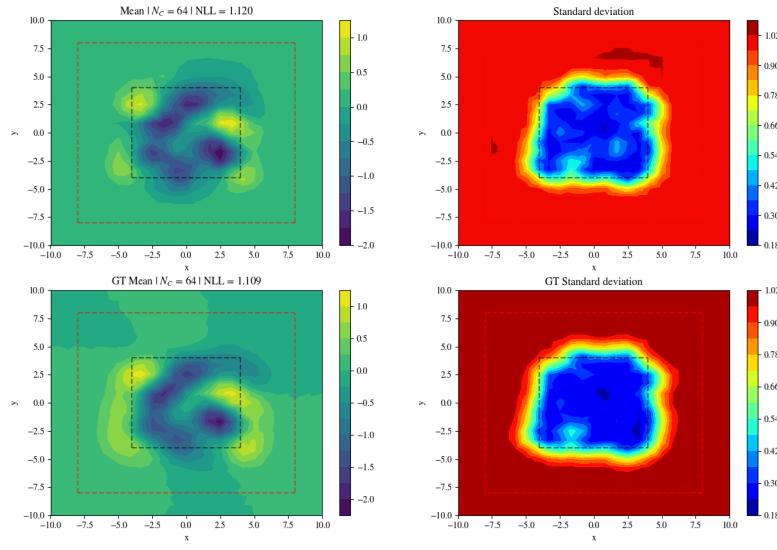
6.3.1 Gaussian Process

As mentioned previously, the Gaussian Process dataset is not very difficult to learn and the ConvNP and TETNP both perform very well on this dataset with the TETNP outperforming the ConvNP by a small margin as shown in Table 6.3.1.

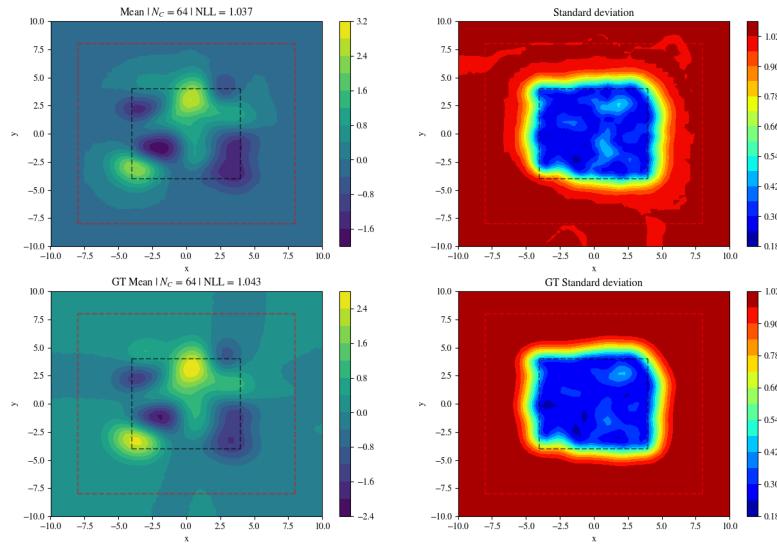
Model	Validation Loss
ConvNP	1.168
TETNP	1.134

Table 6.3.1: Validation Loss of ConvNP and TETNP on the 2D Gaussian Process dataset after training for 3 hours using 1 million parameters models and 64 context points. Lower is better.

Observing the samples from the ConvNP and TETNP for low frequency Figure 6.3.1 and high frequency functions Figure 6.3.2, we can see both models are able to generate very similar predictions to the ground truth GP. The TETNP performs slightly better than the ConvNP in the low frequency case, but the ConvNP performs better in the high frequency case for these samples. However overall the TETNP performs better than the ConvNP on the Gaussian Process dataset.

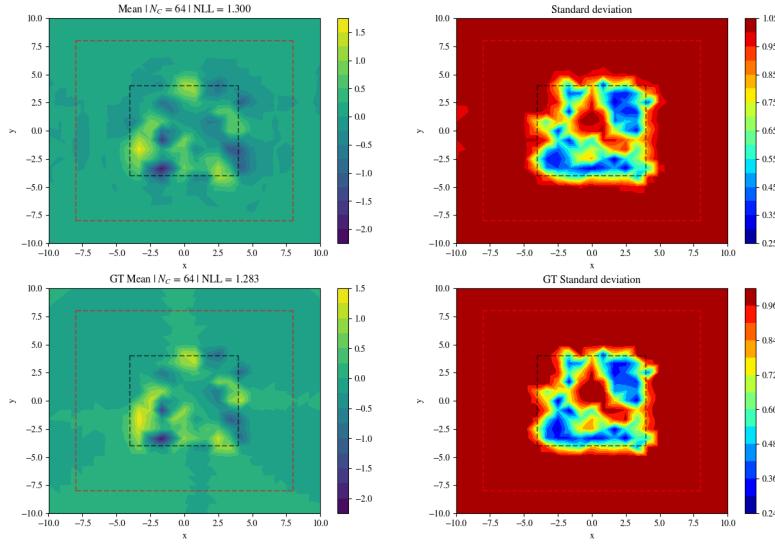


(a) ConvNP (top plot is the model prediction and bottom is the ground truth GP)

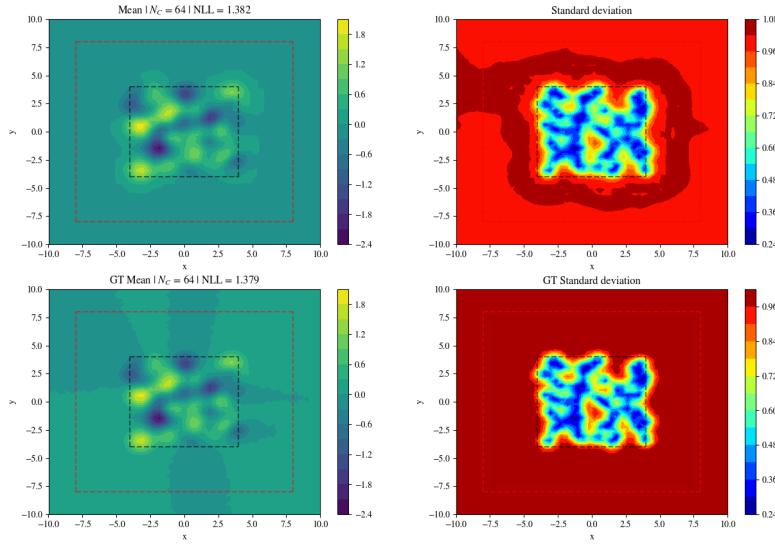


(b) TETNP (top plot is the model prediction and bottom is the ground truth GP)

Figure 6.3.1: Samples from ConvNP and TETNP on a low frequency 2D Gaussian Process.



(a) ConvNP (top plot is the model prediction and bottom is the ground truth GP)

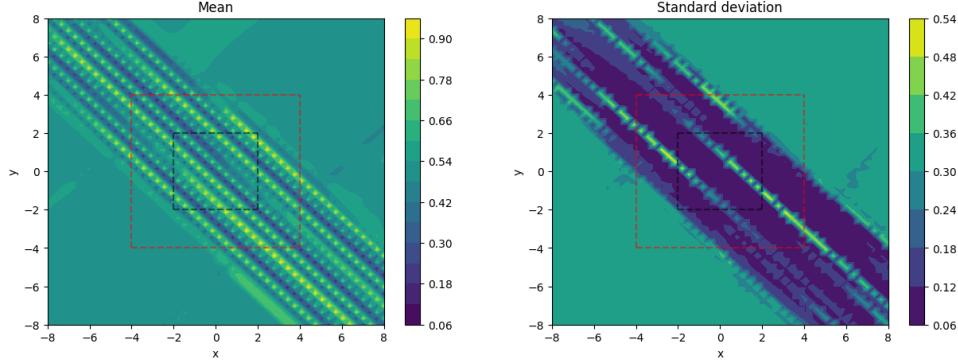


(b) TETNP (top plot is the model prediction and bottom is the ground truth GP)

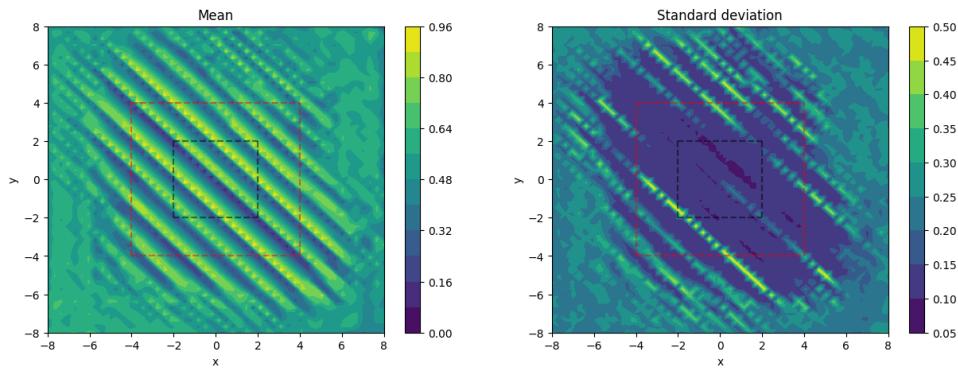
Figure 6.3.2: Samples from ConvNP and TETNP on a high frequency 2D Gaussian Process.

6.3.2 Restricted Sawtooth and Rotational Equivariance

As previously stated the restricted sawtooth dataset is a subset of the full sawtooth dataset which restricts the ‘direction of travel’ of the sawtooth function to the line of $x_1 = x_2$ or $x_1 = -x_2$. By training both models on this dataset we can see how well the models can generalize.



(a) TETNP



(b) ConvNP

Figure 6.3.3: Samples from TETNP and ConvNP on the Restricted Sawtooth dataset. The region inside the black dotted box is the region the models were trained on and the region outside the black dotted box is the region the models were not trained on.

Figure 6.3.3 shows that the ConvNP performs excellently on this dataset, which is expected as CNNs have filters which learn features and patterns in the data explicitly, thus performs well on extrapolation tasks (outside the black dotted box region). The TETNP on the other hand struggles to generalize fully within the target region (red dotted box) and outside the target region. Instead, it learns to extrapolate the sawtooth along one axis, but not the other. This clearly highlights a limitation of the Transformer architecture which has very little interpretability hence producing weird results.

Is this TETNP able to generalize to the full sawtooth dataset? To answer this question we will simply run the TETNP on a rotated version of the restricted sawtooth dataset which is the full sawtooth dataset.

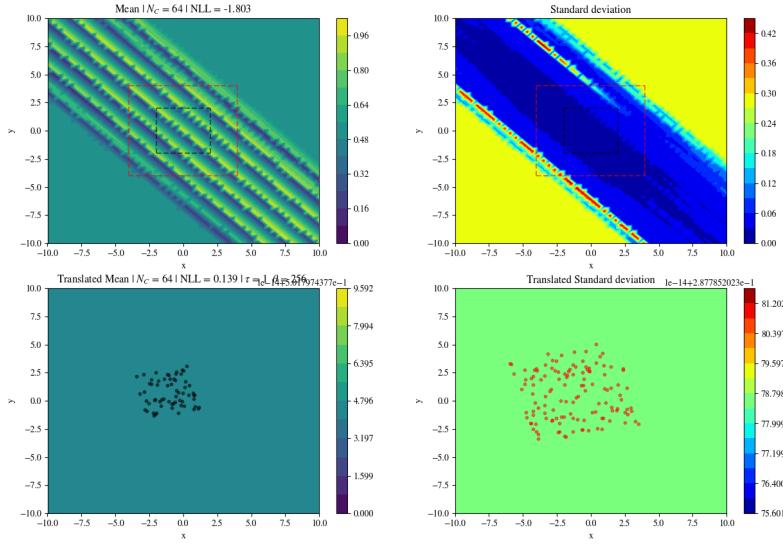


Figure 6.3.4: Samples from TETNP on the full Sawtooth dataset. The top plot is the model prediction on the restricted sawtooth dataset and the bottom plot is the model prediction on the full sawtooth dataset which is a rotated version of the restricted sawtooth dataset. The region inside the black dotted box is the region the model was trained on and the region outside the black dotted box is the region the model was not trained on.

In Figure 6.3.4 we explicitly reduced the target and context region size to allow for the TETNP to cover the full target region. The bottom plot shows the predictions when we rotate the context points by 256 degrees, clearly the TETNP completely fails, it just predicts a constant mean and standard deviation. *Could introducing rotational equivariance to the TETNP help it generalize to the full sawtooth dataset?*

Rotational Equivariance

To introduce rotational equivariance is fairly simple, in our formulation of the Translation Equivariance Attention, we use a matrix Δ which is the difference between the \mathbf{x} values of all the data points.

$$\text{Not RE : } \Delta_{ij} = \mathbf{x}_i - \mathbf{x}_j \quad (6.3.1)$$

$$\text{RE : } \Delta_{ij} = \|\mathbf{x}_i - \mathbf{x}_j\|_2 \quad (6.3.2)$$

To introduce rotational equivariance we can simply take the L2 norm of the Δ matrix which will give us the distance between all the data points. Distances are invariant to rotation, hence the TETNP should be rotationally equivariant. Using this new Δ matrix we can train the TETNP on the restricted sawtooth dataset and see if it can generalize to the full sawtooth dataset.

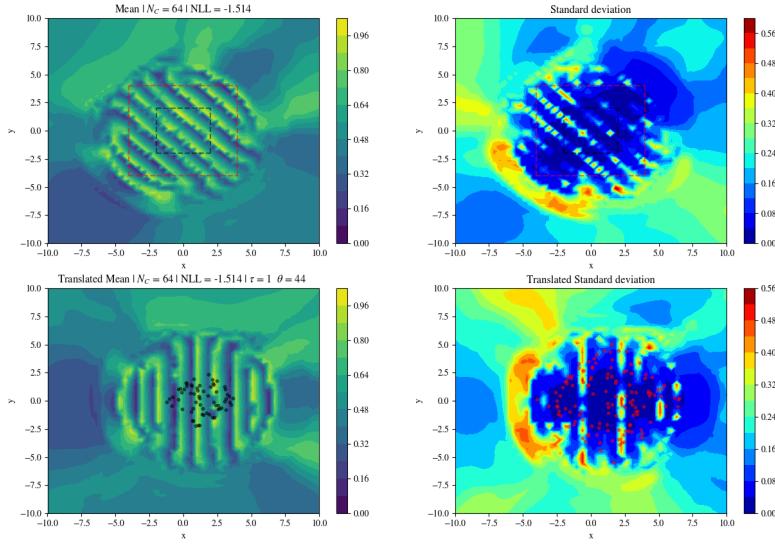


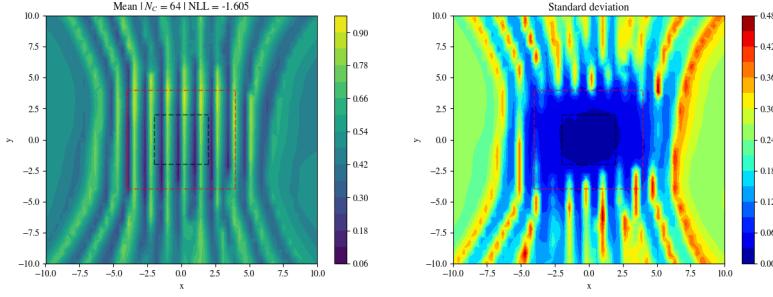
Figure 6.3.5: Samples from TETNP on the full Sawtooth dataset with Rotational The region inside the black dotted box is the region the model was trained on and the region outside the black dotted box is the region the model was not trained on.

Figure 6.3.5 demonstrates a massive improvement in the TETNP's ability to generalize to the full sawtooth dataset. Hence, we can conclude that RE is beneficial for the case when the inherent structure of the data is rotationally invariant. This illustrates a massive benefit of the Transformer architecture, as **it is very easy to introduce inductive biases to the Transformer model**, which is not the case for CNNs.

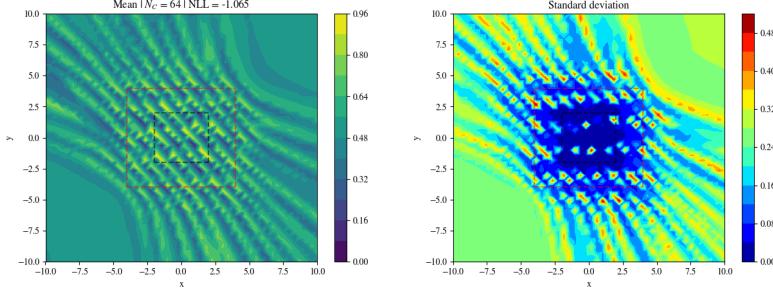
However, as we will see in the next section, if the data given to the model contains samples from many directions, the model will learn to be RE.

6.3.3 Full Sawtooth

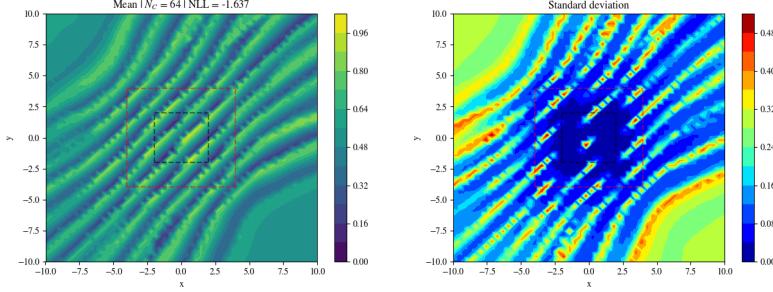
The full sawtooth dataset is the full sawtooth function which is not restricted to any direction of travel. We observe that the TETNP is able to generalize to the full sawtooth dataset without the need for rotational equivariance Figure 6.3.6.



(a)



(b)



(c)

Figure 6.3.6: Samples from TETNP on the full Sawtooth dataset. The region inside the black dotted box is the region the model was trained on and the region outside the black dotted box is the region the model was not trained on.

TODO

GET PLOTS OF CONVNP ON FULL SAWTOOTH AND COMPARE RESULTS!!!!!!!!!!!!

!
!
!
!

N_c	N_t	ConvNP Memory (MB)	TETNP Memory (MB)
10	10	24	13
100	10	29	23
1000	10	257	984
5000	10	1273	24082
<hr/>			
10	1000	268	26
100	1000	268	113
1000	1000	268	985
5000	1000	1273	24083

Table 6.3.2: Memory usage of ConvNP and TETNP on the 2D Sawtooth dataset.

6.4 Computational Complexity

TODO

Run the code for this

!
!
!
!

Chapter 7

Linear Runtime Models

7.1 Introduction

From Table 5.4.1 and Table 6.3.2, it is clear that the Transformer models are memory-intensive due to their quadratic complexity in terms of the dataset size. This is a significant bottleneck for scaling up the Transformer models to larger datasets with limited compute resources. To address this issue, several methods have been proposed to reduce the memory complexity of the Transformer models which we will explore in this chapter.

7.2 Pseudotokens

Pseudotokens (also known as Inducing Points) are a set of tokens that are used to approximate the full set of tokens, it can be thought of as a lower dimensional representation of the data set. They have been widely used in the context of Gaussian Processes (GPs) to reduce the complexity of the model with great success [Hensman, Fusi, and Lawrence 2013]. The original tokens $\mathbf{X} \in \mathbb{R}^{N \times D}$ are projected onto into a lower-dimensional space $\mathbf{I} \in \mathbb{R}^{M \times D}$ where $M \ll N$ through some translation equivariant network [Ashman et al. 2024] giving us the pseudotokens \mathbf{I} which are translation equivariant to the original tokens \mathbf{X} . The pseudotokens $\mathbf{I} \in \mathbb{R}^{M \times D}$ can then be used to perform cross-attention with the original tokens \mathbf{X} thus reducing the memory complexity to $\mathcal{O}(MN_c + MM_t)$ making the model linear with respect to context and target set size.

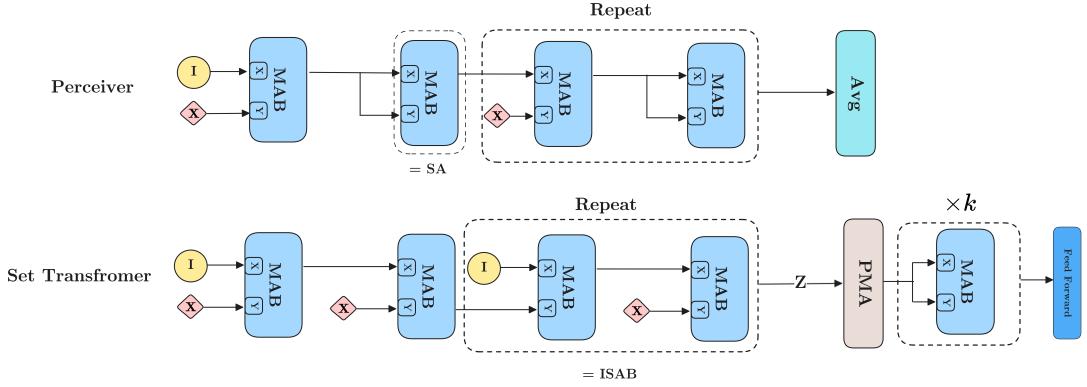


Figure 7.2.1: Perceiver vs Set Transformer. MAB is equivalent to Multi-Head Cross Attention, refer to [Ashman et al. 2024] for more details.

We consider two implementations of a pseudotokens based Transformer, one is the Set Transformer [Lee et al. 2019] and the other is the Perceiver [Jaegle et al. 2021]. Both models are very similar and only really differ in the way they implement the cross-attention mechanism between the original and pseudo tokens. [Feng et al. 2023] implemented the Perceiver model in the context of NPs creating the ‘Latent Bottled Attention Neural Process’ (LBANP) model. [Ashman et al. 2024] implemented the Set Transformer model into a NP creating the ‘Inducing Set Transformer’ (IST) model. Due to the similarity of the models, we expect both to have very similar performance on the tasks they are evaluated on.

7.3 Linear Transformer

[Katharopoulos et al. 2020] introduces a kernelized form of the self-attention mechanism that allows the model to be linear in the number of tokens by avoiding the softmax operation. The output of an attention head \mathbf{H} is computed as follows:

$$\mathbf{H}_i = \frac{\phi(\mathbf{Q}_i)^T \sum_{j=1}^N \phi(\mathbf{K}_j) \mathbf{V}_j^T}{\phi(\mathbf{Q}_i)^T \sum_{j=1}^N \phi(\mathbf{K}_j)} \quad (7.3.1)$$

Where ϕ is a function that introduces non-linearities, the authors use the ELU function [Clevert, Unterthiner, and Hochreiter 2016]. It is clear that we only compute $\sum_{j=1}^N \phi(\mathbf{K}_j)$ and $\sum_{j=1}^N \phi(\mathbf{K}_j) \mathbf{V}_j^T$ once for all the queries \mathbf{Q}_i , thus reducing the complexity to $\mathcal{O}(N)$.

This can simply replace the transformer in the original TNP model, we will refer to this model as the ‘Linear Transformer NP’ (LinearTNP).

7.4 HyperMixer

Is self-attention required for a Transformer model to be effective? The majority of parameters in Transformers are in the MLPs and not the self-attention mechanism. [Tolstikhin

[et al. 2021] proposes a Transformer model that removes the self-attention mechanism and replaces with MLPs across rows and columns of the input, effectively ‘mixing’ the feature and data points dimensions in the input to learn patterns in a way akin to self-attention. This model is called the ‘MLP-Mixer’ and is shown to perform on par with the original Transformer models on image classification tasks. However, the ‘MLP-Mixer’ requires a **known fixed input size** which breaks the flexibility of the model if we apply it to a neural process.

To overcome this limitation in flexibility, the ‘HyperMixer’ [Mai et al. 2023] model was proposed which is a variant of the MLP-Mixer that is designed to work with variable input sizes. The model uses hypernetworks [Ha, Dai, and Le 2016] to generate the weights of the MLPs by using hypernetworks (h_1, h_2) on the queries \mathbf{Q} and keys \mathbf{K} of the input (row wise).

$$\mathbf{W}_1 = h_1(\mathbf{K}) = \begin{bmatrix} \text{MLP}_1(\mathbf{k}_1) \\ \vdots \\ \text{MLP}_1(\mathbf{k}_N) \end{bmatrix} \quad \mathbf{W}_2 = h_2(\mathbf{Q}) = \begin{bmatrix} \text{MLP}_2(\mathbf{q}_1) \\ \vdots \\ \text{MLP}_2(\mathbf{q}_N) \end{bmatrix} \quad (7.4.1)$$

Where $\text{MLP}_1, \text{MLP}_2 : \mathbb{R}^d \rightarrow \mathbb{R}^p$ transforms the dimension of the input to a lower dimension p . The output of the model is computed as follows:

$$\mathbf{H} = \mathbf{W}_1 \sigma(\mathbf{W}_2^T \mathbf{V}) \quad (7.4.2)$$

where σ is the activation function, the authors use the GELU function [hendrycks2016gaussian]. Cross attention can be performed by generating the queries from one input and the keys and values from another input. We simply replace attention in the original TNP model with the HyperMixer model to create the ‘HyperMixNP’ model.

Lack of Translation Equivariance

Both LinearTNP and HyperMixNP models are not translation equivariant. To introduce TE without using pseudotokens, we require the use of the pairwise differences’ matrix Δ as described in 4.3.3, this would increase the complexity of the model to $\mathcal{O}(N^2)$. It is possible to use pseudotokens in conjunction with these model to achieve translation equivariance however this was not explored in this work.

7.5 Experimental Results

7.5.1 2D Gaussian Process

7.5.2 2D Sawtooth

7.5.3 Memory Complexity

Chapter 8

Conclusion

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