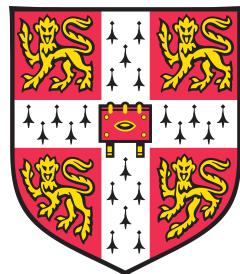


Uses of Complex Wavelets in Deep Convolutional Neural Networks



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Declaration

I hereby declare that except where specific reference is made to the work of others, the contents of this dissertation are original and have not been submitted in whole or in part for consideration for any other degree or qualification in this, or any other university. This dissertation is my own work and contains nothing which is the outcome of work done in collaboration with others, except as specified in the text and Acknowledgements. This dissertation contains fewer than 65,000 words including appendices, bibliography, footnotes, tables and equations and has fewer than 150 figures.

Fergal Cotter
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Abstract

Image understanding has long been a goal for computer vision. It has proved to be an exceptionally difficult task due to the large amounts of variability that are inherent to objects in a scene. Recent advances in supervised learning methods, particularly convolutional neural networks (CNNs), have pushed forth the frontier of what we have been able to train computers to do.

Despite their successes, the mechanics of how these networks are able to recognize objects are little understood, and the networks themselves are often very difficult and time-consuming to train. It is very important that we improve our current approaches in every way possible.

A CNN is built from connecting many learned convolutional layers in series. These convolutional layers are fairly crude in terms of signal processing - they are arbitrary taps of a finite impulse response filter, learned through stochastic gradient descent from random initial conditions. We believe that if we reformulate the problem, we may achieve many insights and benefits in training CNNs. Noting that modern CNNs are mostly viewed from and analyzed in the spatial domain, this thesis aims to view the convolutional layers in the frequency domain (viewing things in the frequency domain has proved useful in the past for denoising, filter design, compression and many other tasks). In particular, we use *complex wavelets* (rather than the Fourier transform or the discrete wavelet transform) as basis functions to reformulate image understanding with deep networks.

In this thesis, we explore the most popular and well-developed form of using complex wavelets in deep learning, the ScatterNet from Stephane Mallat. We explore its current limitations by building a DeScatterNet and found that while it has many nice properties, it may not be sensitive to the most appropriate shapes for understanding natural images.

We then develop a *locally invariant* convolutional layer, a combination of a complex wavelet transform, a modulus operation, and a learned mixing. To do this, we derive backpropagation equations and allow gradients to flow back through the (previously fixed) ScatterNet front end. Connecting several such locally invariant layers allows us to build *learnable ScatterNet*, a more flexible and general form of the ScatterNet (while still maintaining its desired properties).

We show that the learnable ScatterNet can provide significant improvements over the regular ScatterNet when being used as a front end for a learning system. Additionally, we show that the locally invariant convolutional layer can directly replace convolutional layers in a deep CNN (and not just at the front-end). The locally invariant convolutional layers

naturally downsample the input (because of the complex modulus) while increasing the channel dimension (because of the multiple wavelet orientations used). This is an operation that often happens in a CNN by a combination of a pooling and convolutional layer. It was at these locations in a CNN where the learnable ScatterNet performed best, implying it may be useful as learnable pooling layer.

Finally, we develop a system to learn complex weights that act directly on the wavelet coefficients of signals, in place of a convolutional layer. We call this layer the *wavelet gain layer* and show it can be used alongside convolutional layers. The network designer may then choose to learn in the pixel *or* wavelet domains. This layer shows a lot of promise and affords more control over what regions of the frequency space we want our layer to learn from. Our experiments show that it can improve on learning in the pixel domain for early layers of a CNN.

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Chapter 1

Introduction

It has long been the goal of computer vision researchers to be able to develop systems that can reliably recognize objects in a scene. Achieving this unlocks a huge range of applications that can benefit society as a whole: fully autonomous vehicles, automatic labelling of uploaded videos/images for searching, interpretation and screening of security video feeds, and many more, all far-reaching and extremely valuable. Many of these tasks are very tedious for humans and would be done much better by machines if the missed-detection rate can be kept low enough. The challenge does not lie in finding the right application, but in the difficulty of training a computer to see.

Some of the difficulties associated with vision are the presence of nuisance variables such as changes in lighting condition, changes in viewpoint, and background clutter. These variables do not affect the scene but can drastically change the pixel representation of it. Humans, even at early stages of their lives, have little difficulty filtering out these nuisance variables and are excellent at extracting the necessary information from a scene. To design a robust system, it makes sense to take account of how our brains see and understand scenes.

Unfortunately, biological vision is also a complex system. It has more to it than simply collecting photons in the eye. An excerpt from a recent Neurology paper [1] sums up the problem well:

It might surprise some to learn that visual information is significantly degraded as it passes from the eye to the visual cortex. Thus, of the unlimited information available from the environment, only about 10^{10} bits/sec are deposited in the retina ... only $\sim 6 \times 10^6$ bits/sec leave the retina and only 10^4 make it to layer IV of V1 [2], [3]. These data clearly leave the impression that visual cortex receives an impoverished representation of the world ... it should be noted that estimates of the bandwidth of conscious awareness itself (i.e., what we 'see') are in the range of 100 bits/sec or less[2], [3].

Current video cameras somewhat act as a combination of the first and second stage of this system, collecting photons in photosensitive sensors and then converting this to a stream of images. Standard definition digital television typically has a bit rate between 3×10^6 and 10^7 bits/sec (slightly larger but comparable to the 10^6 bits/sec travelling through the optic nerve).

If we are to build effective vision systems, it makes sense to emulate this compression of information between the optic nerve and the later stages of the visual cortex. Hubel and Wiesel revolutionized our understanding of the V1 cortex in their Nobel prize-winning work (awarded in 1981 in Physiology/Medicine) by studying cats [4], [5], macaques and spider monkeys [6]. They found that neurons in the V1 cortex fired most strongly when edges of a particular (i.e., neuron-dependent) orientation were presented to the animal, so long as the edge was inside the receptive field of this neuron. Continuing on this work, Blakemore and Cooper [7] analysed the perception of kittens that had restricted visual information presented to them. In one of their experiments, the kittens were kept in darkness and then exposed for a few hours a day to only horizontal or vertical lines. After five months, they were taken into natural environments and their reactions were monitored. The two groups of cats would only play with objects when presented in an orientation that matched the orientation of their original environment. This suggest that these early layers of perception are *learned*.

The current state of the art in image understanding systems are Convolutional Neural Networks (CNNs). These are a learned model that cascades many convolutional filters serially in layers, separated by nonlinearities. They are seemingly inspired by the visual cortex in the way that they are hierarchically connected, progressively compressing the information into a richer representation.

Figure 1.1 shows an example architecture for the famous AlexNet [8]. Inputs are resized to a manageable size, in this case, 224×224 pixels. Then multiple convolutional filters of size 11×11 are convolved over this input to give 96 output *channels* (or *activation maps*). In the figure, these are split onto two graphics cards or GPUs for memory purposes. These are then passed through a pointwise nonlinear function, or *nonlinearity*. The activations are then pooled (a form of downsampling) and convolved with more filters to give 256 new channels at the second stage. This is repeated 3 more times until the 13×13 output with 256 channels is unravelled and passed through a fully connected neural network to classify the image as one of 1000 possible classes.

CNNs have garnered lots of attention since 2012 when the previously mentioned AlexNet nearly halved the top-5 classification error rate (from 26% to 16%) in the ImageNet Large Scale Visual Recognition Competition (ILSVRC) [9]¹. In the years since then, the complexity of CNNs has grown significantly. AlexNet had only 5 convolutional layers, whereas the 2015

¹The previous state of the art classifiers had been built by combining keypoint extractors like SIFT[10] and HOG[11] with classifiers such as Support Vector Machines[12] and Fisher Vectors[13], for example [14].

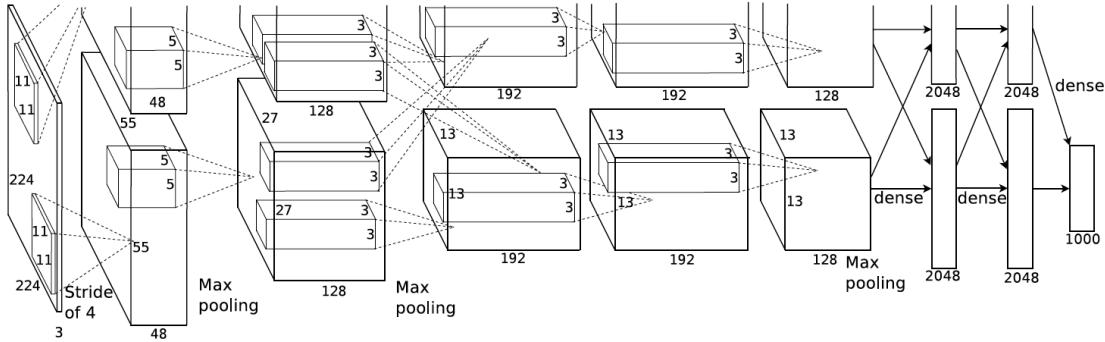


Figure 1.1: Convolutional Architecture example. The previous layer’s activations are combined with a learned convolutional filter. Note that while the activation maps are 3-D arrays, the convolution is only a 2-D operation. This means the filters have the same number of channels as the input and produce only one output channel. Multiple channels are made by convolving with multiple filters. Not shown here are the nonlinearities that happen in between convolution operations. Image taken from [8].

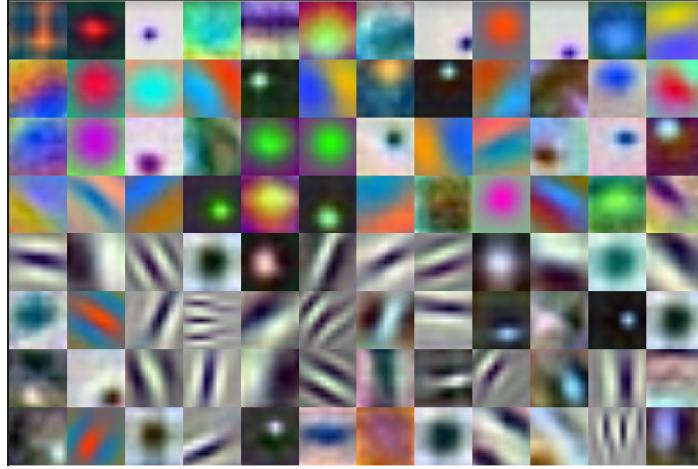
ILSVRC winner ResNet [15] achieved 3.57% top-5 error with 151 convolutional layers (and had some experiments with 1000 layer networks).

1.1 Motivation

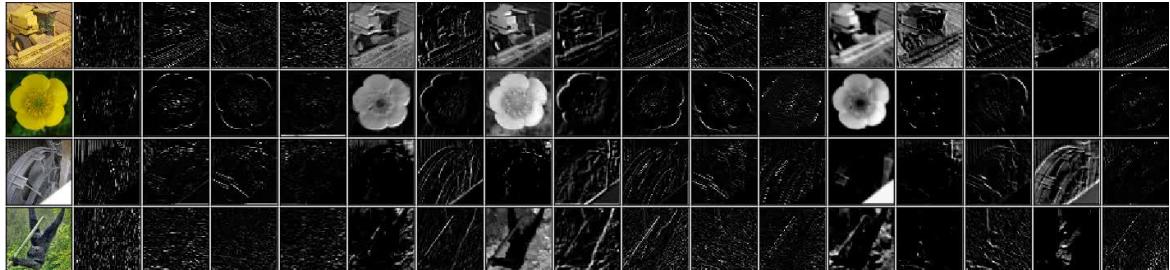
Despite their success, CNNs are often criticized for being *black-box* methods. You can view the first layer of filters quite easily (see Figure 1.2a) as they exist in RGB space, but beyond that things get trickier as the filters have a third, *channel* dimension, typically much larger than the two spatial dimensions. Additionally, it is not clear what the input channels themselves correspond to. For illustration purposes, we have also shown some example activations from the first three convolutional layers for AlexNet in Figure 1.2(b)-(d)². We can see in Figure 1.2b that in the conv1 activations, some of the first layer channels are responding to edges or colour information, but as we go deeper to conv2 and conv3, it becomes less and less clear what each activation is responding to.

This has started to become a problem, and while we are happy to trust modern CNNs for isolated tasks, we are less likely to be comfortable with them driving cars through crowded cities, or making executive decisions that affect people directly. In a commonly used contrived example, it is not hard to imagine a deep network that could be used to assess whether giving a bank loan to an applicant is a safe investment. Trusting a black box solution is deeply unsatisfactory in this situation. Not only from the customer’s perspective, who, if declined, has the right to know why [16], but also from the bank’s — before lending large

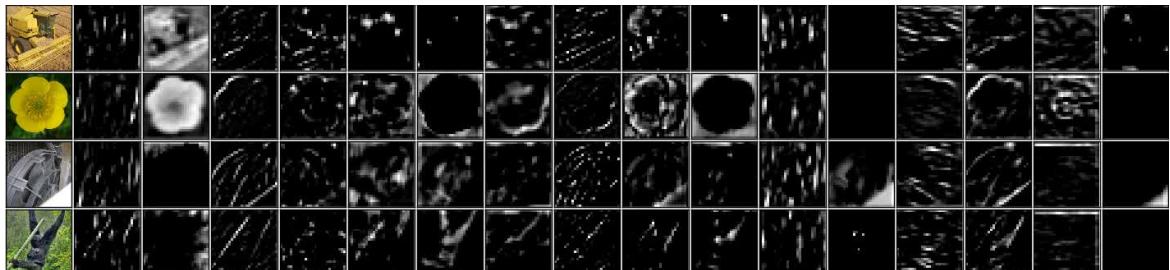
²These activations are taken after a specific nonlinearity that sets negative values to 0, hence the large black regions.



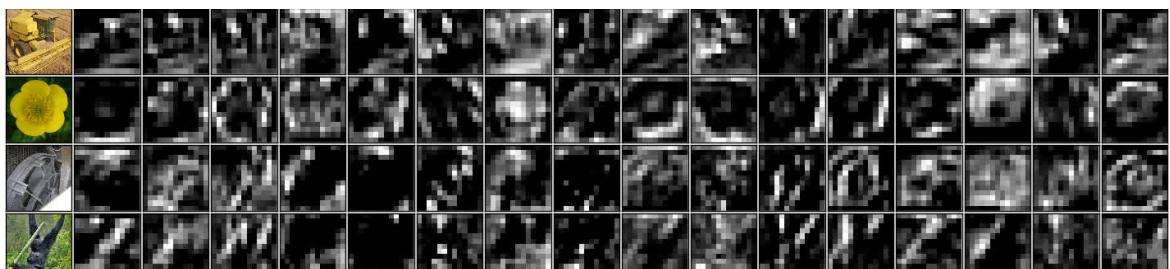
(a) conv1 filters



(b) conv1 activations



(c) conv2 activations



(d) conv3 activations

Figure 1.2: Example first layer filters and the first three layer's outputs. (a) The 11×11 filters for the first stage of AlexNet. Of the 96 filters, 48 were learned on one GPU and another 48 on another GPU. Interestingly, one GPU has learned mostly lowpass/colour filters and the other has learned oriented bandpass filters. (b) - (d) Randomly chosen activations from the output of the first, second and third convolutional layers of AlexNet (see Figure 1.1) with negative values set to 0. Filters and activation images taken from supplementary material of [8].

sums of money, most banks would like to know why the network has given the all clear. ‘It has worked well before’ is a poor rule to live by.

Aside from their lack of interpretability, it takes a long time and a lot of effort to train state-of-the-art CNNs. Typical networks that have won ILSVRC since 2012 have had roughly 100 million parameters and take up to a week to train. This is optimistic and assumes that you already know the necessary optimization or architecture hyperparameters, which you often have to find out by trial and error. In a conversation the author had with Yann LeCun, the attributed father of CNNs, at a Computer Vision Summer School (ICVSS 2016), LeCun highlighted this problem himself:

There are certain recipes (for building CNNs) that work and certain recipes that don’t, and we don’t know why.

Considering the recent success of CNNs, it is becoming more and more important to understand *how* and *what* a network learns, so we can interrogate what in the input has contributed to it making its classification or regression choice. Without this information, the use of these incredibly powerful tools could be restricted to research and proprietary applications.

1.2 Approach

The structure of convolutional layers is fairly crude in terms of signal processing - arbitrary taps of an FIR filter are learned typically via stochastic gradient descent from random starting states to minimize either a mean-squared error or cross-entropy loss.

This leads us to ask a motivating question:

Is it possible to learn convolutional filters as combinations of basis functions rather than individual filter taps?

In achieving this, it is important to find ways to have an adequate richness of filtering while reducing the number of parameters needed to specify resulting filters. We want to contract the space of learning to a subspace or manifold that is more useful. In much the same way, the convolutional layer in a CNN is a restricted version of a fully connected layer in a multi-layer perceptron, yet adding this restriction allowed us to train more powerful networks.

The intuition that we explore in this thesis is that *complex wavelets* are good (possibly optimal) basis functions for convolutional filtering in CNNs.

1.2.1 Why Complex Wavelets?

Most modern approaches to CNNs are framed entirely in the spatial domain; our choice of complex wavelets as the basis function to explore comes from the deeper intuition that it

may be helpful to rethink about CNNs in the *frequency domain*. Historically, the frequency domain has been an excellent space for solving many signal processing problems such as noise removal, filter design, edge detection and data compression. We believe it may prove to have advantages for CNNs too (beyond just an efficient space to do convolution in).

The Fourier transform, which uses complex sinusoids as its basis function, is perhaps the most ubiquitous tool to use for frequency domain analysis. The problem with these complex sinusoids is that they have *infinite* support. This means that small changes in one part of an image affect every Fourier coefficient. Additionally, they are not stable to small deformations, as small changes can produce unbounded changes in the representation [17].

The common remedy to this problem is to use the localized, and more stable, short-time Fourier transform (STFT). The STFT (or the Gabor transform) is a natural extension of the Fourier transform, windowing the complex sinusoids with a Gaussian(or similar) function. The STFT has the undesirable property that all frequencies are sampled with the same resolution. A close relative of the STFT is the continuous wavelet transform (CWT). The shorter duration of the wavelet basis functions as the frequency increases means that their time resolving power improves with centre-frequency. Another commonly used wavelet transform is the discrete wavelet transform (or the DWT) often favoured over the CWT because of its speed of computation. It can use many different finite support basis functions, all with different frequency localization properties, but it is usually limited to using real filters. As such, it suffers from many problems such as shift-dependence and lack of directionality in two dimensions (2-D). These problems can be remedied by using the slower CWT with complex basis functions, but we choose instead to use the dual-tree complex wavelet transform, or DT \mathbb{C} WT [18] with q-shift filters [19].

The DT \mathbb{C} WT allows for complex basis functions that have shift-invariance and directionality but is fast, like the DWT (in 2-D it can be thought of as the application of 4 DWTs in parallel). It is also more easily invertible than the CWT, forming a tight frame [20], which we believe may prove to be a very important property for visualizing what a CNN is responding to.

We revisit the properties of the Fourier transform, STFT, CWT, DWT and DT \mathbb{C} WT and expand on the properties behind our choice of basis functions in the literature review section 2.6.

On top of the intuition that the wavelet domain is a good space in which to frame CNNs, there are some experimental motivating factors too. Firstly, the wavelet transform has had much success in image and video compression [21]. Good compression performance implies an ability of the basis functions to represent the input data sparsely (as seems to happen in the brain). Secondly, the filters from the first layer of AlexNet (Figure 1.2) look like oriented wavelets. Given that there was no prior placed on the filters to make them have this similarity to wavelets, this result is noteworthy. And finally, the aforementioned work of Hubel and Wiesel suggests that the early layers of the visual system act like a Gabor transform.

These experimental observations imply that complex wavelets would do well in replacing the first layer of a CNN, but we should also ask if they can be used at deeper layers. Their well-understood and well-defined behaviour would help us to answer the above *how* and *why* questions. Additionally, they allow us to enforce a certain amount of smoothness and near orthogonality; smoothness seems to be important to avoid sensitivity to adversarial or spoofing attacks [22] and near orthogonality allows you to cover a large space with fewer coefficients.

But first, we must find out *if* it is possible to get the same or nearly the same performance by using wavelets as the building blocks for CNNs, and this is the core goal of this thesis.

1.3 Method

1.3.1 ScatterNets

To explore the uses of complex wavelets in CNNs, we begin by looking at one of the most popular current uses of wavelets in image recognition tasks, the Scattering Transform.

The Scattering Transform, or the *ScatterNet*, was introduced in [17], [23] at the same time as AlexNet. It is a non-black-box network that can be thought of as a restricted complex-valued CNN [24]. Unlike a CNN, it has predefined convolutional kernels, set to complex wavelet (and scaling) functions and uses the complex magnitude as its nonlinearity. Due to its well-defined structure, it can be analyzed and bounds on its stability to shifts, noise and deformations are found in [17].

For a simple task like identifying small handwritten digits, the variabilities in the data are simple and small and the ScatterNet can easily reduce the problem into a space which a Gaussian Support Vector Machine (or SVM [12]) can easily solve [23]. For a more complex task like identifying real-world objects, the ScatterNet can somewhat reduce the variabilities and get good results with an SVM, but there is a significant performance gap between this and what a CNN can achieve. For example, in [25] a second-order ScatterNet can achieve 82.3% top-1 classification accuracy on CIFAR-10, a commonly used dataset, whereas modern CNNs such as [15] can achieve 93.4%.

1.3.2 Learnable ScatterNets

To start to address the performance gap between ScatterNet front ends and CNNs we first investigate the properties of current ScatterNets. Inspired by the visualization work of Zeiler and Fergus [26] we build a DeScatterNet. The DeScatterNet leverages the perfect reconstruction properties of the DT^CWT and allows us to investigate what in the input image the ScatterNet is responding to.

The DeScatterNet shows that the ScatterNet may be limiting itself by not combining the filtering of different wavelet orientations (it does not mix the channels as a CNN does).

Inspired by the work of [27], we propose the learnable ScatterNet, which includes this mixing while keeping the desirable properties of the ScatterNet.

The learnable ScatterNet can be thought of as using the scattering outputs as the *basis functions*³ for our convolutional layers. We show that this improves greatly on the ScatterNet design, and under certain constraints can improve on the performance of CNNs too.

1.3.3 Wavelet Domain Filtering

We find that the complex modulus of the ScatterNet design to be useful for some operations in a CNN, but it has a demodulating effect on the frequency energy (all the outputs have significantly more energy in lower frequencies). This limits repeated application of it as the demodulating effect compounds.

We develop a system that does not use the complex modulus; instead, it learns *complex* gains in the wavelet domain. Rather than mixing subbands together, we keep them independent and only learn to mix across the channel dimension. This is important, as it allows us to then use the inverse DTCWT to return to the pixel domain. The shift-invariant properties of the DTCWT mean the reconstructed outputs are (mostly) free from aliasing effects, despite much of the processing being carried out at significantly reduced sample rates in the wavelet domain.

We show that our layer can be used alongside regular convolutional layers. I.e., it becomes possible to ‘step’ into the wavelet domain to do wavelet filtering for one layer, before ‘stepping’ back into the pixel domain to do pixel filtering for the next layer.

1.4 Thesis Layout and Contributions to Knowledge

This thesis has one literature review chapter and four novel-work chapters:

- Chapter 2 explores some of the background necessary for starting to develop image understanding models. In particular, it covers the inspiration for CNNs and the workings of CNNs themselves, as well as covering the basics of wavelets and ScatterNets.
- Chapter 3 proposes a change to the core of the ScatterNet. In addition to performance issues with ScatterNets, they are slow and both memory-intensive and compute-intensive to calculate. This in itself is enough of an issue to make it unlikely that they would be used as part of deep networks. To overcome this, we change the computation to use the DTCWT [18] instead of Morlet wavelets, achieving a 20 to 30 times speed-up while achieving a small improvement in classification performance.

³Although they are not true basis functions as they are the combination of a complex wavelet with a modulus nonlinearity, and are thus data-dependent.

- Chapter 4 describes our *DeScatterNet*, a tool used to interrogate the structure of ScatterNets. We also perform tests to determine the usefulness of the different scattered outputs finding that many of them are not useful for image classification.
- Chapter 5 describes the *Learnable ScatterNet* we have developed to address some of the issues found from the interrogation in chapter 4. We find that a learnable ScatterNet layer performs better than a regular ScatterNet, and can improve on the performance of a CNN if used instead of pooling layers. We also find that scattering works well not just on RGB images, but can also be useful when used after one layer of learning.
- In chapter 6, we step away from ScatterNets and present the *Wavelet Gain Layer*. The gain layer uses the wavelet space as a latent space to learn representations. We find possible nonlinearities and describe how to learn in both the pixel and wavelet domain. This work showed that there may well be benefits to learning in the wavelet domain for earlier layers of CNNs, but we have not yet found advantages to using the wavelet gain layer for deeper layers.

1.4.1 Contributions and Publications

The key contributions of this thesis are:

- Software for wavelets and DT \mathbb{C} WT based ScatterNet (described in chapter 3) and publicly available at [28].
- ScatterNet analysis and visualizations (described in chapter 4). This chapter expands on the paper we presented at MLSP2017 [29].
- Invariant Layer/Learnable ScatterNet (described in chapter 5)). This chapter expands on the paper accepted at ICIP2019 [30]. Software available at [31].
- Learning convolutions in the wavelet domain (described in chapter 6). We have published preliminary results on this work to arXiv [32] but have expanded on this paper in the chapter. Software available at [33].

1.4.2 Related Research

Readers may also be interested in the theses [34] and [35].

In [34] Singh looks at using the ScatterNet as a fixed front end and combining it with well-known machine learning methods such as SVMs, Autoencoders and Restricted Boltzmann Machines. By combining frameworks in a defined way he creates unsupervised feature extractors which can then be used with simple classifiers. Some relevant papers that makeup this thesis are [36]–[38]. In [36] Singh shows that the DT \mathbb{C} WT-ScatterNet outperforms a Morlet-ScatterNet when used as a front end for an SVM, which is similar to the work we do

in chapter 3 where we show the DTcCWT-ScatterNet outperforms a Morlet-ScatterNet when used as a front end for CNNs. He then expands on this work by testing other backends in [37], [38].

In [35] Oyallon looks at ScatterNets as front ends to deeper learning systems, such as CNNs. Some relevant papers that makeup Oyallon's thesis are [25], [39], [40]. [39], [40] are particularly relevant as he uses a ScatterNet as a feature extractor for a CNN. We do similar research in chapter 5, but allow for learned weights in the ScatterNet in our design.

Chapter 2

Background

2.1 Chapter Layout

This thesis combines work in several fields. We provide a background for the most important and relevant fields in this chapter. We first introduce the basics of deep learning, looking at general supervised learning in [section 2.2](#) before more specifically examining the structure of Neural Networks in [section 2.3](#) and CNNs in [section 2.4](#).

We then defining the properties of Wavelet Transforms, looking at the difference between the discrete WT and the complex Morlet and DTCWT in [section 2.6](#). Finally, we introduce the Scattering Transform or ScatterNet, the original inspiration for this thesis in [section 2.7](#).

2.2 Supervised Machine Learning

While this subject is general and covered in many places, we take inspiration from [41] (chapters 1, 2, 7, 8) and [42] (chapter 5-10). Consider a sample space over inputs and targets $\mathcal{X} \times \mathcal{Y}$ and a data generating distribution p_{data} . Given a dataset of input-target pairs $\mathcal{D} = \{(x^{(n)}, y^{(n)})\}_{n=1}^N$ we would like to make predictions about $p_{data}(y|x)$ that generalize well to unseen data. A common way to do this is to build a parametric model to directly estimate this conditional probability. For example, regression asserts the data are distributed according to a function of the inputs plus a noise term ϵ :

$$y = f(x, \theta) + \epsilon \tag{2.2.1}$$

This noise is often modelled as a zero-mean Gaussian random variable, $\epsilon \sim \mathcal{N}(0, \sigma^2 I)$, which means we can write:

$$p_{model}(y|x, \theta, \sigma^2) = \mathcal{N}(y; f(x, \theta), \sigma^2 I) \tag{2.2.2}$$

where (θ, σ^2) are the parameters of the model.

We can find point estimates of the parameters by maximizing the likelihood of $p_{model}(y|x, \theta)$ (or equivalently, minimizing $KL(p_{model}||p_{data})$, the KL-divergence between p_{model} and p_{data}). As the data are all assumed to be i.i.d., we can multiply individual likelihoods, and solve for θ :

$$\theta_{MLE} = \arg \max_{\theta} p_{model}(y|x, \theta) \quad (2.2.3)$$

$$= \arg \max_{\theta} \prod_{n=1}^N p_{model}(y^{(n)}|x^{(n)}, \theta) \quad (2.2.4)$$

$$= \arg \max_{\theta} \sum_{n=1}^N \log p_{model}(y^{(n)}|x^{(n)}, \theta) \quad (2.2.5)$$

Using the Gaussian regression model from above, this becomes:

$$\theta_{MLE} = \arg \max_{\theta} \sum_{n=1}^N \log p_{model}(y^{(n)}|x^{(n)}, \theta) \quad (2.2.6)$$

$$= \arg \max_{\theta} \left(-N \log \sigma - \frac{N}{2} \log(2\pi) - \sum_{n=1}^N \frac{(y^{(n)} - f(x^{(n)}, \theta))^2}{2\sigma^2} \right) \quad (2.2.7)$$

$$= \arg \min_{\theta} \frac{1}{N} \sum_{n=1}^N \frac{(y^{(n)} - f(x^{(n)}, \theta))^2}{2} \quad (2.2.8)$$

which gives us the well-known result that we would like to find parameters that minimize the mean squared error (MSE) between targets y and predictions $\hat{y} = f(x, \theta)$.

For binary classification $y \in \{0, 1\}$ and instead of the model in (2.2.2), we have:

$$p_{model}(y|x, \theta) = \text{Ber}(y; \sigma(f(x, \theta))) \quad (2.2.9)$$

where $\sigma(x)$ is the sigmoid function and Ber is the Bernoulli distribution. Note that we have used σ to refer to noise standard deviation thus far but now use $\sigma(x)$ to refer to the sigmoid and softmax functions, a confusing but common practice. $\sigma(x)$ and $\text{Ber}(y; p)$ are defined as:

$$\sigma(z) = \frac{1}{1 + e^{-z}} \quad (2.2.10)$$

$$\text{Ber}(y; p) = p^{\mathbb{I}(y=1)}(1-p)^{\mathbb{I}(y=0)} \quad (2.2.11)$$

where $\mathbb{I}(x)$ is the indicator function. The sigmoid function is useful here as it can convert a real output $f(x, \theta)$ into a probability estimate. In particular, large positive values get mapped to 1, large negative values to 0, and values near 0 get mapped to 0.5 [42, Chapter 6].

This expands naturally to multi-class classification by making y a 1-hot vector in $\{0, 1\}^C$. We must also swap the Bernoulli distribution for the Multinoulli or Categorical distribution,

and the sigmoid function for a softmax. The softmax function for vector \mathbf{z} is defined as:

$$\sigma_i(\mathbf{z}) = \frac{e^{z_i}}{\sum_{k=1}^C e^{z_k}} \quad (2.2.12)$$

which has the nice properties that $0 \leq \sigma_i \leq 1$ and $\sum_i \sigma_i = 1$. The categorical distribution uses the softmax output:

$$\text{Cat}(y; \boldsymbol{\sigma}) = \prod_{c=1}^C \sigma_c^{\mathbb{I}(y_c=1)} \quad (2.2.13)$$

If we let $\hat{y}_c = \sigma_c(f(x, \theta))$, this makes (2.2.9):

$$p_{model}(y|x, \theta) = \text{Cat}(y; \sigma(f(x, \theta))) \quad (2.2.14)$$

$$= \prod_{c=1}^C \prod_{n=1}^N \left(\hat{y}_c^{(n)} \right)^{\mathbb{I}(y_c^{(n)}=1)} \quad (2.2.15)$$

As $y_c^{(n)}$ is either 0 or 1, we remove the indicator function. Maximizing this likelihood to find the ML estimate for θ :

$$\theta_{MLE} = \arg \max_{\theta} \prod_{c=1}^C \prod_{n=1}^N \left(\hat{y}_c^{(n)} \right)^{y_c^{(n)}} \quad (2.2.16)$$

$$= \arg \max_{\theta} \frac{1}{N} \sum_{n=1}^N \sum_{c=1}^C y_c^{(n)} \log \hat{y}_c^{(n)} \quad (2.2.17)$$

which we recognize as the cross-entropy between y and \hat{y} .

2.2.1 Priors on Parameters and Regularization

Maximum likelihood estimates for parameters, while straightforward, can often lead to overfitting. A common practice is to regularize learnt parameters θ by putting a prior over them. If we do not have any prior information about what we expect them to be, it may still be useful to put an uninformative prior over them. For example, if our parameters are in the reals, a commonly used uninformative prior is a Gaussian.

Let us extend the regression example from above by saying we would like the prior on the parameters θ to be a Gaussian, i.e. $p(\theta) = \mathcal{N}(0, \tau^2 I_D)$. The corresponding maximum a

posteriori (MAP) estimate is then obtained by finding:

$$\theta_{MAP} = \arg \max_{\theta} p(\theta | \mathcal{D}, \sigma^2) \quad (2.2.18)$$

$$= \arg \max_{\theta} \frac{p(y|x, \theta, \sigma^2)p(\theta)}{p(y|x)} \quad (2.2.19)$$

$$= \arg \max_{\theta} \log p(y|x, \theta, \sigma^2) + \log p(\theta) \quad (2.2.20)$$

$$= \arg \min_{\theta} \frac{1}{N} \sum_{n=1}^N \frac{(y^{(n)} - f(x^{(n)}, \theta))^2}{2} + \frac{\lambda}{2} \|\theta\|_2^2 \quad (2.2.21)$$

where $\lambda = \sigma^2/\tau^2$ is the ratio of the observation noise to the strength of the prior [41, Chapter 7]. This is equivalent to minimizing the MSE with an ℓ_2 penalty on the parameters, also known as ridge regression or penalized least squares. λ is often called *weight decay* in the neural network literature, which we will also use in this thesis.

2.2.2 Loss Functions and Minimizing the Objective

It may be useful to rewrite (2.2.18) as an objective function on the parameters $J(\theta)$:

$$J(\theta) = \frac{1}{N} \sum_{n=1}^N \frac{(y^{(n)} - f(x^{(n)}, \theta))^2}{2} + \frac{\lambda}{2} \|\theta\|_2^2 \quad (2.2.22)$$

$$= L_{data}(y, f(x, \theta)) + L_{reg}(\theta) \quad (2.2.23)$$

where L_{data} is the data loss such as MSE or cross-entropy and L_{reg} is the regularization, such as ℓ_2 or ℓ_1 penalized loss. Now $\theta_{MAP} = \arg \min_{\theta} J(\theta)$.

Finding the global minimum of the objective function is task-dependent and is often not straightforward. One commonly used technique is called *gradient descent* (GD). This is not difficult to do as it only involves calculating the gradient at a given point and taking a small step in the direction of steepest descent. The update equation for GD is:

$$\theta_{t+1} = \theta_t - \eta \frac{\partial J}{\partial \theta} \quad (2.2.24)$$

Unsurprisingly, such a simple technique has limitations. In particular, it is sensitive to the choice of step size and has a slow convergence rate when the condition number (ratio of largest to smallest eigenvalues) of the Hessian around the optimal point is large [43]. An example of this is shown in Figure 2.1. In this figure, the step size is chosen with exact line search, i.e.

$$\eta = \arg \min_s f(x + s \frac{\partial f}{\partial x}) \quad (2.2.25)$$

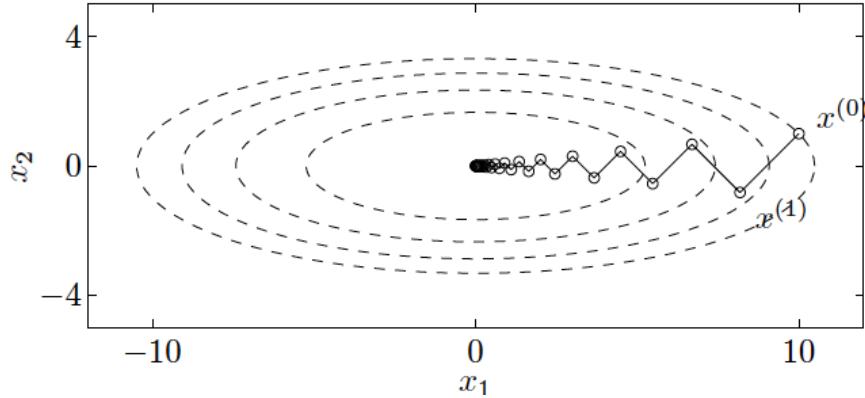


Figure 2.1: **Trajectory of gradient descent in an ellipsoidal parabola.** Some contour lines of the function $f(x) = 1/2(x_1^2 + 10x_2^2)$ and the trajectory of GD optimization using exact line search. This space has condition number 10, and shows the slow convergence of GD in spaces with largely different eigenvalues. Image taken from [43] Figure 9.2.

To truly overcome this problem, we must know the curvature of the objective function $\frac{\partial^2 J}{\partial \theta^2}$. An example optimization technique that uses the second-order information is Newton's method [43, Chapter 9]. Such techniques sadly do not scale with size, as computing the Hessian is proportional to the number of parameters squared, and many neural networks have hundreds of thousands, if not millions of parameters. In this thesis, we only consider *first-order* optimization algorithms.

2.2.3 Stochastic Gradient Descent

Aside from the problems associated the curvature of the function $J(\theta)$, another common issue faced with the gradient descent of (2.2.24) is the cost of computing $\frac{\partial J}{\partial \theta}$. In particular, the first term:

$$L_{data}(y, f(x, \theta)) = \mathbb{E}_{x, y \sim p_{data}} [L_{data}(y, f(x, \theta))] \quad (2.2.26)$$

$$= \frac{1}{N} \sum_{n=1}^N L_{data}\left(y^{(n)}, f(x^{(n)}, \theta)\right) \quad (2.2.27)$$

involves evaluating the entire dataset at the current values of θ . As the training set size grows into the thousands or millions of examples, this approach becomes prohibitively slow.

(2.2.26) writes the data loss as an expectation, hinting at the fact that we can remedy this problem by using fewer samples $N_b < N$ to evaluate L_{data} . This variation is called Stochastic Gradient Descent (SGD).

Choosing the batch size is a hyperparameter choice that we must think carefully about. Setting the value very low, e.g. $N_b = 1$ can be advantageous as the noisy estimates for the gradient have a regularizing effect on the network [44]. Increasing the batch size to larger

values allows you to easily parallelize computation as well as increasing your accuracy for the gradient, allowing you to take larger step sizes [45]. A good initial starting point is to set the batch size to 128 samples and increase/decrease from there [42].

2.2.4 Gradient Descent and Learning Rate

The step size parameter, η in (2.2.24) is commonly referred to as the learning rate. Choosing the right value for the learning rate is key. Unfortunately, the line search algorithm in (2.2.25) would be too expensive to compute for neural networks (as it would involve evaluating the function several times at different values), each of which takes about as long as calculating the gradients themselves. Additionally, as the gradients are typically estimated over a mini-batch and are hence noisy there may be little added benefit in optimizing the step sizes in the estimated direction.

Figure 2.2 illustrates the effect the learning rate can have over a contrived convex example. Optimizing over more complex loss surfaces only exacerbates the problem. Sadly, choosing the initial learning rate is ‘more of an art than a science’ [42], but [46], [47] have some tips on what to how to set it. We have found in our work that searching for a large learning rate that causes the network to diverge and reducing it from there can be a good search strategy. This agrees with Section 1.5 of [48] which states that for regions of the loss space which are roughly quadratic, $\eta_{max} = 2\eta_{opt}$ and any learning rate above $2\eta_{opt}$ causes divergence.

On top of the initial learning rate, the convergence of SGD methods require [46]:

$$\sum_{t=1}^{\infty} \eta_t \rightarrow \infty \quad (2.2.28)$$

$$\sum_{t=1}^{\infty} \eta_t^2 < \infty \quad (2.2.29)$$

Choosing how to do this also contains a good amount of artistry, and there is no one scheme that works best. A commonly used greedy method is to keep the learning rate constant until the training loss stabilizes and then to enter the next phase of training by setting $\eta_{k+1} = \gamma \eta_k$ where γ is a decay factor. Setting γ and the thresholds for triggering a step however must be chosen by monitoring the training loss curve and trial and error [46].

2.2.5 Momentum and Adam

One simple and very popular modification to SGD is to add *momentum*. Momentum accumulates past gradients with an exponential-decay moving average and continues to move in their direction. The name comes from the comparison of finding minima to rolling a ball over a surface – any new force (newly computed gradients) must overcome the current momentum of the ball. This has a smoothing effect on noisy gradients.

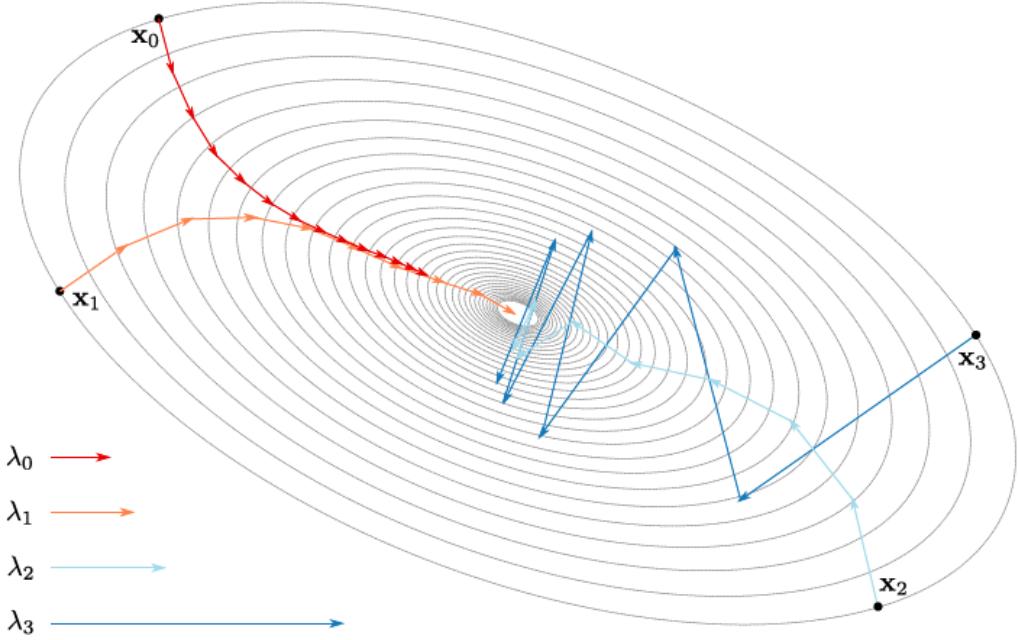


Figure 2.2: **Trajectories of SGD with different initial learning rates.** This figure illustrates the effect the step size has over the optimization process by showing the trajectory for $\eta = \lambda_i$ from equivalent starting points on a symmetric loss surface. Increasing the step size beyond λ_3 can cause the optimization procedure to diverge. Image taken from [49] Figure 2.7.

We can achieve momentum by creating a *velocity* variable v_t and modify (2.2.24) to be:

$$v_{t+1} = \alpha v_t - \eta_k \frac{\partial J}{\partial \theta} \quad (2.2.30)$$

$$\theta_{t+1} = \theta_t + v_{t+1} \quad (2.2.31)$$

where $0 \leq \alpha < 1$ is the momentum term indicating how quickly to ‘forget’ past gradients.

Another popular modification to SGD is the adaptive learning rate technique Adam [50]. There are several other adaptive schemes such as AdaGrad [51] and AdaDelta [52], but they are all quite similar, and Adam is often considered the most robust of the three [42]. The goal of all of these adaptive schemes is to take larger update steps in directions of low variance, helping to minimize the effect of large condition numbers we saw in Figure 2.1. Adam does this by keeping track of the first m_t and second v_t moments of the gradients:

$$g_{t+1} = \frac{\partial J}{\partial \theta} \quad (2.2.32)$$

$$m_{t+1} = \beta_1 m_t + (1 - \beta_1) g_{t+1} \quad (2.2.33)$$

$$v_{t+1} = \beta_2 v_t + (1 - \beta_2) g_{t+1}^2 \quad (2.2.34)$$

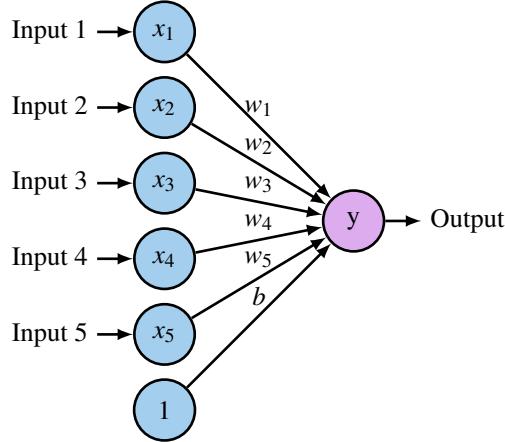


Figure 2.3: **A single neuron.** The neuron is composed of inputs x_i , weights w_i (and a bias term), as well as an activation function. Typical activation functions include the sigmoid function, tanh function and the ReLU

where $0 \leq \beta_1, \beta_2 < 1$. Note the similarity between updating the mean estimate in (2.2.33) and the velocity term in (2.2.30)¹. The parameters are then updated with:

$$\theta_{t+1} = \theta_t - \eta \frac{m_{t+1}}{\sqrt{v_{t+1}} + \epsilon} \quad (2.2.35)$$

where ϵ is a small value to avoid dividing by zero.

2.3 Neural Networks

2.3.1 The Neuron and Single Layer Neural Networks

The neuron, shown in Figure 2.3 is the core building block of Neural Networks. It takes the dot product between an input vector $\mathbf{x} \in \mathbb{R}^D$ and a weight vector \mathbf{w} , before applying a chosen nonlinearity. Historically, the sigmoid nonlinearity was the most popular but today other functions have become more popular. Still, the convention has remained to name this generic nonlinearity σ . I.e.

$$y = \sigma(\langle \mathbf{x}, \mathbf{w} \rangle) = \sigma \left(\sum_{i=0}^D x_i w_i \right) \quad (2.3.1)$$

where we have used the shorthand $b = w_0$ and $x_0 = 1$. Also, note that we will use the common practice in the neural network literature to call the parameters *weights* denoted by w .

¹The m_{t+1} and v_{t+1} terms are then bias-corrected as they are biased towards zero at the beginning of training. We do not include this for conciseness.

Some of the other popular nonlinear functions σ are the tanh and ReLU:

$$\tanh(x) = \frac{e^x - e^{-x}}{e^x + e^{-x}} \quad (2.3.2)$$

$$\text{ReLU}(x) = \max(x, 0) \quad (2.3.3)$$

See [Figure 2.4](#) for plots of these. The original Rosenblatt perceptron [53] also used the Heaviside function $H(x) = \mathbb{I}(x > 0)$.

Note that if $\langle \mathbf{w}, \mathbf{w} \rangle = 1$ then $\langle \mathbf{x}, \mathbf{w} \rangle$ is the distance from the point \mathbf{x} to the hyperplane with normal \mathbf{w} (with non unit-norm \mathbf{w} , this can be thought of as a scaled distance). Thus, the weight vector \mathbf{w} defines a hyperplane in \mathbb{R}^D which splits the space into two. The choice of nonlinearity then affects how points on each side of the plane are treated. For a sigmoid, points far below the plane get mapped to 0 and points far above the plane get mapped to 1 (with points near the plane having a value of 0.5). For tanh nonlinearities, these points get mapped to -1 and 1. For ReLU nonlinearities, every point below the plane ($\langle \mathbf{x}, \mathbf{w} \rangle < 0$) gets mapped to zero and every point above the plane keeps its inner product value.

Nearly all modern neural networks use the ReLU nonlinearity and it has been credited with being a key reason for the recent surge in deep learning success [54], [55]. In particular:

1. It is less sensitive to initial conditions as the gradients that backpropagate through it will be large even if x is large. A common observation of sigmoid and tanh nonlinearities was that their learning would be slow for quite some time until the neurons came out of saturation, and then their accuracy would increase rapidly before levelling out again at a minimum [56]. The ReLU, on the other hand, has constant gradient.
2. It promotes sparsity in outputs, by setting them to a hard 0. Studies on brain energy expenditure suggest that neurons encode information in a sparse manner. [57] estimates the percentage of neurons active at the same time to be between 1 and 4%. Sigmoid and tanh functions will typically have *all* neurons firing, while the ReLU can allow neurons to fully turn off.

2.3.2 Multilayer Perceptrons

As mentioned in the previous section, a single neuron can be thought of as a separating hyperplane with an activation that maps the two halves of the space to different values. Such a linear separator is limited and famously cannot solve the XOR problem [58]. Fortunately, adding a single hidden layer like the one shown in [Figure 2.5](#) can change this, and it is proveable that with an infinitely wide hidden layer, a neural network can approximate any function [59], [60]. This extension is called a multilayer perceptron, or MLP.

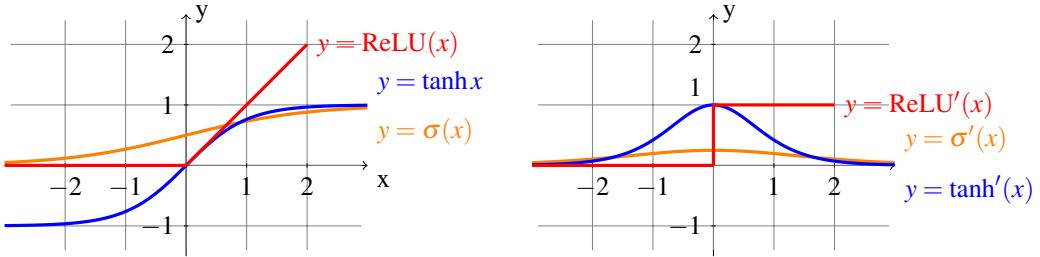


Figure 2.4: **Common Neural Network nonlinearities and their gradients.** The sigmoid, tanh and ReLU nonlinearities are commonly used activation functions for neurons. The tanh and sigmoid have the nice property of being smooth but can have saturation when the input is either largely positive or largely negative, causing little gradient to flow back through it. The ReLU does not suffer from this problem, and has the additional nice property of setting values to exactly 0, making a sparser output activation.

The forward pass of such a network with one hidden layer of H units is:

$$h_i = \sigma \left(\sum_{j=0}^D x_j w_{ij}^{(1)} \right) \quad (2.3.4)$$

$$y = \sum_{k=0}^H h_k w_k^{(2)} \quad (2.3.5)$$

where $w^{(l)}$ denotes the weights for the l -th layer, of which Figure 2.5 has 2. Note that these individual layers are often called *fully connected* as each node in the previous layer affects every node in the next.

If we were to expand this network to have N_l such fully connected layers, we could rewrite the action of each layer in a recursive fashion:

$$y^{(l+1)} = W^{(l+1)} x^{(l)} \quad (2.3.6)$$

$$x^{(l+1)} = \sigma \left(y^{(l+1)} \right) \quad (2.3.7)$$

where W is now a weight matrix, acting on the vector of previous layer's outputs $x^{(l)}$. As we are now considering every layer to be an input to the next stage, we have removed the h notation and added the superscript (l) to define the depth. $x^{(0)}$ is the network input and $y^{(N_l)}$ is the network output. Let us say that the output has C nodes, and a hidden layer $x^{(l)}$ has C_l nodes.

2.3.3 Backpropagation

It is important to truly understand backpropagation when designing neural networks, so we describe the core concepts now for a neural network with N_l layers.

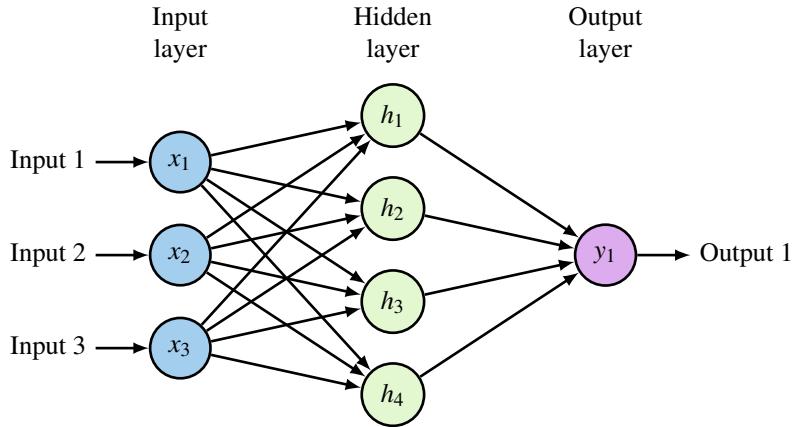


Figure 2.5: **Multi-layer perceptron.** Expanding the single neuron from Figure 2.3 to a network of neurons. The internal representation units are often referred to as the *hidden layer* as they are an intermediary between the input and output.

The delta rule, initially designed for networks with no hidden layers [61], was expanded to what we now consider *backpropagation* in [62]. While backpropagation is conceptually just the application of the chain rule, Rumelhart, Hinton, and Williams successfully updated the delta rule to networks with hidden layers, laying a key foundational step for deeper networks.

With a deep network, calculating $\frac{\partial L_{data}}{\partial w}$ may not seem easy, particularly if w is a weight in one of the earlier layers. We need to define a rule for updating the weights in all N_l layers of the network, $W^{(1)}, W^{(2)}, \dots, W^{(N_l)}$ however, only the final set $W^{(N_l)}$ are connected to the data loss function L_{data} .

2.3.3.1 Regression Loss

Let us start with writing down the derivative of L (dropping the *data* subscript for now) with respect to the network output $\hat{y} = y^{(N_l)}$ using the regression objective function (2.2.8).

$$\frac{\partial L}{\partial y^{(N_l)}} = \frac{\partial}{\partial \hat{y}} \left(\frac{1}{N} \sum_{n=1}^{N_b} \frac{1}{2} (y^{(n)} - \hat{y}^{(n)})^2 \right) \quad (2.3.8)$$

$$= \frac{1}{N} \sum_{n=1}^{N_b} (\hat{y}^{(n)} - y^{(n)}) \quad (2.3.9)$$

$$= e \in \mathbb{R} \quad (2.3.10)$$

where we have used the fact that for the regression case, $y^{(n)}, \hat{y}^{(n)} \in \mathbb{R}$.

2.3.3.2 Classification Loss

For the classification case (2.2.17), let us keep the output of the network as $y^{(N_l)} \in \mathbb{R}^C$ and define \hat{y} as the softmax applied to this vector $\hat{y}_c^{(n)} = \sigma_c(y^{(N_l, n)})$. As we now have two

superscripts, one for the sample number and one for the layer number, we combine them into a tuple of superscripts.

Note that the softmax is a vector-valued function going from $\mathbb{R}^C \rightarrow \mathbb{R}^C$ so has a Jacobian matrix S with values:

$$S_{ij} = \frac{\partial \hat{y}_i}{\partial y_j^{(N_l)}} = \begin{cases} \sigma_i(1 - \sigma_j) & \text{if } i = j \\ -\sigma_i \sigma_j & \text{if } i \neq j \end{cases} \quad (2.3.11)$$

Now, let us return to (2.2.17) and find the derivative of the objective function to this output value \hat{y} :

$$\frac{\partial L}{\partial \hat{y}} = \frac{\partial}{\partial \hat{y}} \left(\frac{1}{N} \sum_{n=1}^{N_b} \sum_{c=1}^C y_c^{(n)} \log \hat{y}_c^{(n)} \right) \quad (2.3.12)$$

$$= \frac{1}{N} \sum_{n=1}^{N_b} \sum_{c=1}^C \frac{y_c^{(n)}}{\hat{y}_c^{(n)}} \quad (2.3.13)$$

$$= d \in \mathbb{R}^C \quad (2.3.14)$$

Note that unlike (2.3.10), this derivative is vector-valued. To find $\frac{\partial L}{\partial y_j^{(N_l)}}$ we use the chain rule. It is easier to find the partial derivative with respect to one node in the output first, and then expand from here. I.e.:

$$\frac{\partial L}{\partial y_j^{(N_l)}} = \sum_{i=1}^C \frac{\partial L}{\partial \hat{y}_i} \frac{\partial \hat{y}_i}{\partial y_j^{(N_l)}} \quad (2.3.15)$$

$$= S_j^T d \quad (2.3.16)$$

where S_j is the j th column of the Jacobian matrix S . It becomes clear now that to get the entire vector derivative for all nodes in $y^{(N_l)}$, we must multiply the transpose of the Jacobian matrix with the error term from (2.3.14):

$$\frac{\partial L}{\partial y^{(N_l)}} = S^T d \quad (2.3.17)$$

2.3.3.3 Final Layer Weight Gradient

Let us continue by assuming $\frac{\partial L}{\partial y^{(N_l)}}$ is vector-valued as was the case with classification. For regression, it is easy to set $C = 1$ in the following to get the necessary results. For clarity we drop the layer superscript in the intermediate calculations.

We call the gradient for the final layer weights the *update* gradient. It can be computed by the chain rule again:

$$\frac{\partial J}{\partial W_{ij}} = \frac{\partial L_{data}}{\partial y_i} \frac{\partial y_i}{\partial W_{ij}} + \lambda W_{ij} \quad (2.3.18)$$

$$= \frac{\partial L_{data}}{\partial y_i} x_j + \lambda W_{ij} \quad (2.3.19)$$

where the second term in the above two equations comes from the regularization loss that is added to the objective. The update gradient of the entire weight matrix is then:

$$\frac{\partial J}{\partial W^{(N_l)}} = \frac{\partial L_{data}}{\partial \hat{y}} x^T + 2\lambda W \quad (2.3.20)$$

$$= S^T d(x^{(N_l-1)})^T + 2\lambda W^{(N_l)} \in \mathbb{R}^{C \times C_{N_l-1}} \quad (2.3.21)$$

2.3.3.4 Final Layer Passthrough Gradient

We also want to find the *passthrough* gradients of the final layer $\frac{\partial L}{\partial x^{(N_l-1)}}$ (these are not affected by the regularization terms, so we only need to find the gradient w.r.t. the data loss L). In a similar fashion, we first find the gradient with respect to individual elements in $x^{(N_l-1)}$ before generalizing to the entire vector:

$$\frac{\partial L}{\partial x_i} = \sum_{j=1}^C \frac{\partial L}{\partial y_j} \frac{\partial y_j}{\partial x_i} \quad (2.3.22)$$

$$= \sum_{j=1}^C \frac{\partial L}{\partial y_j} W_{j,i} \quad (2.3.23)$$

$$= W_i^T \frac{\partial L}{\partial y} \quad (2.3.24)$$

$$(2.3.25)$$

where W_i is the i th column of W . Thus

$$\frac{\partial L}{\partial x^{(N_l-1)}} = (W^{(N_l)})^T \frac{\partial L}{\partial y^{(N_l)}} \quad (2.3.26)$$

$$= (W^{(N_l)})^T S^T d \quad (2.3.27)$$

This passthrough gradient then can be used to update the next layer's weights by repeating [subsubsection 2.3.3.3](#) and [subsubsection 2.3.3.4](#).

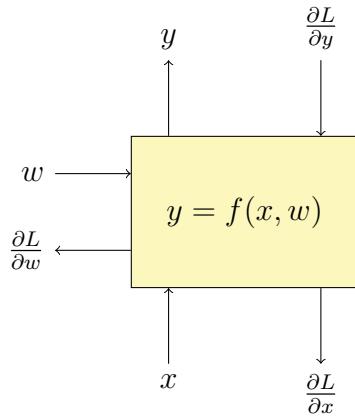


Figure 2.6: **General block form for autograd.** All neural network functions need to be able to calculate the forward pass $y = f(x, w)$ as well as the update and passthrough gradients $\frac{\partial L}{\partial w}, \frac{\partial L}{\partial x}$. Backpropagation is then easily done by allowing data to flow backwards through these blocks from the loss.

2.3.3.5 General Layer Update

The easiest way to handle this flow of gradients, and the basis for most automatic differentiation packages is the block definition shown in Figure 2.6. For all neural network components (even if they do not have weights), the operation must not only be able to calculate the forward pass $y = f(x, w)$ given weights w and inputs x , but also calculate the *update* and *passthrough* gradients $\frac{\partial L}{\partial w}, \frac{\partial L}{\partial x}$ given an input gradient $\frac{\partial L}{\partial y}$. The input gradient will have the same shape as y as will the update and passthrough gradients match the shape of w and x . This way, gradients for the entire network can be computed in an iterative fashion starting at the loss function and moving backwards.

2.4 Convolutional Neural Networks

Convolutional Neural Networks (CNNs) are a special type of Neural Network built mainly from *convolutional layers* rather than fully connected layers. A convolutional layer is one where the weights are shared spatially across the layer. In this way, a neuron at is only affected by nodes from the previous layer in a given neighbourhood, rather than from every node.

First popularized in 1998 by LeCun et. al. in [63], the convolutional layer was introduced to build invariance with respect to translations, as well as reduce the parameter size of early neural networks for pattern recognition. The idea of having a locally receptive field had already been shown to be a naturally occurring phenomenon by Hubel and Wiesel [5]. They did not become popular immediately, and another spatially based keypoint extractor, SIFT [10], was the mainstay of detection systems until the AlexNet CNN [8] won the 2012

ImageNet challenge [9] by a large margin over the next competitors, many of who used SIFT. This CNN had 5 convolutional layers followed by 3 fully connected layers.

We now briefly describe the convolutional layer, as well as many other layers used in CNNs that have become popular in the past few years.

2.4.1 Convolutional Layers

In the analysis of neural networks so far, we have considered column vectors $x^{(l)}, y^{(l)} \in \mathbb{R}^{C_l}$. Convolutional layers for image analysis have a different format, with the spatial component of the input is preserved.

Let us first consider the definition of 2-D convolution for single-channel images:

$$y[\mathbf{n}] = (x * h)[\mathbf{n}] = \sum_{\mathbf{k}} x[\mathbf{k}]h[\mathbf{n} - \mathbf{k}] \quad (2.4.1)$$

$$= \sum_{k_1, k_2} x[k_1, k_2]h[n_1 - k_1, n_2 - k_2] \quad (2.4.2)$$

where the sum is done over the support of h . For an input $x \in \mathbb{R}^{H \times W}$ and filter $h \in \mathbb{R}^{K_H \times K_W}$ the output has spatial support $y \in \mathbb{R}^{H+K_H-1 \times W+K_W-1}$.

The filter h can also be thought of as a *matched filter* that gives its largest output when the input contains the mirror of h , \tilde{h} . If the input has shapes similar to \tilde{h} in many locations, each of these locations in y will also have large outputs.

If we stack red, green and blue input channels on top of each other², we have a 3-dimensional input $x \in \mathbb{R}^{C \times H \times W}$ with $C = 3$. This third dimension is often called the *depth* dimension, to distinguish it from the two spatial dimensions. In a CNN layer each filter h is 3 dimensional with depth exactly equal to C . The convolution is done only over the two spatial dimensions and the C outputs are summed at each pixel location. This makes (2.4.1):

$$y[\mathbf{n}] = \sum_{c=0}^{C-1} \sum_{\mathbf{k}} x[c, \mathbf{k}]h[c, \mathbf{n} - \mathbf{k}] \quad (2.4.3)$$

It is not enough to only have a single matched filter and often we would like to have a bank of them, each one sensitive to a different shape. For example, if one filter is sensitive to horizontal edges, we may also want to detect vertical, and diagonal edges. Let us rename the number of channels in the input layer as C_l and specify that we would like to have C_{l+1} different matched filters. We then stack each of the single channel outputs from (2.4.3) to

²In deep learning literature, there is not a consensus about whether to stack the outputs with the channel first ($\mathbb{R}^{C \times H \times W}$) or last ($\mathbb{R}^{H \times W \times C}$). The latter is more common in Image Processing for colour and spectral images but the former is the standard for many deep learning frameworks, including the one we use – PyTorch [64]. For this reason, we stack channels along the first dimension of our tensors.

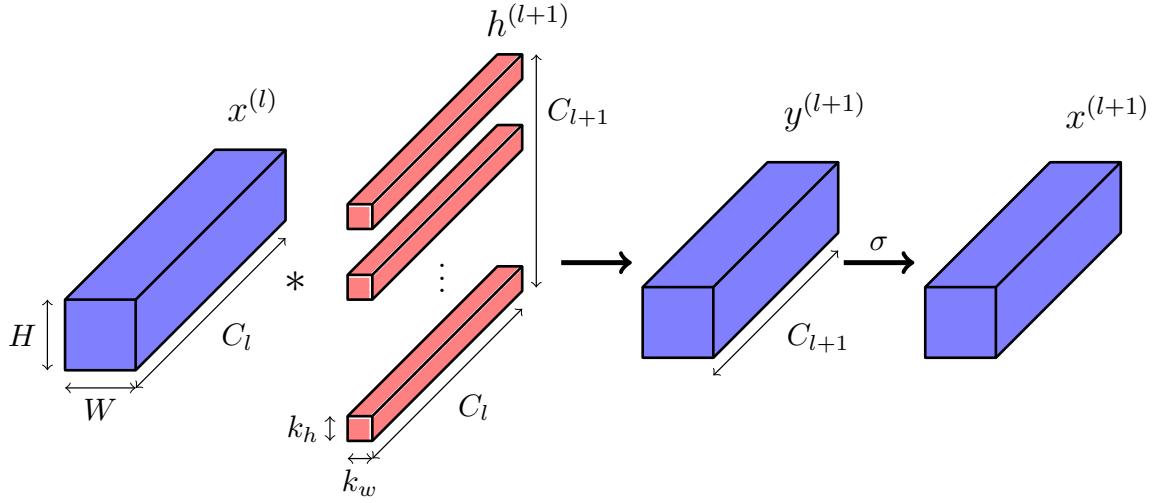


Figure 2.7: **A convolutional layer.** A convolutional layer followed by a nonlinearity σ . The previous layer’s activations are convolved with a bank of C_{l+1} filters, each of which has spatial size $k_h \times k_w$ and depth C_l . Note that there is no convolution across the channel dimension. Each filter produces one output channel in $y^{(l+1)}$.

give the output $y \in \mathbb{R}^{C_{l+1} \times H \times W}$:

$$y[f, \mathbf{n}] = \sum_{c=0}^{C-1} \sum_{\mathbf{k}} x[c, \mathbf{k}] h[f, c, \mathbf{n} - \mathbf{k}] \quad (2.4.4)$$

After a convolutional layer, we can then apply a pointwise nonlinearity to each output location in y . Like multilayer perceptrons, this was typically the sigmoid function σ , but is now more commonly the ReLU. Revisiting (2.3.6) and (2.3.7), we can rewrite this for a convolutional layer at depth l with C_l input and C_{l+1} output channels:

$$y^{(l+1)}[f, \mathbf{n}] = \sum_{c=0}^{C_l-1} x^{(l)}[c, \mathbf{n}] * h^{(l)}[f, c, \mathbf{n}] \quad (2.4.5)$$

$$x^{(l+1)}[f, \mathbf{n}] = \sigma(y^{(l+1)}[f, \mathbf{n}]) \quad (2.4.6)$$

where $f \in \{0, 1, \dots, C_{l+1}-1\}$ indexes the filter number/output channel. A diagram representing this operation is shown in Figure 2.7.

2.4.1.1 Padding and Stride

Regular 2-D convolution expands the input from size $H \times W$ to $(H + K_H - 1) \times (W + K_W - 1)$. In neural networks, this is called *full convolution*. It is often desirable (and common) to have the same output size as input size, which can be achieved by taking the central $H \times W$ outputs of full convolution. This is often called *same-size convolution*. Another option

commonly used is to only evaluate the kernels where they fully overlap the input signal, causing a reduction in the output size to $(H - K_H + 1) \times (W - K_W + 1)$. This is called *valid convolution* and was used in the original LeNet-5 [63].

Signal extension for full and same-size convolution is by default *zero padding*, and most deep learning frameworks have no ability to choose other padding schemes as part of their convolution functions. Other padding such as *symmetric padding* can be achieved by expanding the input signal before doing a valid convolution.

Stride is a commonly used term in deep learning literature. A stride of 2 means that we evaluate the filter kernel at every other input location. In signal processing, this is simply called decimation by 2.

2.4.1.2 Gradients

To get the update and passthrough gradients for the convolutional layer we will need to expand (2.4.5) (again we will drop the layer superscripts for clarity):

$$y[f, n_1, n_2] = \sum_{c=0}^{C-1} \sum_{k_1} \sum_{k_2} x[c, k_1, k_2] h[f, c, n_1 - k_1, n_2 - k_2] \quad (2.4.7)$$

The derivative for a single output $y[f, n_1, n_2]$ w.r.t. a single input $x[c, k_1, k_2]$ is then simply:

$$\frac{\partial y_{f,n_1,n_2}}{\partial x_{c,k_1,k_2}} = h[f, c, n_1 - k_1, n_2 - k_2] \quad (2.4.8)$$

It is clear from (2.4.7) that a single activation $x[c, k_1, k_2]$ affects many output values. Thus, the derivative for the loss function to this single point in x is the sum of the chain rule applied to all output positions:

$$\frac{\partial L}{\partial x_{c,k_1,k_2}} = \sum_f \sum_{n_1} \sum_{n_2} \frac{\partial L}{\partial y_{f,n_1,n_2}} \frac{\partial y_{f,n_1,n_2}}{\partial x_{c,k_1,k_2}} \quad (2.4.9)$$

$$= \sum_f \sum_{n_1} \sum_{n_2} \frac{\partial L}{\partial y_{f,n_1,n_2}} h[f, c, n_1 - k_1, n_2 - k_2] \quad (2.4.10)$$

Now we let $\Delta y[f, n_1, n_2] = \frac{\partial L}{\partial y_{f,n_1,n_2}}$ be the passthrough gradient signal from the next layer, and $\tilde{h}[\alpha, \beta, \gamma, \delta] = h[\beta, \alpha, -\gamma, -\delta]$ be a set of filters that have been mirror-imaged in the spatial domain and had their filter and channel dimensions transposed. Combining these two and

substituting into (2.4.10), we get the passthrough gradient for the convolutional layer:

$$\frac{\partial L}{\partial x_{c,k_1,k_2}} = \sum_f \sum_{n_1} \sum_{n_2} \Delta y[f, n_1, n_2] \tilde{h}[c, f, k_1 - n_1, k_2 - n_2] \quad (2.4.11)$$

$$= \sum_f \Delta y[f, \mathbf{k}] * \tilde{h}[c, f, \mathbf{k}] \quad (2.4.12)$$

which is the same as (2.4.5). I.e. we can backpropagate the gradients through a convolutional layer by mirror-imaging the filters spatially, transposing them in the channel and filter dimensions, and doing a forward convolutional layer with \tilde{h} applied to Δy . Similarly, we find the update gradients to be:

$$\frac{\partial L}{\partial h_{f,c,k_1,k_2}} = \sum_{n_1} \sum_{n_2} \frac{\partial L}{\partial y_{f,n_1,n_2}} \frac{\partial y_{f,n_1,n_2}}{\partial h_{f,c,k_1,k_2}} \quad (2.4.13)$$

$$= \sum_{n_1} \sum_{n_2} \Delta y[f, n_1, n_2] x[c, n_1 - k_1, n_2 - k_2] \quad (2.4.14)$$

$$= (\Delta y[f] * x[c]) [k_1, k_2] \quad (2.4.15)$$

where $*$ is the cross-correlation operation.

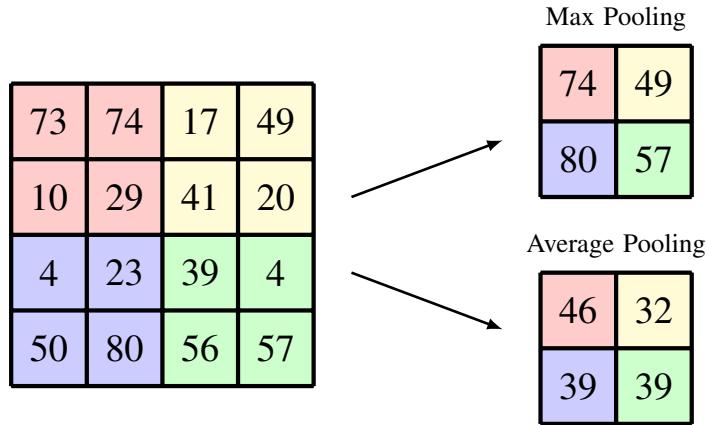
2.4.2 Pooling

Pooling layers are common in CNNs where we want to reduce the spatial size. As we go deeper into a CNN, it is common for the spatial size of the activation to decrease, and the channel dimension to increase. The C_l values at a given spatial location can then be thought of as a feature vector describing the presence of shapes in a given area in the input image.

Pooling is useful to add some invariance to smaller shifts when downsampling. It is often done over small spatial sizes, such as 2×2 or 3×3 . Invariance to larger shifts can be built up with multiple pooling (and convolutional) layers.

Two of the most common pooling techniques are *max pooling* and *average pooling*. Max pooling takes the largest value in its spatial area, whereas average pooling takes the mean. A visual explanation is shown in Figure 2.8. Note that pooling is typically a spatial operation, and only in rare cases is done over the channel/depth dimension.

A review of pooling methods in [65] found both max and average pooling to perform similarly well. While max pooling was the most popular in earlier state of the art networks [8], [66], there has been a recent trend towards using average pooling [67] or even to do away with pooling altogether in favour of strided convolutions (this idea was originally proposed in [68] and used notably in [15], [69], [70]).

Figure 2.8: Max vs Average 2×2 pooling.

2.4.3 Dropout

Dropout is a particularly strong regularization scheme that randomly turns off, or ‘zeros out’, neurons in a neural network [71], [72]. Each neuron has probability p of having its value set to 0 (independently of other neurons) during training time, forcing the network to be more general and preventing ‘co-adaption’ of neurons [72].

During test time, dropout is typically turned off, but can still be used to get an estimate on the uncertainty of the network by averaging over several runs [73].

2.4.4 Batch Normalization

Batch normalization proposed in [74] is a conceptually simple technique which rescales activations of a neural network. Despite its simplicity it has become very popular and has been found to be very useful to train deeper CNNs.

First let us define $\mu_c^{(l)}$ and $\sigma_c^{(l)}$ as the mean and standard deviations for a channel in a given activation at layer l . This mean and standard deviation is taken by averaging across the entire dataset (with N samples), and at every spatial location.

$$\mu_c^{(l)} = \frac{1}{N} \sum_{i=0}^{N-1} \sum_{\mathbf{n}} x^{(l,i)}[c, \mathbf{n}] \quad (2.4.16)$$

$$(\sigma_c^{(l)})^2 = \frac{1}{N} \sum_{i=0}^{N-1} \sum_{\mathbf{n}} \left(x^{(l,i)}[c, \mathbf{n}] \right)^2 - (\mu_c^{(l)})^2 \quad (2.4.17)$$

where $\mu, \sigma \in \mathbb{R}^C$. Batch normalization removes the mean and variance of the data, scales the data by a learnable gain γ and shifts the data to a learnable mean β , with $\gamma, \beta \in \mathbb{R}^C$. If we

drop the layer superscripts, this means the action of batch normalization is defined by:

$$y[c, \mathbf{n}] = \frac{x[c, \mathbf{n}] - \mu_c}{\sigma_c + \epsilon} \gamma_c + \beta_c \quad (2.4.18)$$

where ϵ is a small value to avoid dividing by 0.

Of course, during training, we do not have access to the dataset statistics μ, σ so these values are estimated from the batch statistics. A typical practice is to keep an exponential moving average estimate of these values.

The passthrough and update gradients are:

$$\frac{\partial L}{\partial x_{c,n_1,n_2}} = \frac{\partial L}{\partial y_{c,n_1,n_2}} \frac{\gamma}{\sigma + \epsilon} \quad (2.4.19)$$

$$\frac{\partial L}{\partial \beta_c} = \sum_{\mathbf{n}} \frac{\partial L}{\partial y_{c,\mathbf{n}}} \quad (2.4.20)$$

$$\frac{\partial L}{\partial \gamma_c} = \sum_{\mathbf{n}} \frac{\partial L}{\partial y_{c,\mathbf{n}}} \frac{x_{c,\mathbf{n}} - \mu_c}{\gamma_c + \epsilon} \quad (2.4.21)$$

Batch normalization layers are typically placed *between* convolutional layers and nonlinearities.

Consider a linear operation such as convolution with weights W acting on the previous layer's output X , defined by $Y = WX$. Batch normalization removes the sensitivity of our network to initial scaling of the weights, as $BN(aWX) = BN(WX)$. It is also particularly useful for backpropagation as an increase in weight scale does not change the passthrough gradients and leads to *smaller* update gradients [74], making the network more resilient to the problems of vanishing and exploding gradients:

$$\begin{aligned} \frac{\partial BN((aW)X)}{\partial X} &= \frac{\partial BN(WX)}{\partial X} \\ \frac{\partial BN((aW)X)}{\partial (aW)} &= \frac{1}{a} \cdot \frac{\partial BN(WX)}{\partial W} \end{aligned} \quad (2.4.22)$$

2.5 Relevant Architectures and Datasets

In this section we briefly review some relevant CNN architectures that will be helpful to refer back to in this thesis.

2.5.1 Datasets

When doing image analysis tasks it is important to know comparatively how well different networks perform on the same challenge. To achieve this, the community has developed several datasets that are commonly used to report metrics. For image classification there are five such datasets, listed here in increasing order of difficulty:

1. **MNIST**: 10 classes, 6000 images per class, 28×28 pixels per image. The images contain the digits 0–9 in greyscale on a blank background. The digits have been size normalized and centred. Dataset description and files can be obtained at [75].
2. **CIFAR-10**: 10 classes, 5000 images per class, 32×32 pixels per image. The images contain classes of everyday objects like cars, dogs, planes etc. The images are colour and have little clutter or background. Dataset description can be found in [76] and files at [77].
3. **CIFAR-100**: 100 classes, 500 images per class, 32×32 pixels per image. Similar to CIFAR-10, but now with fewer images per class and ten times as many classes. Dataset description can be found in [76] and files at [77].
4. **Tiny ImageNet**: 200 classes, 500 images per class, 64×64 pixels per image. A more recently introduced dataset that bridges the gap between CIFAR and ImageNet. Images are larger than CIFAR and there are more categories. Dataset description and files can be obtained at [78].
5. **ImageNet CLS**: There are multiple types of challenges in ImageNet, but CLS is the classification challenge and is most commonly reported in papers. It has 1000 classes of objects with a varying amount of images per class. Most classes have 1300 examples in the training set, but a few have less than 1000. The images have variable size, typically a couple of hundred pixels wide and a couple of hundred pixels high. The images can have varying amounts of clutter and can be at different scales, making it a particularly difficult challenge. Dataset description is in [9] and the most reliable source of the data can be found at [79].

Several other classification datasets do exist but are not commonly used, such as PASCAL VOC [80] and Caltech-101 and Caltech-256 [81]³.

2.5.2 LeNet

LeNet-5 [63] is a good network to start with: it is simple yet contains many of the layers used in modern CNNs. Shown in Figure 2.9 it has two convolutional and three fully connected layers. The outputs of the convolutional layers are passed through a sigmoid nonlinearity and downsampled with average pooling. The first two fully-connected layers also have sigmoid nonlinearities. The loss function used is a combination of tanh functions and MSE loss.

³Tiny ImageNet is also not commonly used as it is quite new. We have included it the main list as we have found it to be quite a useful step up from CIFAR without requiring the weeks to train experimental configurations on ImageNet.

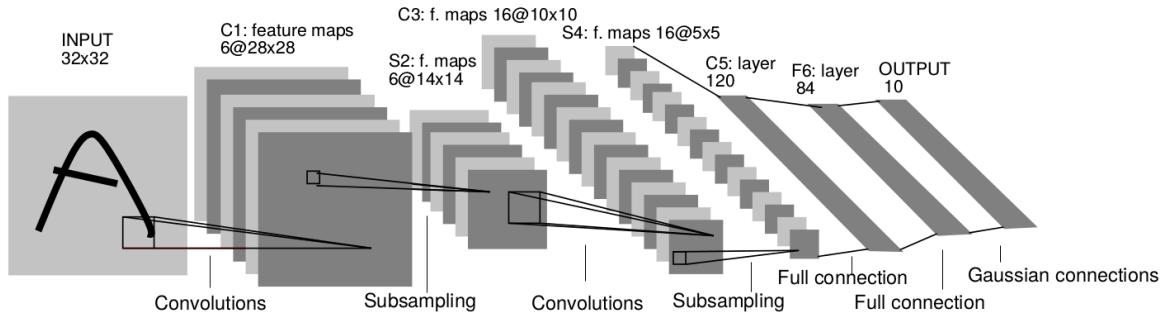


Figure 2.9: **LeNet-5 architecture.** The ‘original’ CNN architecture used for handwriting recognition. LeNet has 2 convolutional and 3 fully connected layers making 5 parameterized layers. After the second convolutional layer, the $16 \times 5 \times 5$ pixel output is unravelled to a 400 long vector. Image taken from [63].

2.5.3 AlexNet

We have already seen AlexNet [8] in chapter 1. It is arguably one of the most important architectures in the development in CNNs as it was able to experimentally prove that CNNs can be used for complex tasks. This required a few innovations: they used multiple GPUs to do fast processing on large images, they used the ReLU to avoid saturation, and also added dropout to aid generalization. Training of AlexNet on 2 GPUs available in 2012 takes roughly a week.

The first layer uses convolutions with spatial support of 11×11 , followed by 5×5 and 3×3 for the final three layers.

2.5.4 VGGnet

The Visual Geometry Group (VGG) at Oxford came second in the ILSVRC challenge in 2014 with their VGG-nets [66], but remains an important network for some of the design choices it inspired. Their optimal network was much deeper than AlexNet, with 19 convolutional layers on top of each other before 3 fully connected layers. These convolutional layers all used the smaller 3×3 seen only at the back of AlexNet.

This network is particularly attractive due its simplicity, compared to the more complex Inception Network [82] which won the 2014 ILSVRC challenge. VGG-16, the 16 layer variant of VGG stacks two or three convolutional layers (and ReLUs) on top of each other before reducing spatial size with max pooling. After processing at five scales, the resulting $512 \times 14 \times 14$ activation is unravelled and passed through a fully connected layer.

These VGG networks also marked the start of a trend that has since become common, where channel depth is doubled after pooling layers. The doubling of channels and quartering the spatial size still causes a net reduction in the number of activations.

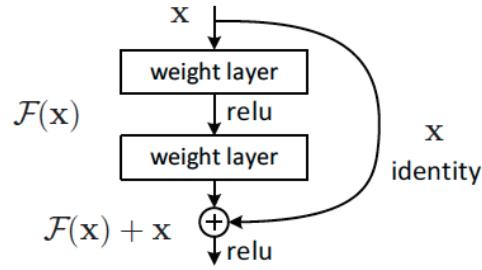


Figure 2.10: **The residual unit from ResNet.** A residual unit. The identity mapping is always present, and the network learns the difference from the identity mapping, $\mathcal{F}(x)$. Taken from [15].

2.5.5 The All Convolutional Network

The All Convolutional Network [68] introduced two popular modifications to the VGG networks:

- They argued for the removal of max pooling layers, saying that a 3×3 convolutional layer with stride 2 works just as well.
- They removed the fully connected layers at the end of the network, replacing them with 1×1 convolutions. Note that a 1×1 convolution still has shared weights across all spatial locations. The output layer then has size $C_L \times H \times W$, where H, W are many times smaller than the input image size, and the vector of C_L coefficients at each spatial location can be interpreted as a vector of scores marking the presence/absence of C_L different shapes. For classification, the output can be averaged over all spatial locations, whereas for localization it may be useful to keep this spatial information.

The new network was able to achieve state of the art results on CIFAR-10 and CIFAR-100 and competitive performance on ImageNet, while only use a fraction of the parameters of other networks.

2.5.6 Residual Networks

Residual Networks or ResNets won the 2015 ILSVRC challenge, introducing the residual layer. Most state of the art models today use this residual mapping in some way [69], [70].

The inspiration for the residual layer came from the difficulties experienced in training deeper networks. Often, adding an extra layer would *decrease* network performance. This is counter-intuitive as the deeper layers could simply learn the identity mapping, and achieve the same performance.

To promote the chance of learning the identity mapping, they define a residual unit, shown in Figure 2.10. If a desired mapping is denoted $\mathcal{H}(x)$, instead of trying to learn this,



Figure 2.11: **Importance of phase over magnitude for images.** The phase of the Fourier transform of the first image is combined with the magnitude of the Fourier transform of the second image and reconstructed. Note that the first image has entirely won out and nothing is left visible of the cameraman.

they instead learn $\mathcal{F}(x) = \mathcal{H}(x) - x$. Doing this promotes a strong diagonal in the Jacobian matrix which improves conditioning for gradient descent.

Recent analysis of a ResNet without nonlinearities [83], [84] proves that SGD fails to converge for deep networks when the network mapping is far away from the identity, suggesting that a residual mapping is a good thing to do.

2.6 The Fourier and Wavelet Transforms

Computer vision is an extremely difficult task. Pixel intensities in an image are typically not very informative in understanding what is in that image. These values are sensitive to lighting conditions and camera configurations. It would be easy to take two photos of the same scene and get two vectors x_1 and x_2 that have a very large Euclidean distance, but to a human, would represent the same objects. What is most important in defining an image is difficult to define, however, some things are notably more important than others. For example, the location or phase of the waves that make up an image is much more important than the magnitude of these waves, something that is not necessarily true for audio processing. A simple experiment to demonstrate this is shown in Figure 2.11.

2.6.1 The Fourier Transform

For a signal $f(t) \in L_2(\mathbb{R})$ (square summable signals), the *Fourier transform* is defined as:

$$F(\omega) = \int_{-\infty}^{\infty} f(t)e^{-j\omega t} dt \quad (2.6.1)$$

This can be extended to two dimensions for signals $f(\mathbf{u}) \in L_2(\mathbb{R}^2)$:

$$F(\boldsymbol{\omega}) = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} f(\mathbf{u}) e^{-j\boldsymbol{\omega}^t \mathbf{u}} d\mathbf{u} = \langle f(\mathbf{u}), e^{j\boldsymbol{\omega}^t \mathbf{u}} \rangle \quad (2.6.2)$$

The Fourier transform is an invaluable signal expansion, as viewing a signal in the frequency space offers many insights, as well as affording many very useful properties (most notably the efficiency of convolution as a product of Fourier transforms). While it is a mainstay in signal processing, it can be a poor feature descriptor due to the infinite support of its basis functions - the complex sinusoids $e^{j\boldsymbol{\omega}^t \mathbf{u}}$. If a single pixel changes in the input it can change all of the Fourier coefficients. As natural images are generally non-stationary, we need to be able to isolate frequency components in local regions of an image, and not have this property of global dependence. To achieve a more local Fourier transform we can use the short time (or short space) Fourier Transform (STFT) or the continuous wavelet transform (CWT). The two are very similar and mainly differ in the way they handle the concept of ‘scale’. We will only discuss the CWT in this review, but for an excellent comparison of the two, we recommend [85, Chapter 1].

2.6.2 The Continuous Wavelet Transform

The *continuous wavelet transform*, like the Fourier Transform, can be used to decompose a signal into its frequency components. Unlike the Fourier transform, these frequency components can be localized in space. To achieve this, we need a bandpass filter, or *mother wavelet* ψ^4 such that:

$$\int_{-\infty}^{\infty} \psi(\mathbf{u}) d\mathbf{u} = \Psi(0) = 0 \quad (2.6.3)$$

Any function that has sufficient decay of energy with frequency and satisfies (2.6.3), is said to satisfy the *admissibility condition*.

As we are working in 2-D for image processing, consider rotations, dilations, and shifts of this function by $\theta \in [0, 2\pi]$, $a > 0$, $\mathbf{b} \in \mathbb{R}^2$ respectively, where

$$\text{Rotation: } R_\theta x(\mathbf{u}) = x(r_{-\theta}\mathbf{u}) \quad (2.6.4)$$

$$\text{Dilation: } D_a x(\mathbf{u}) = \frac{1}{a} x\left(\frac{\mathbf{u}}{a}\right) \quad (2.6.5)$$

$$\text{Translation: } T_{\mathbf{b}} x(\mathbf{u}) = x(\mathbf{u} - \mathbf{b}) \quad (2.6.6)$$

where r_θ is the 2-D rotation matrix. Now consider shifts, scales and rotations of our bandpass filter

$$\psi_{\mathbf{b}, a, \theta}(\mathbf{u}) = \frac{1}{a} \psi\left(\frac{r_{-\theta}(\mathbf{u} - \mathbf{b})}{a}\right) \quad (2.6.7)$$

⁴We use upright ψ, ϕ to distinguish 1-D wavelets from their 2-D counterparts ψ, ϕ

which are called the *daughter wavelets*. The 2D CWT of a signal $x(\mathbf{u})$ is defined as

$$CWT_x(\mathbf{b}, a, \theta) = \int_{-\infty}^{\infty} \psi_{\mathbf{b}, a, \theta}^*(\mathbf{u}) x(\mathbf{u}) d\mathbf{u} = \langle \psi_{\mathbf{b}, a, \theta}(\mathbf{u}), x(\mathbf{u}) \rangle \quad (2.6.8)$$

2.6.2.1 Properties

The CWT has some particularly nice properties, such as *covariance* under the three transformations (2.6.6)-(2.6.4):

$$R_{\theta_0}x \rightarrow CWT_x(r_{-\theta_0}\mathbf{b}, a, \theta + \theta_0) \quad (2.6.9)$$

$$D_{a_0}x \rightarrow CWT_x(\mathbf{b}/a_0, a/a_0, \theta) \quad (2.6.10)$$

$$T_{\mathbf{b}_0}x \rightarrow CWT_x(\mathbf{b} - \mathbf{b}_0, a, \theta) \quad (2.6.11)$$

Most importantly, the CWT is now localized in space, which distinguishes it from the Fourier transform. This means that changes in one part of the image will not affect the wavelet coefficients in another part of the image, so long as the distance between the two parts is much larger than the support region of the wavelets you are examining.

2.6.2.2 Inverse

The CWT can be inverted by using a *dual* function $\tilde{\psi}$. There are restrictions on what dual function we can use, namely the dual-wavelet pair must have an admissible constant C_ψ that satisfies the cross-admissibility constraint [86]. Assuming these constraints are satisfied, we can recover x from CWT_x .

2.6.2.3 Interpretation

As the CWT is a convolution with a zero mean function, the wavelet coefficients are only large in the regions of the parameter space (\mathbf{b}, a, θ) where $\psi_{\mathbf{b}, a, \theta}$ ‘match’ the features of the signal. As the wavelet ψ is well localized, the energy of the coefficients CWT_x will be concentrated on the significant parts of the signal.

For an excellent description of the properties of the CWT in 1-D we recommend [87] and in 2-D we recommend [85].

2.6.3 Discretization and Frames

The CWT is highly redundant. We have taken a 2-D signal and expressed it in 4 dimensions (2 offset, 1 scale and 1 rotation). In reality, we would like to sample the space of the CWT in an efficient manner. We would ideally like to fully retain all information in x (be able to reconstruct x from the samples) while sampling over (\mathbf{b}, a, θ) as little as possible to avoid redundancy. To understand how to do this we must briefly talk about frames.

A set of vectors $\phi = \{\varphi_i\}_{i \in I}$ in a hilbert space \mathbb{H} is a *frame* if there exist two constants $0 < A \leq B < \infty$ such that for all $x \in \mathbb{H}$:

$$A\|x\|^2 \leq \sum_{i \in I} |\langle x, \varphi_i \rangle|^2 \leq B\|x\|^2 \quad (2.6.12)$$

with A, B called the *frame bounds* [20]. The frame bounds relate to the issue of stable reconstruction. In particular, no vector x with $\|x\| > 0$ should be mapped to 0, as this would violate the bound on A from below. This can be interpreted as ensuring our set ϕ covers the entire frequency space. The upper bound ensures that the transform coefficients are bounded.

Any finite set of vectors that spans the space is a frame. An orthonormal basis is a commonly known non-redundant frame where $A = B = 1$ and $|\varphi_i| = 1$ (e.g. the Discrete Wavelet Transform or the Fourier Transform). Tight frames are frames where $A = B$ and Parseval tight frames have the special case $A = B = 1$. It is possible to have frames that have more vectors than dimensions, and this will be the case with many expansions we explore in this thesis.

If $A = B$ and $|\varphi_i| = 1$, then A is the measure of the redundancy of the frame. Of course, for the orthogonal basis, $A = 1$ when $|\varphi_i| = 1$ so there is no redundancy. For the 2-D DTCTWT which we will see shortly, the redundancy is 4.

2.6.3.1 Inversion and Tightness

(2.6.12) specify the constraints that make a frame representation invertible. The tighter the frame bounds, the more easily it is to invert the signal. This gives us some guide to choosing the sampling grid for the CWT.

One particular inverse operator is the *canonical dual frame*. If we define the frame operator $S = \Phi\Phi^*$ then the canonical dual of Φ is defined as $\tilde{\Phi} = \{\tilde{\varphi}_i\}_{i \in I}$ where:

$$\tilde{\varphi}_i = S^{-1}\varphi_i \quad (2.6.13)$$

then from [20] we have:

$$x = \sum_{i \in I} \langle x, \varphi_i \rangle \tilde{\varphi}_i = \sum_{i \in I} \langle x, \tilde{\varphi}_i \rangle \varphi_i \quad (2.6.14)$$

If a frame is tight, then so is its dual.

2.6.4 Discrete Wavelet Transform

(2.6.7) gave the equation for the daughter wavelets in 2-D, in 1-D at scales $a = 2^j, j \in \mathbb{Z}$, and translations $b = 2^j l, l \in \mathbb{Z}$, this is simply:

$$\psi_{j,l}(t) = \frac{1}{\sqrt{a}} \psi \left(\frac{t-b}{a} \right) = 2^{-j/2} \psi(2^{-j}t - l) \quad (2.6.15)$$

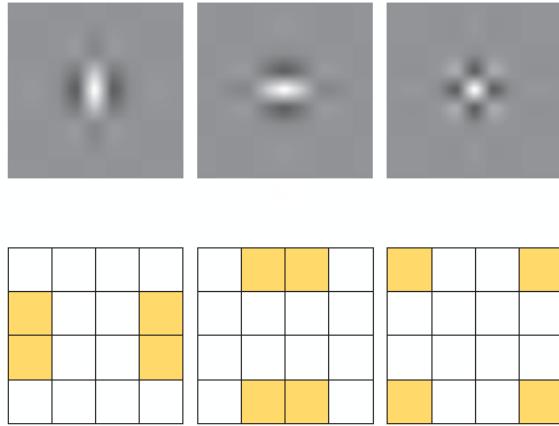


Figure 2.12: **Typical wavelets from the 2D separable DWT.** Top: Wavelet point spread functions for ψ^v (low-high), ψ^h (high-low), and ψ^d (high-high) wavelets. High-high wavelets are in a checkerboard pattern, with no favoured orientation. Bottom: Idealized support of the spectra of each of the wavelets. Image taken from [18].

where we have chosen to redefine the dilation parameter a in terms of a scale factor j . As j increases, we move to *coarser* scales.

The 2-D DWT has one scaling function and three wavelet functions, composed of the product of 1-D wavelets in the horizontal (u_1) and vertical (u_2) directions:

$$\phi(\mathbf{u}) = \phi(u_1)\phi(u_2) \quad (2.6.16)$$

$$\psi^h(\mathbf{u}) = \phi(u_1)\psi(u_2) \quad (2.6.17)$$

$$\psi^v(\mathbf{u}) = \psi(u_1)\phi(u_2) \quad (2.6.18)$$

$$\psi^d(\mathbf{u}) = \psi(u_1)\psi(u_2) \quad (2.6.19)$$

with h, v, d indicating the sensitivity to horizontal, vertical and diagonal edges. The point spread functions for the wavelet functions are shown in Figure 2.12.

For the four equations above (2.6.16) – (2.6.19), define the daughter wavelets as:

$$\phi_{lm}^j(\mathbf{u}) = \phi_{j,l}(u_1)\phi_{j,m}(u_2) \quad (2.6.20)$$

$$\psi_{lm}^{h,j}(\mathbf{u}) = \phi_{j,l}(u_1)\psi_{j,m}(u_2) \quad (2.6.21)$$

$$\psi_{lm}^{v,j}(\mathbf{u}) = \psi_{j,l}(u_1)\phi_{j,m}(u_2) \quad (2.6.22)$$

$$\psi_{lm}^{d,j}(\mathbf{u}) = \psi_{j,l}(u_1)\psi_{j,m}(u_2) \quad (2.6.23)$$

for $l, m \in \mathbb{Z}$ where l, m define horizontal and vertical translation. We can then get an orthonormal basis with the set $\{\phi_{lm}^j, \psi_{lm}^{h,j}, \psi_{lm}^{v,j}, \psi_{lm}^{d,j}\}_{j,l,m}$. The wavelet coefficients at a chosen

scale and location can then be found by taking the inner product of the signal x with the daughter wavelets.

2.6.4.1 Shortcomings

The Discrete Wavelet Transform (DWT) is an orthogonal basis. It is a natural first signal expansion to consider when frustrated with the limitations of the Fourier transform. It is also a good example of the limitations of non-redundant transforms, as it suffers from several drawbacks:

- The DWT is sensitive to the zero crossings of its wavelets. We would like singularities in the input to yield large wavelet coefficients, but this may not always be the case. See [Figure 2.13](#).
- They have poor directional selectivity. As the wavelets are purely real, they have passbands in all four quadrants of the frequency plane. While they can pick out edges aligned with the frequency axis, they are not specific to other orientations. See [Figure 2.12](#).
- They are not shift-invariant - small shifts greatly perturb the wavelet coefficients. [Figure 2.13](#) shows this for the centre-left and centre-right images.

The lack of shift-invariance and the possibility of low outputs at singularities is a price to pay for the critically sampled property of the transform. This shortcoming can be overcome with the undecimated DWT [88], [89], but it comes with a heavy computational and memory cost.

2.6.5 Complex Wavelets

Fortunately, we can improve on the DWT with complex wavelets, as they can solve these new shortcomings while maintaining the desired localization properties the Fourier transform lacked.

The Fourier transform does not suffer from a lack of directional selectivity and shift-variance because its basis functions are derived from complex sinusoids⁵:

$$e^{j\omega t} = \cos(\omega t) + j \sin(\omega t) \quad (2.6.24)$$

whereas the DWT's basis functions are based on only the real sinusoid, $\cos(\omega t)$. As t moves along the real line, the phase of the Fourier coefficients change linearly, while their magnitude remains constant. In contrast, as t moves along the real line, the sign of the real coefficient flips between -1 and 1, and its magnitude is a rectified sinusoid.

⁵We have temporarily switched to 1D notation here as it is clearer and easier to use, but the results still hold for 2D.

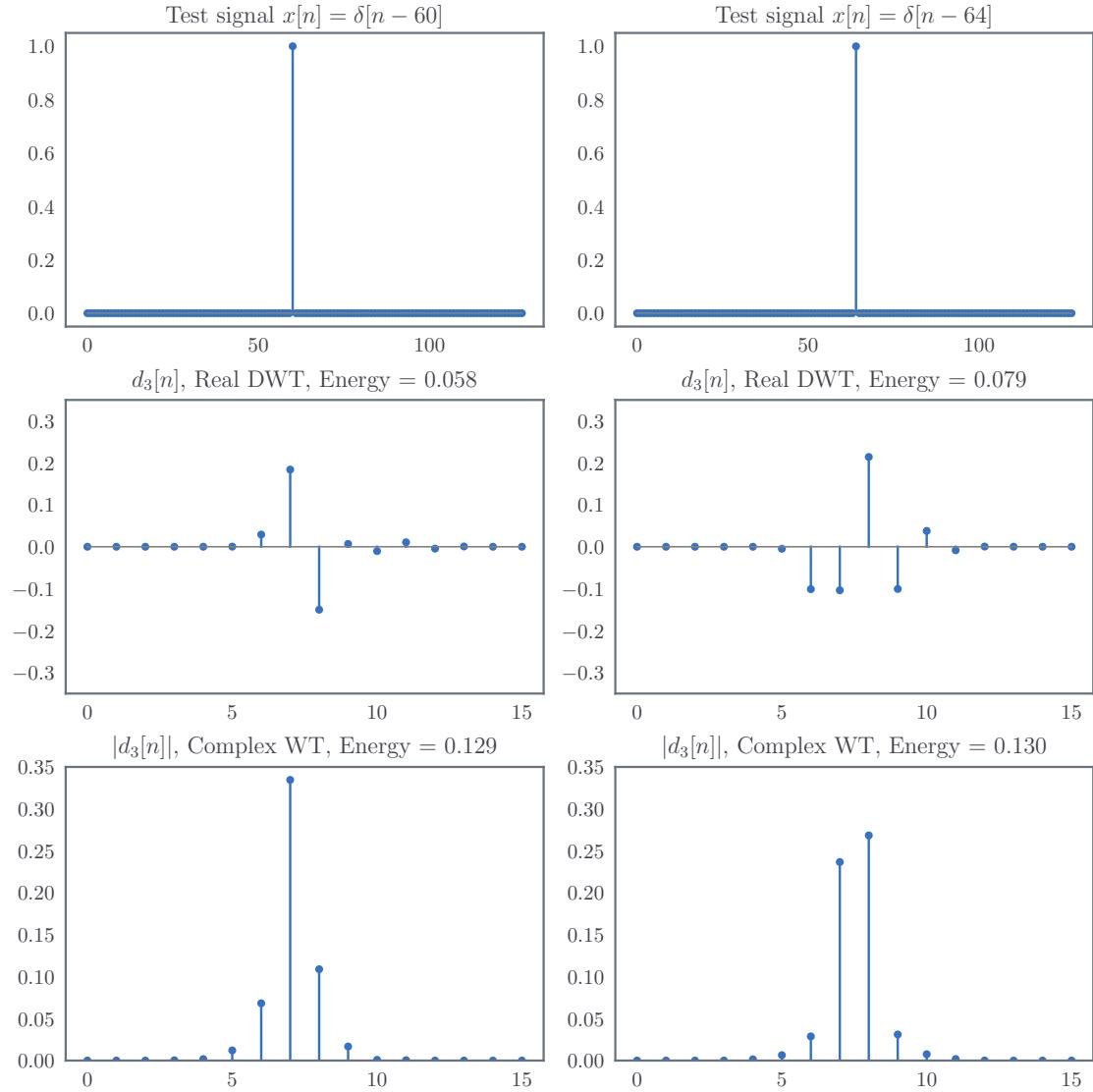


Figure 2.13: **Sensitivity of DWT coefficients to zero crossings and small shifts.** Two impulse signals $\delta(n - 60)$ and $\delta(n - 64)$ are shown (top), as well as the wavelet coefficients for scale $j = 3$ for the DWT (middle) and for the DTCWT (bottom). In the middle row, not only are the coefficients very different from a shifted input, but the energy has almost doubled. As the DWT is an orthonormal transform, this means that this extra energy has come from other scales. In comparison, the energy of the magnitude of the DTCWT coefficients has remained far more constant, as has the shape of the envelope of the output. Image adapted from [18].

The nice properties of the complex sinusoids come from the fact that the cosine and sine functions of the Fourier transform form a Hilbert pair and together constitute an analytic signal.

We can achieve these nice properties if the mother wavelet for our wavelet transform is analytic:

$$\psi_c(t) = \psi_r(t) + j\psi_i(t) \quad (2.6.25)$$

where $\psi_r(t)$ and $\psi_i(t)$ form a Hilbert pair (i.e., they are 90° out of phase with each other).

There are a number of possible ways to do a wavelet transform with complex wavelets. We examine two in particular: a Fourier-based, sampled CWT using Morlet wavelets, and the Dual-Tree Complex Wavelet Transform (DTCWT) developed by Kingsbury [18], [19], [90]–[95].

We look at the Morlet wavelet transform because it is used by Mallat et. al. in their scattering transform [23], [25], [96]–[102]. We believe the DTCWT several advantages over the Morlet based implementation, and has been the basis for most of our work.

Let us write the wavelet transform of an input x as

$$\mathcal{W}x = \{x * \phi_J, x * \psi_\lambda\}_\lambda \quad (2.6.26)$$

where $\lambda = (j, k)$ with $j \in \{1, 2, \dots, J\}$ indexing the scale⁶ and $k \in \{0, 1, \dots, K - 1\}$ indexing the orientations of the chosen wavelet transform, whether it be the DTCWT or Morlet transform.

2.6.6 Sampled Morlet Wavelets

The wavelet transform used by Mallat et. al. in their scattering transform is an efficient implementation of the Gabor Transform. While the Gabor wavelets have the best theoretical trade-off between spatial and frequency localization, they have a (usually small) non-zero mean. This violates (2.6.3) making them inadmissible as wavelets. Instead, the Morlet wavelet has the same shape, but with an extra degree of freedom chosen to set $\int \psi(\mathbf{u}) d\mathbf{u} = 0$. This wavelet has equation (in 2D):

$$\psi(\mathbf{u}) = \frac{1}{2\pi\sigma^2} (e^{i\mathbf{u}^t \xi} - \beta) e^{-\frac{|\mathbf{u}|^2}{2\sigma^2}} \quad (2.6.27)$$

where β is this extra degree of freedom, and usually $\beta \ll 1$. σ is the size of the gaussian window and ξ is the approximate location of the peak frequency response in the Fourier plane, with $-\pi \leq \xi_1, \xi_2 \leq \pi$.

Bruna and Mallat [23] add a further additional degree of freedom in their original design by allowing for a non-circular Gaussian window over the complex sinusoid, which gives control

⁶We try to number everything from zero in this thesis, but number j from 1 as is the practice in [18], [19].



Figure 2.14: **Single Morlet filter with varying slants and window sizes.** Top left — 45° plane wave (real part only). Top right — plane wave with $\sigma = 3, \gamma = 1$. Bottom left — plane wave with $\sigma = 3, \gamma = 0.5$. Bottom right — plane wave with $\sigma = 2, \gamma = 0.5$.

over the angular resolution of the final wavelet. This makes (2.6.27):

$$\psi(\mathbf{u}) = \frac{\gamma}{2\pi\sigma^2} (e^{i\mathbf{u}^t \xi} - \beta) e^{-\mathbf{u}^t \Sigma^{-1} \mathbf{u}} \quad (2.6.28)$$

Where

$$\Sigma^{-1} = \begin{bmatrix} \frac{1}{2\sigma^2} & 0 \\ 0 & \frac{\gamma^2}{2\sigma^2} \end{bmatrix}$$

The effects of modifying the eccentricity parameter γ and the window size σ are shown in Figure 2.14. A full family of Morlet wavelets at varying scales and orientations is shown in Figure 2.15.

2.6.6.1 Tightness and Invertibility

Recall our definition of the wavelet transform \mathcal{W} from (2.6.26). Assuming the transform is bounded, we can always scale it so that it satisfies Plancherel's equality

$$\|\mathcal{W}x\| = \|x\| \quad (2.6.29)$$

which is a nice property to have for invertibility, as well as for analysing how different signals get transformed (e.g. white noise versus standard images). Scaling the transform changes the upper bound B in (2.6.12) to 1 and makes the lower bound $A = 1 - \alpha$, where α is a measure of how non-tight a frame is.

Using the capital notation to denote the Fourier transform, define the function $A(\omega)$ to be the coverage each wavelet family has over the frequency plane:

$$A(\omega) = |\Phi_J(\omega)|^2 + \sum_{\lambda} |\Psi_{\lambda}(\omega)|^2 \quad (2.6.30)$$

For a unit norm input $\|x\|^2 = 1$ and scaled wavelets, we can now change (2.6.12) to be:

$$1 - \alpha \leq A(\omega) \leq 1 \quad (2.6.31)$$

If $A(\omega)$ is ever close to 0, then there is not a good coverage of the frequency plane at that location. Figure 2.15 shows the frequency coverage of a few sample grids over the CWT parameters used by Mallat. Invertibility is possible, but not guaranteed for all configurations.

The tightness of the frame is determined by the sampling grid of our wavelets parameters (\mathbf{b}, a, θ) . Common choices for sampling grids for 2-D wavelets are [85, Section 2.2]:

- For dilations, $a = 2^{j/Q}$ for $j \in \mathbb{Z}$ controlling the scale and Q , the number of scales per octave.
- For rotations, subdivide the interval $[0, \pi)$ into K sections, and choose $\theta_k = \frac{k\pi}{K}$, $k = \{0, 1, \dots, K-1\}$.
- For the translations, set the offsets $\mathbf{b} = (l2^{j/Q}, m2^{j/Q})$, $l, m \in \mathbb{Z}$.

2.6.7 The DT \mathbb{C} WT

The DT \mathbb{C} WT was first proposed by Kingsbury in [91], [92] as a way to combat many of the shortcomings of the DWT such as its poor directional selectivity and its poor shift-invariance. A thorough analysis of the properties and benefits of the DT \mathbb{C} WT is done in [18], [93]. Building on these properties, it has been used successfully for denoising and inverse problems [103]–[106], texture classification [107], [108], image registration [109], [110] and SIFT-style keypoint generation matching [111]–[115] amongst many other applications. Compared to Gabor (or Morlet) image analysis, the authors of [18] sum up the dangers as:

A typical Gabor image analysis is either expensive to compute, is noninvertible, or both.

This nicely summarises the difference between this method and the Fourier based method outlined in subsection 2.6.6. The DT \mathbb{C} WT is a filter bank (FB) based wavelet transform. It is faster to implement than the Morlet analysis, as well as being more readily invertible.

2.6.7.1 Design Criteria for the DT \mathbb{C} WT

As in subsection 2.6.5, we want to have a complex mother wavelet $\psi_c = \psi_r + j\psi_i$ and complex scaling function $\phi_c = \phi_r + j\phi_i$, but now achieved with filter banks. The complex component

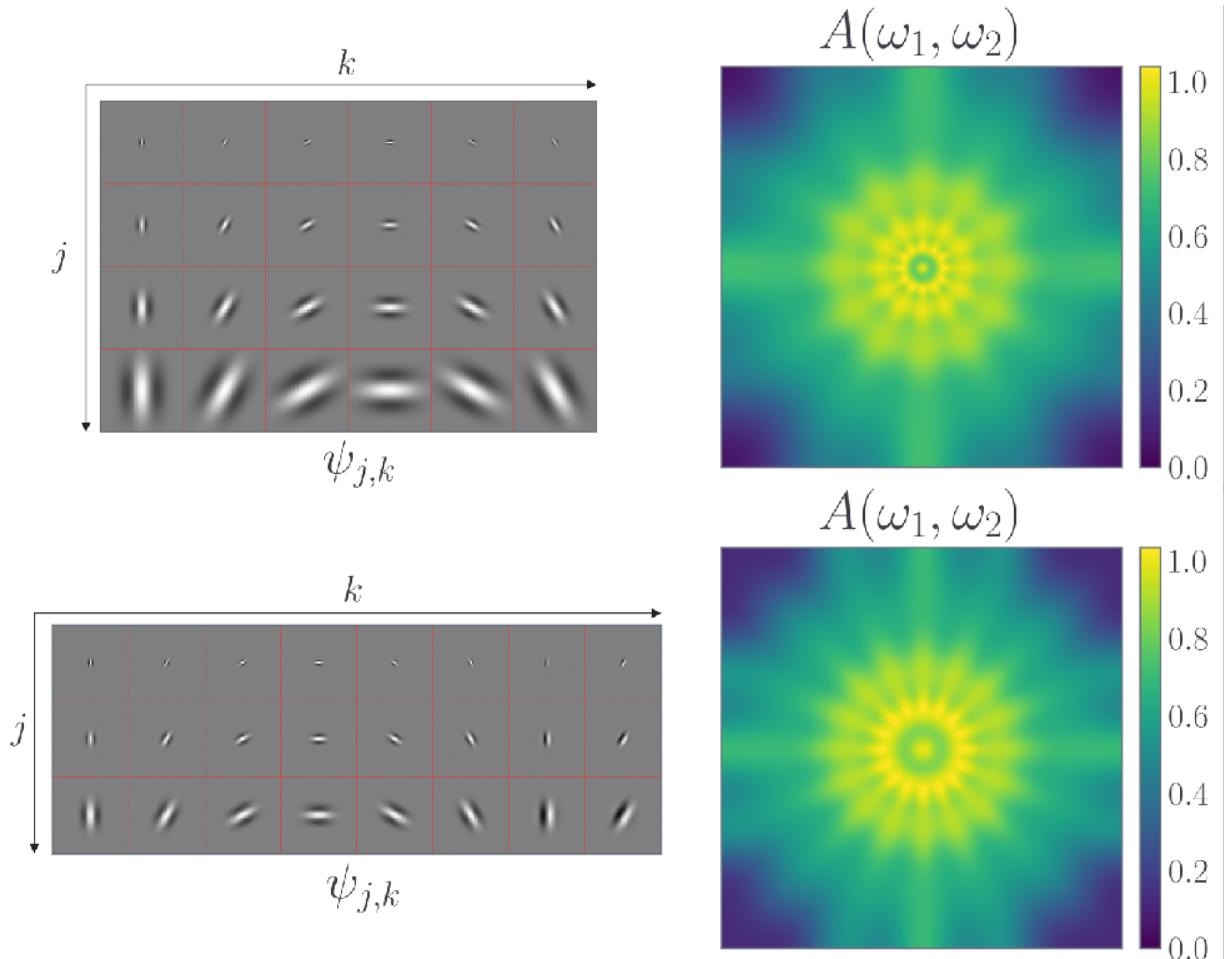


Figure 2.15: Two Morlet Wavelet families and their tiling of the frequency plane. For each set of parameters, the point spread functions of the *real* wavelet bases are shown, next to their covering of the frequency plane $A(\boldsymbol{\omega})$. Each square is 45×45 pixels. Top: $J = 3, K = 6, Q = 1$, Bottom: $J = 4, K = 8, Q = 1$. None of the configurations cover the corners of the frequency plane, but this is often mostly noise. Increasing J, K or Q gives better frequency localization but at the cost of spatial localization and added complexity. Image adapted from [101].

allows for support of both the wavelet and scaling functions on only one half of the frequency plane.

The dual-tree framework shown in [Figure 2.16](#) can achieve this by making the real and imaginary components with their own DWT. We define:

- h_0, h_1 the low and high-pass analysis filters for ϕ_r, ψ_r
- g_0, g_1 the low and high-pass analysis filters for ϕ_i, ψ_i
- \tilde{h}_0, \tilde{h}_1 the low and high pass synthesis filters for $\tilde{\phi}_r, \tilde{\psi}_r$.
- \tilde{g}_0, \tilde{g}_1 the low and high pass synthesis filters for $\tilde{\phi}_i, \tilde{\psi}_i$.

The dilation and wavelet equations for a 1D filter bank implementation are [18]:

$$\phi_r(t) = \sqrt{2} \sum_n h_0(n) \phi_r(2t - n) \quad (2.6.32)$$

$$\psi_r(t) = \sqrt{2} \sum_n h_1(n) \phi_r(2t - n) \quad (2.6.33)$$

$$\phi_i(t) = \sqrt{2} \sum_n g_0(n) \phi_i(2t - n) \quad (2.6.34)$$

$$\psi_i(t) = \sqrt{2} \sum_n g_1(n) \phi_i(2t - n) \quad (2.6.35)$$

Designing a filter bank implementation that results in Hilbert-symmetric wavelets does not appear to be an easy task. However, it was shown by Kingsbury in [93] (and later proved by Selesnick in [116]) that the necessary conditions are conceptually very simple. One low-pass filter must be a *half-sample shift* of the other. I.e., if $g_0(n) = h_0(n - 1/2)$ then the corresponding wavelets are a Hilbert transform pair

$$\psi_g(t) \approx \mathcal{H}\{\psi_h(t)\} \quad (2.6.36)$$

As the DTCWT is designed as an invertible filter bank implementation, this is only one of the constraints. As with conventional (real) discrete wavelets, there are also perfect reconstruction, finite support, linear phase and vanishing moment constraints to consider in the filter bank design.

The derivation of the filters that meet these conditions is covered in detail in [19], [117], and in general in [18]. The result is the option of three main families of filters: biorthogonal filters ($h_0[n] = h_0[N - 1 - n]$ and $g_0[n] = g_0[N - n]$), q-shift filters ($g_0[n] = h_0[N - 1 - n]$), and common-factor filters.

2.6.7.2 2-D DTCWT and its Properties

While analytic wavelets in 1D are useful for their shift-invariance, the real beauty of the DTCWT lies in its ability to make a separable 2D wavelet transform with oriented wavelets.

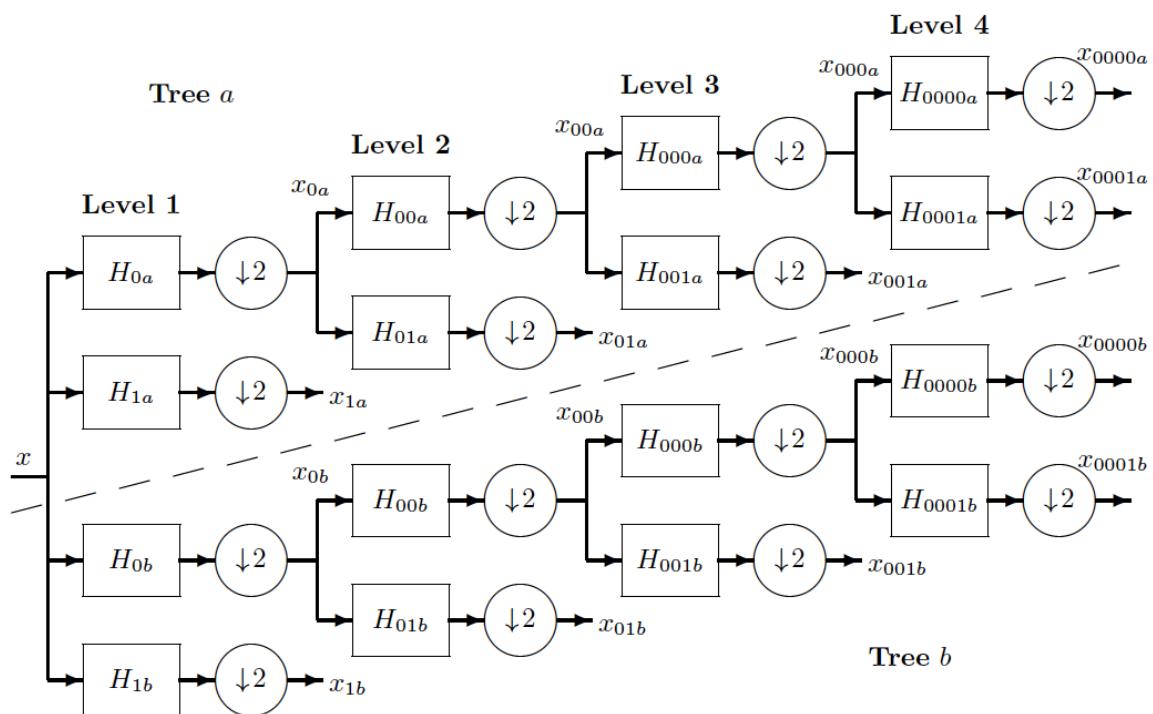


Figure 2.16: **Analysis FBs for the 1-D DTcWT.** Top ‘tree’ forms the real component of the complex wavelet ψ_r , and the bottom tree forms the imaginary (Hilbert pair) component ψ_i . Image taken from [93].

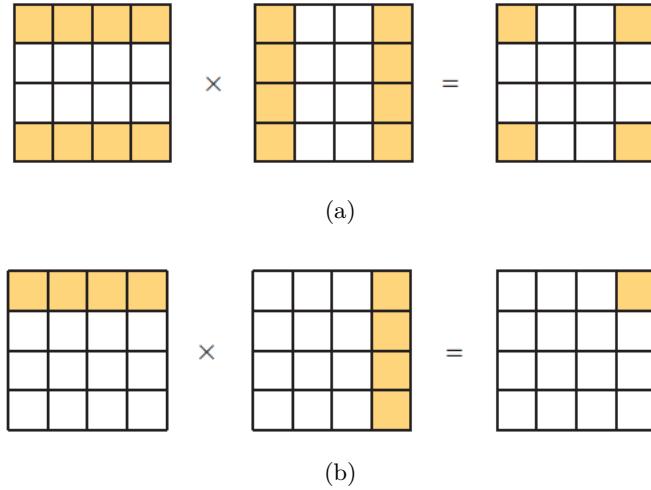


Figure 2.17: The DWT high-high vs the DT-CWT high-high frequency support. (a) The high-high DWT wavelet having a passband in all 4 corners of the frequency plane vs (b) the high-high DT-CWT wavelet frequency support only existing in one quadrant. Taken from [18]

Figure 2.17a shows the spectrum of the wavelet when the separable product uses purely real wavelets, as is the case with the DWT. Figure 2.17b however, shows the separable product of two complex, analytic wavelets resulting in a localized and oriented 2D wavelet.

Note that in this thesis, we name the wavelets by the direction of the edge that they are most sensitive to. For example, the 135° is the second image in Figure 2.18 and can be obtained by the separable product:

$$\psi(\mathbf{u}) = \psi_c(u_1)\psi_c^*(u_2) \quad (2.6.37)$$

$$= (\psi_r(u_1) + j\psi_i(u_1))(\psi_r(u_2) - j\psi_i(u_2)) \quad (2.6.38)$$

$$= (\psi_r(u_1)\psi_r(u_2) + \psi_i(u_1)\psi_i(u_2)) + j(\psi_r(u_1)\psi_i(u_2) - \psi_i(u_1)\psi_r(u_2)) \quad (2.6.39)$$

Similar equations can be obtained for the other five wavelets and the scaling function, by replacing ψ with ϕ for each direction in turn (but not both together), and not taking the complex conjugate in (2.6.37) to get the filters in the right-hand half of the frequency plane. The 2-D DT-CWT requires four 2-D DWTs to calculate the four possible combinations of real and imaginary components. The high and lowpass outputs from these DWTs can then be summed in different ways as in (2.6.39) to get the complex bandpass wavelets. Figure 2.18 shows the resulting wavelets both in the spatial domain and their idealized support in the frequency domain.

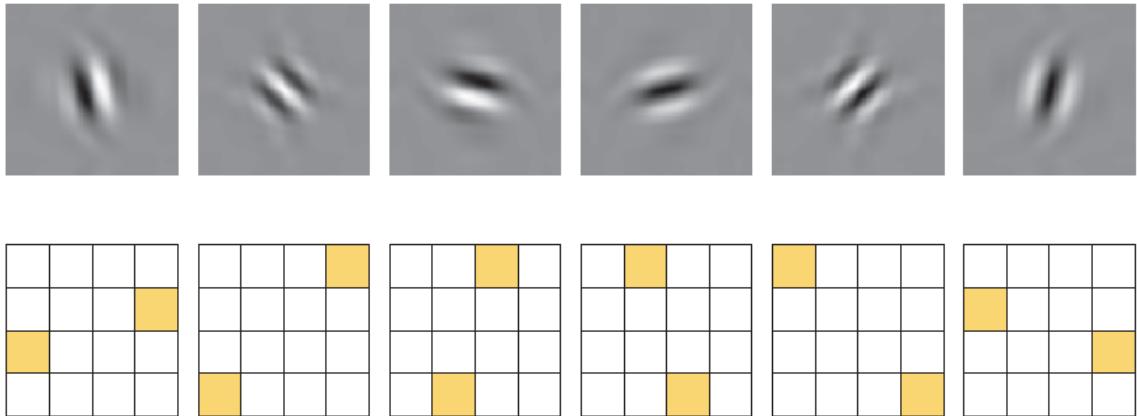


Figure 2.18: **Wavelets from the 2D DTcwt.** **Top:** The six oriented filters in the space domain (only the real wavelets are shown). From left to right these are the $105^\circ, 135^\circ, 165^\circ, 15^\circ, 45^\circ, 75^\circ$ wavelets. **Bottom:** Idealized support of the Fourier spectrum of each wavelet in the 2D frequency plane. Spectra of the the real wavelets are shown — the spectra of the complex wavelets ($\psi_r + j\psi_i$) only has support in the top half of the plane. Image taken from [18].

2.6.7.3 Tightness and Invertibility

We analysed the coverage of the frequency plane for the Morlet wavelet family and saw what areas of the spectrum were better covered than others. How about for the DTcwt?

It is important to note that in the case of the q-shift DTcwt, the wavelet transform is also approximately unitary, i.e.,

$$\|x\|^2 \approx \|\mathcal{W}x\|^2 \quad (2.6.40)$$

and the implementation is perfectly invertible as $A(\omega)$ from (2.6.30) function is unity (or very near unity) $\forall \omega \in [-\pi, \pi] \times [-\pi, pi]$. This is not a surprise, as it is a design constraint in choosing the filters, but nonetheless is important to note.

2.6.8 Summary of Methods

One final comparison to make between the DTcwt and the Morlet wavelets is their frequency coverage. The Morlet wavelets have flexibility at the cost of computational expense and can be made to have tighter angular resolution than the DTcwt. However it is not always better to keep using finer and finer resolutions, indeed the Fourier transform gives the ultimate in angular resolution but as mentioned, this makes it less stable to shifts and deformations. We will explore this in more depth in Chapter 3.

2.7 ScatterNets

ScatterNets have been a very large influence on our work, as well as being quite distinct from the previous discussions on learned methods. They were first introduced by Bruna and Mallat in their work [96], and then were rigorously defined by Mallat in [17]. Perhaps the clearest explanation of them, and the most relevant to our work is in [23].

While CNNs have the ability to learn invariances to nuisance variabilities, their properties and optimal configurations are not well understood. It typically takes multiple trials by an expert to find the correct hyperparameters for these networks. A scattering transform instead builds well understood and well-defined invariances.

We first review some of the desirable invariances before describing how a ScatterNet achieves them.

2.7.1 Desirable Properties

2.7.1.1 Translation-Invariance

Translation is often said to be uninformative for classification — an object appearing in the centre of the image *should* be treated the same way as a similar object appearing near the corner of an image. This can be quantified by saying a representation Φx is invariant to global translations $x_c(\mathbf{u}) = x(\mathbf{u} - \mathbf{c})$ by $\mathbf{c} = (c_1, c_2) \in \mathbb{R}^2$ if:

$$\|\Phi x_c - \Phi x\| \leq C \quad (2.7.1)$$

for some small constant $C > 0$. Note that we may instead want only local translation-invariance and restrict the distance $|\mathbf{c}|$ for which (2.7.1) is true.

Convolutional filters are naturally covariant to translations in the pixel space, so $\Phi x_c = (\Phi x)_c$, $\mathbf{c} \in \mathbb{Z}^2$. Of course, natural objects exist in continuous space and are sampled, and any two images of the same scene taken with small camera disturbances are unlikely to be at integer pixel shifts of each other.

2.7.1.2 Stability to Noise

Stability to additive noise is another useful invariance to incorporate, as it is a common feature in sampled signals. Stability is defined in terms of Lipschitz continuity, which is a strong form of uniform continuity for functions, which we briefly introduce here.

Formally, a Lipschitz continuous function is limited in how fast it can change; there exists an upper bound on the gradient the function can take, although it doesn't necessarily need to be differentiable everywhere. The modulus operator $|x|$ is a good example of a function that has a bounded derivative and so is Lipschitz continuous, but isn't differentiable everywhere.

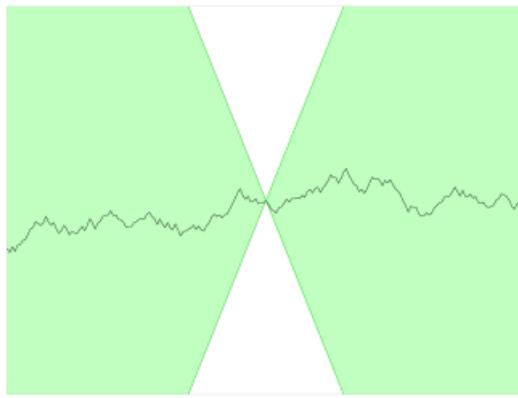


Figure 2.19: **A Lipschitz continuous function.** There is a cone for this function (shown in white) such that the graph always remains entirely outside the cone as it is shifted across. The minimum gradient needed for this to hold is called the ‘best Lipschitz constant’.

Alternatively, the modulus squared has derivative everywhere but is not Lipschitz continuous as its gradient grows with x .

To be stable to additive noise, we require that for a new signal $x'(\mathbf{u}) = x(\mathbf{u}) + \epsilon(\mathbf{u})$, there must exist a bounded $C > 0$ s.t.

$$\|\Phi x' - \Phi x\| \leq C \|x' - x\| \quad (2.7.2)$$

2.7.1.3 Stability to Deformations

Small deformations are important to be invariant to but this must be limited. It is important to keep intra-class variations small but not be so invariant that an object can morph into another (in the case of MNIST for example, we do not want to be so stable to deformations that 7s can map to 1s).

Formally, for a new signal $x_\tau(\mathbf{u}) = x(\mathbf{u} - \tau(\mathbf{u}))$, where $\tau(\mathbf{u})$ is a non constant displacement field (i.e., not just a translation) that deforms the image, we require a $C_\tau > 0$ s.t.

$$\|\Phi x_\tau - \Phi x\| \leq C_\tau \|x\| \sup_{\mathbf{u}} |\nabla \tau(\mathbf{u})| \quad (2.7.3)$$

The term on the right $|\nabla \tau(\mathbf{u})|$ measures the deformation amplitude, so the supremum of it is a limit on the global deformation amplitude.

2.7.2 Definition

A Fourier modulus satisfies the first two of these requirements, in that it is both translation-invariant and stable to additive noise, but it is unstable to deformations due to the large support (infinite in theory) of the sinusoid basis functions it uses. It also loses too much

information — very different signals can all have the same Fourier modulus, e.g. a chirp, white noise and the Dirac delta function all have flat spectra.

Another translation-invariant and stable operator is the averaging kernel which Mallat et. al. use to make the zeroth scattering coefficient:

$$S[\emptyset]x \triangleq x * \phi_J(2^J \mathbf{u}) \quad (2.7.4)$$

which is translation-invariant to shifts less than 2^J . It, unfortunately, results in a loss of information due to the removal of high-frequency content. This is easy to see as the wavelet operator from (2.6.26) $\mathcal{W}x = \{x * \phi_J, x * \psi_\lambda\}_\lambda$ contains all the information of x , whereas the zeroth scattering coefficient is simply the lowpass portion of \mathcal{W} .

This high-frequency content can be ‘recovered’ by keeping the wavelet coefficients. The wavelet terms, like a convolutional layer in a CNN, are only covariant to shifts rather than invariant. This covariance happens in the real and imaginary parts which both vary rapidly. Fortunately, its modulus is much smoother and gives a good measure for the frequency-localized energy content at a given spatial location⁷. Unlike the Fourier modulus, the complex wavelet modulus is stable to deformations due to the grouping together of frequencies into dyadic packets [17].

We combine the wavelet transform and modulus operators into one operator $\widetilde{\mathcal{W}}$:

$$\widetilde{\mathcal{W}}x = \{x * \phi_J, |x * \psi_\lambda|\}_\lambda \quad (2.7.5)$$

$$= \{x * \phi_J, U[\lambda]x\}_\lambda \quad (2.7.6)$$

where the U terms are called the *propagated* signals. These U terms are approximately invariant for shifts of up to 2^j . Mallat et. al. choose to keep the same level of invariance as the zeroth-order coefficients (2^J) by further averaging. This makes the first ordering scattering coefficients:

$$S[\lambda_1]x \triangleq U[\lambda_1]x * \phi_J = |x * \psi_{\lambda_1}| * \phi_J \quad (2.7.7)$$

Again this averaging comes at a cost of discarding high-frequency information, this time about the wavelet sparsity signal $U[\lambda] = |x * \psi_\lambda|$ instead of the input signal x . We can recover this information by repeating the above process.

$$S[\lambda_1, \lambda_2]x \triangleq U[\lambda_2]U[\lambda_1]x \quad (2.7.8)$$

$$= ||x * \psi_{\lambda_1}| * \psi_{\lambda_2}| * \phi_J \quad (2.7.9)$$

⁷Interestingly, the modulus operator can often still be inverted due to the redundancies of the complex wavelet transform [118], hence it does not lose any information.

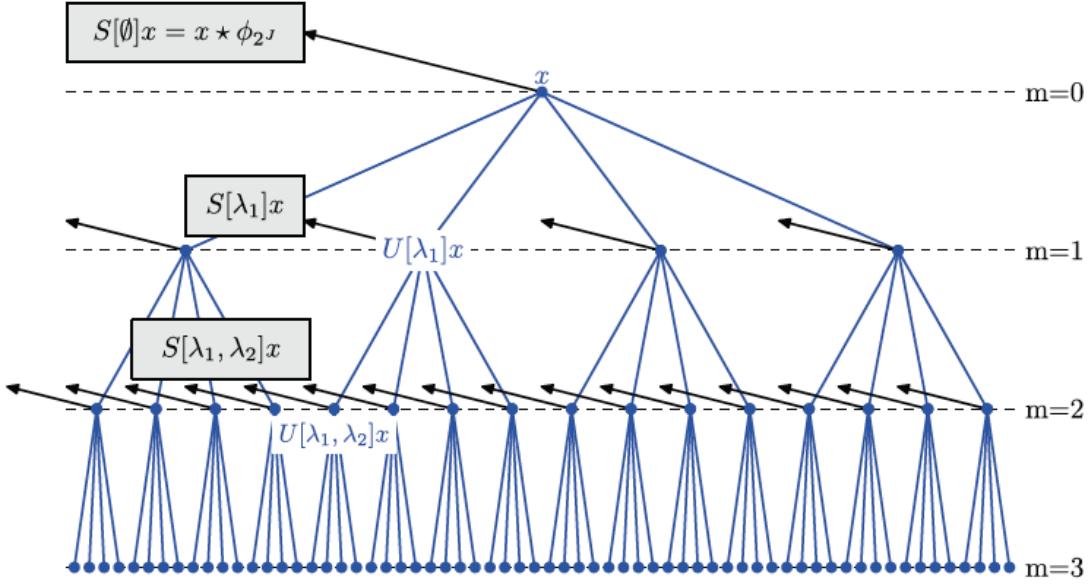


Figure 2.20: **The Scattering Transform.** Scattering outputs are the leftward pointing arrows $S[p]x$, and the intermediate coefficients $U[p]x$ are the centre nodes of the tree. Taken from [23].

In general, let $p = (\lambda_1, \lambda_2, \dots, \lambda_m)$ be a path of length m describing the order of application of wavelets, and define:

$$U[p]x = U[\lambda_m]U[\lambda_{m-1}] \cdots U[\lambda_1]x \quad (2.7.10)$$

$$= ||\cdots|x * \psi_{\lambda_1}| * \psi_{\lambda_2}| \cdots * \psi_{\lambda_m}| \quad (2.7.11)$$

and the m th order scattering coefficient along the path p is $S[p]x = U[p]x * \phi_J$. Further, let $p + \lambda = (\lambda_1, \lambda_2, \dots, \lambda_m, \lambda)$. This allows us to recursively define the next set of *propagated* and *scattering* coefficients by using $\widetilde{\mathcal{W}}$:

$$\widetilde{\mathcal{W}}U[p]x = \{S[p]x, U[p + \lambda]x\}_{\lambda} \quad (2.7.12)$$

which is shown in Figure 2.20

2.7.3 Resulting Properties

For ease, let us define the ‘ m th order scattering coefficients’ as S_m which is the set of all coefficients with path length m . Further, let S be the set of all scattering coefficients of any path length. The energy $\|Sx\|^2$ we then define as

$$\|Sx\|^2 = \sum_p \|S[p]x\|^2 \quad (2.7.13)$$

We can make \mathcal{W} non-expansive with appropriate scaling. Define the energy $\|\mathcal{W}x\|^2$ as

$$\|\mathcal{W}x\|^2 = \|x * \phi\|^2 + \sum_{\lambda} \|x * \psi_{\lambda}\|^2 \quad (2.7.14)$$

then by Plancherel's formula

$$(1 - \epsilon) \|x\|^2 \leq \|\mathcal{W}x\|^2 \leq \|x\|^2 \quad (2.7.15)$$

For the Morlet wavelets originally used in [23], $\epsilon = 0.25$, for the DT \mathbb{C} WT $\epsilon \approx 0$ (for the q-shift DT \mathbb{C} WT it is 0, but for the biorthogonal DT \mathbb{C} WT it is close to but not exactly 0).

2.7.3.1 Translation-Invariance

This is proven in section 2.4 of [17]. We have so far described the Scattering representation as being ‘translation-invariant for shifts up to 2^J ’. We formalize this statement here.

For a 2-D averaging filter based on a father wavelet ϕ , $\phi_J = 2^{-J}\phi(2^{-J}\mathbf{u})$ it is proven in Appendix B of [17] that shifting it by \mathbf{c} , which we denote as \mathcal{L}_c , is Lipschitz continuous:

$$\|\mathcal{L}_c\phi_J - \phi_J\| \leq 2^{-J+2}\|\nabla\phi\|_1|\mathbf{c}| \quad (2.7.16)$$

where $\|\nabla\phi\|_1$ is the ℓ_1 norm of the grad of ϕ .

For simplicity, let us define $A_Jx = \phi_J * x$ and $Sx = A_JUx$. Then we get:

$$\|S\mathcal{L}_c x - Sx\| = \|\mathcal{L}_c A_J Ux - A_J Ux\| \quad (2.7.17)$$

$$\leq \|\mathcal{L}_c A_J - A_J\| \|Ux\| \quad (2.7.18)$$

$$\leq 2^{-J+2}\|\nabla\phi\|_1|\mathbf{c}|\|x\| \quad (2.7.19)$$

2.7.3.2 Stability to Noise

As \mathcal{W} is non-expansive and the complex modulus is also non-expansive

$$\|\widetilde{\mathcal{W}}x - \widetilde{\mathcal{W}}y\| \leq \|x - y\| \quad (2.7.20)$$

We have already shown that S is the repeated application of $\widetilde{\mathcal{W}}$ in (2.7.12), we can then say

$$\|Sx - Sy\| \leq \|x - y\| \quad (2.7.21)$$

making scattering non-expansive and stable to noise. With the DT \mathbb{C} WT, both (2.7.20) and (2.7.21) are nearly inequalities as ϵ is close to 0.

2.7.3.3 Stability to deformations

If $\mathcal{L}_\tau x = x(\mathbf{u} - \tau(\mathbf{u}))$ is an image deformed by a diffeomorphism τ with $\|\tau\|_\infty = \sup_u |\tau(\mathbf{u})|$ and $\|\nabla\tau\|_\infty = \sup_u |\nabla\tau(\mathbf{u})| < 1$ then it is proven in [17] that:

$$\|S\mathcal{L}_\tau x - Sx\| \leq CP \|x\| (2^{-J} \|\tau\|_\infty + \|\nabla\tau\|_\infty) \quad (2.7.22)$$

where $P = \text{length}(p)$ is the scattering order and C is a constant (dependent on J). For deformations with small absolute displacement relative to 2^J , the first term disappears and we have:

$$\|S\mathcal{L}_\tau x - Sx\| \leq CP \|x\| \|\nabla\tau\|_\infty \quad (2.7.23)$$

This theorem shows that S is locally Lipschitz stable to diffeomorphisms and in the case of small deformations, it linearizes them.

2.7.3.4 Energy Decay

As $m \rightarrow \infty$ the invariant coefficients of path length m , U_m , decay towards zero [17]:

$$\lim_{m \rightarrow \infty} U_m = 0 \quad (2.7.24)$$

This is an important property that suggests that we can stop scattering beyond a certain point. Experimental results [23] for image sizes on the order of a few hundred pixels by a few hundred pixels, $m = 3$ captures about 99% of the input energy. For many works using scattering transforms after [23] such as [25], [39], [40], setting $m = 2$ was found to be sufficient.

2.7.3.5 Number of Coefficients

While we have so far talked about non-sampled signals $x(\mathbf{u})$, $\mathbf{u} \in \mathbb{R}^2$, in practice, we want to apply scattering to sampled signals $x[\mathbf{n}]$, $\mathbf{n} \in \mathbb{Z}^2$. The averaging by ϕ_J means that we can subsample Sx by 2^J in each direction. However, now we need also need to index all the paths p that can be used to create the scattering coefficients. Limiting ourselves to $m = 2$ and using a wavelet transform with J scales and K discrete orientations the number of paths for each S_m is the cardinality of the set p_m :

$$n(p_0) = 1 \quad (2.7.25)$$

$$n(p_1) = JK \quad (2.7.26)$$

$$n(p_2) = (J-1)K^2 + (J-2)K^2 + \dots + K^2 \quad (2.7.27)$$

$$= \frac{1}{2}J(J-1)K^2 \quad (2.7.28)$$

Table 2.1: **Redundancy of Scattering Transform.** Shows the number of output channels C_{out} and number of pixels per channel N_{out} for different scattering orders m , scales J , and orientations K for a single channel input image with N pixels.

| m | J | K | C_{out} | N_{out} |
|-----|-----|-----|-----------|-----------|
| 1 | 2 | 6 | 13 | $N/16$ |
| 1 | 2 | 8 | 17 | $N/16$ |
| 2 | 2 | 6 | 49 | $N/16$ |
| 2 | 2 | 8 | 81 | $N/16$ |
| 2 | 3 | 6 | 127 | $N/64$ |
| 2 | 3 | 8 | 217 | $N/64$ |
| 3 | 3 | 6 | 343 | $N/64$ |
| 3 | 3 | 8 | 729 | $N/64$ |

The reason $n(p_2) \neq J^2 K^2$ is due to the demodulating effect of the complex modulus. As $|x * \psi_\lambda|$ is more regular than $x * \psi_\lambda$, $|x * \psi_\lambda| * \psi_{\lambda'}$ is only non-negligible if $\psi_{\lambda'}$ is located at lower frequencies than ψ_λ . This means we can discard over half of the scattering paths as their value will be near zero.

Summing up the above three equations and factoring in the reduced sample rate allowable due to averaging, for an input with N pixels, a second-order scattering representation will have $N 2^{-2J} (1 + JK + \frac{1}{2}J(J-1)K^2)$ pixels. Table 2.1 shows some example values of the ScatterNet redundancy for different J, K and scattering order m .

Chapter 3

A Faster ScatterNet

The drive of this thesis is in exploring if wavelet theory, in particular the DT \mathbb{C} WT, has any place in deep learning and if it does, quantifying how beneficial it can be. The introduction of more powerful GPUs and fast and popular deep learning frameworks such as PyTorch, Tensorflow and Caffe in the past few years has helped the field of deep learning grow very rapidly. Never before has it been so possible and so accessible to test new designs and ideas for a machine learning algorithm than today. Despite this rapid growth, there has been little interest in building wavelet analysis software in modern frameworks.

This poses a challenge and an opportunity. To pave the way for more detailed investigation (both in the rest of this thesis and by other researchers who want to explore wavelets applied to deep learning), we must have the right foundation and tools to facilitate research.

A good example of this is the current implementation of the ScatterNet. While ScatterNets have been the most promising start in using wavelets in a deep learning system, they have tended to be orders of magnitude slower, and significantly more difficult to run than a standard convolutional network.

Additionally, any researchers wanting to explore the DWT in a deep learning system have had to rewrite the filter bank implementation themselves, ensuring they correctly handle boundary conditions and ensure correct filter tap alignment to achieve perfect reconstruction.

3.1 Chapter Layout

This chapter describes how we have built a fast ScatterNet implementation in PyTorch with the DT \mathbb{C} WT as its wavelet transform. First we describe how to do an efficient DWT in PyTorch in [section 3.4](#) before showing how to expand this to an efficient DT \mathbb{C} WT in [section 3.5](#). We then use the DT \mathbb{C} WT to define our own ScatterNet in [section 3.6](#) (in particular, see [Algorithm 3.5](#)). All of the code is available as an open-source library at *PyTorch Wavelets* [28].

In parallel with our efforts, the original authors of the ScatterNet have improved their implementation, making a new package called KyMatIO[119]. We compare the speed and classification performance of our package to KyMatIO in section 3.7 as this provides some interesting insights into the choice of complex wavelet for a ScatterNet. This is similar to the work of [36], where Singh and Kingsbury show that a DT^CWT-ScatterNet outperforms a Morlet-ScatterNet when used as a front end to an SVM for some simpler classification tasks. We find that our proposed DT^CWT-ScatterNet is 7 to 15 times faster than KyMatIO (depending on the padding style and wavelet length), as well as giving a small improvement in performance when used as a front end to a CNN.

3.2 The Design Constraints

3.2.1 Original Design

The original authors implemented their ScatterNet in Matlab [25] using a Fourier-domain based Morlet wavelet transform. The standard procedure for using ScatterNets in a deep learning framework up until recently has been to:

1. Pre-scatter a dataset using conventional CPU-based hardware and software and store the features to disk. This can take several hours to several days depending on the size of the dataset and the number of CPU cores available.
2. Build a network in another framework, usually Tensorflow [120] or Pytorch [64].
3. Load the scattered data from disk and train on it.

We saw that this approach was suboptimal for a number of reasons:

- It is slow and must run on CPUs.
- It is inflexible to any changes you wanted to investigate in the Scattering design; you would have to re-scatter all the data and save elsewhere on disk.
- You can not easily do preprocessing techniques like random shifts and flips, as each of these would change the scattered data.
- The scattered features are often larger than the original images and require you to store entire datasets twice (or more) times.
- The features are fixed and can only be used as a front end to any deep learning system.

3.2.2 Improvements

To address these shortcomings, all of the above limitations become design constraints. In particular, the new software should be:

- Able to run on GPUs (ideally on multiple GPUs in parallel).
- Flexible and fast so that it can run as part of the forward pass of a neural network (allowing preprocessing techniques like random shifts and flips).
- Able to pass gradients through, so that it can be part of a larger network and have learning stages before scattering.

To achieve all of these goals, we choose to build our software on PyTorch using the DTCWT. PyTorch is a popular open-source deep learning framework that can do many operations on GPUs with native support for automatic differentiation. PyTorch uses the CUDA and cuDNN libraries for its GPU-accelerated primitives. Its popularity is of key importance, as it means users can build complex networks involving ScatterNets without having to use or learn extra software.

3.3 A Brief Description of Autograd

As part of a modern deep learning framework, we need to define functional units like the one shown in [Figure 2.6](#). Not only must we be able to calculate the forward evaluation of a block $Y = f(X, h)$ given an input X and (possibly) some learnable weights h , we must also be able to calculate the passthrough and update gradients $\frac{\partial \mathcal{L}}{\partial X}$, $\frac{\partial \mathcal{L}}{\partial h}$ given $\frac{\partial \mathcal{L}}{\partial Y}$. This typically involves saving $\frac{\partial Y}{\partial X}$ and $\frac{\partial Y}{\partial h}$ evaluated at the current values of X and h when we calculate the forward pass.

For example, the simple ReLU: $Y = \max(0, X)$, is not memory-less. On the forward pass, we need to put a 1 in all the positions where $X > 0$, and a 0 elsewhere. For a convolutional layer, we need to save X and h to correlate with $\frac{\partial \mathcal{L}}{\partial Y}$ on the backwards pass. It is up to the block designer to manually calculate the gradients and design the most efficient way of programming them. For clarity and repeatability, we give pseudocode for all the core operations developed in our package *PyTorch Wavelets*.

3.4 Fast Calculation of the DWT and IDWT

To have a fast implementation of the Scattering Transform, we need a fast implementation of the DTCWT. Similarly, for a fast implementation of the DTCWT we need a fast implementation of the DWT. Future work may also explore the DWT as a basis for learning, so having an implementation that is fast and can pass gradients through will prove beneficial in and of itself.

3.4.1 The Input

Recall from [subsection 2.4.1](#) that our input is a 3-D array with channel dimension first, followed by the two spatial dimensions. For a minibatch of images, this becomes a 4-D array, with the sample number in the first dimension. I.e., for a minibatch of N images with C channels and $H \times W$ spatial support, the input has shape: $N \times C \times H \times W$.

3.4.2 Preliminary Notes

There has been much research into the best way to do the DWT on a GPU, in particular comparing the speed of Lifting [\[121\]](#), or second-generation wavelets, to the direct convolutional methods. [\[122\]](#), [\[123\]](#) are two notable such publications, both of which find that the convolution-based implementations are better suited for the massively parallel architecture found in modern GPUs. For this reason, we implement a convolutional-based DWT.

As the DWT and IDWT uses fixed filters, we do not need to calculate $\frac{\partial L}{\partial h}$ on the backwards pass. This means on the forward pass we can save on memory by not storing $\frac{\partial Y}{\partial h} = X$ (see [subsubsection 2.4.1.2](#)). We found it easiest to explicitly rewrite forward and backward passes for the DWT, as the autograd mechanics would often unnecessarily save intermediate activations.

For example, consider an input $X \in \mathbb{R}^{128 \times 256 \times 64 \times 64}$. This is not an unrealistically large input for a CNN, as often minibatch sizes are 128, channels can be in the hundreds and 64×64 pixels is a relatively common input size. If we were to represent these numbers as floats, it would require 512MB of space on a GPU (modern GPUs have about 10GB of memory, so this single activation already requires 5%). If we perform a DWT on this input, we would require another 512MB of space for the output. Using naive autograd, PyTorch (or another framework) saves an extra 512MB of memory for the backwards pass to calculate $\frac{\partial L}{\partial h}$, even if we specify it as not requiring a gradient. This means we can save 50% of the memory cost by being explicit about what is, and what is not needed for backpropagation.

3.4.3 Primitives

We start with the commonly known property that for a convolutional block, the passthrough gradient is the gradient with respect to the output convolved with the time reverse of the filter. More formally, if $Y(z) = H(z)X(z)$, then:

$$\Delta X(z) = H(z^{-1})\Delta Y(z) \tag{3.4.1}$$

where $H(z^{-1})$ is the Z -transform of the time/space reverse of $H(z)$, $\Delta Y(z) \triangleq \frac{\partial L}{\partial Y}(z)$ is the gradient of the loss with respect to the output, and $\Delta X(z) \triangleq \frac{\partial L}{\partial X}(z)$ is the gradient of the loss with respect to the input. This was proved in [subsubsection 2.4.1.2](#), we have just rewritten it in terms of Z -transforms here.

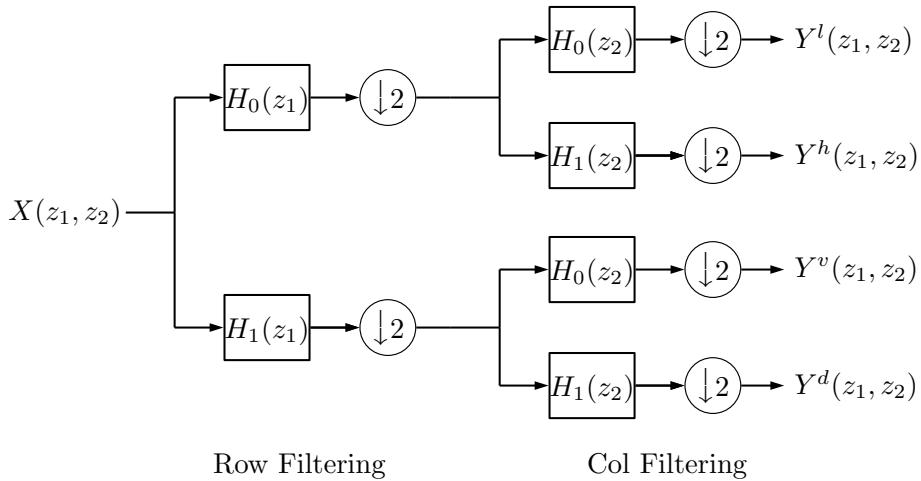


Figure 3.1: 2-D DWT Filter Bank Layout. The components of a filter bank DWT in two dimensions. Y^l, Y^h, Y^v, Y^d represent the lowpass, horizontal, vertical, and diagonal components.

Additionally, for a decimation by two block, the passthrough gradient is interpolation by two (see [section B.1](#) for a proof of this).

[Figure 3.1](#) shows the block diagram for performing the forward pass of a DWT. We can draw a similar figure for the backwards pass by replacing all the forward operations with their passthrough gradients. Using the above two properties, we see that the backwards pass of a DWT has the same form as the inverse wavelet transform, with the time reverse of the analysis filters used as the synthesis filters. For orthogonal wavelet transforms, the synthesis filters are the time reverse of the analysis filters so backpropagation can be done by doing the inverse wavelet transform with $\Delta Y^l(z), \Delta Y^h(z), \Delta Y^v(z), \Delta Y^d(z)$. See [section B.2](#) for more information on this and the equivalent figure for the backwards pass.

Like Matlab, most deep learning frameworks have an efficient function for doing convolution followed by downsampling. Similarly, there is an efficient function to do upsampling followed by convolution (for the inverse transform). Let us call these `conv2d_down` and `conv2d_up` respectively¹. These functions, in turn, call the cuDNN low-level functions which can only support zero padding or no padding (we called this *valid convolution* in [subsubsection 2.4.1.1](#)). If another padding type is desired, it must be done beforehand and valid convolution used.

With both `conv2d_down` and `conv2d_up`, we want to apply the same filter to each channel of the input and *not* sum them up (recall [\(2.4.5\)](#) sums over the channel dimension after doing convolution). There are options available to prevent the summing in most frameworks, related

¹E.g. In PyTorch, convolution followed by downsampling is done with a call to `torch.nn.functional.conv2d` with the stride parameter set to 2. Upsampling by 2 followed by convolution is done by calling `torch.nn.functional.conv2d_transpose`.

Algorithm 3.1 1-D analysis and synthesis stages of a DWT

```

1: function AFB1D( $x, h_0, h_1, mode, axis$ )
2:   if  $axis = 3$  then
3:      $h_0, h_1 \leftarrow h_0^t, h_1^t$                                  $\triangleright$  row filtering
4:   end if
5:    $p \leftarrow \lfloor (len(x) + len(h_0) - 1)/2 \rfloor$            $\triangleright$  calculate output size
6:    $b \leftarrow \lfloor p/2 \rfloor$                                  $\triangleright$  calculate pad size before
7:    $a \leftarrow \lceil p/2 \rceil$                                  $\triangleright$  calculate pad size after
8:    $x \leftarrow \text{pad}(x, b, a, mode)$                           $\triangleright$  pre pad the signal with selected mode
9:    $lo \leftarrow \text{conv2d\_down}(x, h_0)$ 
10:   $hi \leftarrow \text{conv2d\_down}(x, h_1)$ 
11:  return  $lo, hi$ 
12: end function

1: function SFB1D( $lo, hi, g_0, g_1, axis$ )
2:   if  $axis = 3$  then
3:      $g_0, g_1 \leftarrow g_0^t, g_1^t$                                  $\triangleright$  row filtering
4:   end if
5:    $p \leftarrow \text{len}(g_0) - 2$                                  $\triangleright$  calculate output size
6:    $lo \leftarrow \text{pad}(lo, p, p, 'zero')$                        $\triangleright$  pre pad the signal with zeros
7:    $hi \leftarrow \text{pad}(hi, p, p, 'zero')$                        $\triangleright$  pre pad the signal with zeros
8:    $x \leftarrow \text{conv2d\_up}(lo, g_0) + \text{conv2d\_up}(hi, g_1)$ 
9:   return  $x$ 
10: end function

```

to doing *grouped convolution*. For a good explanation on grouped convolution, we recommend Chapter 5 of [49]. In our code below, we assume that `conv2d_down` and `conv2d_up` act independently on the C channels and do not sum across them, giving an output with the same shape as the input - $N \times C \times H \times W$.

3.4.4 1-D Filter Banks

Let us assume that the analysis (h_0, h_1) and synthesis (g_0, g_1) filters are already in the form needed to do *column* filtering. The necessary steps to do the 1-D analysis and synthesis filtering are described in Algorithm 3.1. We do not need to define backpropagation functions for the `afb1d` (analysis filter bank 1-d) and `sfb1d` (synthesis filter bank 1-D) functions as they are each other's backwards pass.

3.4.5 2-D Transforms and Their Gradients

Having built the 1-D filter banks, we can easily generalize them to 2-D. Furthermore, we can define the backwards steps of both the forward 2-D DWT and the inverse 2-D DWT using

Algorithm 3.2 2-D DWT and its gradient

```

1: function DWT.FORWARD( $x, h_0^r, h_1^r, h_0^c, h_1^c, mode$ )
2:   save  $h_0^r, h_1^r, h_0^c, h_1^c, mode$                                  $\triangleright$  For the backwards pass
3:    $lo, hi \leftarrow \text{afb1d}(x, h_0^r, h_1^r, mode, axis = 3)$            $\triangleright$  row filter
4:    $ll, lh \leftarrow \text{afb1d}(lo, h_0^c, h_1^c, mode, axis = 2)$            $\triangleright$  column filter
5:    $hl, hh \leftarrow \text{afb1d}(hi, h_0^c, h_1^c, mode, axis = 2)$            $\triangleright$  column filter
6:   return  $ll, lh, hl, hh$ 
7: end function

1: function DWT.BACKWARD( $\Delta ll, \Delta lh, \Delta hl, \Delta hh$ )
2:   load  $h_0^r, h_1^r, h_0^c, h_1^c, mode$ 
3:    $\Delta lo \leftarrow \text{sfb1d}(\Delta ll, \Delta lh, h_0^c, h_1^c, mode, axis = 2)$ 
4:    $\Delta hi \leftarrow \text{sfb1d}(\Delta hl, \Delta hh, h_0^r, h_1^r, mode, axis = 2)$ 
5:    $\Delta x \leftarrow \text{sfb1d}(\Delta lo, \Delta hi, h_0^r, h_1^r, mode, axis = 3)$ 
6:   return  $\Delta x$ 
7: end function

```

these filter banks. We show how to do this in [Algorithm 3.2](#). The inverse transform logic is moved to the appendix - in [Algorithm B.1](#).

Note that we have allowed for different row and column filters in [Algorithm 3.2](#). Most commonly used wavelets will use the same filter for both directions (e.g. the orthogonal Daubechies family), but later when we use the DT \mathbb{C} WT, we will want to have different horizontal and vertical filters.

Further, note that the only things that need to be saved are the filters, as seen in [Algorithm 3.2.2](#). These are typically only a few floats, giving us a large saving over relying on autograd.

A multiscale DWT can easily be made by calling [Algorithm 3.2](#) multiple times on the lowpass output. Again, no intermediate activations need to be saved, giving this implementation almost no memory overhead.

3.5 Fast Calculation of the DT \mathbb{C} WT

We have built upon previous implementations of the DT \mathbb{C} WT [124]–[126]. The ‘dual-tree’ part of the DT \mathbb{C} WT comes from it having two trees, each with their own set of filters: $a = \{h_0^a, h_1^a, g_0^a, g_1^a\}$ and $b = \{h_0^b, h_1^b, g_0^b, g_1^b\}$. In one dimension, this can be done with two DWTs, one using the a filters and the other using the b filters. In two dimensions, we must instead do four multiscale DWTs. We follow the notation in [93] and maintain separate imaginary operators j_1 and j_2 for the row and column processing. We call our four DWTs r, j_1, j_2, j_1j_2 and show how the filter bank system is connected in [Figure 3.2](#). The output of these four DWTs can be interpreted as a four-element ‘complex’ vector $\{a, b, c, d\} = a + bj_1 + cj_2 + dj_1j_2$. This 4-vector can be converted into a pair of complex 2-vectors by letting $j_1 = j_2 = j$ in one

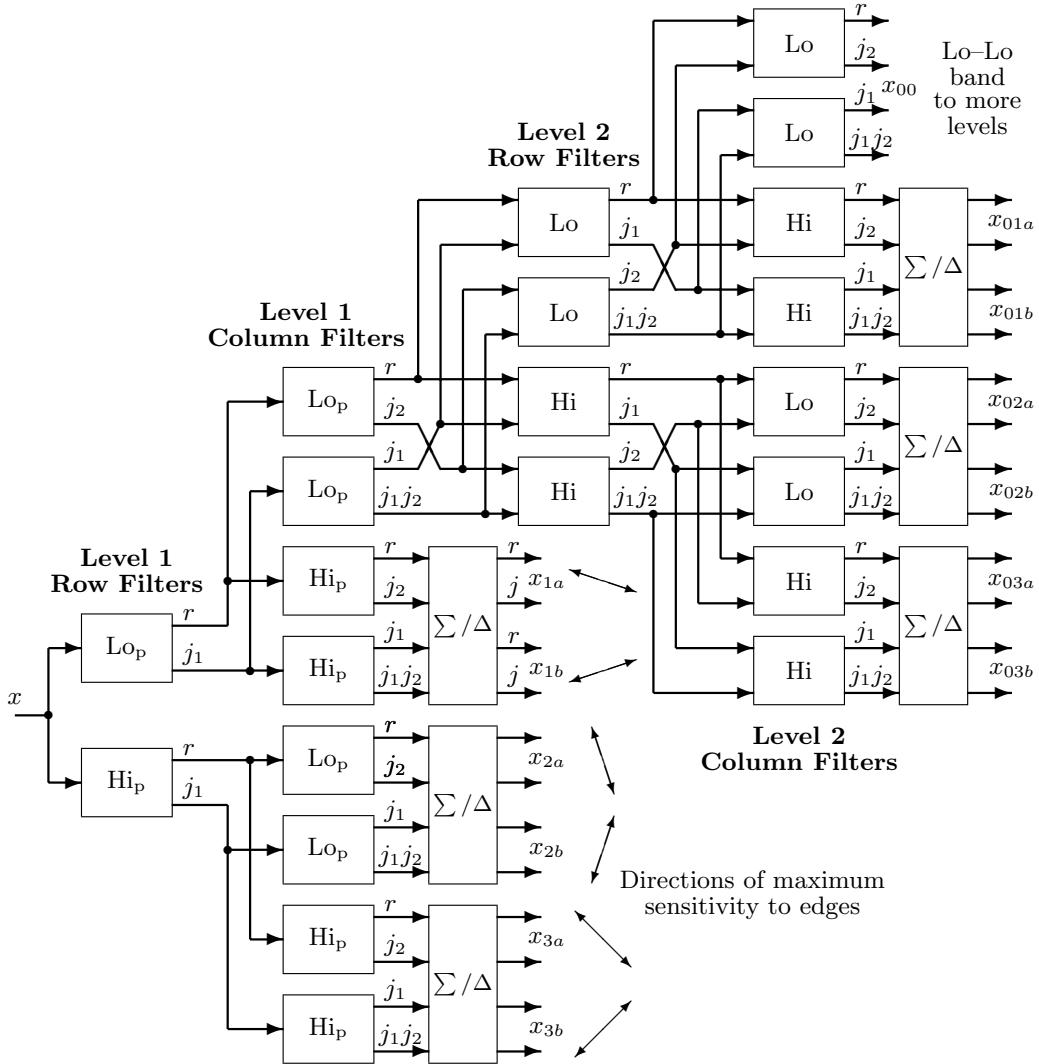


Figure 3.2: **2-D DTcCWT filter bank layout.** Two levels of the complex wavelet tree for a real 2-D input image x . The four DWTs involved in the DTcCWT are labelled with r , j_1 , j_2 and j_1j_2 . Image taken from [93].

Algorithm 3.3 2-D DTcWT.

```

1: function DTcWT( $x$ ,  $J$ , mode)
2:   load  $h_0^a$ ,  $h_1^a$ ,  $h_0^b$ ,  $h_1^b$                                  $\triangleright$  Load from memory
3:    $ll_r$ ,  $ll_{j_1}$ ,  $ll_{j_2}$ ,  $ll_{j_1j_2} \leftarrow x/2$ 
4:   for  $1 \leq j \leq J$  do
5:      $ll_r$ ,  $lh_r$ ,  $hl_r$ ,  $hh_r \leftarrow \text{DWT}_r(ll_r, h_0^a, h_1^a, h_0^a, h_1^a, mode)$ 
6:      $ll_{j_1}$ ,  $lh_{j_1}$ ,  $hl_{j_1}$ ,  $hh_{j_1} \leftarrow \text{DWT}_{j_1}(ll_{j_1}, h_0^b, h_1^b, h_0^a, h_1^a, mode)$ 
7:      $ll_{j_2}$ ,  $lh_{j_2}$ ,  $hl_{j_2}$ ,  $hh_{j_2} \leftarrow \text{DWT}_{j_2}(ll_{j_2}, h_0^a, h_1^a, h_0^b, h_1^b, mode)$ 
8:      $ll_{j_1j_2}$ ,  $lh_{j_1j_2}$ ,  $hl_{j_1j_2}$ ,  $hh_{j_1j_2} \leftarrow \text{DWT}_{j_1j_2}(ll_{j_1j_2}, h_0^b, h_1^b, h_0^b, h_1^b, mode)$ 
9:      $x_{1a} \leftarrow (lh_r - lh_{j_1j_2}) + j(lh_{j_1} + lh_{j_2})$ 
10:     $x_{1b} \leftarrow (lh_r + lh_{j_1j_2}) + j(-lh_{j_1} + lh_{j_2})$ 
11:     $x_{2a} \leftarrow (hl_r - hl_{j_1j_2}) + j(hl_{j_1} + hl_{j_2})$ 
12:     $x_{2b} \leftarrow (hl_r + hl_{j_1j_2}) + j(-hl_{j_1} + hl_{j_2})$ 
13:     $x_{3a} \leftarrow (hh_r - hh_{j_1j_2}) + j(hh_{j_1} + hh_{j_2})$ 
14:     $x_{3b} \leftarrow (hh_r + hh_{j_1j_2}) + j(-hh_{j_1} + hh_{j_2})$ 
15:     $yh[j] \leftarrow (x_{1b}, x_{3b}, x_{2b}, x_{2a}, x_{3a}, x_{1a})$ 
16:   end for
17:    $yl \leftarrow \text{interleave}(ll_r, ll_{j_1}, ll_{j_2}, ll_{j_1j_2})$ 
18:   return  $yl$ ,  $yh$ 
19: end function

```

case and $j_1 = -j_2 = -j$ in the other case, producing the two outputs:

$$(a - d) + (b + c)j \quad (3.5.1)$$

$$(a + d) + (-b + c)j \quad (3.5.2)$$

corresponding to the first and second quadrant directional filters respectively. The sum and difference blocks Σ/Δ in Figure 3.2 do this operation [93].

The four lowpass coefficients from each scale are used for the next scale DWTs. At the final scale, they are interleaved to get four times the expected lowpass output area expected from a single decimated DWT. See Algorithm 3.3 for the code on how to do the full DTcWT.

A requirement of the DTcWT is the need to use different filters for the first scale to all subsequent scales [18]. We have not shown this in Algorithm 3.3 for simplicity, but it would simply mean we would have to handle the $j = 1$ case separately.

We have moved the inverse DTcWT algorithm to Algorithm B.3. Note that for both the DTcWT and inverse DTcWT, we rely on autograd to calculate the backwards pass by calling our defined DWT.Forward and DWT.Backward methods.

Algorithm 3.4 Magnitude forward and backward steps

```

1: function MAG.FORWARD( $x, y$ )
2:    $r \leftarrow \sqrt{x^2 + y^2}$ 
3:    $\theta \leftarrow \arctan 2(y, x)$                                  $\triangleright \arctan 2$  handles  $x = 0$ 
4:   save  $\theta$ 
5:   return  $r$ 
6: end function

1: function MAG.BACKWARD( $\Delta r$ )
2:   load  $\theta$ 
3:    $\Delta x \leftarrow \Delta r \cos \theta$                              $\triangleright$  Reinsert phase
4:    $\Delta y \leftarrow \Delta r \sin \theta$                              $\triangleright$  Reinsert phase
5:   return  $\Delta x, \Delta y$ 
6: end function

```

3.6 The DT^CWT ScatterNet

3.6.1 The Magnitude Operation

Now that we have a forward and backward pass for the DT^CWT, the final missing piece is the magnitude operation. If $z = x + jy$, then:

$$r = |z| = \sqrt{x^2 + y^2} \quad (3.6.1)$$

This has two partial derivatives, $\frac{\partial r}{\partial x}, \frac{\partial r}{\partial y}$:

$$\frac{\partial r}{\partial x} = \frac{x}{\sqrt{x^2 + y^2}} = \frac{x}{r} \quad (3.6.2)$$

$$\frac{\partial r}{\partial y} = \frac{y}{\sqrt{x^2 + y^2}} = \frac{y}{r} \quad (3.6.3)$$

Given an input gradient, Δr , the passthrough gradient is:

$$\Delta z = \Delta r \frac{\partial r}{\partial x} + j \Delta r \frac{\partial r}{\partial y} \quad (3.6.4)$$

$$= \Delta r \frac{x}{r} + j \Delta r \frac{y}{r} \quad (3.6.5)$$

$$= \Delta r e^{j\theta} \quad (3.6.6)$$

where $\theta = \arctan \frac{y}{x}$. This has a nice interpretation to it as well, as the backwards pass is simply reinserting the discarded phase information. The pseudocode for this operation is shown in [Algorithm 3.4](#). These partial derivatives are restricted to be in the range $[-1, 1]$

but have a singularity at the origin. In particular:

$$\lim_{x \rightarrow 0^-, y \rightarrow 0} \frac{\partial r}{\partial x} = -1 \quad (3.6.7)$$

$$\lim_{x \rightarrow 0^+, y \rightarrow 0} \frac{\partial r}{\partial x} = +1 \quad (3.6.8)$$

$$\lim_{x \rightarrow 0, y \rightarrow 0^-} \frac{\partial r}{\partial y} = -1 \quad (3.6.9)$$

$$\lim_{x \rightarrow 0, y \rightarrow 0^+} \frac{\partial r}{\partial y} = +1 \quad (3.6.10)$$

Rather than using the subgradient method [127] which is commonly used to handle the magnitude operation, we propose to smooth the magnitude operator:

$$r_s = \sqrt{x^2 + y^2 + b^2} - b \quad (3.6.11)$$

This keeps the magnitude near zero for small x, y but does slightly shrink larger values, however our gradient is smoother and we no longer have to worry about dividing by zero issues when calculating gradients. We can choose the size of b as a hyperparameter in optimization, which we revisit in [subsection 3.6.3](#). The partial derivatives now become:

$$\frac{\partial r_s}{\partial x} = \frac{x}{\sqrt{x^2 + y^2 + b^2}} = \frac{x}{r_s} \quad (3.6.12)$$

$$\frac{\partial r_s}{\partial y} = \frac{y}{\sqrt{x^2 + y^2 + b^2}} = \frac{y}{r_s} \quad (3.6.13)$$

There is a memory cost associated with this, as we will now need to save both $\frac{\partial r_s}{\partial x}$ and $\frac{\partial r_s}{\partial y}$ as opposed to saving only the phase. [Algorithm B.2](#) has the pseudocode for the smooth magnitude.

3.6.2 Putting it all Together

Now that we have the DTcWT and the magnitude operation, it is straightforward to get a DTcWT scattering layer, shown in [Algorithm 3.5](#).

For a second-order ScatterNet, the first C channels of Z from [Algorithm 3.5](#) are the S_0 coefficients, the next JKC are the S_1 coefficients and the final $\frac{1}{2}(J-1)JK^2C$ channels are the S_2 coefficients.

For ease in handling the different sample rates of the lowpass and the bandpass, we have averaged the lowpass over a 2×2 window and downsampled it by 2 in each direction. This slightly affects the higher-order scattering coefficients, as the true DTcWT needs the doubly sampled lowpass for the second scale, however we noticed little difference in performance from doing the true DTcWT and this slightly modified one

Algorithm 3.5 DTCWT ScatterNet Layer

```

1: function DTCWT_SCAT( $x$ ,  $J = 2$ ,  $M = 2$ )
2:    $Z \leftarrow x$ 
3:   for  $1 \leq m \leq M$  do
4:      $yl, yh \leftarrow \text{DTCWT}(Z, J = 1, mode = \text{'symmetric'})$ 
5:      $S \leftarrow \text{avg\_pool}(yl, 2)$ 
6:      $U \leftarrow \text{mag}(\text{Re}(yh), \text{Im}(yh))$ 
7:      $Z \leftarrow \text{concatenate}(S, U, axis = 1)$             $\triangleright$  stack 1 lowpass with 6 magnitudes
8:   end for
9:   if  $J > M$  then
10:     $Z \leftarrow \text{avg\_pool}(Z, 2^{J-M})$ 
11:   end if
12:   return  $Z$ 
13: end function

```

Table 3.1: Hyperparameter settings for the DT^CWT scatternet.

| Hyperparameter | Values |
|-------------------------|---------------------------------|
| Wavelet | near_sym_a 5,7 tap filters |
| | near_sym_b 13,19 tap filters |
| | near_sym_b_bp 13,19 tap filters |
| Padding Scheme | symmetric |
| | zero |
| Magnitude Smoothing b | 0 |
| | 1e-3 |
| | 1e-2 |
| | 1e-1 |

3.6.3 DT^CWT ScatterNet Hyperparameter Choice

Before comparing to the Morlet-based ScatterNet, we test different padding schemes, wavelet lengths and magnitude smoothing parameters (see (3.6.11)) for the DT^CWT ScatterNet. We test these over a grid of values described in Table 3.1.

The different wavelets have different lengths and hence different frequency responses. Additionally, the ‘near_sym_b_bp’ wavelet is a rotationally symmetric wavelet with diagonal passband brought in by a factor of $\sqrt{2}$, making it have similar $|\omega|$ to the horizontal and vertical bandpasses.

The results of these experiments are shown in Figure 3.3. The choice of options can have a significant impact on classification accuracy. Symmetric padding appears to work marginally better than zero padding. Surprisingly, the shorter filters (near_sym_a) fare better than their longer counterparts and bringing in the diagonal subbands (near_sym_b_bp) does not help.

Additionally, the smoothing bias may indeed be a good idea for the forward pass as well as aiding the backwards pass, so long as it is less than 0.1.

However, it is difficult to tell just how much ‘better’ the different filter lengths and padding schemes are, as there are many other factors to take into consideration when training a CNN. The standard deviation on the results from [Figure 3.3](#) were in the range from 0.1 to 0.4% for the two CIFAR datasets, which is smaller than the 2% range seen from the best option to the worst. However, for the larger Tiny ImageNet dataset the standard deviations were in the range from 0.2 to 0.6%, which is comparable to the 1% difference between the ‘best’ and ‘worst’ configurations.

We proceed by using the ‘near_sym_a’ filters, with symmetric padding and a smoothing factor of 0.01, but leave all configurations in our code in [\[28\]](#) and note that more experimentation needs to be done in analysing and comparing the choice of options.

3.7 Comparisons

Now that we have the ability to do a DT^CWT based ScatterNet, how does this compare with the original Matlab implementation [\[25\]](#) and the newly developed KyMatIO [\[119\]](#)? [Table 3.2](#) lists the different properties and options of the competing packages.

3.7.1 Speed

We test the speed of the various packages on our reference architecture (see [appendix A](#)) with a moderately large input with $128 \times 3 \times 256 \times 256$ pixels. The CPU experiments used all cores available, whereas the GPU experiments ran on a single GPU. We include two permutations of our proposed ScatterNet with different length filters and different padding schemes. Type A uses the long ‘near_sym_b’ filters and has symmetric padding, and Type B uses shorter ‘near_sym_a’ filters and the faster zero padding scheme. We compare it to a Morlet based implementation with $K = 6$ orientations (the same number of orientations as in the DT^CWT).

See [Table 3.3](#) for the execution time results. Type A is ~ 7 times faster than the Fourier-based KyMatIO on GPUs and Type B has a ~ 14 times speedup over the Morlet backend.

Additionally, when compared with version 0.1.0 of KyMatIO, the memory footprint of the DT^CWT based implementation is only 2% of KyMatIO’s, highlighting the importance of being explicit in not saving unnecessary information for backpropagation.

3.7.2 Performance

To confirm that changing the ScatterNet core has not impeded the performance of the ScatterNet as a feature extractor, we build a simple Hybrid ScatterNet, similar to [\[39\]](#), [\[40\]](#). Our net has a second order scattering transform before four convolutional layers. See

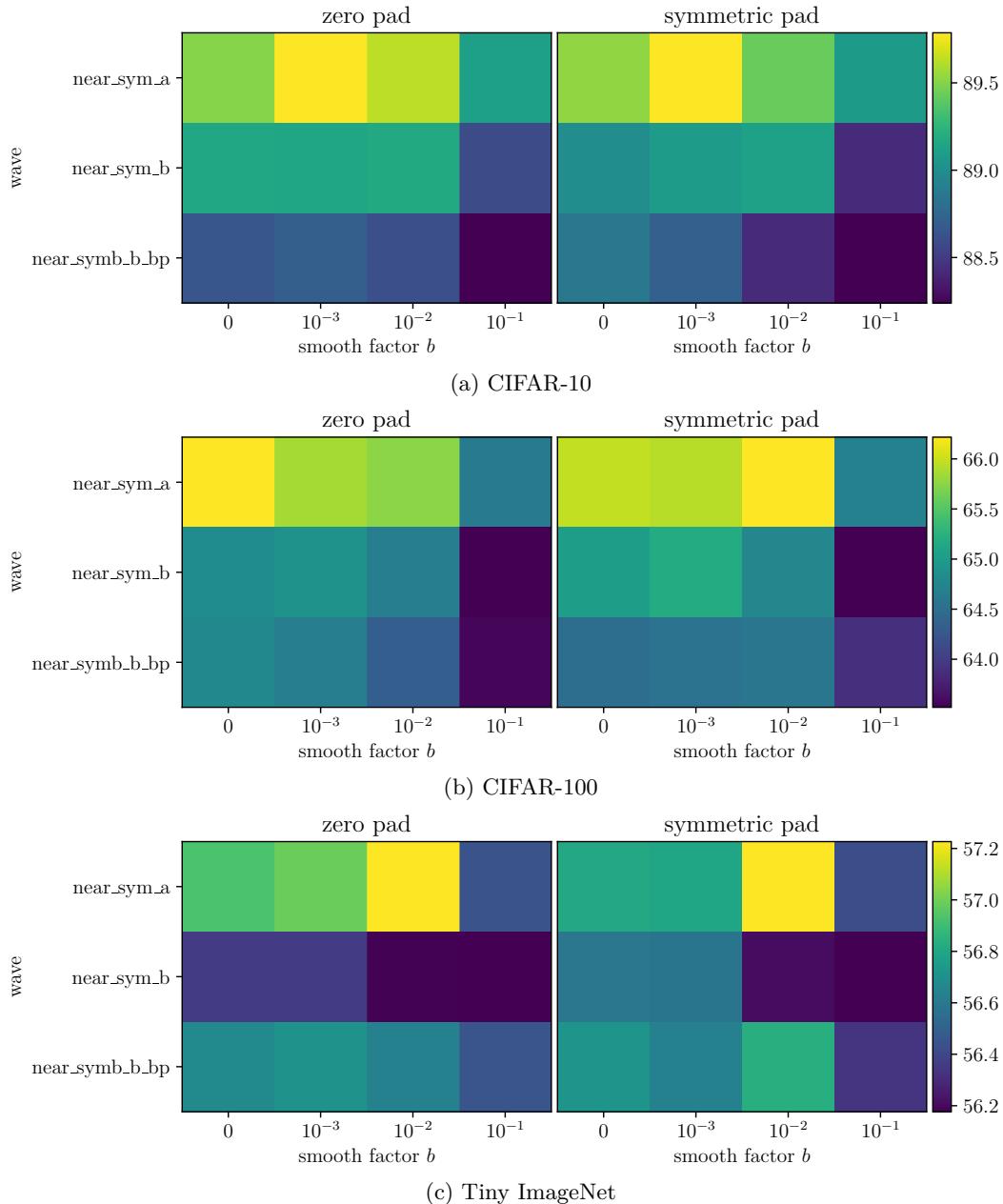


Figure 3.3: Hyperparameter results for the DTCWT scatternet on various datasets. Image showing relative top-1 accuracies (in %) on the given datasets using architecture described in [Table 3.1](#). Each subfigure is a new dataset and therefore has a new colour range (shown on the right). Results are averaged over 3 runs from different starting positions. The choice of options can have a very large impact on the classification accuracy. Symmetric padding is marginally better than zero padding. Surprisingly, the shorter filter (near_sym_a) fares better than its longer counterparts, and bringing in the diagonal subbands (near_symb_b_bp) does not help. Additionally, the smoothing bias may indeed be a good idea for the forward pass as well as aiding the backwards pass, so long as it is less than 0.1.

Table 3.2: **Comparison of properties of different ScatterNet packages.** In particular the wavelet backend used, the number of orientations available, the available boundary extension methods, whether it has GPU support and whether it supports backpropagation.

| Package | Backend | Orientations | Boundary Ext. | GPU | Backprop |
|------------------|------------------------|--------------|---------------|-----|----------|
| ScatNetLight[25] | FFT-based | Flexible | Periodic | No | No |
| KyMatIO[119] | FFT-based | Flexible | Periodic | Yes | Yes |
| DTCWT Scat | Separable filter banks | 6 | Flexible | Yes | Yes |

Table 3.3: **Comparison of execution time for the forward and backward passes of the competing ScatterNet Implementations.** Tests were run on the reference architecture described in [appendix A](#). The input for these experiments is a batch of images of size $128 \times 3 \times 256 \times 256$ in 4 byte floating precision. We list two different types of options for our ScatterNet. Type A uses 16 tap filters and has symmetric padding, whereas type B uses 6 tap filters and uses zero padding at the image boundaries. Values are given to 2 significant figures, averaged over 5 runs.

| Package | CPU | | GPU | |
|-------------------|----------|---------|---------|---------|
| | Fwd (s) | Bwd (s) | Fwd (s) | Bwd (s) |
| ScatNetLight[25] | > 200.00 | n/a | n/a | n/a |
| KyMatIO[119] | 95.0 | 130.0 | 1.44 | 2.5 |
| DTCWT Scat Type A | 3.3 | 3.6 | 0.21 | 0.27 |
| DTCWT Scat Type B | 2.8 | 3.2 | 0.10 | 0.16 |

[Table 3.4](#) for the network layout. We use the optimal hyperparameter choices from the previous section, and compare these to Morlet based ScatterNet with 6 and 8 orientations.

We run tests on the following datasets:

- CIFAR-10: 10 classes, 5000 images per class, 32×32 pixels per image.
- CIFAR-100: 100 classes, 500 images per class, 32×32 pixels per image.
- Tiny ImageNet[78]: 200 classes, 500 images per class, 64×64 pixels per image.

The images in Tiny ImageNet are four times the size, so the output after scattering is 16×16 . We add a max pooling layer after conv4, followed by two more convolutional layers conv5 and conv6, before average pooling.

These networks are optimized with SGD with momentum. The initial learning rate is 0.5, momentum is 0.85, batch size $N = 128$ and weight decay is 10^{-4} . For CIFAR-10/CIFAR-100 we scale the learning rate by a factor of 0.2 after 60, 80 and 100 epochs, training for 120

epochs in total. For Tiny ImageNet, the rate change is at 18, 30 and 40 epochs (training for 45 in total). Our experiment code is available at [31].

The results of this experiment are shown in Table 3.5. It is promising to see that the DT \mathbb{C} WT based ScatterNet has not only sped up, but slightly improved upon the Morlet based ScatterNet as a frontend. Interestingly, both with Morlet and DT \mathbb{C} WT wavelets, 6 orientations performed better than 8, despite having fewer parameters in conv1.

3.8 Conclusion

In this chapter we have proposed changing the wavelet used in Scattering transforms from the Morlet wavelet to the spatially-separable, biorthogonal DT \mathbb{C} WT. This was originally inspired by the need to speed up the slow Matlab scattering package, as well as to provide GPU accelerated code that could do wavelet transforms as part of a deep learning program.

We have derived the forward and backpropagation functions necessary to do fast and memory-efficient DWTs, DT \mathbb{C} WTs, and Scattering based on the DT \mathbb{C} WT, and have made this code publically available at [28]. We hope that this will reduce some of the barriers we initially faced in using wavelets and Scattering in deep learning.

In parallel with our efforts, the original ScatterNet authors rewrote their package to do faster Scattering. In theory, a spatially separable wavelet transform acting on N pixels has order $\mathcal{O}(N)$ whereas an FFT based implementation has order $\mathcal{O}(N \log N)$. We have shown experimentally that on modern GPUs, the difference is far larger than this, with the DT \mathbb{C} WT backend an order of magnitude faster than Fourier-based Morlet implementation [119].

We have experimentally verified that using a different complex wavelet slightly improves the performance of a ScatterNet front end to a CNN.

Table 3.4: **Hybrid architectures for performance comparison.** Comparison of Morlet based ScatterNets (Morlet6 and Morlet8) to the DT \mathbb{C} WT based ScatterNet on CIFAR. The output after scattering has $3(K+1)^2$ channels (243 for 8 orientations or 147 for 6 orientations) of spatial size 8×8 . This is passed to 4 convolutional layers of width $C = 192$ before being average pooled and fed to a single fully connected classifier. $N_c = 10$ for CIFAR-10 and 100 for CIFAR-100. In the DT \mathbb{C} WT architecture, we test different padding schemes and wavelet lengths.

| Morlet8 | Morlet6 | DTCWT |
|---|---|---|
| Scat $J = 2, K = 8, m = 2$ $y \in \mathbb{R}^{243 \times 8 \times 8}$ | Scat $J = 2, K = 6, m = 2$ $y \in \mathbb{R}^{147 \times 8 \times 8}$ | Scat $J = 2, K = 6, m = 2$ $y \in \mathbb{R}^{147 \times 8 \times 8}$ |
| conv1, $w \in \mathbb{R}^{C \times 243 \times 3 \times 3}$ | | conv1, $w \in \mathbb{R}^{C \times 147 \times 3 \times 3}$ |
| | conv2, $w \in \mathbb{R}^{C \times C \times 3 \times 3}$ | |
| | conv3, $w \in \mathbb{R}^{2C \times C \times 3 \times 3}$ | |
| | conv4, $w \in \mathbb{R}^{2C \times 2C \times 3 \times 3}$ | |
| | avg pool, 8×8 | |
| | fc, $w \in \mathbb{R}^{2C \times N_c}$ | |

Table 3.5: **Performance comparison for a DT \mathbb{C} WT based ScatterNet vs Morlet based ScatterNet.** We report top-1 classification accuracy for the 3 listed datasets as well as training time for each model in hours.

| Type | CIFAR-10 | | CIFAR-100 | | Tiny ImgNet | |
|--------------------|----------|----------|-----------|----------|-------------|----------|
| | Acc. (%) | Time (h) | Acc. (%) | Time (h) | Acc. (%) | Time (h) |
| Morlet8 | 88.6 | 3.4 | 65.3 | 3.4 | 57.6 | 5.6 |
| Morlet6 | 89.1 | 2.4 | 65.7 | 2.4 | 57.5 | 4.4 |
| DT \mathbb{C} WT | 89.8 | 1.1 | 66.2 | 1.1 | 57.3 | 2.7 |

Chapter 4

Visualizing and Improving Scattering Networks

Despite the success of CNNs, deep nets are often criticized for being ‘black box’ methods. Once you train a CNN, you can view the first layer of filters quite easily (see [Figure 1.2a](#)) as they exist in RGB space. Beyond that things get trickier, as the filters have a third, *channel* dimension, typically much larger than its two spatial dimensions. Representing these dimensions with different colours becomes tricky and uninformative, so we must do something else.

A recent paper titled ‘Why Should I Trust You?’ [\[128\]](#) explored the consequences of interpretability on human trust for machine learning models. Unsurprisingly, a model with an interpretable methodology was trusted more than those which did not have one, even if it had a lower prediction accuracy on the test set. To build trust and to aid training, we need to probe these networks and visualize how and why they are making decisions.

Some good work has been done in this area. In particular, Zeiler and Fergus [\[26\]](#) design a DeConvNet to visualize what input patterns a filter in a given layer is mostly highly activated by. In [\[129\]](#), Mahendran and Vedaldi learn to invert representations by updating a noisy input via GD until its latent feature vector matches a desired target. Simonyan, Vedaldi, and Zisserman [\[130\]](#) develop saliency maps by projecting gradients back to the input space and measuring where they have largest magnitude.

We introduced ScatterNets in [section 2.7](#) and looked at making them faster in [chapter 3](#). They have been one of the main successes in applying wavelets to deep learning systems, and are particularly inspiring due to their well-defined properties. They are typically used as unsupervised feature extractors [\[23\], \[25\], \[36\], \[131\]](#) and can outperform CNNs for classification tasks with reduced training set sizes, e.g. in CIFAR-10 and CIFAR-100 (Table 6 from [\[39\]](#) and Table 4 from [\[131\]](#)). They are also near state-of-the-art for Texture Discrimination tasks (Tables 1–3 from [\[99\]](#)). Despite this, there still exists a considerable gap

between Scatternets and CNNs on challenges like CIFAR-10 with the full training set (83% vs. > 90%). Even considering the benefits of ScatterNets, this gap must be addressed.

While ScatterNets have good theoretical foundations and properties [17], it is difficult to understand the second-order scattering. In particular, how useful are these coefficients for training and how similar are the scattered features to a modern state of the art convolutional network? To answer these questions, this chapter interrogates ScatterNet frontends. Taking inspiration from the work done for CNNs, we build a DeScatterNet to visualize what the second-order features are. We also heuristically probe a trained hybrid network (ScatterNet front end + CNN backend) and quantify the importance of the individual features.

4.1 Chapter Layout

We first redefine the operations that form a ScatterNet in section 4.3 before introducing our DeScatterNet in section 4.4, and show how we can use it to examine the layers of ScatterNets (using a similar technique to the CNN visualization in [26]). We use this analysis tool to highlight what patterns a ScatterNet is sensitive to (section 4.5), showing that they are very different from what their CNN counterparts are sensitive to, and possibly less useful for discriminative tasks.

We then measure the ScatterNet channel saliency by performing occlusion tests on a trained hybrid network ScatterNet-CNN, iteratively switching off individual Scattering channels and measuring the effect this has on the validation accuracy in section 4.6. The results from the occlusion tests strengthen the idea that some of the ScatterNet patterns may not be well suited for deep learning systems.

We use these observations to propose an architectural change to ScatterNets, which have not changed much since their inception in [17], and show that it is possible to get visually more appealing shapes by filtering across the orientations of the ScatterNet. We present this in section 4.7.

4.2 Related Work

Zeiler, Taylor, and Fergus first attempted to use ‘deconvolution’ to improve their learning [132], then later for purely visualization purposes [26]. Their method involves monitoring network nodes, seeing what input image causes the largest activity and mapping activations at different layers of the network back to the pixel space using meta-information from these images.

Figure 4.1 shows the block diagram for how deconvolution is done. They invert a convolutional layer by taking the 2D transpose of each slice of the filter. Inverting a ReLU is done by simply applying a ReLU again (ensuring only positive values can flow back through the network). Inverting a max pooling step is a little trickier, as max pooling is quite a lossy

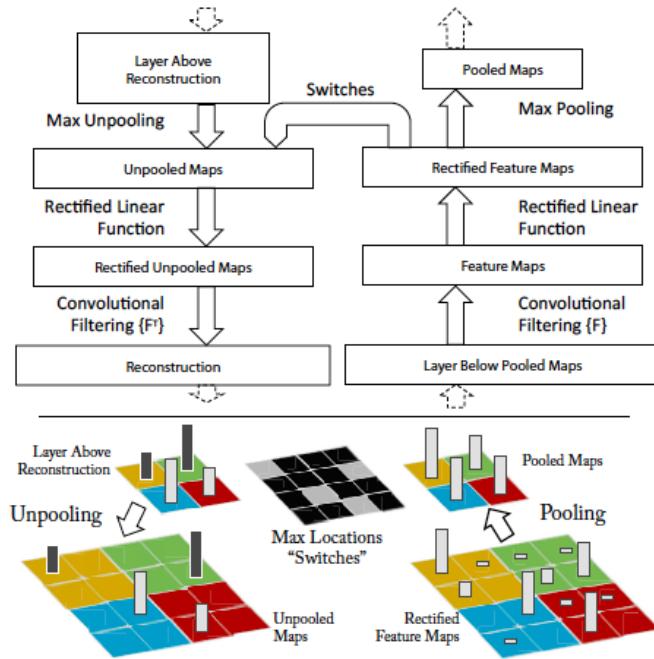


Figure 4.1: Deconvolution Network Block Diagram. Switches are saved alongside pooled features, and the filters used for deconvolution are the transpose of the filters used for a forward pass. Taken from [26].

operation. Zeiler, Taylor, and Fergus get around this by saving extra information on the forward pass of the model — switches that store the location of the input that caused the maximum value. This way, on the backwards pass, it is trivial to store activations to the right position in the larger feature map. The positions that did not contribute to the max pooling operation remain as zero on the backwards pass. This is shown in the bottom half of Figure 4.1.

Mahendran and Vedaldi take a slightly different route on deconvolution networks [129]. They do not store this extra information but instead define a cost function to maximize. This results in visualization images that look very surreal, and can be quite different from the input.

In [68], the authors design an *all convolutional* model, where they replace the max pooling layers commonly seen in CNNs with a convolutional block with stride 2, i.e. a convolution followed by decimation. They did this by first adding an extra convolutional layer before the max pooling layers, then taking away the max pooling and adding decimation after convolution, noting that removing the max pooling had little effect.

The benefit of this is that they can now reconstruct images as Zeiler and Fergus did, but without having to save switches from the max pooling operation. Additionally, they modify

the handling of the ReLU in the backwards pass to combine the regular backpropagation action and the ReLU action from deconvolution. They call this ‘guided backprop’.

Another interrogation tool commonly used is occlusion or perturbation. In [26], Zeiler and Fergus occlude regions in the input image and measure the impact this has on classification score. In [133], Fong and Vedaldi use gradients to find the minimal mask to apply to the input image that causes misclassification.

4.3 The Scattering Transform

While we have introduced the scattering transform before, we clarify the format we use for this chapter’s analysis.

We use the DTCWT based ScatterNet introduced in the previous chapter – [Algorithm 3.5](#) as a front end, with $K = 6$ orientations, $J = 2$ scales and $M = 2$ orders.

Consider a single-channel input signal $x(\mathbf{u})$, $\mathbf{u} \in \mathbb{R}^2$. The zeroth-order scatter coefficient is the lowpass output of a J level filter bank:

$$S_0 x(\mathbf{u}) \triangleq x(\mathbf{u}) * \phi_J(\mathbf{u}) \quad (4.3.1)$$

This is approximately invariant to translations of up to 2^J pixels¹. In exchange for gaining invariance, the S_0 coefficients have lost information (contained in the rest of the frequency space). The remaining energy of x is contained within the first-order *wavelet* coefficients:

$$W_1 x(\lambda_1, \mathbf{u}) \triangleq x * \psi_{\lambda_1} \quad (4.3.2)$$

for $\lambda_1 = (j_1, \theta_1)$, $j_1 \in \{1, 2\}$, $\theta_1 = \frac{\pi+2k\pi}{12}$ with $k \in \{0, 1, \dots, 5\}$.

Taking the magnitude of W_1 gives us the first-order *propagated* signals:

$$U_1 x(\lambda_1, \mathbf{u}) \triangleq |x * \psi_{\lambda_1}| = \sqrt{(x * \psi_{\lambda_1}^r)^2 + (x * \psi_{\lambda_1}^i)^2} \quad (4.3.3)$$

The first-order scattering coefficient make U_1 invariant up to the coarsest scale J by averaging it:

$$S_1 x(\lambda_1, \mathbf{u}) \triangleq |x * \psi_{\lambda_1}| * \phi_J \quad (4.3.4)$$

This has $KJ = 6 \times 2 = 12$ output channels for each input channel. Later in this chapter we will want to distinguish between the first and second scale coefficients of the S_1 terms, which we will do by moving the j index to a superscript. I.e., S_1^1 and S_1^2 refer to the set of 6 S_1 terms at the first and second scales.

¹From here on, we drop the \mathbf{u} notation when indexing x , for clarity.

The second-order scattering coefficients are defined only on paths of decreasing frequency (i.e. $J_2 < J_1$) [23]

$$S_2x(\lambda_1, \lambda_2, \mathbf{u}) \triangleq |U_1x * \psi_{\lambda_2}| * \phi_J \quad (4.3.5)$$

Previous work shows that for natural images we get diminishing returns after $m = 2$. Our output is then a stack of these 3 outputs:

$$Sx = \{S_0x, S_1x, S_2x\} \quad (4.3.6)$$

with $1 + 12 + 36 = 49$ channels per input channel.

4.3.1 Scattering Colour Images

A wavelet transform like the DTcWT accepts single-channel input yet we often work on RGB images. This leaves us with a choice. We can either:

1. Apply the wavelet transform (and the subsequent scattering operations) on each channel independently. This would triple the output size to $3C$.
2. Define a frequency threshold below which we keep colour information, and above which, we combine the three channels.

The second option uses the well known fact that the human eye is far less sensitive to higher spatial frequencies in colour channels than in luminance channels. This also fits in with the first layer filters seen in the well known Convolutional Neural Network, AlexNet. Roughly one half of the filters were low frequency colour ‘blobs’, while the other half were higher frequency, greyscale, oriented wavelets.

For this reason, we choose the second option for the architecture described in this chapter. We keep the 3 colour channels in our S_0 coefficients but work only on greyscale for high orders (the S_0 coefficients are the lowpass bands of a J-scale wavelet transform, so we have effectively chosen a colour cut-off frequency of $2^{-J} \frac{f_s}{2}$).

We combine the three channels by modifying our magnitude operation from (3.6.11) to now be:

$$r_s = \sqrt{x_r^2 + y_r^2 + x_g^2 + y_g^2 + x_b^2 + y_b^2 + b^2 - b} \quad (4.3.7)$$

Where x_r, x_g, x_b are the real parts of the wavelet response for the red, green and blue channels, and y is the corresponding imaginary part. This only affects the S_1 coefficients and the S_2 coefficients then are calculated as per (4.3.5).

An alternative to (4.3.7) is to combine the colours *before* scattering into a luminance channel. However, we choose to use (4.3.7) instead as this has the ability to detect colour edges with constant luminance.

With $J = 2$ the resulting scattering output now has $3 + 12 + 36 = 51$ channels at $1/16$ the spatial input size.

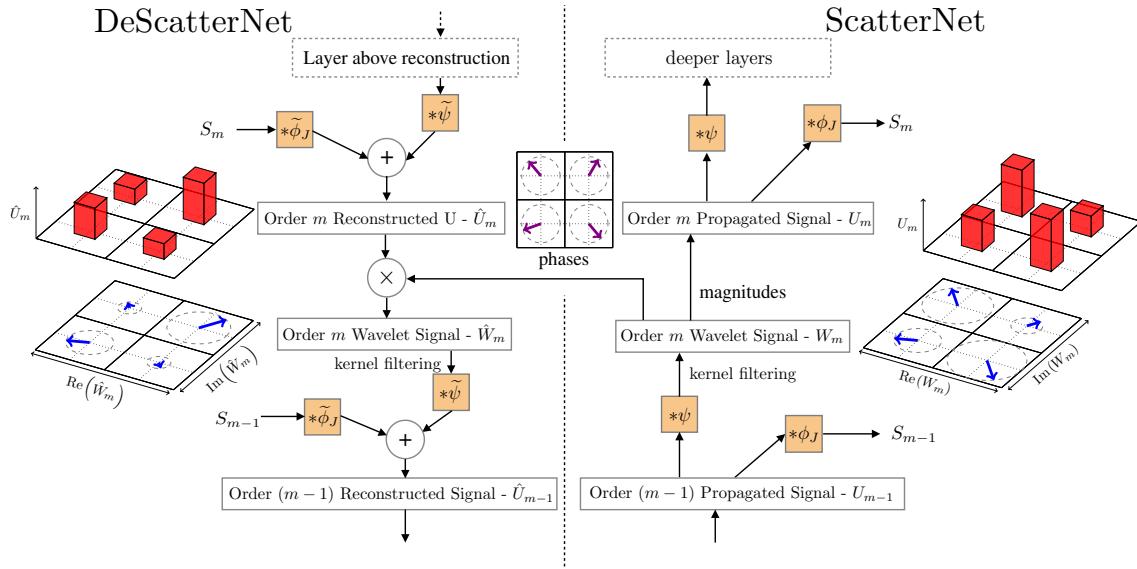


Figure 4.2: **The Descattering Network.** Comprises a DeScattering layer (left) attached to a Scattering layer (right). We are using the same convention as [26] Figure 1 - i.e., the input signal starts in the bottom right-hand corner, passes forwards through the ScatterNet (up the right half), and then is reconstructed in the DeScatterNet (downwards on the left half). The DeScattering layer will reconstruct an approximate version of the previous order’s propagated signal. The 2×2 grids shown around the image are either Argand diagrams representing the magnitude and phase of small regions of *complex* (De)ScatterNet coefficients or bar charts showing the magnitude of the *real* (De)ScatterNet coefficients (after applying the modulus nonlinearity). For reconstruction, we need to save the discarded phase information and reintroduce it by multiplying it with the reconstructed magnitudes.

4.4 The Inverse Scatter Network

We now introduce our inverse scattering network. This allows us to back-project scattering coefficients to the image space; it is inspired by the DeconvNet used by Zeiler and Fergus in [26] to look into the deeper layers of CNNs. Like the DeConvNet, the inverse scattering network is similar to backpropagating a single strong activation (rather than usual gradient terms).

We emphasize that instead of thinking about perfectly reconstructing x from $S \in \mathbb{R}^{C \times H' \times W'}$, we want to see what signal/pattern in the input image caused a large activation in each channel. This can then give us a good idea of what each output channel is sensitive to, or what the ScatterNet is ‘extracting’ from the input. Note that we do not use any of the log normalization layers described in [25], [131].

4.4.1 Inverting the Low-Pass Filtering

Going from the U coefficients to the S coefficients in the forward pass involves convolving by a low pass filter ϕ_J , possibly followed by decimation to make the output $(H \times 2^{-J}) \times (W \times 2^{-J})$.

ϕ_J is a purely real filter, and we can ‘invert’ this operation by interpolating S to the same spatial size as U and convolving with the mirror image of ϕ_J , $\tilde{\phi}_J$ (this is equivalent to the transpose convolution described in [26]).

$$\hat{S}_m x = (S_m x) * \tilde{\phi}_J \quad (4.4.1)$$

We note that interpolation usually involves lowpass smoothing of the signal, so this can all be one operation.

4.4.2 Inverting the Magnitude Operation

In the same vein as [26], we face a difficult task in inverting the nonlinearity in our system. We lend inspiration from the switches introduced in the DeconvNet; the switches in a DeconvNet save the location of maximal activations so that on the backwards pass activation layers could be unpooled trivially. We do an equivalent operation by saving the *phase* of the complex activations. On the backwards pass we reinsert the phase to give our recovered W^2

$$\hat{W}_m x = (\hat{U}_m x) e^{j\theta_m} \quad (4.4.2)$$

4.4.3 Inverting the Wavelet Decomposition

Using the DT \mathbb{C} WT makes inverting the wavelet transform simple, as we can simply feed the coefficients through the synthesis filter banks to regenerate the signal. For complex ψ , this is convolving with the conjugate transpose $\tilde{\psi}$:

$$\hat{U}_{m-1} x = \hat{S}_{m-1} x + \hat{W}_m x \quad (4.4.3)$$

$$= (S_{m-1} x) * \tilde{\phi}_J + \sum_{\lambda_m} (\hat{U}_m x) e^{j\theta_m} * \tilde{\psi}_{\lambda_m} \quad (4.4.4)$$

4.4.4 The DT \mathbb{C} WT ScatterNet

The combination of the above three stages can be repeated for higher orders. The resulting DeScatterNet is shown in Figure 4.2.

For the DT \mathbb{C} WT ScatterNet (from Algorithm 3.5), this is the same as finding the *gradient* from the corresponding channels in Z .

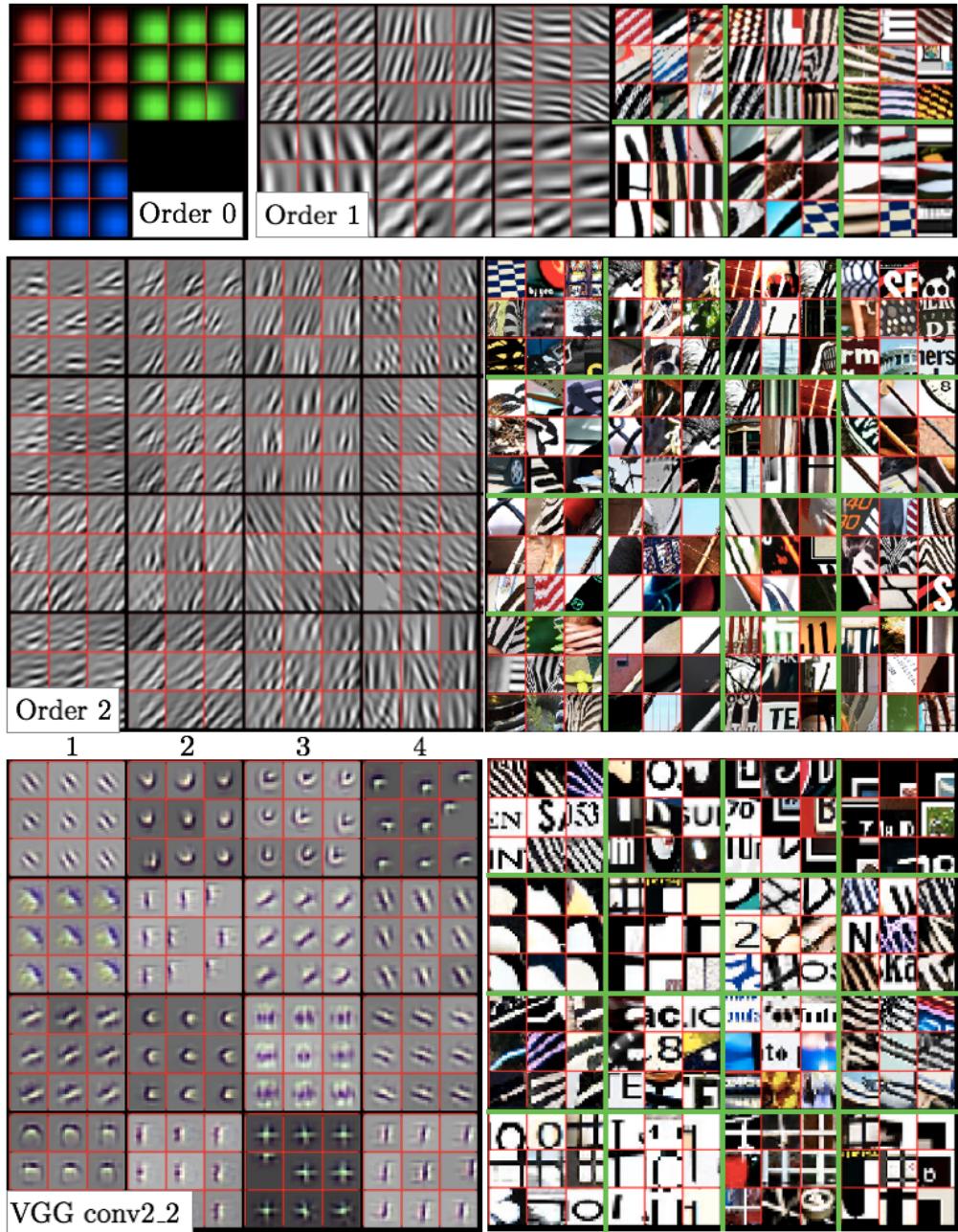


Figure 4.3: Comparison of Scattering to Convolutional features. Visualization of a random subset of features from S_0 (all 3), S_1 (6 from the 12) and S_2 (16 from the 36) scattering outputs. We record the top 9 activations for the chosen features and project them back to the pixel space. We show them alongside the input image patches which caused the large activations. We also include reconstructions from layer conv2_2 of VGG Net [66] (a popular CNN, often used for feature extraction) for reference — here we display 16 of the 128 channels. The VGG reconstructions were made with a CNN DeconvNet based on [26]. Image best viewed digitally.

4.5 Visualization with Inverse Scattering

We scatter all of the images from ImageNet’s validation set and record the top 9 images which most highly activate each of the C channels in the ScatterNet. This is the *identification* phase (in which no inverse scattering is performed).

Then, in the *reconstruction* phase, we load in the $9 \times C$ images and scatter them one by one. We take the resulting 52 channel output vector and mask all but the largest value in the channel we are currently examining and all values in the other channels. This 1-sparse tensor is then presented to the inverse scattering network from [Figure 4.2](#) and projected back to the image space.

Some results of this are shown in [Figure 4.3](#) for the first and second-order coefficients. For a given output channel, we show the top 9 activations projected independently to pixel space. We also show the patch of pixels in the input image which cause this large output. As there are 12 S_1 coefficients, we randomly choose 3 orientations from S_1^1 and 3 S_1^2 . Similarly, there are 36 S_2 coefficients, so we randomly choose 16 of these.

The order 0 and order 1 scattering (labelled with ‘Order 1’ in [Figure 4.3](#)) coefficients look quite similar to the first layer filters from the well-known AlexNet CNN [8]. This is not too surprising, as the first-order scattering coefficients are simply a wavelet transform followed by average pooling. They are responding to images with strong edges aligned with the wavelet orientation.

The second-order coefficients (labelled with ‘Order 2’ in [Figure 4.3](#)) appear very similar to the order 1 coefficients at first glance. They too are sensitive to edge-like features, and some of them (e.g. third row, third column and fourth row, second column) are mostly just that. These are features that have the same oriented wavelet applied at both the first and second-order ($\theta_1 = \theta_2$). Others, such as the nine in the top left square (first row, first column), and top right square (first row, fourth column) are more sensitive to checker-board like patterns. These are activations where the orientation of the wavelet for the first and second-order scattering were far from each other (15° and 105° for the first row, first column and 105° and 45° for the first row, fourth column).

For comparison, we include reconstructions from the second layer of the well-known VGG CNN (labelled with ‘VGG conv2_2’, in [Figure 4.3](#)). These were made with a DeconvNet, following the same method as [26]. Note that while some of the features are edge-like, we also see higher order shapes like corners, crosses and curves.

These reconstructions show that the features extracted from ScatterNets vary significantly from those learned in CNNs after the first-order. In many respects, the features extracted from a CNN like VGGNet look preferable for use as part of a classification system.

²We note that this is equivalent to finding the gradient through the magnitude operation, just as the ‘switches’ from [26] are equivalent to taking the gradient of the max pooling layer.

4.6 Channel Saliency

To get another heuristic on the importance of the ScatterNet channels, let us examine the effect on inference scores observed when zeroing out Scattering channels. Zeiler and Fergus [26] and Zhou, Khosla, Lapedriza, *et al.* [134] have done similar studies but over patches of the input image.

We must be careful to occlude with a sensible mask, the S_0x , S_1x and S_2x all have very different probability densities. [26] occlude with a patch of grey values whereas [134] use random values. Assuming $x \sim \mathcal{N}(0, \sigma^2 I)$ (already a fairly weak assumption), the pdf of S_0x will also be a zero-mean gaussian. However, the distributions of S_1x and S_2x are more complex - the real and imaginary parts of the DT \mathbb{C} WT are sparse but are strongly correlated in energy. Further, after the modulus operation, there is a strong positive bias to all the pdfs until the signal passes through another bandpass filter. Choosing a sensible random mask is therefore difficult, so we instead use a constant mask. Analysis of the datasets showed that zero is very close to the maximum likelihood value for each channel so we occlude channels by simply setting them to zero at every spatial location.

4.6.1 Experiment Setup

We take a network similar to the one from [Table 3.4](#) (2 layers of ScatterNet followed by 4 convolutional layers). Unlike the previous chapter, we use the colour operation described in [subsection 4.3.1](#) so the scattering output has 51 output channels. Further, we set the first convolutional layer after the ScatterNet to have 100 channels for display purposes later.

We train this network on the same 3 datasets - CIFAR-10, CIFAR-100 and Tiny ImageNet, and report the drop in classification scores on the validation set after removing one channel at a time.

We additionally display the weight matrix for the first learned layer of the network trained on Tiny ImageNet. As the scattering output has 51 channels and the first layer of the CNN has 100 channels, this weight matrix has shape: $w \in \mathbb{R}^{100 \times 51 \times 3 \times 3}$ (it is a 3×3 convolution over the 51 channels with 100 filters). This can give us a second perspective on the channel importance by looking at the relative weights of the ScatterNet channels across the 100 filters. We define:

$$A_{c,f}^{rms} = \sqrt{\frac{\sum_{i,j} w[f, c, i, j]^2}{\sum_f \sum_{i,j} w[f, c, i, j]^2}} \quad (4.6.1)$$

This gives us a matrix A^{rms} (for root mean squared) which has columns of unit energy representing the different output channels after conv1. The row values then show how much each scattering channel contributes to each output channel. This is shown in [Figure 4.5](#).

4.6.2 Results

First, we look at Tiny ImageNet in [Figure 4.4](#). Note that when any of the S_0 channels are removed, the validation accuracy drops sharply for all 3 colours. A similar result happens when any of the S_1 channels are zeroed out.

For both the first and second scales of the first-order coefficients, S_1^1 and S_1^2 , there are two channels that seem less important - the second and fifth channels, corresponding to the 45° and 135° wavelets. Often the high-high portion of the first scale coefficients are considered mostly noise, but this does not explain why the 45° and 135° channels for the second scale coefficients are also less important. A possible interesting conclusion to be drawn from this is that the dataset does not have as many important diagonal edges in it as horizontal and vertical edges, and the network has learned this difference in importance.

To test this, we retrain the network but this time rotate the input images randomly 30° clockwise or anti-clockwise in both training and validation. We then rerun the occlusion experiment for all channels and plot the resulting changes in [Figure 4.4b](#). Interestingly, for this network, the 45° and 135° wavelets for S_1^2 are now the most important of the 6, which validates our assumption. The corresponding wavelets for S_1^1 have become more important, but it is likely that they remain less salient because of the effects of the higher bandwidth for the diagonal wavelets.

Comparatively, the S_2 channels have little effect on the classification score when individually masked. The four largest drops in accuracy for S_2 are happening when $\theta_1 = \theta_2 \in \{15^\circ, 75^\circ, 105^\circ, 165^\circ\}$. When we drop channels in S_2 with $\theta_1 \neq \theta_2$, the network performance is not affected very much. Recall that $\theta_1 \neq \theta_2$ corresponds to the ripple-like patterns in [Figure 4.3](#).

[Figure 4.5](#) shows the size of $A_{c,f}^{rms}$. The columns of the matrix all have unit-norm, so each entry represents how much relative energy comes from each scattering channel (brighter values indicating more energy). Looking across the rows we see how often a scattering output is used for the CNN next layer. Most of the filters are heavily dependent on S_0 , many are dependent on S_1 and only a few take information from S_2 .

We include the same occlusion results for the two CIFAR datasets in [Figure 4.6](#) for completeness, although the insight gained here is the same - the S_2 coefficients are the least important. One notable difference to [Figure 4.4](#) is in the S_1^2 coefficients which have reduced importance in CIFAR. As CIFAR has a smaller input spatial size than Tiny ImageNet this comes as no surprise.

4.7 Corners, Crosses and Curves

As a final part of this chapter, we would like to highlight some of the filters possible by making small modifications to the ScatterNet design. The visualizations shown here are

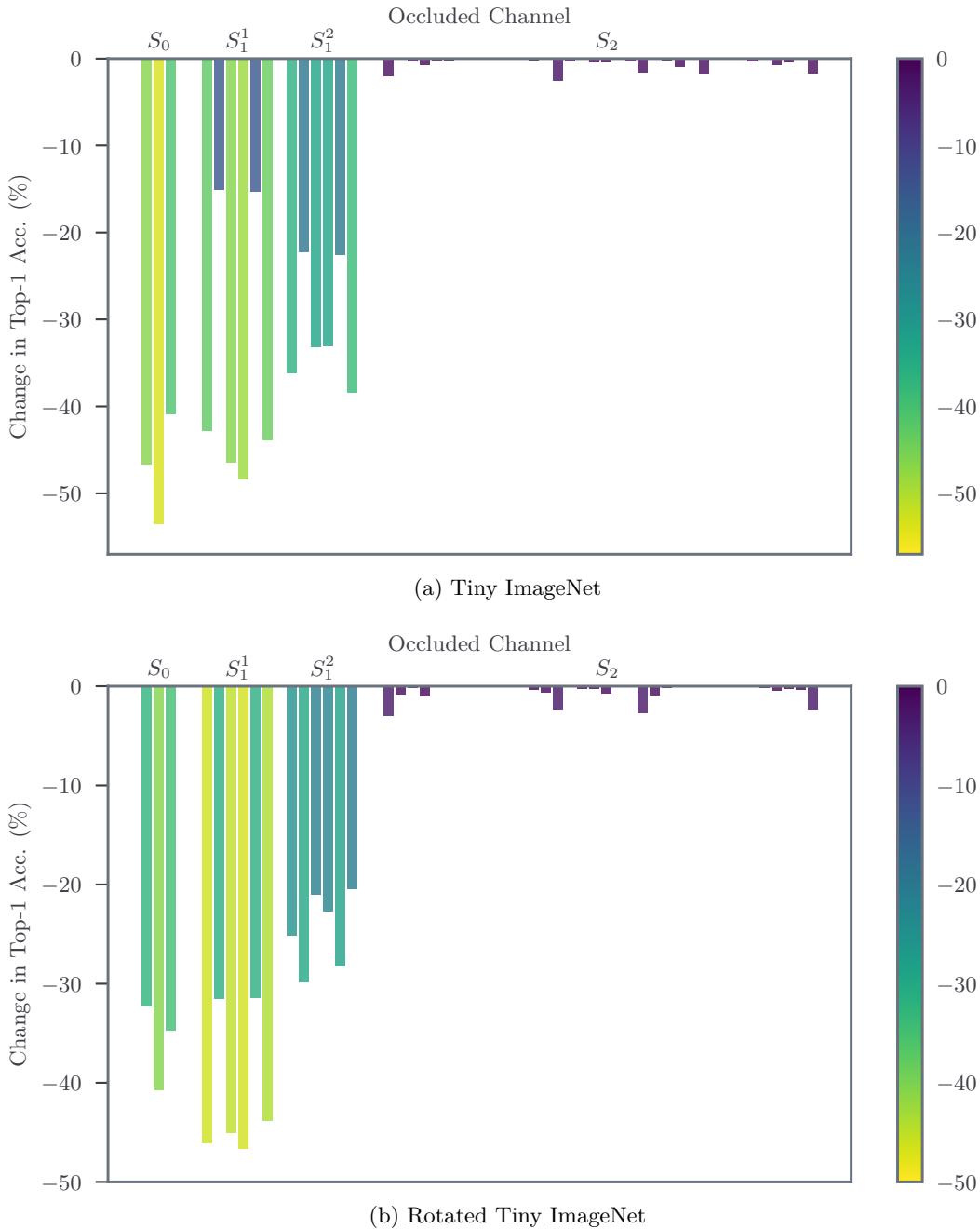


Figure 4.4: **Tiny ImageNet changes in accuracy from channel occlusion.** Numbers reported are the drop in final classification accuracy when a channel is set to zero. The bars are coloured relative to their magnitude to aid seeing the differences for the S_1 coefficients. (a) When any of the lowpass channels S_0 are removed, the classification accuracy drops sharply, note that the middle channel, corresponding to green, is unsurprisingly the most important of the three colours. The first scale, first-order scattering coefficients S_1^1 are slightly more important than the second scale coefficients. The 36 S_2 coefficients have little individual effect on the validation score when removed. (b) The same network trained with input samples rotated by $\pm 30^\circ$. In (a) the second and fifth orientations for both S_1^1 and S_1^2 , corresponding to the 45° and 135° wavelets, are comparatively less important than other orientations at the same scale. This suggests that perhaps the dataset does not have much diagonal information. When rotated this trend changes and the diagonal wavelets at both scales become more important.

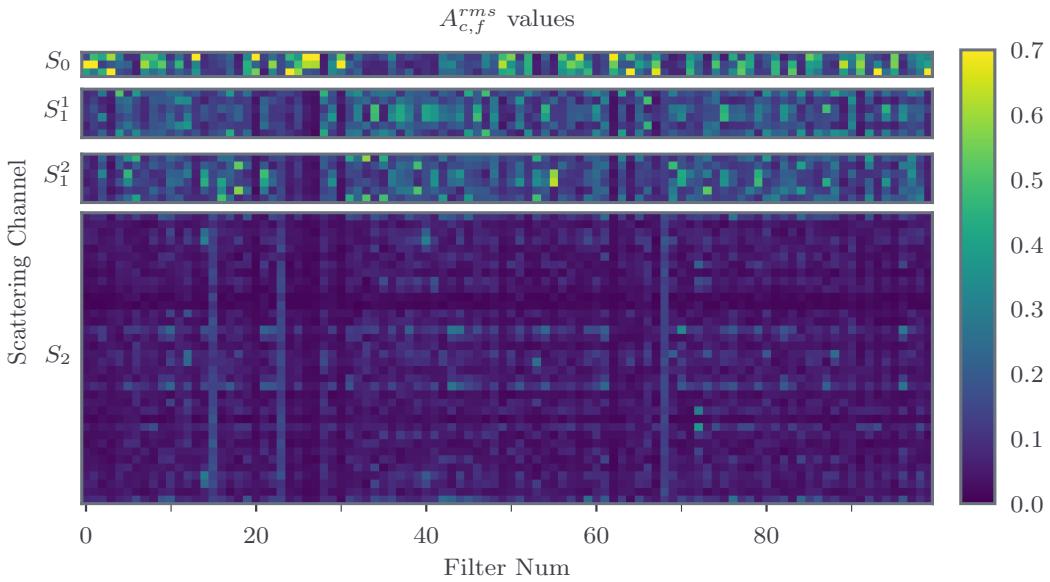


Figure 4.5: Channel weights for first learned layer in a hybrid ScatterNet-CNN. A visualization of the matrix A^{rms} from (4.6.1) for a network trained on Tiny ImageNet. The columns of the matrix all have unit norm and represent how much relative energy comes from each scattering output channel. Most of the filters are heavily dependent on S_0 , many are dependent on S_1 and only a few take information from S_2 .

mostly inspirational, as we did not see any marked improvement in using them as a fixed front end for the ScatterNet system. However, they are the basis for the next chapter of work in adding learning in between Scattering layers.

Sifre and Mallat introduced the idea of a ‘Roto-Translation’ ScatterNet in [99]. Invariance to rotation could be made by applying averaging (and bandpass) filters across the K orientations from the wavelet transform *before* applying the complex modulus. Let us call the averaging and bandpass filters they use $h \in \mathbb{C}^K$. We can think of this stage as stacking the K outputs of a complex wavelet transform on top of each other and convolving a filter h_α over all spatial locations of the wavelet coefficients $W_m x$. Let us call the output from filtering across the channel dimension V_m :

$$V_m x(j, \alpha, \mathbf{u}) = \sum_{\theta} W_m x(j, \theta, \mathbf{u}) h_{j, \alpha}(\theta) \quad (4.7.1)$$

We present a variation on this idea by filtering with a more general $h \in \mathbb{C}^{12 \times H \times W}$. We use 12 channels rather than 6, as we use the $K = 6$ orientations and their complex conjugates; each wavelet is a 30° rotation of the previous, so with 12 rotations, we can cover the full 360° .

Figure 4.7 shows some reconstructions from these V coefficients with hand-designed h 's using the above DeScatterNet. All shapes are shown in real and imaginary Hilbert-like pairs;

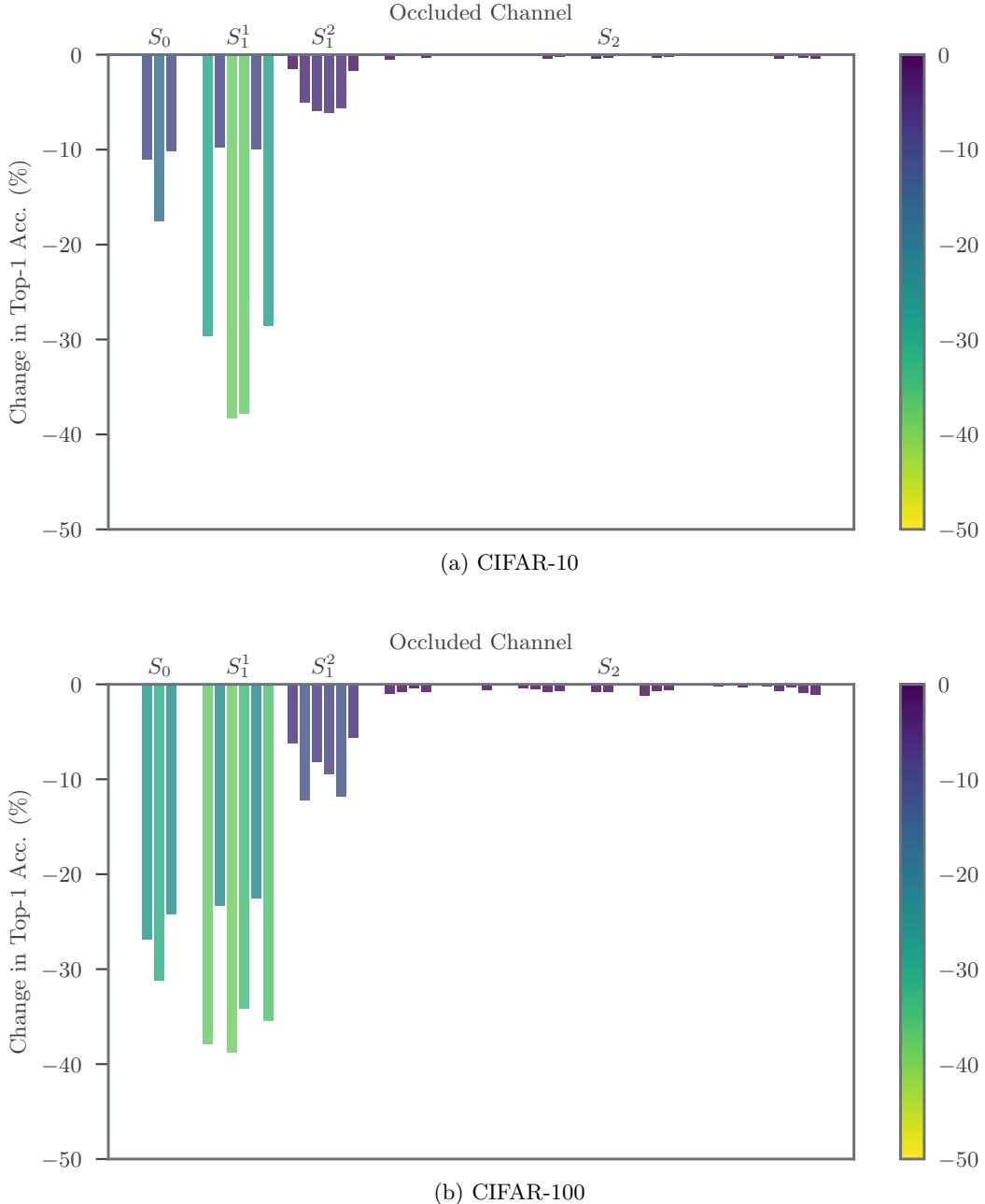


Figure 4.6: **CIFAR changes in accuracy from channel occlusion.** Numbers reported are the drop in final classification accuracy when a channel is set to zero. The bars are coloured relative to their magnitude to aid seeing the differences for the S_1 coefficients. Unlike Figure 4.4 the S_1^2 coefficients are less important. CIFAR has a smaller image size than Tiny ImageNet so this is not surprising.

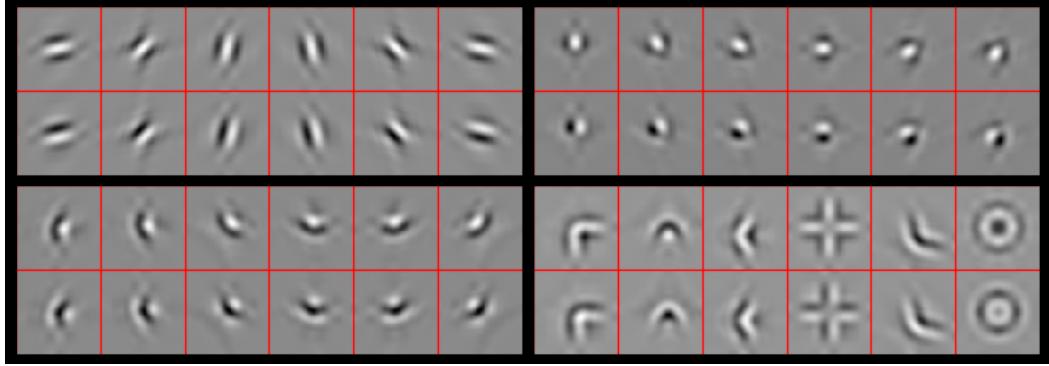


Figure 4.7: Shapes possible by filtering across the wavelet orientations with complex coefficients. All shapes are shown in pairs: the top image is reconstructed from a purely real output, and the bottom image from a purely imaginary output. These ‘real’ and ‘imaginary’ shapes are nearly orthogonal in the pixel space (normalized dot product < 0.01 for all but the doughnut shape in the bottom right, which has 0.15) but produce the same U' , something that would not be possible without the complex filters of a ScatterNet. Top left - reconstructions from U_1 (i.e. no cross-orientation filtering). Top right- reconstructions from U'_1 using a $12 \times 1 \times 1$ Morlet Wavelet, similar to what was done in the ‘Roto-Translation’ ScatterNet described in [25], [99]. Bottom left - reconstructions from U'_1 made with a more general $12 \times 1 \times 1$ filter, described in Equation 4.7.2. Bottom right - some reconstructions possible by filtering a general $12 \times 3 \times 3$ filter.

the top row of images in each quadrant are reconstructed from a purely real V , while the bottom row are reconstructed from a purely imaginary V .

In the top left, we display the 6 wavelet filters for reference. In the top right of the figure we see some of the shapes made by using the h ’s from the Roto-Translation ScatterNet [25], [99]. The bottom left is where we present some of our novel kernels. These are simple corner-like shapes made by filtering with $h \in \mathbb{C}^{12 \times 1 \times 1}$ where h is set to

$$h = [1, j, j, 1, 0, 0, 0, 0, 0, 0, 0, 0] \quad (4.7.2)$$

The six orientations are made by rolling the coefficients in h along one sample (i.e. $[0, 1, j, j, 1, 0, \dots]$, $[0, 0, 1, j, j, 1, 0, \dots]$, $[0, 0, 0, 1, j, j, 1, 0, \dots]$ …). Coefficients roll back around (like circular convolution) when they reach the end. The canonical filter is then $[1, j, j, 1]$ where the 90° phase offset of the middle two weights from the outer two allows for nicely continuous ridges of similar intensity around the centre of the corner.

Finally, in the bottom right we see shapes made by $h \in \mathbb{C}^{12 \times 3 \times 3}$. Note that with the exception of the ring-like shape which has 12 non-zero coefficients, all of these shapes were reconstructed with h ’s that have 4 to 8 non-zero coefficients of a possible 64. These shapes are now beginning to more closely resemble the more complex shapes seen in the middle stages of CNNs.

The shapes in the bottom row of the bottom right quadrant appear very similar to those in the top row but have nearly zero inner product. This shows one level of invariance of this system, as after taking the complex magnitude, both the top and the bottom shape will activate the network with the same strength. In comparison, for the purely real filters of a CNN if the top shape would cause a large output then the bottom shape would cause near 0 activity.

4.8 Conclusion

This chapter presents a way to investigate what the higher orders of a ScatterNet are responding to - the DeScatterNet described in [section 4.4](#). Using this, we have shown that the second-order of a ScatterNet responds strongly to patterns that are very different to those that highly activate the second layer of a CNN. As well as being dissimilar to CNNs, visual inspection of the ScatterNet's patterns reveal that they may be less useful for discriminative tasks, and we believe this may be causing the current gaps in state-of-the-art performance between the two.

Additionally, we performed occlusion tests to heuristically measure the importance of the individual scattering channels. The results of this test reaffirmed the suspicions raised from the visualizations. In particular, many of the second-order Scattering coefficients may not be very useful in a deep classifier. Those that were more useful were typically when the second-order wavelet had the same orientation as the first. We also noted that diagonal orientations appear less important than horizontal and vertical ones, even at coarser scales. This appears to be an artefact of the CIFAR and Tiny ImageNet datasets, as rotating the images by 30° made diagonal edges more important.

Finally, we demonstrated the possible shapes attainable when we filter across orientations with complex mixing coefficients. We believe that this mixing is a key step in the development of improved ScatterNets and wavelets in deep learning systems.

Chapter 5

A Learnable ScatterNet: Locally Invariant Convolutional Layers

In this chapter, we explore tying together the ideas from Scattering Transforms and Convolutional Neural Networks (CNN) for Image Analysis by proposing a learnable ScatterNet. The work presented in [chapter 4](#) implies that while the Scattering Transform has been a promising start in using complex wavelets in image understanding tasks, there is something missing from them. To address this, we propose a learnable ScatterNet by building it with our proposed “Locally Invariant Convolutional Layers”.

Previous attempts at combining ScatterNets with CNNs in hybrid networks [34], [39], [40] have tended to keep the two parts separate, with the ScatterNet forming a fixed front end and the CNN forming a learned backend. We instead look at adding learning between scattering orders, as well as adding learned layers before the ScatterNet.

We do this by adding a second stage after a scattering order, which mixes output activations together in a learnable way. The flexibility of the mixing we introduce allows us to build a layer that acts as a Scattering Layer with no learning, or as one that acts more as a convolutional layer with a controlled number of input and output channels, or more interestingly, as a hybrid between the two.

Our experiments show that these locally invariant layers can improve accuracy when added to either a CNN or a ScatterNet. We also discover some surprising results in that the ScatterNet may be best positioned after one or more layers of learning rather than at the front of a neural network.

5.1 Chapter Layout

In [section 5.2](#) we discuss related work before briefly reviewing the convolutional layer and scattering notation we will use in [section 5.3](#). We introduce our learnable scattering layers in [section 5.4](#) and describe their properties and implementation in [section 5.5](#). Finally, we

present some experiments we have run in ?? and then draw conclusions about how these new ideas might improve neural networks in the future.

5.2 Related Work

There have been several similar works that look into designing new convolutional layers by separating them into two stages — a first stage that performs a non-standard filtering process, and a second stage that combines the first stage into single activations. The inception layer [135] by Szegedy *et al.* does this by filtering with different kernel sizes in the first stage, and then combining with a 1×1 convolution in the second stage. Ioannou *et al.* also do something similar by making a first stage with horizontal and vertical filters, and then combining in the second stage again with a 1×1 convolution[136]. But perhaps the most similar works are those that use a first stage with fixed filters, combining them in a learned way in the second stage. Of particular note are:

- “Local Binary Convolutional Neural Networks” [137]. This paper builds a first stage with a small 3×3 kernel filled with zeros, and randomly insert ± 1 in several locations, keeping a set sparsity level. This builds a very crude spatial differentiator in random directions. The output of the first stage is then passed through a sigmoid nonlinearity before being mixed with a 1×1 convolution. The imposed structure on the first stage was found to be a good regularizer and prevented overfitting, and the combination of the mixing in the second layer allowed for a powerful and expressive layer, with performance near that of a regular CNN layer.
- “DCFNet: Deep Neural Network with Decomposed Convolutional Filters” [27]. This paper decomposes convolutional filters as linear combinations of Fourier Bessel and random bases. The first stage projects the inputs onto the chosen basis, and the second stage learns how to mix these projections with a 1×1 convolution. Unlike [137] this layer is purely linear. The supposed advantage being that the basis can be truncated to save parameters and make the input less susceptible to high frequency variations. The work found that this layer had marginal benefits over regular CNN layers in classification, but had improved stability to noisy inputs.

5.3 Recap of Useful Terms

5.3.1 Convolutional Layers

Let the output of a CNN at layer l be $x^{(l)}(c, \mathbf{u})$, $c \in \{0, \dots, C_l - 1\}$, $\mathbf{u} \in \mathbb{R}^2$ where c indexes the channel dimension, and \mathbf{u} is a vector of coordinates for the spatial position. Of course, \mathbf{u} is typically sampled on a grid, but we keep it continuous to more easily differentiate between

the spatial and channel dimensions. Recall from (2.4.5) and (2.4.6) that a convolutional layer in a standard CNN is defined by the two operations:

$$y^{(l+1)}(f, \mathbf{u}) = \sum_{c=0}^{C_l-1} x^{(l)}(c, \mathbf{u}) * h_f^{(l)}(c, \mathbf{u}) \quad (5.3.1)$$

$$x^{(l+1)}(f, \mathbf{u}) = \sigma\left(y^{(l+1)}(f, \mathbf{u})\right) \quad (5.3.2)$$

where $h_f^{(l)}(c, \mathbf{u})$ is the f th filter of the l th layer with C_l different point spread functions, and $f \in \{0, \dots, C_{l+1}-1\}$. σ is a nonlinearity such as the ReLU, possibly combined with scaling such as batch normalization. The convolution is done independently for each c in the C_l channels and the resulting outputs are summed together to give one activation map.

5.3.2 Wavelet Transforms

Recall from (2.6.26) that:

$$\mathcal{W}x(c, \mathbf{u}) = \{x(c, \mathbf{u}) * \phi_J(\mathbf{u}), x(c, \mathbf{u}) * \psi_\lambda(\mathbf{u})\}_\lambda \quad (5.3.3)$$

Where ψ_λ is a mother wavelet dilated by 2^j , $1 \leq j \leq J$ and rotated by $\theta = \frac{\pi+2k\pi}{12}$, $0 \leq k < 6$:

$$\psi_{j,\theta}(\mathbf{u}) = 2^{-j}\psi(2^{-j}r_{-\theta}\mathbf{u}) \quad (5.3.4)$$

Define the set of all possible λ s as Λ whose size is $|\Lambda| = JK$.

5.3.3 Scattering Transforms

As the real and imaginary parts of complex wavelets are in quadrature with each other, taking the modulus of the resulting transformed coefficients removes the high frequency oscillations of the output signal while preserving the energy of the coefficients over the frequency band covered by ψ_λ . This is crucial to ensure that the scattering energy is concentrated towards zero-frequency as the scattering order increases, allowing sub-sampling. We define the wavelet modulus propagator to be:

$$\widetilde{\mathcal{W}}x(c, \mathbf{u}) = \{x(c, \mathbf{u}) * \phi_J(\mathbf{u}), |x(c, \mathbf{u}) * \psi_\lambda(\mathbf{u})|\}_{\lambda \in \Lambda} \quad (5.3.5)$$

The modulus terms are called $U[\lambda]x = |x * \psi_\lambda|$, and the scattering terms are $S[\lambda]x = U[\lambda]x * \phi_J(\mathbf{u})$.

5.4 Locally Invariant Layer

We propose to mix the terms at the output of each wavelet modulus propagator $\tilde{\mathcal{W}}$. The second term in $\tilde{\mathcal{W}}$, the U terms, are often called ‘covariant’ terms but in this work, we will call them *locally invariant*, as they tend to be invariant up to a scale 2^j . We propose to mix these locally invariant terms U and the lowpass terms S with learned weights $a_{f,\lambda}$ and b_f .

For example, consider the wavelet modulus propagator from (5.3.5), and let the input to it be $x^{(l)}$. Our proposed output is then:

$$y^{(l+1)}(f, \mathbf{u}) = \sum_{\lambda \in \Lambda} \sum_{c=0}^{C-1} a_{f,\lambda}(c) \left| x^{(l)}(c, \mathbf{u}) * \psi_\lambda(\mathbf{u}) \right| + \sum_{c=0}^{C-1} b_f(c) \left(x^{(l)}(c, \mathbf{u}) * \phi_J(\mathbf{u}) \right) \quad (5.4.1)$$

$$= \sum_{c=0}^{C-1} \left(b_f(c) \left(x^{(l)}(c, \mathbf{u}) * \phi_J(\mathbf{u}) \right) + \sum_{\lambda \in \Lambda} a_{f,\lambda}(c) \left| x^{(l)}(c, \mathbf{u}) * \psi_\lambda(\mathbf{u}) \right| \right) \quad (5.4.2)$$

Note that an input to the wavelet modulus propagator $\tilde{\mathcal{W}}$ with C channels has $(JK+1)C$ output channels – C lowpass channels and JKC modulus bandpass channels. Let us define a new output z with index variable $q \in \mathbb{Z}$ such that:

$$z^{(l+1)}(q, \mathbf{u}) = \begin{cases} x^{(l)}(c, \mathbf{u}) * \phi_J(\mathbf{u}) & \text{if } 0 \leq q < C \\ |x^{(l)}(c, \mathbf{u}) * \psi_\lambda(\mathbf{u})| & \text{if } C \leq q < (JK+1)C \end{cases} \quad (5.4.3)$$

i.e. the lowpass channels are the first C channels of z , the modulus of the 15° ($k=0$) wavelet coefficients with $j=1$ are the next C channels, then the modulus coefficients with $k=1$ and $j=1$ are the third C channels, and so on. This is similar to the view of a ScatterNet we introduced in [Algorithm 3.5](#), but restricted to a single order.

We do the same for the weights a, b by defining $\tilde{a}_f = \{b_f, a_{f,\lambda}\}_{\lambda}$ and let:

$$\tilde{a}_f(q) = \begin{cases} b_f(c) & \text{if } 0 \leq q < C \\ a_{f,\lambda}(c) & \text{if } C \leq q < (JK+1)C \end{cases} \quad (5.4.4)$$

we can then use (5.4.3) and (5.4.4) to simplify (5.4.2), giving:

$$y^{(l+1)}(f, \mathbf{u}) = \sum_{q=0}^{(JK+1)C-1} z^{(l+1)}(q, \mathbf{u}) \tilde{a}_f(q) \quad (5.4.5)$$

or in matrix form with $A_{f,q} = \tilde{a}_f(q)$

$$Y^{(l+1)}(\mathbf{u}) = AZ^{(l+1)}(\mathbf{u}) \quad (5.4.6)$$

This equation now looks very similar to the standard convolutional layer from (5.3.1), except we have replaced the previous layer’s x with intermediate coefficients z with $|Q|=$

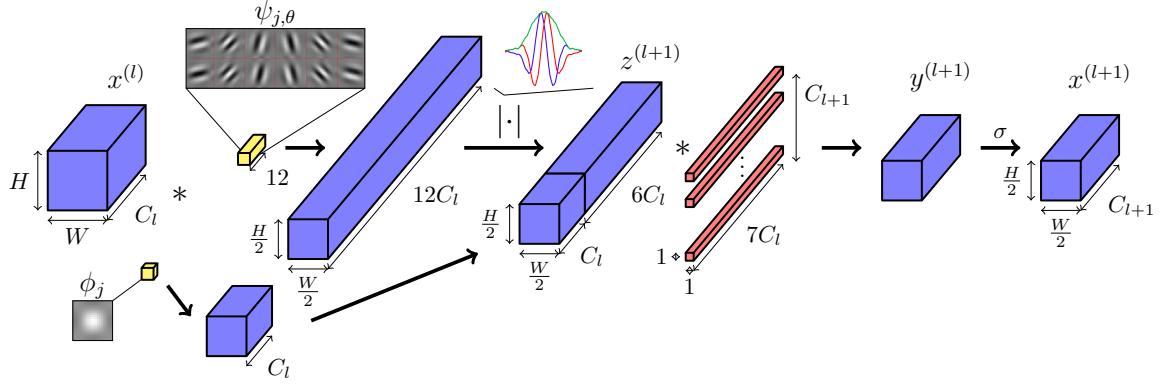


Figure 5.1: **Block Diagram of Proposed Invariant Layer for $j = J = 1$.** Activations are shaded blue, fixed parameters yellow and learned parameters red. Input $x^{(l)} \in \mathbb{R}^{C_l \times H \times W}$ is filtered by real and imaginary oriented wavelets and a lowpass filter and is downsampled. The channel dimension increases from C_l to $(2K + 1)C_l$, where the number of orientations is $K = 6$. The real and imaginary parts are combined by taking their magnitude (an example of what this looks like in 1D is shown above the magnitude operator) - taking the envelope of the components oscillating in quadrature. These are concatenated with the lowpass activations to give $z^{(l+1)}$. The resulting activations z are mixed across the channel dimension and then passed through a nonlinearity σ to give $x^{(l+1)}$. If the desired output spatial size is $H \times W$, $x^{(l+1)}$ can be bilinearly upsampled paying only a few multiplies per pixel.

$(JK + 1)C$ channels and the convolutions of (5.3.1) have been replaced by a matrix multiply (which can also be seen as a 1×1 convolutional layer). We can then apply (5.3.2) to (5.4.5) to get the next layer's output:

$$x^{(l+1)}(f, \mathbf{u}) = \sigma(y^{(l+1)}(f, \mathbf{u})) \quad (5.4.7)$$

Figure 5.1 shows a block diagram for this process.

5.4.1 Properties

5.4.1.1 Recovering the original ScatterNet Design

The first thing to note is that with careful choice of A and σ , we can recover the original translation invariant ScatterNet [23], [39]. If $C_{l+1} = (JK + 1)C_l$ and A is the identity matrix $I_{C_{l+1}}$, there is no mixing and $y^{(l+1)} = \tilde{\mathcal{W}}x$.

Further, if $\sigma = \text{ReLU}$ as is commonly the case in training CNNs, σ has no effect on the positive locally invariant terms U . It will affect the averaging terms if the signal is not positive, but this can be dealt with by adding a channel-dependent bias term α_c to $x^{(l)}$ to ensure it is positive. This bias term will not affect the propagated signals as $\int \alpha_c \psi_\lambda(\mathbf{u}) d\mathbf{u} = 0$. The bias can then be corrected by subtracting $\alpha_c \|\phi_J\|_2$ from the averaging terms after taking the ReLU, then $x^{(l+1)} = \tilde{\mathcal{W}}x$.

This makes one layer of our system equivalent to a first-order scattering transform, giving S_0 and U_1 (invariant to input shifts of 2^1). We can repeat the same process for the next layer (recall the recursive scattering relation we saw in (2.7.12)), giving S_1 and U_2 (invariant to shifts of 2^2). If we want to build higher invariance, we can continue or simply average these outputs with an average pooling layer.

5.4.1.2 Flexibilty of the Layer

Unlike a regular ScatterNet, we are free to choose the size of C_{l+1} . This means we can set $C_{l+1} = C_l$ as is commonly the case in a CNN, and make a convolutional layer from mixing the locally invariant terms. This avoids the exponentially increasing complexity that comes with extra network layers that standard ScatterNets suffer from.

5.4.1.3 Stability to Noise and Deformations

Let us define the action of our layer on the scattering coefficients to be Vx . We would like to find a bound on $\|V\mathcal{L}_\tau x - Vx\|$. To do this, we note that the mixing is a linear operator and hence is Lipschitz continuous. The authors in [27] find constraints on the mixing weights to make them non-expansive (i.e. Lipschitz constant 1). Further, the ReLU is non-expansive meaning the combination of the two is also non-expansive, so $\|V\mathcal{L}_\tau x - Vx\| \leq \|S\mathcal{L}_\tau x - Sx\|$, and (2.7.23) holds.

5.5 Implementation Details

Again, we use the DTCWT [18] for our wavelet filters due to their fast implementation with separable convolutions which we discuss more in subsection 5.5.3. There are two side effects of this choice. The first is that the number of orientations of wavelets is restricted to $K = 6$. The second is that we naturally downsample the output activations by a factor of 2 for each direction for each scale j , giving a 4^j downsampling factor overall. This represents the source of the invariance in our layer. If we do not wish to downsample the output (say to make the layer fit in a larger network), we can bilinearly interpolate the output of our layer. This is computationally cheap to do on its own, but causes the next layer's computation to be higher than necessary (there will be almost no energy for frequencies higher than $f_s/4$).

In all our experiments we set $J = 1$ for each invariant layer, meaning we can mix the lowpass and bandpass coefficients at the same resolution. Figure 5.1 shows how this is done. Note that setting $J = 1$ for a single layer does not restrict us from having $J > 1$ for the entire system, as if we have a second layer with $J = 1$ after the first, including downsampling (\downarrow), we would have:

$$(((x * \phi_1) \downarrow 2) * \psi_{1,\theta}) \downarrow 2 = (x * \psi_{2,\theta}) \downarrow 4 \quad (5.5.1)$$

Algorithm 5.1 Locally Invariant Convolutional Layer forward and backward passes

```

1: procedure INVLAYER.FORWARD( $x, A$ )
2:    $yl, yh \leftarrow \text{DT\textcircled{C}WT}.\text{Forward}(x^l, \text{nlevels} = 1)$ 
3:    $U \leftarrow \text{MAG\_SMOOTH}.\text{Forward}(yh)$                                  $\triangleright$  See Algorithm B.2
4:    $yl \leftarrow \text{AVGPOOL2x2}(yl)$                                           $\triangleright$  Downsample lowpass to match U size
5:    $Z \leftarrow \text{CONCATENATE}(yl, U)$                                           $\triangleright$  Concatenate along the channel dim
6:    $Y \leftarrow AZ$                                                                 $\triangleright$  Mix
7:   save  $A, Z$                                                                 $\triangleright$  For the backwards pass
8:   return  $Y$ 
9: end procedure

1: procedure INVLAYER.BACKWARD( $\frac{\partial L}{\partial Y}$ )
2:   load  $A, Z$ 
3:    $\frac{\partial L}{\partial A} \leftarrow \frac{\partial L}{\partial Y} Z^T$                                 $\triangleright$  Calculate update gradient
4:    $\Delta Z \leftarrow A^T \frac{\partial L}{\partial Y}$ 
5:    $\Delta yl, \Delta U \leftarrow \text{UNSTACK}(\Delta Z)$ 
6:    $\Delta yl \leftarrow \text{AVGPOOL2x2}.\text{Backward}(\Delta yl)$ 
7:    $\Delta yh \leftarrow \text{MAG\_SMOOTH}.\text{Backward}(\Delta U)$ 
8:    $\frac{\partial L}{\partial x} \leftarrow \text{DT\textcircled{C}WT}.\text{Backward}(\Delta yl, \Delta yh)$        $\triangleright$  Calculate passthrough gradient
9:   return  $\frac{\partial L}{\partial x}, \frac{\partial L}{\partial A}$ 
10: end procedure

```

5.5.1 Parameter Memory Cost

A standard convolutional layer with C_l input channels, C_{l+1} output channels and kernel size $L \times L$ has $L^2 C_l C_{l+1}$ parameters.

The number of learnable parameters in each of our proposed invariant layers with $J = 1$ and $K = 6$ orientations is:

$$\# \text{params} = (JK + 1)C_l C_{l+1} = 7C_l C_{l+1} \quad (5.5.2)$$

The spatial support of the wavelet filters is typically 5×5 pixels or more, and we have reduced the number of parameters to fewer than $3 \times 3 = 9$ per filter, while producing filters that are significantly larger than this.

5.5.2 Activation Memory Cost

A standard convolutional layer needs to save the activation $x^{(l)}$ to convolve with the back-propagated gradient $\frac{\partial L}{\partial y^{(l+1)}}$ on the backwards pass (to give $\frac{\partial L}{\partial w^{(l)}}$). For an input with C_l channels of spatial size $H \times W$, this means HWC_l floats must be saved.

Our layer requires us to save the activation $z^{(l+1)}$ for updating the \tilde{a} terms. This has $7C_l$ channels of spatial size $\frac{HW}{4}$. This means that our proposed layer needs to save $\frac{7}{4}HWC_l$ floats, a $\frac{7}{4}$ times memory increase on the standard layer.

5.5.3 Computational Cost

A standard convolutional layer with kernel size $L \times L$ needs $L^2 C_{l+1}$ multiplies per input pixel (of which there are $C_l \times H \times W$).

There is an overhead in doing the wavelet decomposition for each input channel. A separable 2-D discrete wavelet transform (DWT) with 1-D filters of length L will have $2L(1 - 2^{-2J})$ multiplies per input pixel for a J scale decomposition. A DTCWT has 4 DWTs for a 2-D input, so its cost is $8L(1 - 2^{-2J})$, with $L = 6$ a common size for the filters. It is important to note that unlike the filtering operation, this does not scale with C_{l+1} , the end result being that as C_{l+1} grows, the cost of C_l forward transforms is outweighed by that of the mixing process whose cost is proportional to $C_l C_{l+1}$.

Because we are using a decimated wavelet decomposition, the sample rate decreases after each wavelet layer. The benefit of this is that the mixing process then only works on 1/4 the spatial area after the first scale and 1/16 the spatial area after the second scale. Restricting ourselves to $J = 1$ as we mentioned in [section 5.5](#), the computational cost is then:

$$\underbrace{\frac{7}{4}C_{l+1}}_{\text{mixing}} + \underbrace{36}_{\text{DTCWT}} \quad \text{multiplies per input pixel} \quad (5.5.3)$$

In most CNNs, C_{l+1} is several dozen if not several hundred, which makes (5.5.3) significantly smaller than $L^2 C_{l+1} = 9C_{l+1}$ multiplies for 3×3 convolutions.

5.5.4 Forward and Backward Algorithm

There are two layer-hyperparameters to choose:

- The number of output channels C_{l+1} . This may be restricted by the architecture.
- The variance of the weight initialization for the mixing matrix A .

Assuming we have already chosen these values, then the forward and backward algorithms can be computed with Algorithm [Algorithm 5.1](#).

5.6 Layer Introduction with MNIST

To begin experiments on the proposed locally invariant layer, we look at how well a simple system works on MNIST, and compare it to an equivalent system with convolutions. Because of the small size of the MNIST challenge, we can quickly get results, allowing a large number of trials and a broad search over hyperparameters to be done. In this way, we can use the findings from these experiments to guide our work on more difficult tasks like CIFAR and Tiny ImageNet.

Table 5.1: **Architectures for MNIST hyperparameter experiments.** The activation size rows are offset from the layer description rows to convey the input and output shapes. The ‘project’ layers in both architectures are unlearned, so all of the learning has to be done by the first two layers and the reshuffle layer.

| (a) Reference Arch with 3×3 convolutions | | (b) Invariant Architecture | |
|---|--|----------------------------|--|
| Activation Size | Layer Name + Info | Activation Size | Layer Name + Info |
| $1 \times 28 \times 28$ | | $1 \times 28 \times 28$ | |
| $7 \times 28 \times 28$ | conv1, $w \in \mathbb{R}^{7 \times 1 \times 3 \times 3}$ | $7 \times 14 \times 14$ | inv1, $A \in \mathbb{R}^{7 \times 7}$ |
| $7 \times 14 \times 14$ | maxpool1, 2×2 | $49 \times 7 \times 7$ | inv2, $A \in \mathbb{R}^{49 \times 49}$ |
| $49 \times 14 \times 14$ | conv2 $w \in \mathbb{R}^{49 \times 7 \times 3 \times 3}$ | 2401×1 | unravel |
| $49 \times 7 \times 7$ | maxpool2, 2×2 | 10×1 | project, $w \in \mathbb{R}^{2401 \times 10}$ |
| $49 \times 7 \times 7$ | unravel | 10×1 | reshuffle, $w \in \mathbb{R}^{10 \times 10}$ |
| 2401×1 | project, $w \in \mathbb{R}^{2401 \times 10}$ | 10×1 | |
| 10×1 | reshuffle, $w \in \mathbb{R}^{10 \times 10}$ | | |
| 10×1 | | | |

Table 5.2: **Hyperparameter settings for the MNIST experiments.** The weight gain is the term a from [Equation 5.6.1](#). Note that $\log_{10} 3.16 = 0.5$.

| Hyperparameter | Values |
|--------------------|---|
| Learning Rate (lr) | {0.0316, 0.1, 0.316, 1} |
| Momentum (mom) | {0, 0.25, 0.5, 0.9} |
| Weight Decay (wd) | { 10^{-5} , 3.16×10^{-5} , 10^{-4} , 3.16×10^{-4} } |
| Weight Gain (a) | {0.5, 1.0, 1.5, 2.0} |

To minimize the effects of learning from other layers, we build a custom small network, as described in [Table 5.1](#). The first two layers are learned convolutional/invariant layers, followed by a fully connected layer with *fixed* weights that we can use to project down to the number of output classes. Finally, we add a small learned layer that linearly combines the 10 outputs from the random projection, to give 10 new outputs. This is to facilitate reordering of the outputs to the correct class. This simple network is meant to test the limits of our layer, rather than achieve state of the art performance on MNIST.

Given that our layer is quite different to a standard convolutional layer, we must do a full hyperparameter search over optimizer parameters such as the learning rate, momentum, and weight decay, as well as layer hyperparameters such as the variance of the random initialization for the mixing matrix A .

Table 5.3: **Architecture performance comparison.** Numbers reported are the mean and standard deviation of accuracy over 10 runs with the optimal hyperparameters, θ . Note that for both architectures we found that the learning rate, lr, was the most important hyperparameter to choose correctly, and had the largest impact on the performance.

| Architecture | $\theta = \{\text{lr, mom, wd, a}\}$ | accuracy | |
|---------------|---|----------|------|
| | | mean | std |
| Convolutional | $\{0.1, 0.5, 10^{-5}, 1.5\}$ | 97.3 | 0.29 |
| Invariant | $\{0.032, 0.9, 3.2 \times 10^{-5}, 1.0\}$ | 96.6 | 0.26 |

To simplify the weight variance search, we use Glorot Uniform Initialization [56] and only vary the gain value a :

$$A_{ij} \sim U \left[-a \sqrt{\frac{6}{(C_l + C_{l+1})HW}}, a \sqrt{\frac{6}{(C_l + C_{l+1})HW}} \right] \quad (5.6.1)$$

where C_l , C_{l+1} are the number of input and output channels as before, and the kernel size is $H = W = 1$ for an invariant layer and $H = W = 3$ for a convolutional layer.

We do a grid search over these hyperparameters and use Hyperband [138] to schedule early stopping of poorly performing runs. Each run has a grace period of 5 epochs and can train for a maximum of 20 epochs. We do not do any learning rate decay. We found the package Tune [139] was very helpful in organising parallel distributed training runs. The hyperparameter options are described in Table 5.2, note that we test $4^4 = 256$ different options.

Once we find the optimal hyperparameters for each network, we then run the two architectures 10 times with different random seeds and report the mean and variance of the accuracy. The results of these runs are listed in Table 5.3.

5.6.1 Proposed Expansions

The results from the previous section seem to indicate that our proposed invariant layer is a slightly worse substitute for a convolutional layer.

We posit that this is due to the centred nature of the wavelet bases that were used to generate the z and later the y coefficients. As they are all centred at roughly the same location (the phase of the complex wavelet allows for some small spatial separation) 1×1 convolutions may not be capable of building richer shapes by separating the wavelet centres. This effect was seen in the previous chapter when we presented some possible new shapes in Figure 4.7. The bottom right quadrant were shapes made from mixing wavelet coefficients in a 3×3 area, and these were much richer than those attainable by only mixing in a 1×1 area¹

¹This comparison is only a guide however, as in this chapter we are mixing coefficients after taking the modulus, where as Figure 4.7 were generated by mixing with complex gains *before* the modulus.

To get spatial separation, we must replace the 1×1 mixing kernel from (5.4.5) ($\tilde{\alpha}_f(q)$) with a 3×3 filter $g_f(q, \mathbf{u})$. We also change the multiply for convolution, giving us:

$$y^{(l+1)}(f, \mathbf{u}) = \sum_{q \in Q} z^{(l+1)}(q, \mathbf{u}) * g_f(q, \mathbf{u}) \quad (5.6.2)$$

If all of the parameters of $g_f(q, \mathbf{u})$ are learned, then the number of parameters in (5.5.2) increases from $7C_l C_{l+1}$ to $63C_l C_{l+1}$. This is quite a lot of parameters for a single layer, and much more than the $9C_l C_{l+1}$ of a standard 3×3 convolution, or even the $25C_l C_{l+1}$ of a larger 5×5 convolution.

However, we can still get spatial separation without having to learn *all* of the coefficients, if we factor g as:

$$g_f(q, \mathbf{u}) = \tilde{\alpha}_f(q)\alpha_f(q, \mathbf{u}) \quad (5.6.3)$$

where $\alpha_f(q, \mathbf{u})$ is an introduced *fixed* kernel, designed to allow mixing of wavelets from neighbouring spatial locations.

We test a range of possible α 's each with varying complexity/overhead:

- (a) **Random Shifts:** We randomly shift each of the $7C$ subbands horizontally by $\{-1, 0, 1\}$ pixels, and vertically by $\{-1, 0, 1\}$ pixels. This is determined at the beginning of a training session and is consistent between batches. This theoretically is free to do, but practically it involves setting $\alpha_f(q, \mathbf{u})$ to be a 3×3 kernel with a single 1 and eight 0's.
- (b) **Shifted Gaussians:** Instead of shifting impulses as in the previous option, we can shift a gaussian kernel by $\{-1, 0, 1\}$ pixel horizontally and $\{-1, 0, 1\}$ pixel vertically, making a smoother filter than in (a).
- (c) **Random Kernels:** We can set α to be a random 3×3 kernel. This is chosen once at the beginning of training and then is kept fixed between batches.
- (d) **DCT Kernels:** We can learn coefficients for the top three 3×3 discrete cosine transform (DCT) coefficients² and sum them to get g . This is equivalent to having three different α 's, and learning three sets of $\tilde{\alpha}_f(q)$. This is a step between the parameterless kernels of (a) - (c) and the nine-fold parameter increase from learning all of g .

5.6.2 Expanded MNIST experiments

We test if the expansions from the previous section improve the MNIST performance. Again, we search over the hyperparameter space to find the optimal hyperparameters and then run 10 runs at the best set of hyperparameters, and report the results in [Table 5.4](#).

The top two rows are the old results form [Table 5.3](#), the next four rows are the expanded kernels described in the previous section, options (a) - (d). We also include tests for a fully

²The top three DCT coefficients are the constant, the horizontal and the vertical filters.

learned g kernel (despite the expense) in the seventh row under ‘Learned 3×3 ’. To compare this option to an equivalent convolutional system, we modify the convolutional architecture from [Table 5.1](#) to use 5×5 convolutions and $C_1 = 10$, $C_2 = 100$ channels and include these results in the final row of [Table 5.4](#) under ‘Wide Convolutional’. For ease of comparison, we list the computational and parameter cost associated with each option alongside the accuracy results.

As expected, adding in random shifts significantly helps the invariant layer for all but the random kernel (c) option. Two systems of note are the shifted impulse (a) system and the fully learned 3×3 kernel system. The first improves the mean accuracy by 1.3% without any extra learning. The second improves the performance by 2.4% but with a large parameter cost.

Table 5.4: Modified architecture performance comparison. Numbers reported are the mean and standard deviation of accuracy over 10 runs with the optimal hyperparameters, θ . We also list parameter cost and number of multiplies for each layer option, relative to the standard 3×3 convolutional layer to highlight the benefits/drawbacks of each option.

| Architecture | $\theta = \{\text{lr, mom, wd, a}\}$ | cost | | accuracy | |
|-------------------------------------|---|---------------|----------------|-------------|------|
| | | param | mults | mean | std |
| Convolutional | $\{0.1, 0.5, 10^{-5}, 1.5\}$ | 1 | 1 | 97.3 | 0.29 |
| Invariant | $\{0.032, 0.9, 3.2 \times 10^{-5}, 1.0\}$ | $\frac{7}{9}$ | $\frac{7}{36}$ | 96.6 | 0.26 |
| Shifted impulses (a) | $\{0.32, 0.5, 10^{-4}, 1.0\}$ | $\frac{7}{9}$ | $\frac{7}{36}$ | 97.9 | 0.25 |
| Shifted gaussians (b) | $\{1.0, 0.0, 10^{-5}, 1.0\}$ | $\frac{7}{9}$ | $\frac{7}{4}$ | 97.7 | 0.56 |
| Random 3×3 kernel (c) | $\{1.0, 0.9, 10^{-5}, 1.0\}$ | $\frac{7}{9}$ | $\frac{7}{4}$ | 95.8 | 1.01 |
| Learned 3×3 DCT coeffs (d) | $\{1.0, 0.0, 10^{-5}, 1.0\}$ | $\frac{7}{3}$ | $\frac{7}{4}$ | 98.1 | 0.37 |
| Learned 3×3 kernel | $\{0.1, 0.5, 10^{-4}, 1.0\}$ | 7 | $\frac{7}{4}$ | 99.0 | 0.12 |
| Wide Convolutional | $\{0.32, 0.5, 10^{-5}, 1.5\}$ | 7 | 7 | 98.7 | 0.25 |

5.7 Ablation Experiments with CIFAR and Tiny ImageNet

Now we look at expanding our layer to harder datasets, focusing more on the final classification accuracy. We do this again by comparing to a reference architecture. For this task, we choose a VGG-like network as our reference. It has six convolutional layers for CIFAR and eight layers for Tiny ImageNet as shown in [Table 5.5a](#). The initial number of channels C we use is 64. Despite this simple design, this reference architecture achieves near state-of-the-art

Table 5.5: **CIFAR and Tiny ImageNet Base Architecture.** Reference architecture used for experiments on CIFAR-10, CIFAR-100 and Tiny ImageNet. This architecture is based off the VGG[66] architecture. C is a hyperparameter that controls the network width, we use $C = 64$ for our initial tests. The activation size rows are offset from the layer description rows to convey the input and output shapes.

| (a) CIFAR Architecture | | (b) Tiny ImageNet Architecture | |
|-----------------------------|--|--------------------------------|--|
| Activation Size | Layer Name + Info | Activation Size | Layer Name + Info |
| $3 \times 32 \times 32$ | | $3 \times 64 \times 64$ | |
| $C \times 32 \times 32$ | convA, $w \in \mathbb{R}^{C \times 3 \times 3 \times 3}$ | $C \times 64 \times 64$ | convA, $w \in \mathbb{R}^{C \times 3 \times 3 \times 3}$ |
| $C \times 32 \times 32$ | convB, $w \in \mathbb{R}^{C \times C \times 3 \times 3}$ | $C \times 64 \times 64$ | convB, $w \in \mathbb{R}^{C \times C \times 3 \times 3}$ |
| $C \times 32 \times 32$ | pool1, max pooling 2×2 | $C \times 64 \times 64$ | pool1, max pooling 2×2 |
| $C \times 16 \times 16$ | convC, $w \in \mathbb{R}^{2C \times C \times 3 \times 3}$ | $C \times 32 \times 32$ | convC, $w \in \mathbb{R}^{2C \times C \times 3 \times 3}$ |
| $2C \times 16 \times 16$ | convD, $w \in \mathbb{R}^{2C \times 2C \times 3 \times 3}$ | $2C \times 32 \times 32$ | convD, $w \in \mathbb{R}^{2C \times 2C \times 3 \times 3}$ |
| $2C \times 16 \times 16$ | pool2, max pooling 2×2 | $2C \times 32 \times 32$ | pool2, max pooling 2×2 |
| $2C \times 8 \times 8$ | convE, $w \in \mathbb{R}^{4C \times 2C \times 3 \times 3}$ | $2C \times 16 \times 16$ | convE, $w \in \mathbb{R}^{4C \times 2C \times 3 \times 3}$ |
| $4C \times 8 \times 8$ | convF, $w \in \mathbb{R}^{4C \times 4C \times 3 \times 3}$ | $4C \times 16 \times 16$ | convF, $w \in \mathbb{R}^{4C \times 4C \times 3 \times 3}$ |
| $4C \times 8 \times 8$ | avg, 8×8 average pooling | $4C \times 16 \times 16$ | pool3, max pooling 2×2 |
| $4C \times 1 \times 1$ | fc1, fully connected layer | $4C \times 8 \times 8$ | convG, $w \in \mathbb{R}^{8C \times 4C \times 3 \times 3}$ |
| $10 \times 1, 100 \times 1$ | | $8C \times 8 \times 8$ | convH, $w \in \mathbb{R}^{8C \times 8C \times 3 \times 3}$ |
| | | $8C \times 1$ | avg, 8×8 average pooling |
| | | 200×1 | fc1, fully connected layer |

performance for the three datasets (92.6% on CIFAR-10, 72.0% on CIFAR-100 and 59.3% on Tiny ImageNet).

We perform an ablation study where we progressively swap out convolutional layers for invariant layers, keeping the input and output activation sizes the same. As there are 6 layers (or 8 for Tiny ImageNet), there are too many permutations to list the results for swapping out all layers for our locally invariant layer, so we restrict our results to only swapping 1 or 2 layers.

The invariant layer naturally downsamples the input. If we need to keep the output resolution the same as the input (i.e. ‘convA’, ‘convF’), we bilinearly interpolate the output of the invariant layer. If we swap out a layer immediately *before* a pooling layer (i.e. ‘convB’, ‘convD’) then we remove the pooling and do not upsample. Similarly, if we swap out a layer immediately *after* a pooling layer (i.e. ‘convC’, ‘convE’) then we remove the pooling layer and do the downsampling with the invariant layer.

[Figure 5.2](#) and [Figure 5.3](#) reports the top-1 classification accuracies for CIFAR-10, CIFAR-100 and Tiny ImageNet. In the table, ‘invX’ means that the ‘convX’ layer from [Table 5.5a](#) was replaced with an invariant layer. If the convolutional layer is before a pooling layer, then we do not interpolate the output of the invariant layer and we remove the pooling layer.

This network is optimized with stochastic gradient descent with momentum. The initial learning rate is 0.5, momentum is 0.85, batch size $N = 128$ and weight decay is 10^{-4} . For CIFAR-10/CIFAR-100 we scale the learning rate by a factor of 0.2 after 60, 80 and 100 epochs, training for 120 epochs in total. For Tiny ImageNet, the rate change is at 18, 30 and 40 epochs (training for 45 in total).

We see improvements for all three datasets for many possible locations for the invariant layer. The best permutations happen when one or two invariant layers are used near the start of a system, but *not* for the first layer. In particular, the best position for the invariant layer seems to be when we remove the first max pooling layer and use the natural downsampling of the invariant layer (at either the ‘convB’ or ‘convC’ position). This also worked for the second pooling layer (at the ‘convD’ position), but having an invariant layer at both of these positions (i.e. ‘convBD’ and ‘convCE’) performed slightly worse than using only one.

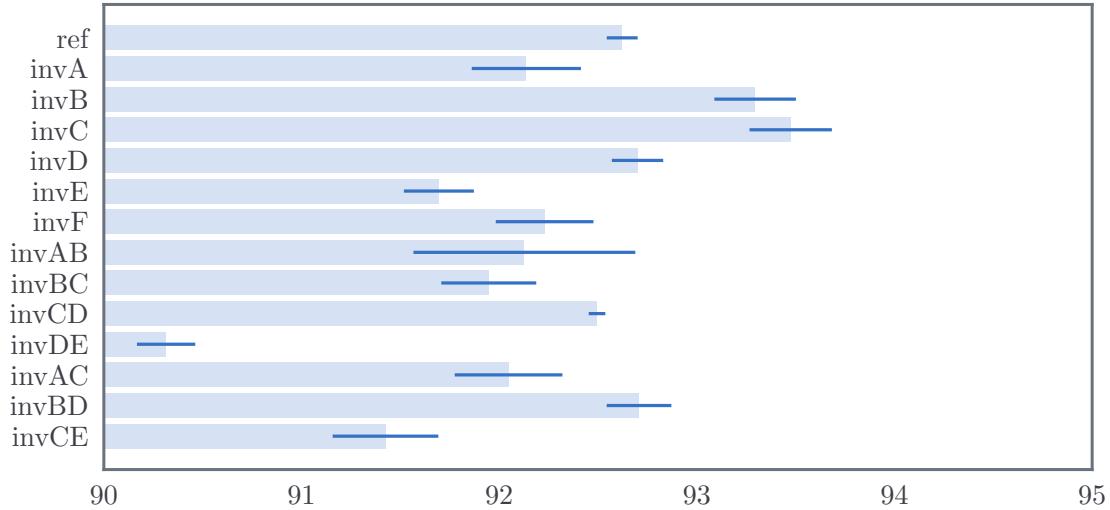
We recall that the magnitude operation in the ScatterNet effectively demodulates the energy from higher spatial frequencies to lower ones. This perhaps may explain why the best place for learnable scattering layers are at positions where you want to downsample in a network.

We also tested the variations from [subsection 5.6.1](#) that added spatial offsets to the gain layer by a fixed kernel and saw similar, if not slightly worse, results than those in [Figure 5.2](#) and [Figure 5.3](#). We did not have time to explore fully as to why this was the case, but we believe that the convolutional backend in the reference architecture ([Table 5.5](#)) provided the necessary spatial separation to build rich shapes, something that we did not have in the MNIST experiments.

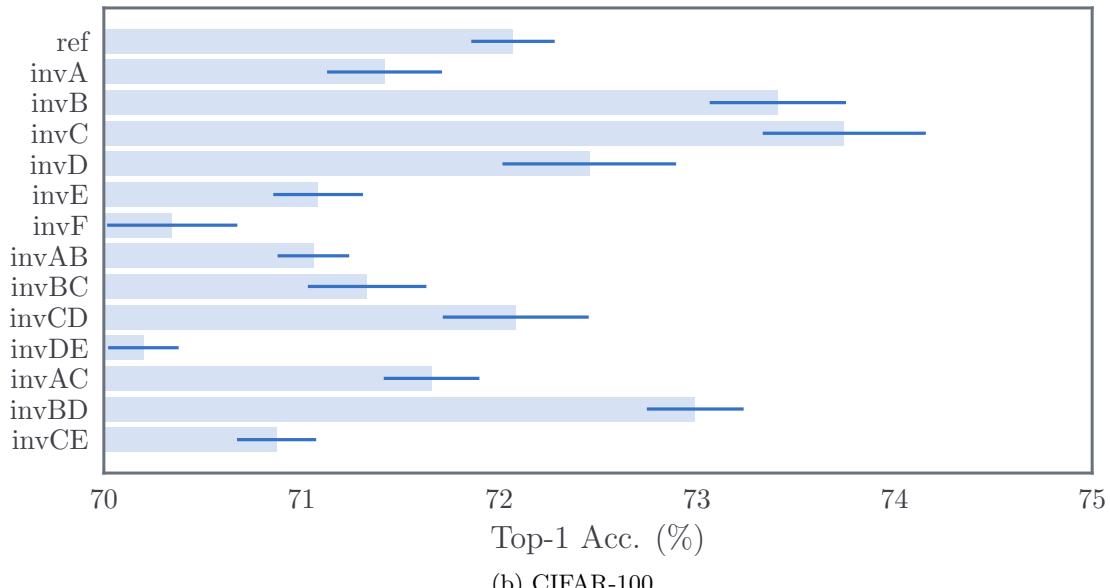
5.8 A New Hybrid ScatterNet

In the previous section, we examined how the locally invariant layer performs when directly swapped for a convolutional layer in an CNN based architecture. In this section, we look at how it performs in a hybrid ScatterNet-like network [39], [40]. Recall that a ScatterNet naturally increases the channel dimension while downsampling the spatial dimensions. A single scattering order for our ScatterNet increases the channel dimension seven-fold, and reduces the spatial area by four. After two layers of a ScatterNet, the output has 49 times more channels than the input.

We allow for this natural channel growth, using a ScatterNet front-end before four convolutional layers similar to ‘convC’ to ‘convF’ from [Table 5.5](#). The input to ‘convC’ already has spatial size 8×8 so we do not need the pooling layer between ‘convD’ and ‘convE’.



(a) CIFAR-10



(b) CIFAR-100

Figure 5.2: **Ablation results for invariant layer.** Results for testing VGG like architecture with convolutional and invariant layers on several datasets. These graphs show the average top-1 accuracy on the validation set, and ± 1 standard deviation results (dark blue lines) for 5 runs. An architecture with ‘invX’ means the equivalent convolutional layer ‘convX’ from Table 5.5 was swapped for our proposed layer. The top row is the reference architecture using all convolutional layers. It appears that using a single invariant layer is the best option, and this will improve performance at many possible locations. The best position for the invariant layer appears to be around ‘convC’, which is just after the first pooling layer.

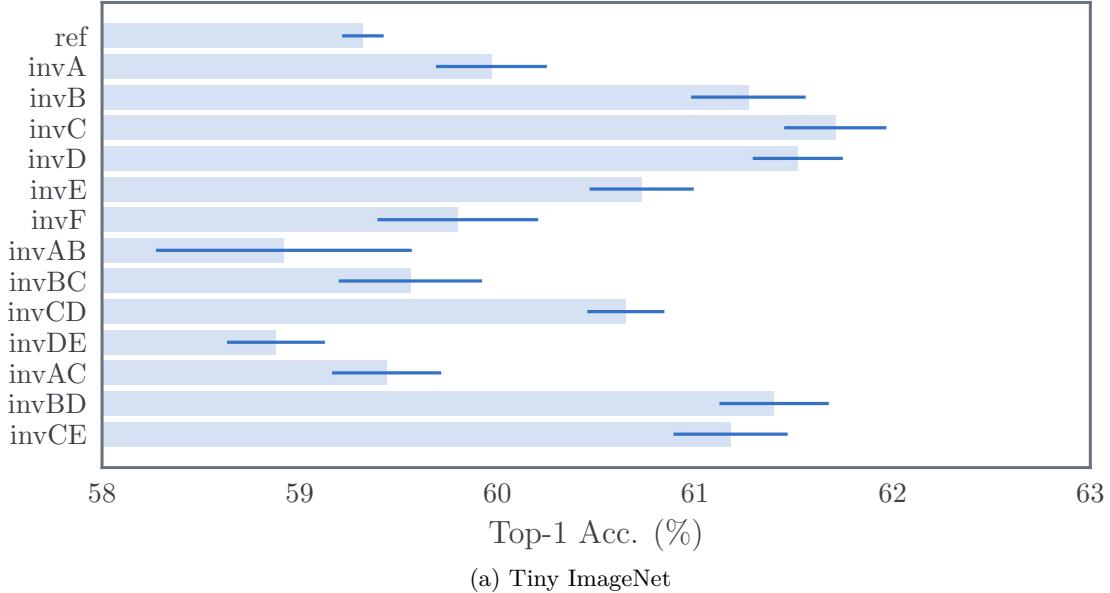


Figure 5.3: **Ablation results for Tiny ImageNet.** These graphs show the average top-1 accuracy on the validation set, and ± 1 standard deviation results (dark blue lines) for 3 runs. Like with CIFAR, it appears gains can be made by using the invariant layer at least once in the network. The best place for it appears to be around the first drop in resolution (which happens between ‘convB’ and ‘convC’).

The final classification layer is the same global 8×8 max pooling followed by a fully connected layer. In addition, we use dropout in these later convolutional layers with drop probability $p = 0.3$.

We compare a ScatterNet with no learning in between scattering orders (ScatNet A in [Table 5.6b](#)) to one with our proposal for a learned mixing matrix (ScatNet B in [Table 5.6c](#)). For completeness, we compare these architectures for a similarly shaped convolutional frontend ([Reference 1](#) in [Table 5.6a](#)).

We also test the hypothesis seen from [section 5.7](#) that scattering layers may work better after the first layer of a network. To do this, we put a small convolutional layer before the learnable ScatterNet front-end that takes the three colour inputs and outputs 16 new channels. Increasing the ScatterNet input to 16 channels means the default output size would be $16 \times 49 = 784$ channels. This is quite large, so we also use a system that uses a non-square A matrix for the two learnable invariant layers, keeping the output to 147 channels. We call this option ScatNet C (see [Table 5.7b](#)), and the full option with 784 output channels ScatNet D (see [Table 5.7c](#)). To compare the added benefit of this extra layer, we include a second reference architecture with the same small convolutional layer before it (see [Reference 2](#) in [Table 5.7a](#)).

Table 5.6: Hybrid ScatterNet models. Hybrid ScatterNet architectures used for experiments on CIFAR-10 and CIFAR-100. For the reference architecture we show the front and back ends. The ScatNet A and ScatNet B architectures use the same back end as the reference. N_c is the number of output classes; in our experiments we set the channel multiplier to be $C = 96$.

| (a) Reference 1 | |
|---|---------------------------|
| Layer | Act. Size |
| convA, $w \in \mathbb{R}^{21 \times 3 \times 3 \times 3}$ | $3 \times 32 \times 32$ |
| pool1, max pool 2×2 | $21 \times 32 \times 32$ |
| convB, $w \in \mathbb{R}^{147 \times 21 \times 3 \times 3}$ | $21 \times 16 \times 16$ |
| pool2, max pool 2×2 | $147 \times 16 \times 16$ |
| | $147 \times 8 \times 8$ |
| convC, $w \in \mathbb{R}^{2C \times 147 \times 3 \times 3}$ | $2C \times 8 \times 8$ |
| convD, $w \in \mathbb{R}^{2C \times 2C \times 3 \times 3}$ | $2C \times 8 \times 8$ |
| convE, $w \in \mathbb{R}^{4C \times 2C \times 3 \times 3}$ | $4C \times 8 \times 8$ |
| convF, $w \in \mathbb{R}^{4C \times 4C \times 3 \times 3}$ | $4C \times 8 \times 8$ |
| avg pool 8×8 | $4C$ |
| fc1, $4C \times N_c$ | N_c |

| (b) ScatNet A Front End | | (c) ScatNet B Front End | |
|-------------------------|--------------------------|---|--------------------------|
| Layer | Act. Size | Layer | Act. Size |
| scat1, no w | $3 \times 32 \times 32$ | inv1, $A \in \mathbb{R}^{21 \times 21}$ | $3 \times 32 \times 32$ |
| scat2, no w | $21 \times 16 \times 16$ | inv2, $A \in \mathbb{R}^{147 \times 147}$ | $21 \times 16 \times 16$ |
| | $147 \times 8 \times 8$ | | $147 \times 8 \times 8$ |

The results of our experiments on our three datasets are shown in [Table 5.8](#). The reference architecture does not perform as well as the reference architecture from [section 5.7](#) in these experiments, but recall that the network is different in these experiments. After two layers, the activation spatial size is already 8×8 .

ScatNet A underperforms the fully convolutional reference architecture with the same shape, suggesting that using a fixed ScatterNet as a front-end may be no better than simply learning the front-end. However, the proposed learnable ScatterNet outperforms both ScatNet A and the reference architecture on all datasets, suggesting that it is important to mix across the Scattering channels.

Table 5.7: **Hybrid ScatterNet models with convolutional layer first.** New front ends for 3 models. The two ScatNet models are the same learnable scatternet from Table 5.6c but with a small convolutional layer ('conv0') before it. ScatNet C ensures the same $147 \times 8 \times 8$ output size as the models in Table 5.6 but ScatNet D has a larger output size, allowing for the natural growth of a second-order ScatterNet model from C input channels to $49C$ output channels.

| (a) Reference 2 | |
|---|---------------------------|
| Layer | Act. Size |
| conv0, $w \in \mathbb{R}^{16 \times 3 \times 3 \times 3}$ | $3 \times 32 \times 32$ |
| convA, $w \in \mathbb{R}^{50 \times 16 \times 3 \times 3}$ | $16 \times 32 \times 32$ |
| pool1, max pooling 2×2 | $50 \times 32 \times 32$ |
| convB, $w \in \mathbb{R}^{147 \times 50 \times 3 \times 3}$ | $50 \times 16 \times 16$ |
| pool2, max pooling 2×2 | $147 \times 16 \times 16$ |
| | $147 \times 8 \times 8$ |

| (b) ScatNet C | | (c) ScatNet D | |
|---|--------------------------|---|---------------------------|
| Layer | Act. Size | Layer | Act. Size |
| conv0, $w \in \mathbb{R}^{16 \times 3 \times 3 \times 3}$ | $3 \times 32 \times 32$ | conv0, $w \in \mathbb{R}^{16 \times 3 \times 3 \times 3}$ | $3 \times 32 \times 32$ |
| inv1, $A \in \mathbb{R}^{50 \times 112}$ | $16 \times 32 \times 32$ | inv1, $A \in \mathbb{R}^{112 \times 112}$ | $16 \times 32 \times 32$ |
| inv2, $A \in \mathbb{R}^{147 \times 350}$ | $50 \times 16 \times 16$ | inv2, $A \in \mathbb{R}^{784 \times 784}$ | $112 \times 16 \times 16$ |
| | $147 \times 8 \times 8$ | | $784 \times 8 \times 8$ |

As we anticipated, adding a learned layer before the learnable ScatterNet improves things further, with ScatNet C and ScatNet D both having big improvements on ScatNets A and B, as well as outperforming the second reference convolutional architecture.

For comparison, we have also listed the performance of other architectures as reported by their authors in order of increasing complexity. Our proposed ScatNet D achieves comparable performance with the All Conv, VGG16 and FitNet architectures. The Deep[140] and Wide[141] ResNets perform best, but with very many more multiplies, parameters and layers.

A final note we would like to make is that the proposed hybrid ScatterNet networks architectures (particularly ScatNets C and ScatNets D) achieve near state-of-the-art performance even though they reduce the spatial size by a factor of 16 after only two layers. This meant that these networks took much less time to train than those in the previous section (20 minutes for ScatNets A and B compared to 2 to 4 hours for the networks in section 5.7).

Table 5.8: **Hybrid ScatterNet top-1 classification accuracies on CIFAR.** N_l is the number of learned convolutional layers, #param is the number of parameters, and #mults is the number of multiplies per $3 \times 32 \times 32$ image. An asterisk indicates that the value was estimated from the architecture description.

| Arch. Name | Arch. Properties | | | Top 1 Accuracies (%) | |
|------------------|------------------|---------|---------|----------------------|-----------|
| | N_l | #Mparam | #Mmults | CIFAR-10 | CIFAR-100 |
| Reference 1 | 6 | 2.65 | 165 | 91.2 | 70.0 |
| ScatNet A | 4 | 2.60 | 165 | 89.5 | 68.2 |
| ScatNet B | 6 | 2.64 | 167 | 91.5 | 70.5 |
| Reference 2 | 7 | 2.69 | 165 | 92.2 | 72.2 |
| ScatNet C | 7 | 2.64 | 251 | 92.6 | 72.7 |
| ScatNet D | 7 | 3.7 | 251 | 93.3 | 73.6 |
| All Conv[68] | 8 | 1.4 | 281* | 92.8 | 66.3 |
| VGG16[142] | 16 | 138* | 313* | 91.6 | - |
| FitNet[143] | 19 | 2.5 | 382 | 91.6 | 65.0 |
| ResNet-1001[140] | 1000 | 10.2 | 4453* | 95.1 | 77.3 |
| WRN-28-10[141] | 28 | 36.5 | 5900* | 96.1 | 81.2 |

Table 5.9: **Hybrid ScatterNet top-1 classification accuracies on Tiny ImageNet.** N_l is the number of learned convolutional layers, #param is the number of parameters, and #mults is the number of multiplies per $32 \times 32 \times 3$ image. An asterisk indicates that the value was estimated from the architecture description.

| Arch. Name | Top-1 Acc (%) |
|-------------|---------------|
| Reference1 | |
| ScatNet A | |
| ScatNet B | |
| Reference 2 | |
| ScatNet C | |
| ScatNet D | |

5.9 Conclusion

In this work we have proposed a new learnable scattering layer, dubbed the locally invariant convolutional layer, tying together ScatterNets and CNNs.

The invariant layer takes a single scale complex wavelet decomposition of the input. The bandpass coefficients are demodulated by using the complex modulus nonlinearity. The resulting magnitude coefficients are then mixed with the lowpass coefficients with a learnt mixing matrix (or 1×1 convolution). We did see some issues with using the invariant layer on its own when we tested it on MNIST, as the 1×1 convolution may not have been large enough to make complex shapes by separating the centres of the wavelets. When we used larger kernels, either in the mixing matrix A (see [subsection 5.6.1](#)) or in a convolutional layer after the invariant layer (see [section 5.7](#)) the performance improved.

Our ablation studies on a VGG-like CNN showed that the invariant layer can easily be shaped to allow it to drop in the place of a convolutional layer, theoretically saving on parameters and computation (see [section 5.7](#)). However, care must be taken when doing this, as our ablation study showed that the layer only improves upon regular convolution at certain depths. Typically, it seems wise to use the invariant layer early in the network, but after the first layer. The invariant layer naturally downsamples the input, so it works well when replacing a convolutional layer followed by pooling, or a pooling layer followed by a convolution.

We also tested the invariant layer on a hybrid ScatterNet architecture (see [section 5.8](#)). We saw that two invariant layers worked well as the first two layers of a deep CNN, but performed even better when used as the second and third layers, with a small learned convolutional before them. These hybrid ScatterNets downsample the input by a factor of 16 and increasing the channel dimension by 49. The reduced spatial size meant the networks were very quick to train, yet were able to achieve near state-of-the-art performance.

Chapter 6

Learning in the Wavelet Domain

In this final section of our work, we move away from the ScatterNet ideas from the previous chapters and instead look at using the wavelet domain as a new space in which to learn. With ScatterNets, complex wavelets are used to scatter the energy into different channels (corresponding to the different wavelet subbands), before the complex modulus demodulates the signal to low frequencies. These channels can then be mixed before scattering again (as we saw in the learnable ScatterNet), but successive use of such layers compounds the demodulation of signal energy towards zero frequency. We saw in the previous chapter that as a result, the modulus-based invariant layer worked best at the location of sample rate changes in a CNN. Most modern CNNs operate at only a handful of spatial resolutions, restricting the number of locations it may be useful in.

In this chapter, we introduce the *wavelet gain layer* which starts in a similar fashion to the ScatterNet – by taking the DTCWT of a multi-channel input. Next, instead of taking a complex modulus, we learn a complex gain for each subband in each input channel. A single value here can amplify or attenuate all the energy in one part of the frequency plane. Then, while still in the wavelet domain, we mix the different input channels *by subband* (e.g. all the 15° wavelet coefficients at a given scale are mixed together, but the 75° and 45° coefficients are not). We can then return to the pixel domain with the inverse wavelet transform. The shift-invariant properties of the DTCWT allow the wavelet coefficients to be changed without introducing sampling artefacts.

We also briefly explore the possibility of doing nonlinearities in the wavelet domain. Our ultimate goal is to connect multiple wavelet gain layers together with nonlinearities before returning to the pixel domain. See [subsection 6.2.4](#) for a more detailed description of this.

The proposed wavelet gain layer can be used in conjunction with regular convolutional layers, with a network moving into the wavelet or pixel space and learning filters in one that would be difficult to learn in the other.

Our experiments so far have shown some promise. We are able to learn complex wavelet gains and have found that the ReLU works well as a wavelet nonlinearity. We have found

that the wavelet gain layer works well at the beginning of a CNN but have not yet seen significant improvements for later layers.

6.1 Chapter Layout

We review some related work and notation in section 6.2 before describing the operation of our layer in section 6.3. In section 6.4, we describe some of the preliminary experiments and results we achieve by learning in the wavelet domain but returning to the pixel domain to perform nonlinearities. Section 6.5 describes expansions on this work to include nonlinearities in the wavelet domain and describes the preliminary results we have achieved so far.

6.2 Background

6.2.1 Related Work

Fujieda et. al. use a DWT in combination with a CNN to do texture classification and image annotation [144], [145]. They take a multiscale wavelet transform of the input image, combine the activations at each scale independently with learned weights, and feed these back into the network at locations where the activation resolution matches the subband resolution. The architecture block diagram is shown in Figure 6.1, taken from the original paper. They found that their ‘Wavelet-CNN’ could outperform competitive non-wavelet-based CNNs on both texture classification and image annotation.

Several works also use wavelets in deep neural networks for super-resolution [146] and for adding detail back into dense pixel-wise segmentation tasks [147]. These typically save wavelet coefficients and use them for the reconstruction phase.

In [148], Rippel, Snoek, and Adams parameterize filters in the DFT domain. Rather than having a pixel domain filter $w \in \mathbb{R}^{F \times C \times K \times K}$, they learn a set of Fourier coefficients $\hat{w} \in \mathbb{C}^{F \times C \times K \times \lceil K/2 \rceil}$ (the reduced spatial size is a result of enforcing that the inverse DFT of their filter to be real, so the parameterization is symmetric). On the forward pass of the neural network, they take the inverse DFT of \hat{w} to obtain w and then convolve this with the input x as a normal CNN would do.¹

Note that an important point should be laboured about reparameterizing filters in either the wavelet or Fourier domains: any invertible linear transform of the parameter space will not change parameter updates if a linear optimization scheme is used (for example standard GD, or SGD with momentum)². This is proved in appendix C. We make this point clear

¹The convolution may be done by taking both the image and filter back into the Fourier space but this is typically decided by the framework, which selects the optimal convolution strategy for the filter and input size. Note that there is not necessarily a saving to be gained by enforcing it to do convolution by product of FFTs, as the FFT size needed for the filter will likely be larger than $K \times K$, which would require resampling the coefficients.

²This fact is something that Rippel, Snoek, and Adams briefly allude to, but do not make clear.

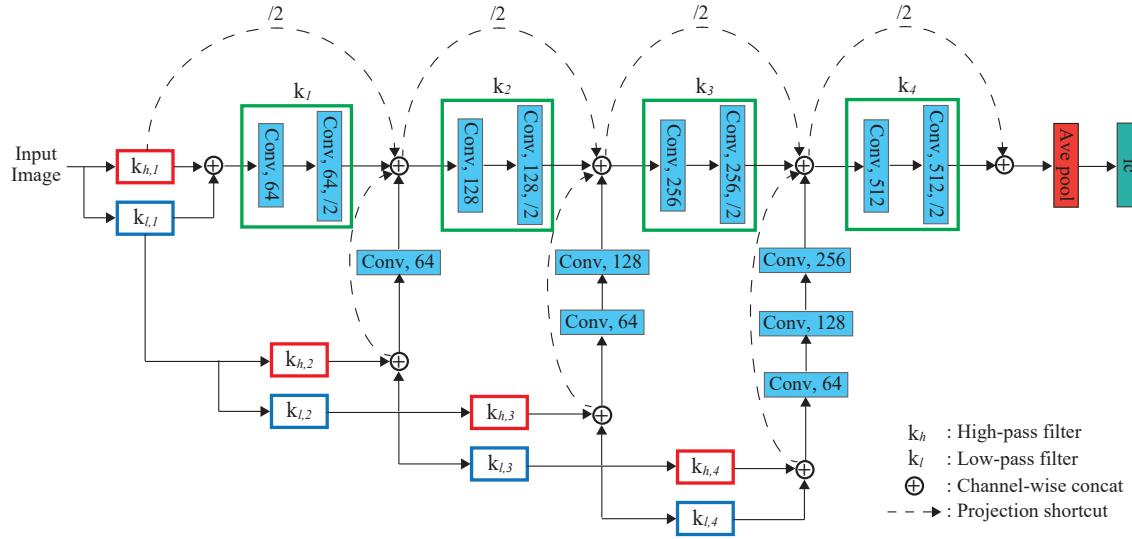


Figure 6.1: **Architecture using the DWT as a frontend to a CNN.** Figure 1 from [145]. Fujieda et. al. take a multiscale wavelet decomposition of the input before passing the input through a standard CNN. They learn convolutional layers independently on each subband and feed these back into the network at different depths, where the resolution of the subband and the network activations match.

as a natural extension to continue the work in [148] would be to parameterize filters in the wavelet domain, taking inverse transforms to the pixel domain and doing normal convolution.

The work presented in this chapter does not learn wavelet coefficients for convolutional filters to be applied in the pixel domain, rather we learn wavelet filters *and* do convolution in the wavelet domain too.

6.2.2 Notation

We make use of the 2-D Z -transform to simplify our analysis:

$$X(\mathbf{z}) = \sum_{n_1} \sum_{n_2} x[n_1, n_2] z_1^{-n_1} z_2^{-n_2} = \sum_{\mathbf{n}} x[\mathbf{n}] \mathbf{z}^{-\mathbf{n}} \quad (6.2.1)$$

As we are working with three-dimensional and four-dimensional arrays but are only doing convolution in two, we introduce a slightly modified 2-D Z -transform which includes the channel index c and the filter number f :

$$X(c, \mathbf{z}) = \sum_{n_1} \sum_{n_2} x[c, n_1, n_2] z_1^{-n_1} z_2^{-n_2} = \sum_{\mathbf{n}} x[c, \mathbf{n}] \mathbf{z}^{-\mathbf{n}} \quad (6.2.2)$$

$$H(f, c, \mathbf{z}) = \sum_{n_1} \sum_{n_2} h[f, c, n_1, n_2] z_1^{-n_1} z_2^{-n_2} = \sum_{\mathbf{n}} h[f, c, \mathbf{n}] \mathbf{z}^{-\mathbf{n}} \quad (6.2.3)$$

We then define the product of these new Z -transform signals to be the channel-wise convolution. E.g. for the 4-D filter $h[f, c, \mathbf{n}]$ with Z -transform $H(f, c, \mathbf{z})$ and the the 3-D signal $x[c, \mathbf{n}]$ with Z -transform $X(c, \mathbf{z})$, let us call the product of the two Z -transforms:

$$X(c, \mathbf{z})H(f, c, \mathbf{z}) = \sum_{\mathbf{n}} \left(\sum_{\mathbf{k}} h[f, c, \mathbf{n} - \mathbf{k}]x[c, \mathbf{k}] \right) \mathbf{z}^{-\mathbf{n}} \quad (6.2.4)$$

Recall from [subsection 2.4.1](#) that a typical convolutional layer in a standard CNN gets the next layer's output in a two-step process:

$$y^{(l+1)}[f, \mathbf{n}] = \sum_{c=0}^{C_l-1} x^{(l)}[c, \mathbf{n}] * h^{(l)}[f, c, \mathbf{n}] \quad (6.2.5)$$

$$x^{(l+1)}[f, \mathbf{n}] = \sigma(y^{(l+1)}[f, \mathbf{n}]) \quad (6.2.6)$$

If we define the nonlinearity σ_z to be the action of σ to each z -coefficient in the polynomial $Y(f, \mathbf{z})$, then we can rewrite (6.2.5) and (6.2.6) as:

$$Y^{(l+1)}(f, \mathbf{z}) = \sum_{c=0}^{C_l-1} X^{(l)}(c, \mathbf{z}) H^{(l)}(f, c, \mathbf{z}) \quad (6.2.7)$$

$$X^{(l+1)}(f, \mathbf{z}) = \sigma_z(Y^{(l+1)}(f, \mathbf{z})) \quad (6.2.8)$$

6.2.3 DT \mathbb{C} WT Notation

For this chapter, we will work with lots of DT \mathbb{C} WT coefficients so we define some slightly new notation here.

A J scale 2-D DT \mathbb{C} WT gives $6J + 1$ coefficients, 6 sets of complex bandpass coefficients for each scale (representing the oriented bands from 15 to 165 degrees) and 1 set of real lowpass (lp) coefficients.

$$\text{DT}\mathbb{C}\text{WT}_J(x) = \{u_{lp}, u_{j,k}\}_{1 \leq j \leq J, 0 \leq k < 6} \quad (6.2.9)$$

Each of these coefficients has size:

$$u_{lp} \in \mathbb{R}^{N \times C \times \frac{H}{2^{J-1}} \times \frac{W}{2^{J-1}}} \quad (6.2.10)$$

$$u_{j,k} \in \mathbb{C}^{N \times C \times \frac{H}{2^J} \times \frac{W}{2^J}} \quad (6.2.11)$$

Recall that the lowpass coefficients are twice as large as in a fully decimated transform due to the interleaving of the four lowpass terms in [Algorithm 3.3](#).

If we ever want to refer to all the subbands at a given scale, we will drop the k subscript and call them u_j . Likewise, u refers to the whole set of DT \mathbb{C} WT coefficients.

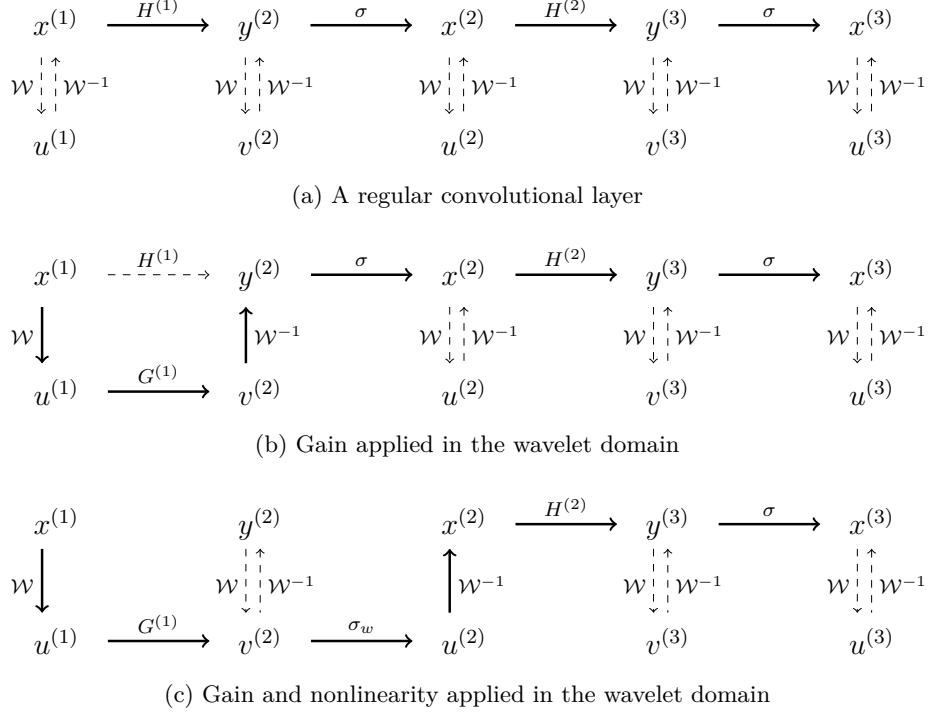


Figure 6.2: Proposed new forward pass in the wavelet domain. Two network layers with some possible options for processing. Solid lines denote the evaluation path and dashed lines indicate relationships. In (a) we see a regular convolutional neural network. We have included the dashed lines to make clear what we are denoting as u and v with respect to their equivalents x and y . In (b) we get to $y^{(2)}$ through a different path. First, we take the wavelet transform of $x^{(1)}$ to give $u^{(1)}$, apply a wavelet gain layer $G^{(1)}$, and take the inverse wavelet transform to give $y^{(2)}$. The dotted line for $H^{(1)}$ indicates that this path is no longer present. Note that there may not be any possible $G^{(1)}$ to make $y^{(2)}$ from (b) equal $y^{(2)}$ from (a). In (c) we have stayed in the wavelet domain longer and applied a wavelet nonlinearity σ_w to give $u^{(2)}$. We then return to the pixel domain to give $x^{(2)}$ and continue on from there in the pixel domain.

6.2.4 Learning in Multiple Spaces

At the beginning of each layer l of a neural network, we have the activations $x^{(l)}$. Naturally, all of these activations have their equivalent wavelet coefficients $u^{(l)}$.

From (6.2.5), convolutional layers also have intermediate activations $y^{(l)}$. Let us discern these from the x coefficients and modify (6.2.9) to say the DTCWT of $y^{(l)}$ gives $v^{(l)}$.

We now propose the *wavelet gain layer* G . The gain layer G can be used instead of a convolutional layer. It is designed to work on the wavelet coefficients of an activation u , to give wavelet domain outputs v .

This can be seen as breaking the convolutional path in Figure 6.2 and taking a new route to get the next layer's coefficients. From here, we can return to the pixel domain by taking the corresponding inverse wavelet transform \mathcal{W}^{-1} . Alternatively, we can stay in the wavelet

domain and apply wavelet-based nonlinearities, σ_{lp} and σ_{bp} for the lowpass and highpass coefficients respectively, to give $u^{(l+1)}$.

Ultimately we would like to explore architecture design with arbitrary sections in the wavelet and pixel domain, but to do this we must first explore:

1. **How effective is a wavelet gain layer G at replacing a standard convolutional layer H ?**
2. **What are effective wavelet nonlinearities σ_{lp} and σ_{bp} ?**

6.3 The DT \mathbb{C} WT Gain Layer

To do the mixing across the C_l channels at each subband, giving C_{l+1} output channels, we introduce the learnable filters g_{lp} , $g_{j,k}$:

$$g_{lp} \in \mathbb{R}^{C_{l+1} \times C_l \times k_{lp} \times k_{lp}} \quad (6.3.1)$$

$$g_{1,1} \in \mathbb{C}^{C_{l+1} \times C_l \times k_1 \times k_1} \quad (6.3.2)$$

$$g_{1,2} \in \mathbb{C}^{C_{l+1} \times C_l \times k_1 \times k_1} \quad (6.3.3)$$

⋮

$$g_{J,6} \in \mathbb{C}^{C_{l+1} \times C_l \times k_J \times k_J} \quad (6.3.4)$$

where the k_j are the sizes of the mixing kernels. These could be 1×1 for simple gain control, or could be larger, say 3×3 , to do more complex filtering on the subbands. Importantly, we can select the support size differently for each subband.

With these gains we define the action of the gain layer $v = Gu$ to be:

$$v_{lp}[f, \mathbf{n}] = \sum_{c=0}^{C_l-1} u_{lp}[c, \mathbf{n}] * g_{lp}[f, c, \mathbf{n}] \quad (6.3.5)$$

$$v_{1,1}[f, \mathbf{n}] = \sum_{c=0}^{C_l-1} u_{1,1}[c, \mathbf{n}] * g_{1,1}[f, c, \mathbf{n}] \quad (6.3.6)$$

$$v_{1,2}[f, \mathbf{n}] = \sum_{c=0}^{C_l-1} u_{1,2}[c, \mathbf{n}] * g_{1,2}[f, c, \mathbf{n}] \quad (6.3.7)$$

⋮

$$v_{J,6}[f, \mathbf{n}] = \sum_{c=0}^{C_l-1} u_{J,6}[c, \mathbf{n}] * g_{J,6}[f, c, \mathbf{n}] \quad (6.3.8)$$

To avoid ambiguity with complex conjugates we remind ourselves that for complex signals a, b the convolution $a * b$ is defined as $(a_r * b_r - a_i * b_i) + j(a_r * b_i + a_i * b_r)$ (see [section E.1](#)).

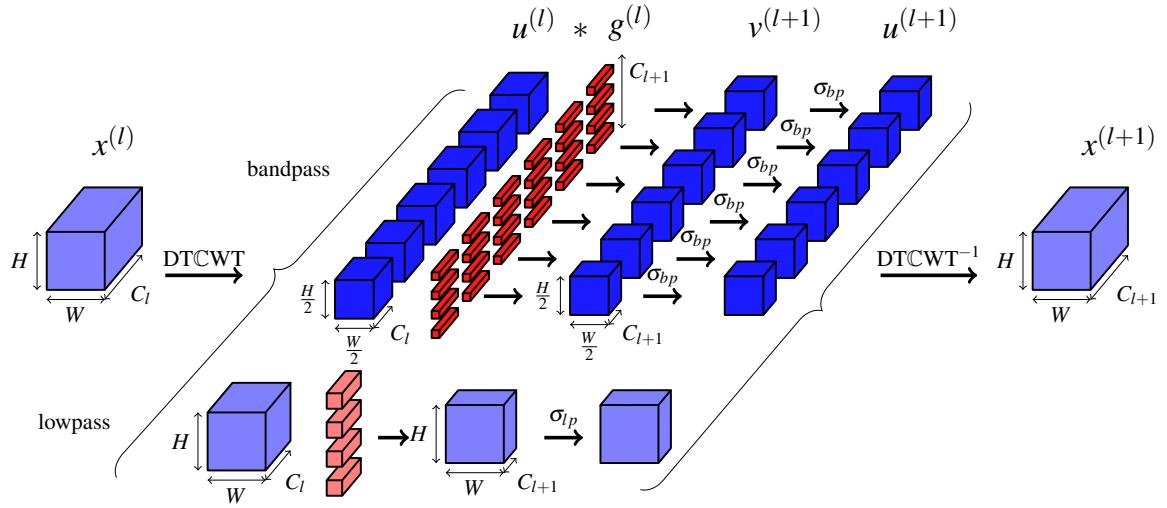


Figure 6.3: Diagram of proposed method to learn in the wavelet domain. Activations are shaded blue and learned parameters red. Deeper shades of blue and red indicate complex valued activations/weights, and lighter values indicate real valued activations/weights. The input $x^{(l)} \in \mathbb{R}^{C_l \times H \times W}$ is taken into the wavelet domain (here $J = 1$) and each subband is mixed independently with C_{l+1} sets of convolutional filters. After mixing, a possible wavelet nonlinearity σ_w is applied to the subbands, before returning to the pixel domain with an inverse wavelet transform. Note the similarity to the regular convolutional layer in [Figure 2.7](#).

The action of the gain layer with only a single-scale wavelet transform, $J = 1$, is shown in [Figure 6.3](#).

6.3.1 The Output

Due to the shift invariant properties of the DTCWT, each wavelet subband has a unique transfer function which is almost free of aliasing (see [Theorem D.1](#)). If we do complex convolution of the wavelet coefficients u with gains g as described in [\(6.3.5\)](#) - [\(6.3.8\)](#), then we preserve the shift invariant properties (see [Theorem D.2](#) and [Theorem D.3](#)) and the inverse DTCWT of the outputs v are free from aliasing.

We can do a complex multiply of the subband coefficients u with gains g by using the filter bank diagram shown in [Figure 6.4a](#). Note that despite the resemblance to many filter bank diagrams for fully decimated DWTS, [Figure 6.4a](#) is different. The top rung corresponds to the real part of a subband and the bottom specifies the imaginary part.

We consider the output from a *single* complex subband. The complex gain for this subband is $G = G_r + jG_i$ where j is the square root of negative one. Let us call the analysis filters $A = A_r + jA_i$ and the synthesis filters $S = S_r - jS_i$ (these are normally called H and G , but we keep those letters reserved for the CNN and gain layer filters). For a single-channel

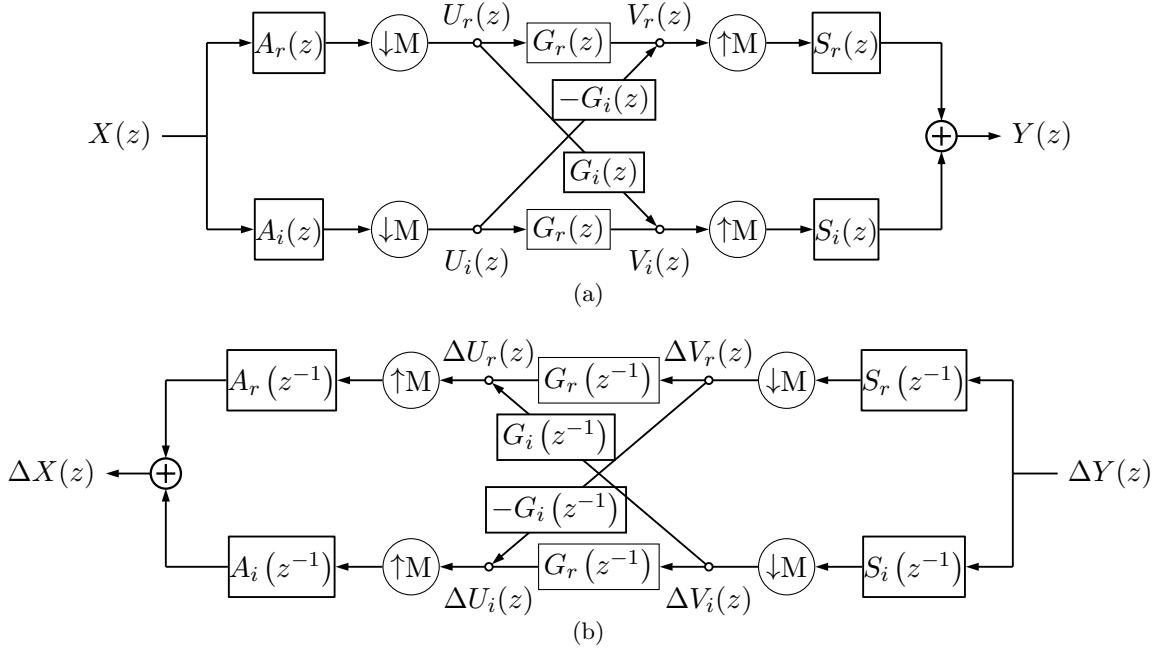


Figure 6.4: **Forward and backward filter bank diagrams for DTCWT gain layer.** Based on Figure 4 in [19]. Ignoring the G gains, the top and bottom paths (through A_r, S_r and A_i, S_i respectively) make up the real and imaginary parts for *one subband* of the dual tree system. Combined, $A_r + jA_i$ and $S_r - jS_i$ make the complex filters necessary to have support on one side of the Fourier domain (see Figure 6.5). Adding in the complex gain $G_r + jG_i$, we can now attenuate/shape the impulse response in each of the subbands. To allow for learning, we need backpropagation. The bottom diagram indicates how to pass gradients $\Delta Y(z)$ through the layer. Note that upsampling has become downsampling, and convolution has become convolution with the time reverse of the filter (represented by z^{-1} terms).

input in 1-D, the output of this layer with decimation and interpolation is:

$$Y(z) = \frac{2}{M} X(z) \left[G_r(z^M) (A_r(z)S_r(z) + A_i(z)S_i(z)) + G_i(z^M) (A_r(z)S_i(z) - A_i(z)S_r(z)) \right] \quad (6.3.9)$$

Again, see [appendix D](#) for the derivation which proves that the aliasing terms caused by the downsampling by M are (largely) eliminated.

This expands to the 2-D case with multiple channels as:

$$Y(f, \mathbf{z}) = \sum_{c=0}^{C_l-1} \frac{2}{M} X(c, \mathbf{z}) \left[G_r(f, c, \mathbf{z}^M) (A_r(\mathbf{z})S_r(\mathbf{z}) + A_i(\mathbf{z})S_i(\mathbf{z})) + G_i(f, c, \mathbf{z}^M) (A_r(\mathbf{z})S_i(\mathbf{z}) - A_i(\mathbf{z})S_r(\mathbf{z})) \right] \quad (6.3.10)$$

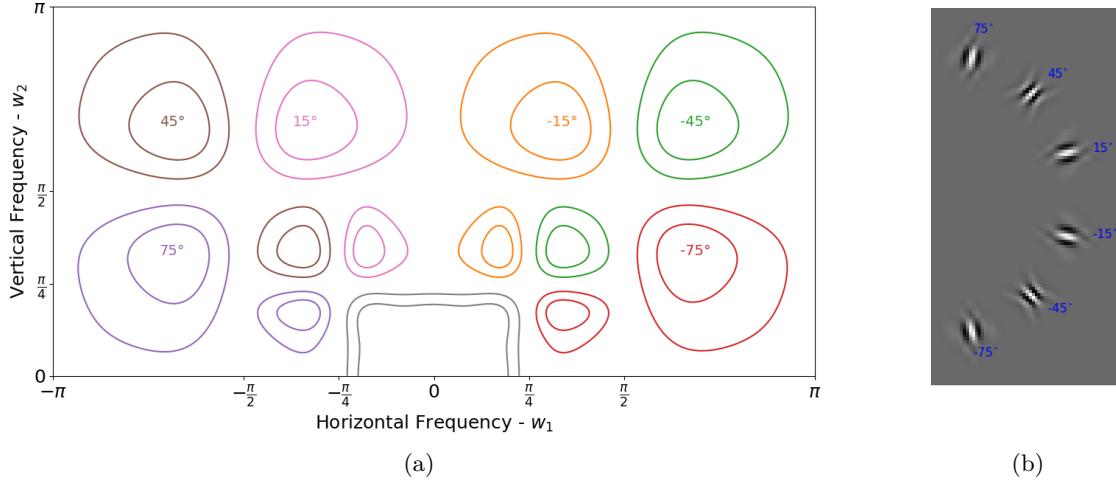


Figure 6.5: DTcCWT subbands. (a) -1dB and -3dB contour plots showing the support in the Fourier domain of the 6 subbands of the DTcCWT at scales 1 and 2, and the scale 2 lowpass. These are the product of the single side-band filters $P(\mathbf{z})$ and $Q(\mathbf{z})$ from Theorem D.1, or half of the support of the double side-band filters $A_r S_r + A_i S_i$ and $A_r S_i - A_i S_r$ from (6.3.10). (b) The pixel domain impulse responses for the second scale wavelets. The Hilbert pair for each wavelet is the underlying sinusoid phase shifted by 90 degrees.

The complex gain G has a real and imaginary part. The real term G_r modifies the subband gain $A_r S_r + A_i S_i$ and the imaginary term G_i modifies its Hilbert Pair $A_r S_i - A_i S_r$. Figure 6.5 shows the contour plots for the frequency support of each of these subbands.

Now we consider the sum of all different subbands. As the operations are all linear, the full output y is simply the sum of all the y 's from individual subbands. The complex gains g adjust the gain and phase of each subband independently. The magnitude of each element controls the amplitude of the frequency response in the region of that subband, while its phase controls the phase of the response and thus modifies the detailed wave-shape (e.g. the locations of its zero crossings).

6.3.2 Backpropagation

Assume we already have access to the quantity $\Delta Y(z)$ (this is the input to the backwards pass). Figure 6.4b illustrates the backpropagation procedure.

We recall the passthrough and update equations for a convolutional block from (2.4.12) and (2.4.15):

$$\frac{\partial L}{\partial x_{c,k_1,k_2}} = \sum_f \Delta y[f, \mathbf{n}] \star h_c[f, \mathbf{n}] \quad (6.3.11)$$

$$\frac{\partial L}{\partial h_{f,c,k_1,k_2}} = \Delta y[f, \mathbf{n}] \star x[c, \mathbf{n}] \quad (6.3.12)$$

In terms of modified Z -transforms, this becomes:

$$\Delta X(c, \mathbf{z}) = \sum_f \Delta Y(f, \mathbf{z}) H'[c, f, \mathbf{z}^{-1}] \quad (6.3.13)$$

$$\Delta H(f, c, \mathbf{z}) = \sum_f \Delta Y(f, \mathbf{z}) X[c, \mathbf{z}^{-1}] \quad (6.3.14)$$

where H' is H transposed in the filter and channel dimensions, and \mathbf{z}^{-1} indicates the filter was mirror-imaged in the spatial dimension. The simplified, single-channel, single filter, 1-D version of this is:

$$\Delta X(z) = \Delta Y(z) H(z^{-1}) \quad (6.3.15)$$

$$\Delta H(z) = \Delta Y(z) X(z^{-1}) \quad (6.3.16)$$

If H were complex, the first term in [Equation 6.3.15](#) would be $\bar{H}(1/\bar{z})$, but as each individual block in the DT \mathbb{C} WT of [Figure 6.4](#) is purely real, we can use the simpler form $H(z^{-1})$.

We calculate $\Delta V_r(z)$ and $\Delta V_i(z)$ by backpropagating $\Delta Y(z)$ through the inverse DT \mathbb{C} WT. This is the same as doing the forward DTCWT on $\Delta Y(z)$ with the synthesis and analysis filters swapped and time-reversed. Then the weight update equations are:

$$\Delta G_r(z) = \Delta V_r(z) U_r(z^{-1}) + \Delta V_i(z) U_i(z^{-1}) \quad (6.3.17)$$

$$\Delta G_i(z) = -\Delta V_r(z) U_i(z^{-1}) + \Delta V_i(z) U_r(z^{-1}) \quad (6.3.18)$$

The passthrough equations have similar form to [\(6.3.10\)](#):

$$\Delta X(z) = \frac{2\Delta Y(z)}{M} [G_r(z^{-M}) (A_r(z) S_r(z) + A_i(z) S_i(z)) + G_i(z^{-M}) (A_r(z) S_i(z) - A_i(z) S_r(z))] \quad (6.3.19)$$

6.3.3 Examples

[Figure 6.6](#) shows example impulse responses of the DTCWT gain layer. For comparison, we also show similar ‘impulse responses’ for a gain layer done in the DWT domain³. The DWT outputs come from three random variables: a 1×1 convolutional weight applied to each of the low-high, high-low and high-high subbands. The DT \mathbb{C} WT outputs come from twelve random variables, again a 1×1 convolutional weight, but now applied to six complex subbands.

To test the space of generated shapes by a vector gain layer gain g_1 , we generate N random vectors of length 12, with each entry taken from a Gaussian distribution with zero

³Modifying DWT coefficients causes a loss of the alias cancellation properties so these are not true impulse responses.

mean and unit variance. We then generate the equivalent point spread functions from (6.3.10) for the N different instances and measure their normalized cross-correlation. We then sort the values and compare the distribution to a set of N random vectors with k degrees of freedom.

Our experiments show that the distribution for the DT^CWT gain layer matches random vectors with roughly 11.5 degrees of freedom (c.f. the 12 variables the layer has). Similarly for the DWT, the normalized cross-correlation matches the distribution for random vectors with roughly 2.8 degrees of freedom (c.f. 3 random variables in the layer). This is particularly reassuring for the DT^CWT as it is showing that there is still representative power despite the redundancy of the transform.

6.3.4 Implementation Details

Before analyzing its performance, we compare the implementation properties of our proposed layer to a standard convolutional layer.

6.3.4.1 Parameter Memory Cost

A standard convolutional layer with C_l input channels, C_{l+1} output channels and kernel size $k \times k$ has $k^2 C_l C_{l+1}$ parameters, with $k = 3$ or $k = 5$ common choices for the spatial size.

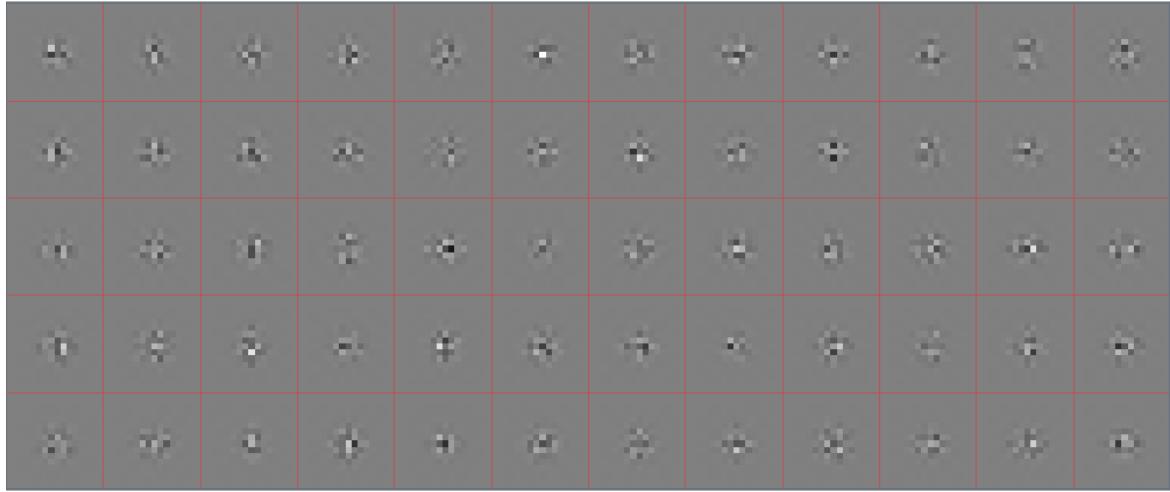
$$\#\text{conv params} = k^2 C_l C_{l+1} \quad (6.3.20)$$

We must choose the spatial sizes of both the lowpass and bandpass mixing kernels. In our work, we are somewhat limited in how large we would like to set the bandpass spatial size, as every extra pixel of support requires $2 \times 6 = 12$ extra parameters. For this reason, we almost always set it to have support 1×1 . The lowpass gains are less costly, and we are free to set them to size $k_{lp} \times k_{lp}$ (with $k_{lp} = 1, 3, 5$ in many of our experiments). Further, due to the size of the datasets we test on, we typically limit ourselves initially to only considering a single scale. If we wish, we can decompose the input into more scales, resulting in a larger net area of effect. In particular, it may be useful to do a two-scale transform and discard the first scale coefficients. This does not increase the number of gains to learn but changes the position of the bands in the frequency space.

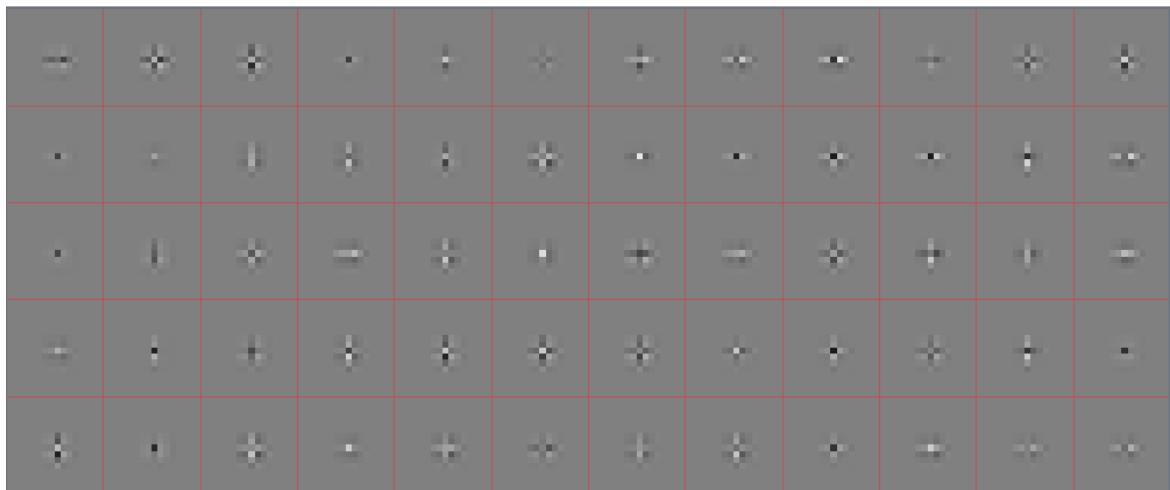
The number of parameters for the gain layer with $k_{lp} = 1$ is:

$$\#\text{params} = (2 \times 6 + 1)C_l C_{l+1} = 13C_l C_{l+1} \quad (6.3.21)$$

This is slightly larger than the $9C_l C_{l+1}$ parameters used in a standard 3×3 convolution, but as Figure 6.6 shows, the spatial support of the full filter is larger than an equivalent one parameterized in the filter domain. If $k_{lp} = 3$ then we would have $21C_l C_{l+1}$ parameters, slightly fewer than the $25C_l C_{l+1}$ of a 5×5 convolution.



(a)



(b)

Figure 6.6: Example outputs from an impulse input for the proposed gain layers. Example outputs $y = \mathcal{W}^{-1}G\mathcal{W}x$ for an impulse x for the DTCWT gain layer and for a similarly designed DWT gain layer. (a) shows the output y for a DTCWT based system. $g_{l_p} = 0$ and g_1 has spatial size 1×1 . The 12 values in g_1 are independently sampled from a random normal of variance 1. The 60 samples come from 60 different random draws of the weights. (b) shows the outputs y when x is an impulse and \mathcal{W} is the DWT with a ‘db2’ wavelet family. Here 3 random numbers are generated for the g_1 coefficients. The strong horizontal and vertical properties of the DWT can clearly be seen in comparison to the much freer DTCWT.

6.3.4.2 Activation Memory Cost

A standard convolutional layer needs to save the activation $x^{(l)}$ to convolve with the back-propagated gradient $\frac{\partial L}{\partial y^{(l+1)}}$ on the backwards pass (to give $\frac{\partial L}{\partial w^{(l)}}$). For an input with C_l channels of spatial size $H \times W$, this means

$$\#\text{conv floats} = HWC_l \quad (6.3.22)$$

Our layer requires us to save the wavelet coefficients u_{lp} and $u_{j,k}$ for updating the g terms as in (6.3.17) and (6.3.18). For the 4 : 1 redundant DT \mathbb{C} WT, this requires:

$$\#\text{DT}\mathbb{C}\text{WT floats} = 4HWC_l \quad (6.3.23)$$

to be saved for the backwards pass. You can see this difference from the difference in the block diagrams in Figure 6.3.

Note that a single scale DT \mathbb{C} WT gain layer requires 16/7 times as many floats to be saved as compared to the invariant layer of the previous chapter. The extra cost of this comes from two things. Firstly, we keep the real and imaginary components for the bandpass (as opposed to only the magnitude), meaning we need $3HWC_l$ floats, rather than $\frac{3}{2}HWC_l$. Additionally, the lowpass was downsampled in the previous chapter, requiring only $\frac{1}{4}HWC_l$, whereas we keep the full sample rate, costing HWC_l .

If memory is an issue and the computation of the DT \mathbb{C} WT is very fast, then we only need to save the $x^{(l)}$ coefficients and can calculate the u 's on the fly during the backwards pass. Note that a two-scale DT \mathbb{C} WT gain layer would still only require $4HWC_l$ floats.

6.3.4.3 Computational Cost

A standard convolutional layer with kernel size $k \times k$ needs k^2C_{l+1} multiplies per input pixel (of which there are $C_l \times H \times W$).

For the DT \mathbb{C} WT, the overhead calculations are the same as in subsection 5.5.3, so we will omit their derivation here. The mixing is however different, requiring complex convolution for the bandpass coefficients, and convolution over a higher resolution lowpass. The bandpass has one quarter spatial resolution at the first scale, but this is offset by the 4 : 1 cost of complex multiplies compared to real multiplies. Again assuming we have set $J = 1$ and $k_{lp} = 1$ then the total cost for the gain layer is:

$$\#\text{mults/pixel} = \underbrace{\frac{6 \times 4}{4}C_{l+1}}_{\text{bandpass}} + \underbrace{C_{l+1}}_{\text{lowpass}} + \underbrace{36}_{\text{DT}\mathbb{C}\text{WT}} + \underbrace{36}_{\text{DT}\mathbb{C}\text{WT}^{-1}} = 7C_{l+1} + 72 \quad (6.3.24)$$

which is marginally smaller than the $9C_{l+1}$ of a 3×3 convolutional layer (if $C_{l+1} > 36$).

6.3.4.4 Parameter Initialization

For both layer types we use the Glorot Initialization scheme [56] with $a = 1$:

$$g_{ij} \sim U \left[-\sqrt{\frac{6}{(C_l + C_{l+1})k^2}}, \sqrt{\frac{6}{(C_l + C_{l+1})k^2}} \right] \quad (6.3.25)$$

where k is the kernel size.

6.4 Gain Layer Experiments

Before we explore the possibilities and performance of using a nonlinearity in the wavelet domain, let us present some experiments and results for the wavelet gain layer where we do nonlinearities purely in the spatial domain, as in a conventional CNN layer. This is the first objective in [subsection 6.2.4](#), comparing G to H .

6.4.1 CNN activation regression

One of the early inspirations for using wavelets in CNNs was the visualizations of the first layer filters learned in AlexNet. These 11×11 colour filters (see [Figure 1.2a](#)) look very much like a 2-D oriented wavelet transform.

So how well can the gain layer emulate the action of this layer? How would it compare to trying to use a reduced size convolutional kernel to learn the action of the layer?

Let us call the action of our target layer H_0 , a learned convolutional layer H and our gain layer G . We assume that we do not have direct access to H_0 but only the convolved outputs $Y = H_0X$. Then, we would like to solve:

$$\arg \min_H (Y - HX)^2, \text{ s.t. } h[c, \mathbf{n}] = 0, \forall \mathbf{n} \notin \mathcal{R} \quad (6.4.1)$$

$$\arg \min_G (Y - \mathcal{W}^{-1}G\mathcal{W}X)^2, \text{ s.t. } g_{j,k}[c, \mathbf{n}] = 0, \forall \mathbf{n} \notin \mathcal{R}' \quad (6.4.2)$$

for some support regions $\mathcal{R}, \mathcal{R}'$. E.g. \mathcal{R} is a 3×3 or 5×5 block, and similarly \mathcal{R}' is a 1×1 region in each subband.

(6.4.1) and (6.4.2) are both regression problems convex in the parameters of H and G , with many possible ways to solve. We are not worried with the optimization procedure chosen here, but of the final distances $\|Y - HX\|$ and $\|Y - \mathcal{W}^{-1}G\mathcal{W}X\|$ (or equivalently, their squares). We choose to find H and G by gradient descent, using the validation set for ImageNet as the data input-output pair (X, Y) . After 3–5 epochs, both H and G typically settle into their global minimum. Because of the large size of the input filters, we allow for both a $J = 1$ and $J = 2$ scale gain layer but only learn weights at the lowest frequency

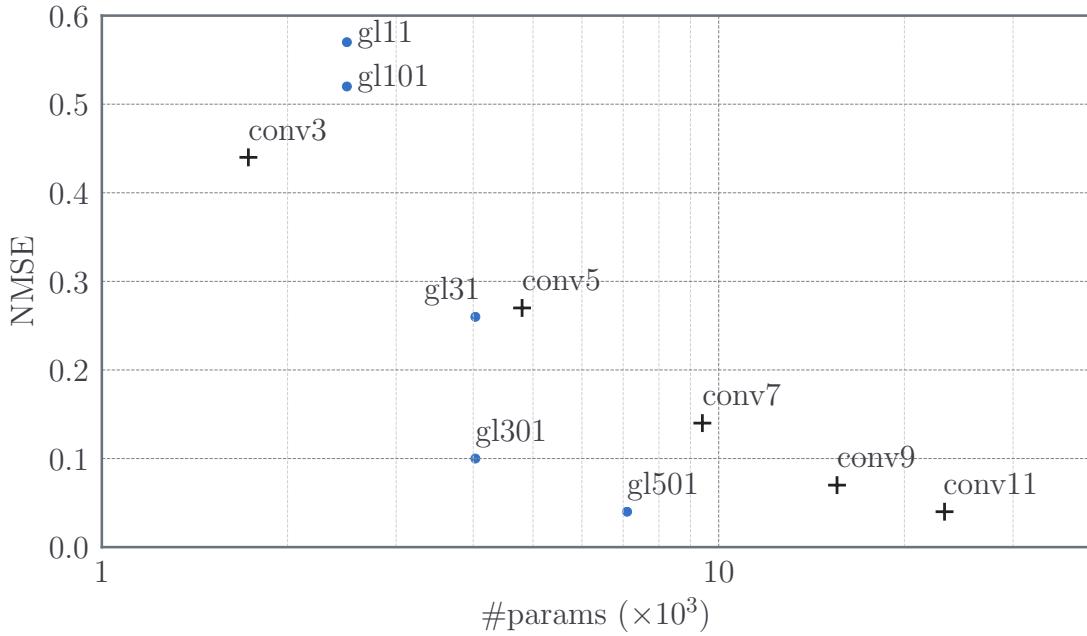


Figure 6.7: Normalized mean squared error for conv layer and wavelet gain layer regression. We minimize (6.4.1) and (6.4.2) on the ImageNet validation set, where the target is made from convolving the input with AlexNet’s first layer filters. This plot shows the final NMSE score compared to the number of learnable parameters. The original conv layer has spatial support 11×11 . The four points labelled ‘convn’ correspond to filters with $n \times n$ spatial support. The four points labelled ‘glabc’ correspond to two-scale gain layers with $a \times a$ support in the lowpass, $b \times b$ spatial support in the first scale, and $c \times c$ spatial support in the second scale. The gain layer can regress to the AlexNet filters quite capably. In this example, it is important to have at least 3×3 lowpass support for the gain layer, and the second scale coefficients are more important than the first scale.

bandpass (i.e. for a 2 scale gain layer, we discard the first scale highpass outputs and only learn g_2).

After training, we report the final normalized mean squared errors between the target values Y and the estimated outputs \hat{Y} , defined as:

$$NMSE = \frac{1}{N} \sum_{n=1}^N \frac{\mathbb{E}[(Y - \hat{Y})^2]}{\mathbb{E}[Y^2]} \quad (6.4.3)$$

The resulting NMSEs are shown in Figure 6.7. A label ‘glab’ indicates a single scale gain layer with $a \times a$ support in the lowpass and $b \times b$ support in the highpass; a label ‘glabc’ indicates a two-scale gain layer with $a \times a$ support in the lowpass, $b \times b$ support in the scale 1 highpass and $c \times c$ support in the scale 2 bandpass gains.

This figure shows several interesting yet unsurprising things. Firstly, bigger lowpass support is very helpful – see the difference between gl101, gl301, and gl501, 3 instances that

only vary in the size of the support of their lowpass filter g_{lp} . Additionally, the second scale coefficients appear more useful than the first scale – see the difference between gl31 and gl301, two instances that have the same number of parameters, but gl31 has g_1 with non-zero support and gl301 has g_2 with non-zero support.

This experiment shows that a gain layer is a good representation for a set of filters like the AlexNet first layer. The next section looks at how well it performs at deeper layers.

6.4.2 Ablation Studies

[Figure 6.7](#) is a useful guide on how the gain layer might be placed in a deep CNN. gl11 (a gain layer with a 1×1 lowpass kernel and a 1×1 bandpass kernel at the first scale), gl101 (same as gl11 but no gain at first scale and 1×1 at second scale), and conv3 all achieve similar NMSEs. Additionally, gl31, gl301, and conv5 all achieve similar NMSEs.

6.4.2.1 Small Kernel Ablation

Most modern CNNs are built with small 3×3 kernels, which we believe are not the best use for the gain layer, built from large support wavelets. For this reason, we deviate from the ablation study done in the previous chapter and build a **shallow**er network with **larger** kernel sizes.

For completeness, we also ran ablation tests on the same deeper network with small kernels used in [chapter 5](#) and include the results in [appendix F](#).

6.4.2.2 Large Kernel Ablation

In this experiment, we build a three-layer CNN with 5×5 convolutional kernels, described in [Table 6.1](#). To help differentiate with the small kernel network introduced in the ablation study of the previous chapter, we have labelled the convolutions here ‘conv1’, ‘conv2’ and ‘conv3’ (as opposed to ‘convA’, ‘convB’, ‘convC’, …).

We test the difference in accuracy achieved by replacing each of the three convolution layers with gl31. Although the gain layers with no gain in the first scale and gain in the second scale (e.g. gl301, gl501) performed better than those with only gain in the first scale (e.g. gl31, gl51) in [subsection 6.4.1](#), we saw them perform consistently worse in the following ablation studies. This is not surprising, as a 5×5 convolutional kernel is too small compared to the width of the central lobe of scale 2 wavelets. For ease of presentation, we only show the results from the single scale gain layer gl31.

On the two CIFAR datasets, we train for 120 epochs, decaying learning rate by a factor of 0.2 at 60, 80 and 100 epochs, and for the Tiny ImageNet dataset, we train for 45 epochs, decaying learning rate at 18, 30 and 40 epochs. We set ℓ_2 weight decay of 10^{-4} for the real gains in the system, and ℓ_1 weight decay of 10^{-5} for the complex gains in the system. See

Table 6.1: **Ablation Base Architecture.** Reference architecture used for experiments on CIFAR-10, CIFAR-100. The activation size rows are offset from the layer description rows to convey the input and output shapes. Unlike Table 5.5, this architecture is shallower and uses 5×5 convolutional kernels as a base. C is a hyperparameter that controls the network width, we use $C = 64$ for our tests. The Tiny ImageNet architecture is very similar but with larger activation sizes and one more convolutional layer ‘conv4’.

| Activation Size | Reference Arch. | Alternate Arch. |
|--------------------------|--|---|
| $3 \times 32 \times 32$ | conv1, $w \in \mathbb{R}^{C \times 3 \times 5 \times 5}$ | or gain1, $g_{lp} \in \mathbb{R}^{C \times 3 \times 3 \times 3}$, $g_1 \in \mathbb{C}^{C \times 6 \times 3 \times 1 \times 1}$ |
| $C \times 32 \times 32$ | pool1, max pool 2×2 | |
| $C \times 16 \times 16$ | conv2, $w \in \mathbb{R}^{2C \times C \times 5 \times 5}$ | or gain2, $g_{lp} \in \mathbb{R}^{2C \times C \times 3 \times 3}$, $g_1 \in \mathbb{C}^{2C \times 6 \times C \times 1 \times 1}$ |
| $2C \times 16 \times 16$ | pool2, max pool 2×2 | |
| $2C \times 8 \times 8$ | conv3, $w \in \mathbb{R}^{4C \times 2C \times 5 \times 5}$ | or gain3, $g_{lp} \in \mathbb{R}^{4C \times 2C \times 3 \times 3}$, $g_1 \in \mathbb{C}^{4C \times 6 \times 2C \times 1 \times 1}$ |
| $4C \times 8 \times 8$ | avg, 8×8 average pool | |
| $4C \times 1 \times 1$ | fc1, fully connected | |
| 10, 100 | | |

section E.2 for information on how we handle regularizing complex gains. The experiment code is available at [33].

The results of various combinations for our three datasets are shown in Figure 6.8. Note that as before, swapping ‘conv1’ with a gain layer is marked by ‘gain1’, and swapping the first two conv layers with two gain layers is marked by ‘gain1_2’ and so forth.

The results are not too promising. Across all three datasets, changing a convolutional layer for a gain layer of similar number of parameters results in a small decrease in accuracy at all depths, and the more layers swapped out the more this degradation compounds.

6.4.3 Network Analysis

It is nonetheless interesting to see that a network with only gain layers (‘gain1_2_3’) can achieve accuracies within a couple of percentage points of a purely convolutional architecture.

In this section, we look at some of the properties of the ‘gain1_2_3’ for CIFAR-100 and compare them to the reference architecture.

6.4.3.1 Bandpass Coefficients

When analyzing the ‘gain1_2_3’ architecture, the most noticeable thing is the distribution of the bandpass gain magnitudes. Figure 6.9a shows these for the second gain layer, gain2. Of the $64 \times 128 = 8192$ complex coefficients most have very small magnitude, in particular, the diagonal wavelet gains. The disparity between the diagonal and horizontal/vertical

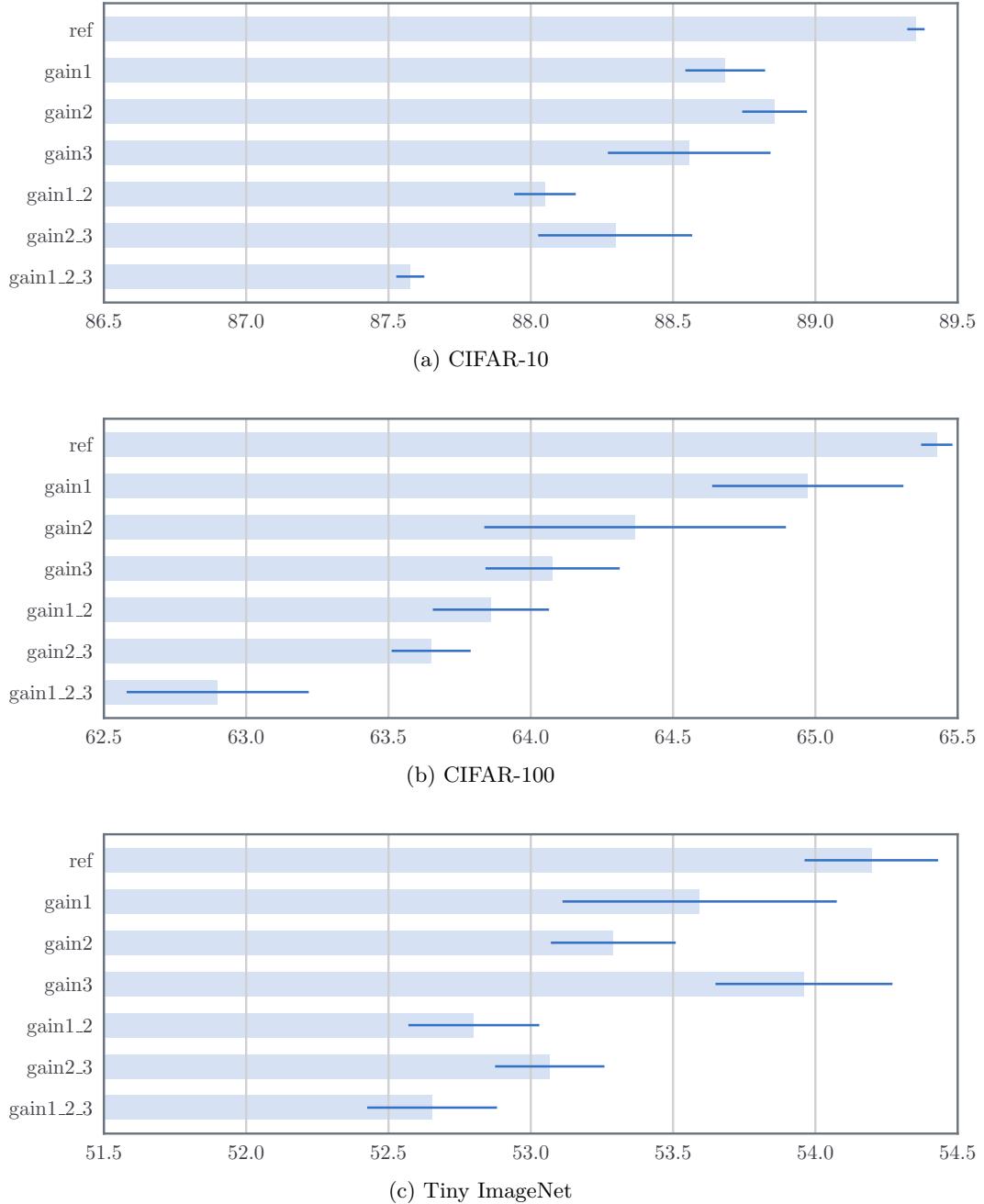


Figure 6.8: **Large kernel ablation results CIFAR and Tiny ImageNet.** Results showing the percentage classification accuracies obtained by swapping combinations of the three conv layers in the reference architecture from Table 6.1 with gain layers. Results shown are averages of 3 runs with the ± 1 standard deviations shown as dark blue lines. These results show that changing a convolutional layer for a gain layer is possible, but comes with a small accuracy cost which compounds as more layers are swapped.

wavelet gain distributions echoes the observation made in [subsection 4.6.2](#), where we saw that occluding diagonal subbands reduced classification performance by less than when occluding horizontal/vertical subbands.

Although the diagonal subbands are particularly small, the horizontal and vertical bands have many coefficients with magnitude very close to zero. This raises an interesting question – how many of these coefficients are important for classification? What if we were to apply a hard thresholding scheme to the weights, would setting some of these values to zero impact the network performance?

We measure the dropoff in accuracy when a hard threshold t is applied to the bandpass gains g_1 for the three gain layers of ‘gain1_2_3’. The resulting sparsity of each layer and the network performance is shown in [Figure 6.9b](#). For example, if we set a hard threshold value $t = 0.4$, only 20% of the gain1 weights, 1% of the gain2 and 0.02% of the gain3 weights are non-zero, yet the classification accuracy has only dropped by 0.5%.

This figure shows that despite the high cost of the bandpass gains – $12C_lC_{l+1}$ for a 1×1 gain, very few of these need to be nonzero.

6.4.3.2 DeConvolution and Filter Sensitivity

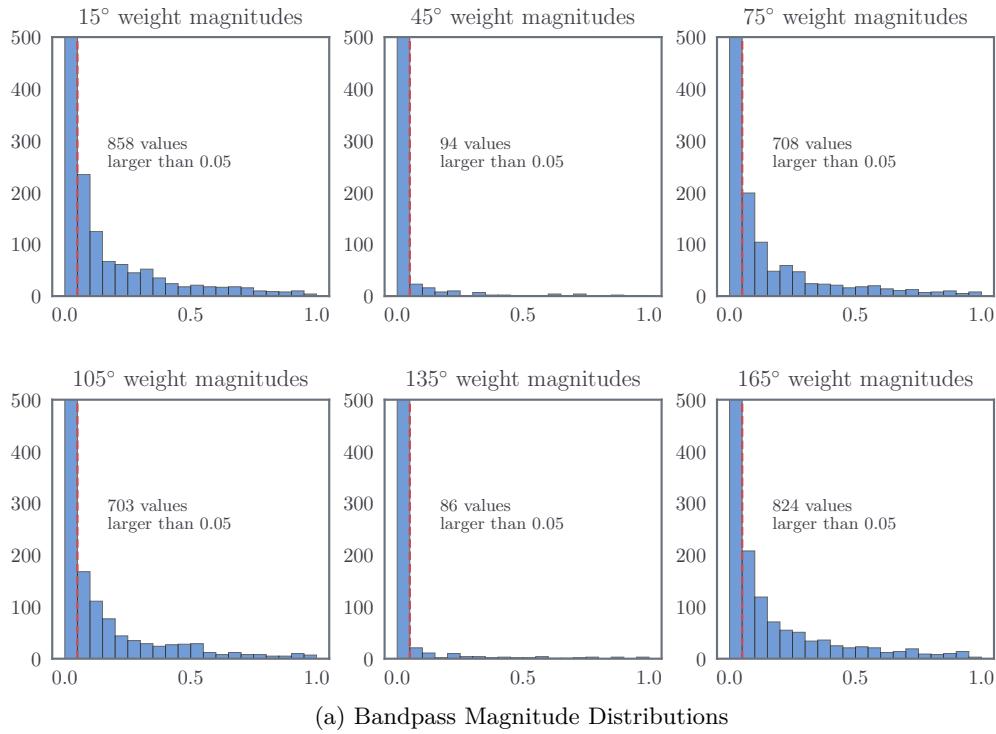
To visualize what the gain layer is responsive to, we build a deconvolutional system similar to the one described in [chapter 4](#). In particular, we present the entire CIFAR-100 validation set to the reference architecture and to the gain1_2_3 architecture, keeping track of what most highly excites each channel. Once we have this information, we present the same image again, storing the ReLU switches and max pooling locations for this same image, then we zero out all but a single value for the given channel, and zero out all other channels, and deconvolve to see the input pattern.

The resulting visualizations for the first two layers are shown in [Figure 6.10](#). We show only the top activation for each filter, rather than the top-9. For the second layer filters, we show only 64 of the 128 filter responses.

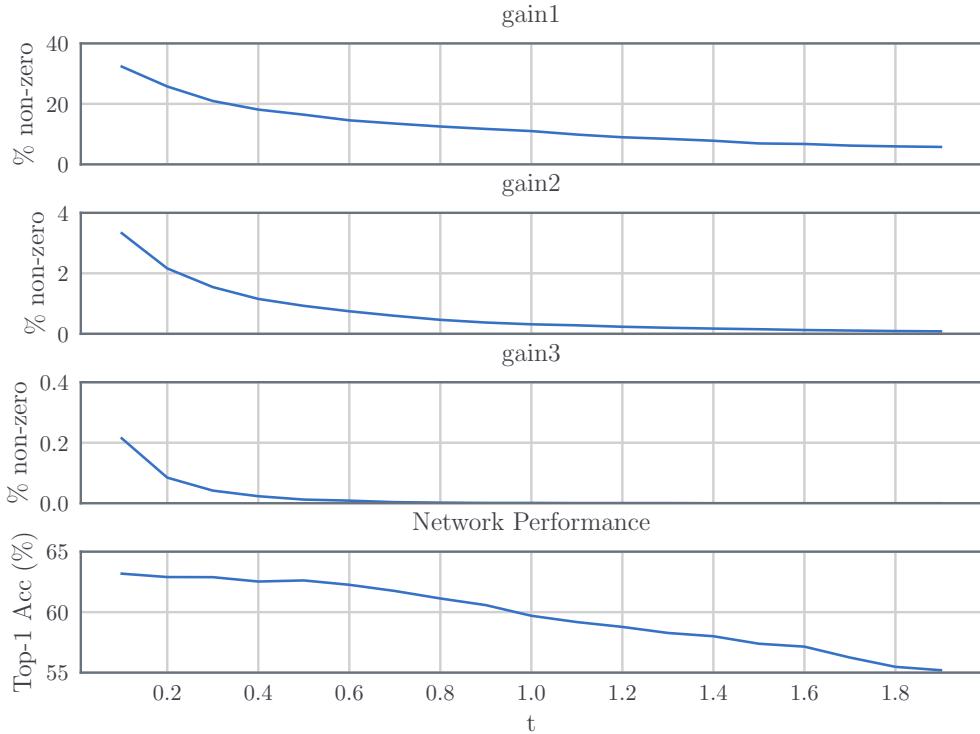
It is reassuring to see that despite the performance difference between the reference architecture and the gain1_2_3 architecture, the filters are responding to similar shapes. Note that for both the first and second layer responses, the gain layer has a smoother roll-off at the edges of the visualization, whereas the convolutional architecture has more square regions of support.

6.5 Wavelet-Based Nonlinearities

Returning to the goals from [subsection 6.2.4](#), the experiments from the previous section have shown that while it is possible to use a wavelet gain layer (G) in place of a convolutional layer (H), this may come with a small performance penalty. Ignoring this effect for the moment,



(a) Bandpass Magnitude Distributions



(b) Accuracy Dropoff from Thresholding

Figure 6.9: Bandpass Gain Properties for all gain layer network. (a) shows the distribution of the magnitudes for bandpass coefficients for the second layer (gain2). Each orientation has $128 \times 64 = 8192$ complex weights, most of which are close to 0 (ℓ_1 regularization was used in training); Y axis for each plot has been clipped to 500. The 45° and 135° weights have many fewer large coefficients. (b) shows the increase in sparsity and dropoff in classification accuracy when the weights are hard-thresholded with value t (same threshold applied to all 3 layers). For a threshold value of $t = 0.4$, 80% of the weights in gain1 are 0, 99.98% of the weights in gain2 are 0, 99.99% of the weights in gain3 are 0, yet classification accuracy is only 0.5% lower than the non-thresholded accuracy.

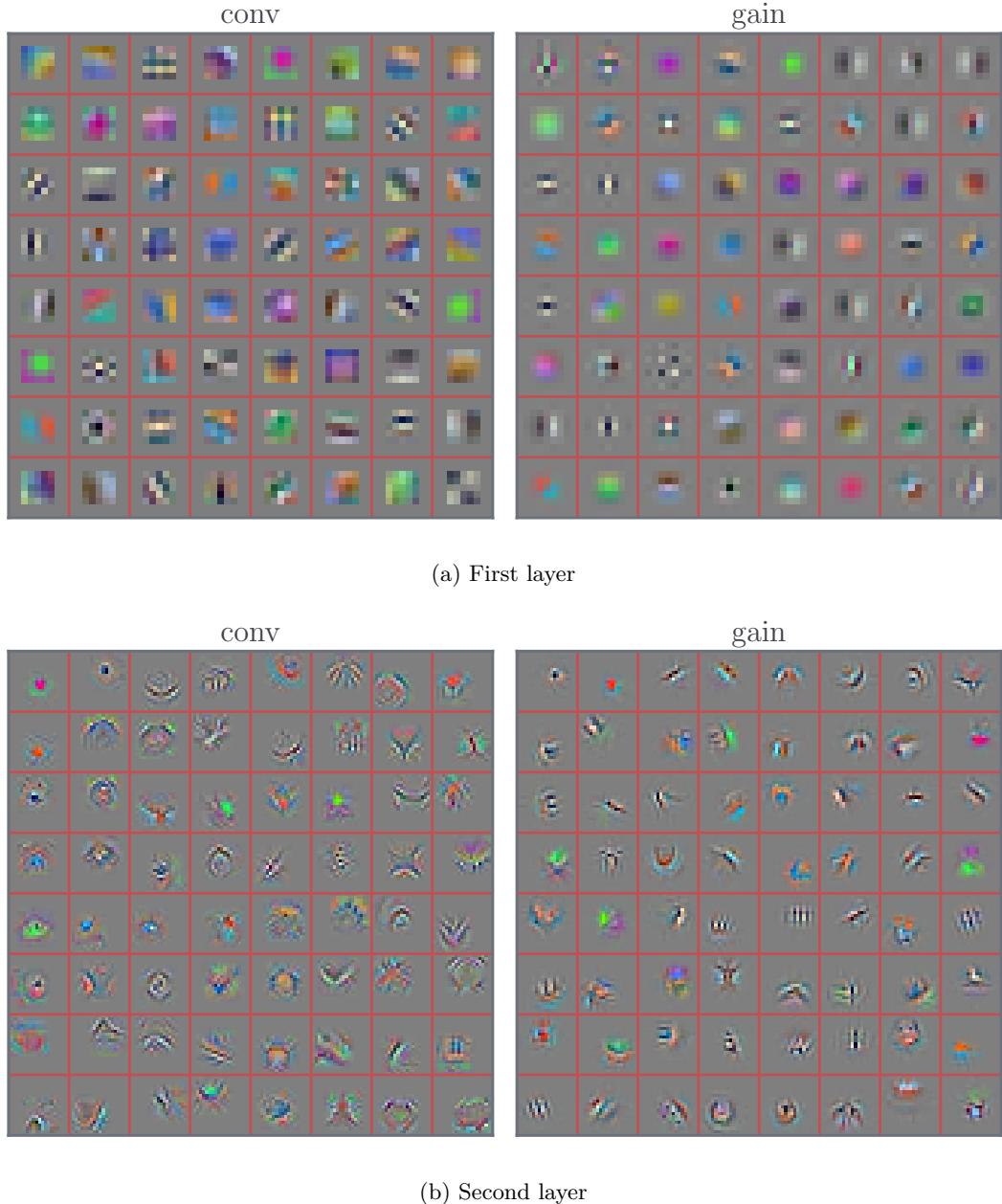


Figure 6.10: Deconvolution reconstructions for the reference architecture and purely gain layer architecture. Visualizations using a DeConvNet method similar to the one described in chapter 4. Here we find the input images in CIFAR-100 validation set that most highly activate each filter. Each image is then re-shown to the network and the meta-information is used to prime the DeConvNet to create the visualizations seen here. The left column has visualizations for the first and second layer filters for the all convolutional method, and the right column has visualizations for the first and second layer filters for the all gain layer method. Note the smoother roll-off at the edge of visualizations in the gain layer compared to the rectangular support regions for the conv layers. Aside from that, the two networks appear to be learning similar shapes.

in this section, we continue with our investigations into learning in the wavelet domain. In particular, is it possible to replace a pixel domain nonlinearity σ with a wavelet-based one σ_w ?

But what sensible nonlinearity to use? Two particular options are good initial candidates:

1. The ReLU: this is a mainstay of most modern neural networks and has proved invaluable in the pixel domain. Perhaps its properties will work well on wavelet coefficients too.
2. Thresholding: a technique commonly applied to wavelet coefficients for denoising and compression. Many proponents of compressed sensing and dictionary learning even like to compare soft thresholding to a two-sided ReLU [149], [150].

In this section, we will look at both and see if they improve the gain layer. If they do, it would be possible to connect multiple layers in the wavelet domain, avoiding the necessity to do inverse wavelet transforms after learning.

6.5.1 ReLUs in the Wavelet Domain

Applying the ReLU to the real lowpass coefficients is not difficult, but it does not generalize so easily to complex coefficients. The simplest option is to apply it independently to the real and imaginary coefficients, effectively only selecting one quadrant of the complex plane:

$$u_{lp} = \max(0, v_{lp}) \quad (6.5.1)$$

$$u_j = \max(0, \operatorname{Re}(v_j)) + j\max(0, \operatorname{Im}(v_j)) \quad (6.5.2)$$

Another option is to apply it to the magnitude of the bandpass coefficients. Of course, these are all strictly positive so the ReLU on its own would not do anything. However, they can be arbitrarily scaled and shifted by using a batch normalization layer. Then the magnitude could shift to (invalid) negative values, which can then be rectified by the ReLU.

Dropping the scale subscript j for clarity (we need it for the square root of negative 1), let a bandpass coefficient at a given scale be $v = r_v e^{j\theta_v}$ and define $\mu_r = \mathbb{E}[r_v]$ and $\sigma_r^2 = \mathbb{E}[(r_v - \mu_r)^2]$, then applying batch-normalization and the ReLU to the magnitude of v means we get:

$$r_u = \operatorname{ReLU}(BN(r_v)) = \max(0, \gamma \frac{r_v - \mu_r}{\sigma_r} + \beta) \quad (6.5.3)$$

$$u = r_u e^{j\theta_v} \quad (6.5.4)$$

This also works equivalently on the lowpass coefficients, although v_{lp} can be negative unlike r_v :

$$u_{lp} = \operatorname{ReLU}(BN(v_{lp})) = \max(0, \gamma' \frac{v_{lp} - \mu_{lp}}{\sigma_{lp}} + \beta') \quad (6.5.5)$$

6.5.2 Thresholding

For $t \in \mathbb{R}$, $z = re^{j\theta} \in \mathbb{C}$ the pointwise hard thresholding is:

$$\mathcal{H}(z, t) = \begin{cases} z & r \geq t \\ 0 & r < t \end{cases} \quad (6.5.6)$$

$$= \mathbb{I}(r > t)z \quad (6.5.7)$$

and the pointwise soft thresholding is:

$$\mathcal{S}(z, t) = \begin{cases} (r - t)e^{j\theta} & r \geq t \\ 0 & r < t \end{cases} \quad (6.5.8)$$

$$= \max(0, r - t)e^{j\theta} \quad (6.5.9)$$

Note that (6.5.9) is very similar to (6.5.3) and (6.5.4). We can rewrite (6.5.3) by taking the strictly positive terms γ, σ outside of the max operator:

$$r_u = \max(0, \gamma \frac{r_v - \mu_r}{\sigma_r} + \beta) \quad (6.5.10)$$

$$= \frac{\gamma}{\sigma_r} \max\left(0, r_v - \left(\mu_r - \frac{\sigma_r \beta}{\gamma}\right)\right) \quad (6.5.11)$$

then if $t' = \mu_v - \frac{\sigma_r \beta}{\gamma} > 0$, **doing batch normalization followed by a ReLU on the magnitude of the complex coefficients is the same as soft shrinkage with threshold t' , scaled by a factor $\frac{\gamma}{\sigma_r}$.**

The same analogy does not apply to the lowpass coefficients, as v_{lp} is not strictly positive.

While soft thresholding is similar to batch normalizations and ReLUs, we would also like to test how well it performs as a wavelet nonlinearity. To do this, we can:

- Learn the threshold t
- Adapt t as a function of the distribution of activations to achieve a desired sparsity level.

In early experiments, we found that trying to set desired sparsity levels by tracking the standard deviation of the statistics and setting a threshold as a function of it performed very poorly (causing a drop in top-1 accuracy of at least 10%). Instead, we choose to learn a threshold t . We make this an unconstrained optimization problem by changing (6.5.9) to:

$$\mathcal{S}(v, t) = \max(0, r - |t|)e^{j\theta} \quad (6.5.12)$$

Learning a threshold is only possible for soft thresholding, as $\frac{\partial L}{\partial t}$ is not defined for hard thresholding. Like batch normalization, we learn independent thresholds t for each channel.

Algorithm 6.1 The *wavelet gain layer* pseudocode

```

1: procedure WAVEGAINLAYER( $x$ )
2:    $u_{lp}, u_1 \leftarrow \text{DT\!CWT}(x, \text{nlevels} = 1)$ 
3:    $v_{lp}, v_1 \leftarrow G(u_{lp}, u_1)$                                  $\triangleright$  the normal wave gain layer
4:    $u_{lp} \leftarrow \sigma_{lp}(v_{lp})$                                  $\triangleright$  lowpass nonlinearity
5:    $u_1 \leftarrow \sigma_{bp}(v_1)$                                  $\triangleright$  bandpass nonlinearity
6:    $y \leftarrow \text{DT\!CWT}^{-1}(u_{lp}, u_1)$ 
7:    $x \leftarrow \sigma_{pixel}(y)$                                  $\triangleright$  pixel nonlinearity
8:   return  $x$ 
9: end procedure

```

6.6 Gain Layer Nonlinearity Experiments

Taking the same ‘gain1_2_3’ architecture used for CIFAR-100, we expand the *wavelet gain layer* by including nonlinearities as described in [Algorithm 6.1](#). In this layer, we have three different nonlinearities: the pixel, the lowpass, and the bandpass nonlinearity.

For these experiments, we test over a grid of possible options for these three functions:

| Nonlinearity | Options | | | |
|--------------|---------|---------|------------|---------------|
| Pixel | None | BN+ReLU | | |
| Lowpass | None | ReLU | BN+ReLU | \mathcal{S} |
| Bandpass | None | ReLU | BN+MagReLU | \mathcal{S} |

Where:

- ‘None’ means no nonlinearity – $\sigma(x) = x$.
- ‘ReLU’ is a ReLU without batch normalization. For real values, is a normal ReLU, for complex values is a ReLU applied independently to real and imaginary parts, i.e. [\(6.5.2\)](#). See [section E.3](#) for equations for the passthrough gradients for this nonlinearity.
- ‘BN+ReLU’ is batch normalization and ReLU (applicable only to real-valued activations) e.g. [\(6.5.5\)](#).
- ‘BN+MagReLU’ applies batch normalization to the magnitude of complex coefficients and then makes them strictly positive with a ReLU. This action is defined in [\(6.5.3\)](#). See [section E.5](#) for information on the passthrough and update equations for this nonlinearity.
- \mathcal{S} is the soft thresholding of [\(6.5.12\)](#) applied to the magnitudes of coefficients with learnable thresholds. See [section E.4](#) for information on the passthrough and update equations for this nonlinearity.

Table 6.2: **Different Nonlinearities in the Gain Layer.** Top-1 Accuracies for the ‘gain1_2_3’ network trained on CIFAR-100 using different wavelet and pixel nonlinearities. The rows of the table correspond to different bandpass nonlinearities and the columns correspond to different lowpass nonlinearities. $\sigma_{pixel} = \sigma_{lp} = \sigma_{bp} = \text{None}$ is a linear system (with max pooling). $\sigma_{pixel} = \text{ReLU}$, $\sigma_{lp} = \sigma_{bp} = \text{None}$ is the system used in earlier experiments, which is linear in the wavelet domain and has a nonlinearity in the pixel domain. The best result is highlighted in bold corresponding to $\sigma_{pixel} = \text{None}$, $\sigma_{lp} = \text{ReLU}$ and $\sigma_{bp} = \text{BN} + \text{ReLU}$.

| | | (a) $\sigma_{pixel} = \text{None}$ | | | | |
|---------------|--|------------------------------------|------|-------------|---------|---------------|
| | | σ_{lp} | None | ReLU | BN+ReLU | \mathcal{S} |
| σ_{bp} | | None | 45.0 | 63.8 | 64.8 | 61.9 |
| None | | ReLU | 42.6 | 62.8 | 63.9 | 61.3 |
| ReLU | | BN+MagReLU | 48.5 | 65.3 | 64.6 | 62.9 |
| BN+MagReLU | | \mathcal{S} | 44.6 | 63.9 | 63.6 | 61.7 |

| | | (b) $\sigma_{pixel} = \text{BN} + \text{ReLU}$ | | | | |
|---------------|--|--|------|------|---------|---------------|
| | | σ_{lp} | None | ReLU | BN+ReLU | \mathcal{S} |
| σ_{bp} | | None | 62.8 | 60.7 | 64.2 | 63.0 |
| None | | ReLU | 62.5 | 60.3 | 64.2 | 63.2 |
| ReLU | | BN+MagReLU | 64.9 | 61.6 | 64.7 | 65.2 |
| BN+MagReLU | | \mathcal{S} | 63.3 | 60.9 | 63.9 | 63.3 |

As the pixel nonlinearity has only two options, the results are best displayed as a pair of tables, firstly for no nonlinearity and secondly for the standard batch normalization and ReLU. See [Table 6.2](#) for these two tables.

Digesting this information gives us some useful insights:

1. It is possible to improve on the gain layer from the previous experiments with the right nonlinearities. The previous section’s gain layer corresponds to $\sigma_{lp} = \sigma_{bp} = \text{None}$ and $\sigma_{pixel} = \text{ReLU}$, or the top left entry of [Table 6.2b](#).
2. Doing a ReLU on the real and imaginary parts of the bandpass coefficients independently (the second row of both tables) almost always performs worse than having no nonlinearity (first row of both tables).
3. The best combination is to have batch normalization and a ReLU applied to the magnitudes of the bandpass coefficients and batch norm and a ReLU applied to either the lowpass or pixel coefficients with no nonlinearity in the pixel domain.

The best accuracy score of 65.3% is now 0.1% lower than the fully convolutional architecture, an improvement from the 62.8% score achieved with only a pixel nonlinearity. This

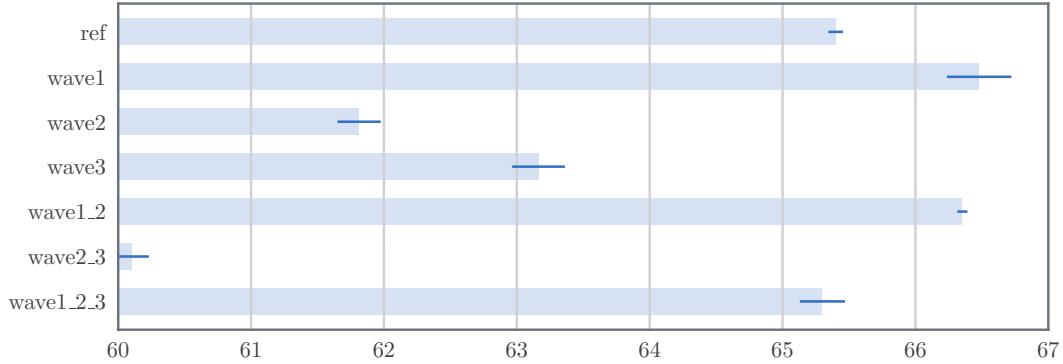


Figure 6.11: **CIFAR-100 Ablation results with the *wave layer*.** We use the same naming scheme from subsection 6.4.2 but to differentiate between the results from Figure 6.8b we call the options ‘waveX’. Results show the mean of 3 runs with ± 1 standard deviation lines in dark blue. When we add nonlinearities in the wavelet domain, the ablation results change dramatically. It appears that learning in the wavelet domain works best for the first layer of the CNN (wave1), and this improves on the purely convolutional method by a whole percentage point. Replacing the second and third layers degrades performance independently of what was used in the first layer. Swapping the first two layers (wave1_2) performs nearly as well and with a slightly narrower spread.

happens when there is no pixel nonlinearity, use a ReLU on the lowpass coefficients and Batch Normalization and a ReLU on the magnitudes of the bandpass coefficients.

6.6.1 Ablation Experiments with Nonlinearities

Now that we have found the best nonlinearity to use for the gain layer, will this improve our ablation study from subsection 6.4.2? To test this, we repeat the same experiment on CIFAR-100 using the newly found nonlinearities in Algorithm 6.1 (i.e., σ_{pixel} = None, σ_{lp} = ReLU, and σ_{bp} = BN + ReLU).

See Figure 6.11 for the results from these experiments. When we use the wavelet-based nonlinearities, the results change considerably. We see an improvement by 1% when the first layer in the CNN is changed for a wave layer, but any other changes degrade performance from this.

6.7 Conclusion

In this chapter we have presented the novel idea of learning filters by taking activations into the wavelet domain. There, we can apply the proposed *wavelet gain layer* G instead of a pixel-wise convolution. We can return to the pixel domain to apply a ReLU, or stay in the wavelet domain and apply wavelet-based nonlinearities σ_{lp}, σ_{bp} . We have considered the

possible challenges this proposes and described how a multi-rate system can learn through backpropagation.

Our experiments have been promising but are still only preliminary. We have shown that the gain layer can learn in an end-to-end system, achieving nearly the same accuracies on CIFAR-10, CIFAR-100 and Tiny ImageNet to the reference system with convolutional layers (Figure 6.8). This is a good start and shows the plausibility of the wavelet gain layer, but more experiments on larger datasets and deeper networks is needed. Despite the slight reduction in performance, we saw some nice properties to the gain layer. Most of the bandpass gains are near zero (Figure 6.9), which does not affect training but could offer speedups for inference. Additionally, doing deconvolution to visualize the sensitivity of the filters in a gain layer showed that the system was still learning sensible shapes with nice spatial roll-off properties (Figure 6.10).

We have searched for good candidates for wavelet nonlinearities, and saw that using a ReLU on the lowpass coefficients, and Batch Normalization and a ReLU on the magnitudes of the bandpass coefficients improved the performance of the gain layer considerably. This is an exciting development and indicates that we may not need to return to the pixel domain at all, possibly eliminating the need for the inverse wavelet transforms used in our experiments (see steps 6 and 7 in Algorithm 6.1). However, one needs to be careful, as taking the inverse transform followed by taking the forward transform does not necessarily give the same wavelet coefficients due to the redundancy of the DT^CWT, instead projecting onto the range space of the transform. Removing the inverse transform is something we did not have time to fully explore and so we have included it in our future work section.

When we added the nonlinearities to the gain layer we saw that we were able to achieve some improvements in performance over a fully convolutional architecture (Figure 6.11). The proposed layer worked best at the beginning of the CNN, which matches the intuition for doing this work described in the introduction to the chapter. More research still needs to be done with the wave layer as part of a deeper system.

Chapter 7

Conclusion and Further Work

In this thesis, we have examined the effectiveness of using complex wavelets as basis functions for deep learning models. We summarize the key results we have found before describing possible future extensions of our work, some natural extensions that we were not able to explore due to time constraints, and some propositions for tying together interesting pieces of work that were developed towards the end of the project.

7.1 Summary of Key Results

Chapter 3 shows how using the separable spatial implementation of the DTCWT as the chosen wavelets for the original ScatterNet design greatly speeds up computation compared with Fourier domain implementations of complex wavelets (see [Table 3.3](#)). We were also able to derive the backpropagation equations for wavelet and scattering layers, aiding the use of ScatterNets as part of deep networks. As part of this, we tested out the performance of the DTCWT based ScatterNet as a front end to a simple CNN for some small classification tasks, and compared its performance to the Morlet based system. We found that as well as being faster, the performance was often better when using the DTCWT wavelets [Table 3.5](#). When doing these tests, we found that of the wavelet choices available for the DTCWT, those with the fewest taps (and hence wider transition bands and worse stopband rejection) performed the best. This is somewhat surprising and while we were not able to investigate why due to time constraints, it may provide some interesting insights as to what the CNN backend is doing. It also may have been a side effect of the small image sizes used in the classification task.

Chapter 4 builds a visualization tool we call the DeScatterNet. We use this to interrogate the input patches that most highly excite the ScatterNet outputs on a subset of images from ImageNet (see [Figure 4.3](#)). We saw that the second order scattering coefficients are most highly excited by ripple and checkerboard like patterns. These were very different to the patterns that most highly excite filters of a CNN. We believe this may explain why ScatterNets

perform well on texture discrimination [23] but less well on natural image classification [25]. We then performed some occlusion tests on a hybrid network with a ScatterNet front end and CNN backend and saw that the CNN was able to operate with little degradation when the second order scattering coefficients were zeroed out (there was only a small drop in classification accuracy), but it suffered greatly when the zeroth or first order coefficients were zeroed out. We also found the surprising result that on the datasets we tested, the filters with diagonal edges were less important than their vertical or horizontal counterparts (Figure 4.4). If the input images were rotated by $\pm 30^\circ$ then the diagonal channels became the most important. This echoes the experiments of Blakemore and Cooper[7] who controlled the orientation of edges exposed to kittens in their development stage. Finally, this chapter showed some ways to expand on the ScatterNet network and shows the features possible with these extensions in Figure 4.7. This last section inspires the work of chapter 5 and chapter 6.

Chapter 5 reworks the ScatterNet into individual layers. This redesign allows us to rethink how we want to use wavelets, and we introduce the *learnable ScatterNet* made up of *locally invariant convolutional layers*. Rather than applying the same layer twice to get a second order ScatterNet, we introduce mixing across the output channels, taken after the magnitude operation. The flexibility of the proposed layer means it can be used in a ScatterNet-like system, where the number of output channels grows exponentially with the number of layers, or in a CNN-like system, where the number of output channels remains mostly constant across layers. We experimented with both possibilities, showing that the extra learnability definitely helps the ScatterNet style system (??) and can help a CNN (??). The demodulation of energy from the complex modulus means that the proposed locally invariant layer can only be used a few times. In particular, we saw that the layer performed best when used where a CNN would naturally downsample (or pool) the input.

Chapter 6 looks at learning in the wavelet space without taking complex magnitudes. We present the wavelet *gain layer* which takes inputs to the wavelet domain, learns complex gains to attenuate/accentuate different subbands, mixes the subbands across the different channels and offers the ability of returning to the pixel domain with the inverse wavelet transform. Our experiments have been promising but are still only preliminary. We show that the *gain layer* can learn in an end-to-end system, achieving nearly the same accuracies on CIFAR-10, CIFAR-100 and Tiny ImageNet to the reference system with only convolutional layers (Figure 6.8). Despite the slight reduction in performance, we saw some nice properties of the gain layer. The bandpass gains were very sparse, needing very few non-zero coefficients and visualizations of what the layers were responding to showed that the gain layer had nice spatial roll-off properties (Figure 6.10). We then found using a ReLU on the lowpass coefficients, and Batch Normalization and a ReLU on the magnitudes of the bandpass coefficients improved the performance of the gain layer considerably (Table 6.2). The gains and nonlinearity together make the *wave layer*. Using this, we saw that we were able to achieve some improvements in performance over a fully convolutional architecture (Figure 6.11). The

wave layer seems to work best at the beginning of the CNN but more research still needs to be done with deeper systems.

7.2 Future Work

This thesis has started to examine ways of using wavelets as basis functions for deep learning models. Our research has found some possible approaches which offer some advantage in terms of number of parameters, interpretability and (theoretical) computation time. But there are many things that we were not able to try, and some of these may show that wavelets have a larger benefit than we were able to find.

7.2.1 Faster Transforms and More Scales

Firstly, despite our best efforts in making a fast wavelet transform, the speed of a DT \mathbb{C} WT in *Pytorch Wavelets* is slower than we believe it ought to be. A 10×10 convolutional filter with 100 multiplies per input pixel is roughly twice as quick to compute as a DT \mathbb{C} WT with 36 multiplies per pixel. We limited our design to use high level cuDNN calls and this was the best we could do with these primitives, and believe that any further speed up would require custom CUDA kernels. The computational time was not a problem for datasets such as CIFAR and Tiny ImageNet, but it did prevent us from testing the wavelet gain layer and invariant layer on ImageNet (see [appendix A](#) for some run times). We believe that these layers would perform better with larger images where the extent of the wavelet is not comparable to the size of the image (hence requiring lots of padding).

Another aspect of testing larger images is the benefit of using multiple scales in any system. Our wavelet gain layer only used the first or second scale in our experiments, but the real benefit of decimated wavelet transforms is the speedup they offer by allowing for multiscale approaches. Little research has been done in splitting the input or mid-level activations into multiple scales and learning different filters for the different scales, but some examples include [\[145\]](#), [\[151\]](#).

7.2.2 Expanding Tests on Invariant Layers

We have shown in [chapter 5](#) that the invariant layer can work quite well in a VGG-like CNN as a replacement for a convolutional layer followed by pooling. We would have liked to test this on larger networks, replacing areas of sample rate change with invariant layers, and see how well this generalizes. One common place that has a large sample rate change is in the first layer of CNNs, with networks like AlexNet and ResNet downsampling by a factor of four in each direction after the first layer. Experiments by Oyallon, Belilovsky, and Zagoruyko [\[39\]](#) have looked at replacing this with a fixed ScatterNet, but it would be interesting to see how well this performs with the *learned* ScatterNet we have developed.

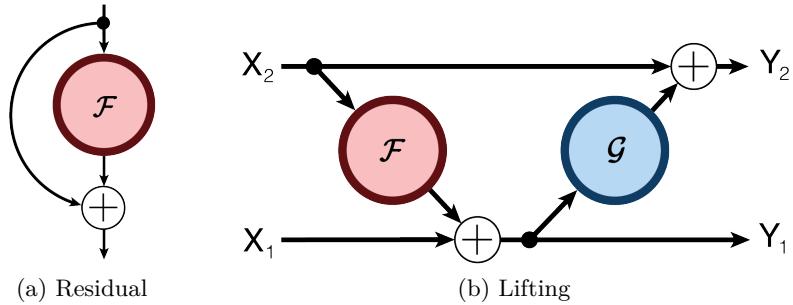


Figure 7.1: Residual vs Lifting Layers. In a residual layer, an input is transformed by a learned function \mathcal{F} and added to itself. Typically $\|\mathcal{F}x\| \ll \|x\|$ meaning the vector y is a small perturbation of the vector x . In a lifting layer, each path is a learned function of the other added to itself. In the original second generation wavelets, \mathcal{F} and \mathcal{G} are FIR filters, but there is no requirement on them to be. Diagrams taken from [152].

7.2.3 Expanding Tests on Gain Layer

We believe that the wavelet gain layer of chapter 6 is the most unique and potentially promising piece of work in this thesis. The results we describe in Table 6.2 were the last experiments we were able to do, and they just start to show some promise for learning in the wavelet domain. There is potentially a lot more research to be done on the wavelet gain layer.

Firstly, the result from Figure 6.9 is very intriguing. This figure shows that a lot of the learned bandpass gains are very near zero, and thresholding them after training does little to affect performance. It would be very interesting to see the performance of the network if the threshold values were set near the end of training, and the nonzero gains were allowed to grow to compensate for this. Also, is this a property unique to the wavelet gain layer? Or do all CNNs learn *mostly* lowpass filters, and need only a few bandpass filters? The parameter cost of the proposed gain layer was dominated by the cost of coding these bandpass gains, despite requiring very few of them. It would be an interesting piece of work to try to redesign the gain layer to allow training with many fewer bandpass parameters.

Secondly, the results from Table 6.2 show that the best performing gain layer had no pixel nonlinearity, a ReLU in the lowpass and a ReLU applied to the magnitude of the bandpass. This is very exciting as it opens up the possibility of staying in the wavelet domain longer. In our experiments, we would take inverse transforms and apply a NoOp in the pixel domain before returning to the wavelet domain. This is marginally different than staying in the wavelet domain, as it involves projecting from a redundant space to a non-redundant one. More experiments could be done to investigate whether this projection process is needed, and hence whether significant DT^CWT computation could be saved.

7.2.4 ResNets and Lifting

We briefly mentioned ResNets in subsection 2.5.6 but did not study them in depth in this thesis. Interestingly, there are many similarities between ResNets and second generation wavelets, or the *lifting* framework [121], [153]. In a residual layer, the output is $y = \mathcal{F}(x) + x$ where for a lifting system, the layer is a two port network defined by:

$$y_1 = \mathcal{F}(x_1) + x_2 \quad (7.2.1)$$

$$y_2 = \mathcal{G}(y_1) + x_1 \quad (7.2.2)$$

Figure 7.1 shows the similarities between the two designs. The works [152], [154] both make the small modifications to the ResNet design to make a lifting style architecture. Gomez, Ren, Urtasun, *et al.* [152] do this to save memory on the activations for backpropagation (you do not need to save meta-information on the forwards pass as you can regenerate activations on the backwards pass). Jacobsen, Smeulders, and Oyallon [154] extend on [152] and also explore merging images with linear interpolation in the activation space, reconstructing from the new latent values. We believe that there are potentially many more benefits to using the lifting design as an extension of our work into learning from basis functions.

7.2.5 Protecting against Attacks

Adversarial examples are starting to become a real concern for CNNs. A classic adversarial example is an image that appears innocuous to a human, but has been corrupted with a low energy signal that can completely convince a CNN that the image is something else. Figure 7.2 shows an example of this taken from [22].

While Szegedy, Zaremba, Sutskever, *et al.* [22] show there is a weakness to CNNs being fooled, these adversarial examples were obtained by gradient ascent to the target class. I.e. they had full access to the parameters of the model. Papernot, McDaniel, Goodfellow, *et al.* [155] expand this to show it is possible to develop an attack without knowing the model, treating it as a *black box*. Protecting against these black box attacks has become an arms race in recent years, with measures and counter-measures constantly being developed. An excellent paper reviewing some attacks and defences is [156].

One such concerning black box attack is described in [157], where Engstrom, Tran, Tsipras, *et al.* show that even simple transformations such as translations and rotations are enough to fool many modern CNNs. This is something our invariant layer may be able to protect against.

Alternatively, a recent defence tactic is described in [158]. In this paper, Cisse, Bojanowski, Grave, *et al.* propose to have non-expansive convolutional and pooling layers (Lipschitz constant 1) and have weight matrices that are approximately Parseval tight frames, a result intimately related to the tight framed DTCWT. In addition to finding these networks more

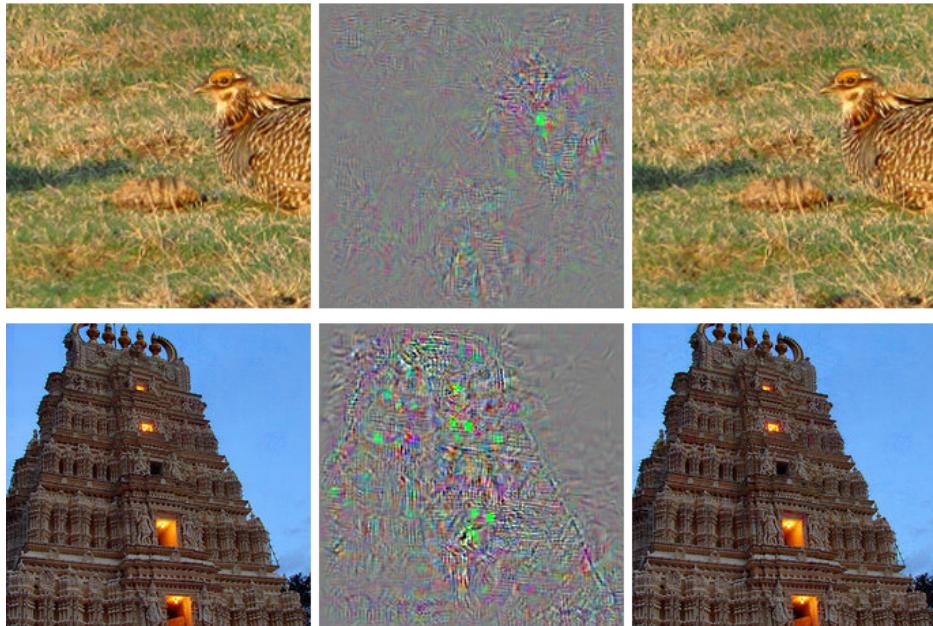


Figure 7.2: **Adversarial examples that can fool AlexNet.** Two examples of images that were correctly classified on the left, with additive signals in the centre (contrast levels were magnified by 10) and the resulting images on the right. Both of the output images are then predicted to be an ostrich. Images taken from [22].

robust to adversarial examples, they show that they can train faster and have a better usage of the capacity of the networks.

We feel that the learning of relatively smooth functions offered by wavelet bases has a potentially large scope for making CNNs more robust to many attacks.

7.2.6 Convolutional Sparse Coding

Recent work on convolutional sparse coding and convolutional dictionary learning [149], [150], [159]–[161] has started to draw many parallels between the structure of modern CNN architectures and the problem of sparse representations. We believe this is a valuable insight and can give a fresh new perspective on how we can better train CNNs.

It would be interesting to try to extend this recent work on convolutional sparse coding to wavelet bases. In [162] Rubinstein, Zibulevsky, and Elad attempted something similar with their double sparsity model, learning to sparsely combine atoms from a fixed base dictionary (they use wavelet, Fourier and DCT base dictionaries), although this was applied to block coding and has not been extended to convolutional sparse coding.

7.2.7 Weight Matrix Properties

Some recent work by Advani and Saxe [163] show that if you ignore the nonlinearities in a neural network, you can analyze the convergence properties by looking at the conditioning of the weight matrix (the ratio of the largest to the smallest singular values). In particular, well conditioned matrices have better convergence, and the size of the *eigengap* (distance between the smallest non-zero eigenvalue and 0) is important to protect against overfitting.

This is something that we have not enforced or considered in the design of our wavelet based layers, and it would be an interesting extension.

7.3 Final Remarks

Need to write something: perhaps emphasising the rich vein of ideas that your work has opened up and the need for substantial improvements in learning efficiency of deep nets that will be required if they are to be widely used in the future.

Appendix A

Architecture Used for Experiments

The experiments for this thesis were run on a single server with 8 GPUs and 14 CPUs. The GPUs were each NVIDIA GeForce GTX 1080 cards released in May 2016. They each have 8GiB of RAM, 2560 CUDA cores and 320 GB/s memory bandwidth. The CPUs were Intel(R) Xeon(R) E5-2660 models.

At the completion of the project, we were running CUDA 10.0 with cuDNN 7.6 and PyTorch version 1.1.

To do hyperparameter search we used the Tune package [139] which we highly recommend, as it makes running trials in parallel very easy.

A.1 Run Times of some of the Proposed Layers

Appendix B

Forward and Backward Algorithms

We have listed some of the forward and backward algorithms here that are not included in the main text for the interested reader. We also have included proofs for the gradients of decimation, interpolation and an entire analysis decomposition.

B.1 Gradients of Sample Rate Changes

Consider 1D decimation and interpolation of a signal X . The results we prove here easily extrapolate to 2D, but for ease we have kept to the 1D case.

Decimation of a signal X by M is defined as:

$$Y[n] = X[Mn] \quad (\text{B.1})$$

and interpolation by M as:

$$Y[n] = \begin{cases} X[\frac{n}{M}] & n = Mk, k \in \mathbb{Z} \\ 0 & \text{otherwise} \end{cases} \quad (\text{B.2})$$

B.1.1 Decimation Gradient

From (B.1) the gradient $\frac{\partial Y_n}{\partial X_k}$ is:

$$\frac{\partial Y_n}{\partial X_k} = \begin{cases} 1 & k = Mn \\ 0 & \text{otherwise} \end{cases} \quad (\text{B.3})$$

By the chain rule, $\frac{\partial J}{\partial X_k}$ is:

$$\frac{\partial J}{\partial X_k} = \frac{\partial J}{\partial Y_n} \frac{\partial Y_n}{\partial X_k} \quad (\text{B.4})$$

$$= \begin{cases} \Delta Y[\frac{k}{M}] & k = Mn \\ 0 & \text{otherwise} \end{cases} \quad (\text{B.5})$$

which is interpolating ΔY by M from (B.2).

B.1.2 Interpolation Gradient

From (B.2) the gradient $\frac{\partial Y_n}{\partial X_k}$ is:

$$\frac{\partial Y_n}{\partial X_k} = \begin{cases} 1 & n = Mk \\ 0 & \text{otherwise} \end{cases} \quad (\text{B.6})$$

and then the gradient $\frac{\partial J}{\partial X_k}$ is:

$$\frac{\partial J}{\partial X_k} = \frac{\partial J}{\partial Y_n} \frac{\partial Y_n}{\partial X_k} \quad (\text{B.7})$$

$$= \begin{cases} \Delta Y[n] & n = Mk \\ 0 & \text{otherwise} \end{cases} \quad (\text{B.8})$$

$$= \Delta Y[Mk] \quad (\text{B.9})$$

which is decimation of ΔY by M .

B.2 Gradient of Wavelet Analysis Decomposition

We mention in subsection 3.4.3 that the gradient of a forward DWT with orthogonal wavelets is just the inverse DWT. This easily follows from applying the chain rule and using the gradients of each of the stages of the DWT (convolution becomes correlation, downsampling becomes upsampling). The equivalent backwards pass of Figure 3.1 is shown in Figure B.1.

For an orthogonal wavelet transform, the synthesis filters are the time reverse of the analysis filters [87, Chapter 3]. This means that the blocks $H_0(z^{-1}), H_1(z^{-1})$ can be replaced with $G_0(z), G_1(z)$ respectively, giving the inverse DWT.

B.3 Extra Algorithms

In the text we refer to the 2-D inverse DWT which we have listed in Algorithm B.1, as well as the smooth magnitude operation (Algorithm B.2) and the inverse DT-CWT Algorithm B.3.

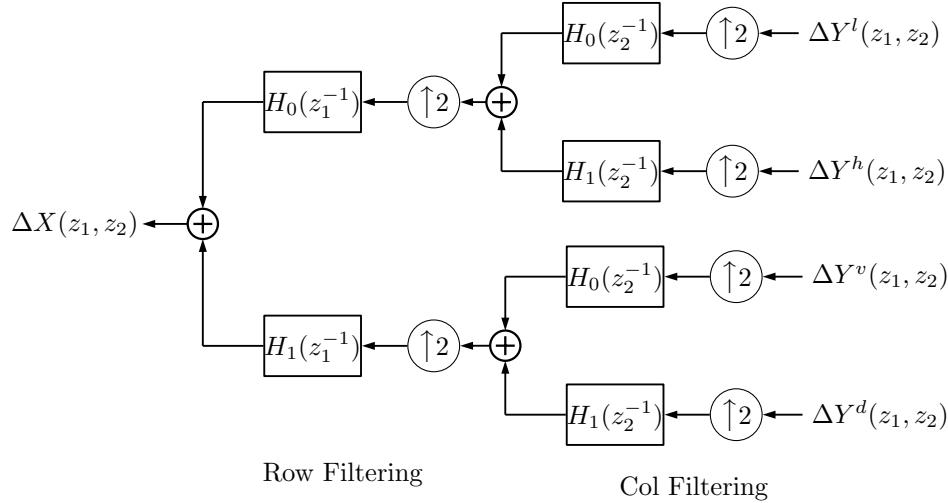


Figure B.1: **Gradient of DWT analysis.** Each block in the forward pass of Figure 3.1 has been swapped for its gradient. The resulting gradient has the same form as the inverse DWT.

Algorithm B.1 2-D Inverse DWT and its gradient

```

1: function IDWT.FORWARD( $ll, lh, hl, hl, g_0^c, g_1^c, g_0^r, g_1^r, mode$ )
2:   save  $g_0^c, g_1^c, g_0^r, g_1^r, mode$                                  $\triangleright$  For the backwards pass
3:    $lo \leftarrow \text{sfb1d}(ll, lh, g_0^c, g_1^c, mode, axis = 2)$            $\triangleright$  See Algorithm 3.1
4:    $hi \leftarrow \text{sfb1d}(hl, hh, g_0^c, g_1^c, mode, axis = 2)$ 
5:    $x \leftarrow \text{sfb1d}(lo, hi, g_0^r, g_1^r, mode, axis = 3)$ 
6:   return  $x$ 
7: end function

1: function IDWT.BACKWARD( $\Delta y$ )
2:   load  $g_0^c, g_1^c, g_0^r, g_1^r, mode$ 
3:    $\Delta lo, \Delta hi \leftarrow \text{afb1d}(\Delta y, g_0^r, g_1^r, mode, axis = 3)$        $\triangleright$  See Algorithm 3.1
4:    $\Delta ll, \Delta lh \leftarrow \text{afb1d}(\Delta lo, g_0^c, g_1^c, mode, axis = 2)$ 
5:    $\Delta hl, \Delta hh \leftarrow \text{afb1d}(\Delta hi, g_0^c, g_1^c, mode, axis = 2)$ 
6:   return  $\Delta ll, \Delta lh, \Delta hl, \Delta hh$ 
7: end function

```

Algorithm B.2 Smooth Magnitude

```

1: function MAG_SMOOTH.FORWARD( $x, y, b$ )
2:    $b \leftarrow \max(b, 0)$ 
3:    $r \leftarrow \sqrt{x^2 + y^2 + b^2}$ 
4:    $\frac{\partial r}{\partial x} \leftarrow \frac{x}{r}$ 
5:    $\frac{\partial r}{\partial y} \leftarrow \frac{y}{r}$ 
6:   save  $\frac{\partial r}{\partial x}, \frac{\partial r}{\partial y}$ 
7:   return  $r - b$ 
8: end function

1: function MAG_SMOOTH.BACKWARD( $\Delta r$ )
2:   load  $\frac{\partial r}{\partial x}, \frac{\partial r}{\partial y}$ 
3:    $\Delta x \leftarrow \Delta r \frac{\partial r}{\partial x}$ 
4:    $\Delta y \leftarrow \Delta r \frac{\partial r}{\partial y}$ 
5:   return  $\Delta x, \Delta y$ 
6: end function

```

Algorithm B.3 2-D Inverse DT^CWT

```

1: function IDTCWT( $yl, yh, mode$ )
2:   load  $g_0^a, g_1^a, g_0^b, g_1^b$                                  $\triangleright$  Load from memory
3:    $ll_r, ll_{j_1}, ll_{j_2}, ll_{j_1j_2} \leftarrow \text{deinterleave}(yl)$ 
4:   for  $j = J; j \geq 1; j = j - 1$  do
5:      $x_{1b}, x_{3b}, x_{2b}, x_{2a}, x_{3a}, x_{1a} \leftarrow yh[j]$ 
6:      $lh_r \leftarrow \text{Re}(x_{1a} + x_{1b})$ 
7:      $lh_{j_1} \leftarrow \text{Im}(x_{1a} - x_{1b})$ 
8:      $lh_{j_2} \leftarrow \text{Im}(x_{1a} + x_{1b})$ 
9:      $lh_{j_1j_2} \leftarrow \text{Re}(x_{1b} - x_{1a})$ 
10:     $hl_r \leftarrow \dots$                                           $\triangleright$  Same procedure for  $hl$  and  $hh$ 
11:     $ll_r \leftarrow \text{IDWT}_r(ll_r, lh_r, hl_r, hh_r, g_0^a, g_1^a, g_0^b, g_1^b, mode)$ 
12:     $ll_{j_1} \leftarrow \text{IDWT}_r(ll_{j_1}, lh_{j_1}, hl_{j_1}, hh_{j_1}, g_0^b, g_1^b, g_0^a, g_1^a, mode)$ 
13:     $ll_{j_2} \leftarrow \text{IDWT}_r(ll_{j_2}, lh_{j_2}, hl_{j_2}, hh_{j_2}, g_0^a, g_1^a, g_0^b, g_1^b, mode)$ 
14:     $ll_{j_1j_2} \leftarrow \text{IDWT}_r(ll_{j_1j_2}, lh_{j_1j_2}, hl_{j_1j_2}, hh_{j_1j_2}, g_0^b, g_1^b, g_0^a, g_1^a, mode)$ 
15:   end for
16:   return  $(ll_r + ll_{j_1} + ll_{j_2} + ll_{j_1j_2}) / 2$ 
17: end function

```

Appendix C

Invertible Transforms and Optimization

This Appendix proves the assertion made in [subsection 6.2.1](#), namely that:

Any invertible linear transform of the parameter space will not change the updates if a linear optimization scheme is used (for example standard gradient descent, or SGD with momentum).

We prove the case for the DFT as the invertible transform, but it can easily be extended to other transforms so long as [\(C.3\)](#) and [\(C.4\)](#) hold.

C.1 Background

We define the DFT as the orthonormal version, i.e.

$$U_{ab} = \frac{1}{\sqrt{N}} \exp\left\{-\frac{2j\pi ab}{N}\right\}$$

then call $X = \text{DFT}\{x\}$. In matrix form the 2-D DFT is then [\[164, Chapter 2\]](#):

$$X = \text{DFT}\{x\} = UxU \tag{C.1}$$

$$x = \text{DFT}^{-1}\{X\} = U^*XU^* \tag{C.2}$$

The gradients of these linear operators are:

$$\frac{\partial L}{\partial X} = U \frac{\partial L}{\partial x} U = \text{DFT}\left\{\frac{\partial L}{\partial x}\right\} \tag{C.3}$$

$$\frac{\partial L}{\partial x} = U^* \frac{\partial L}{\partial X} U^* = \text{DFT}^{-1}\left\{\frac{\partial L}{\partial X}\right\} \tag{C.4}$$

C.2 Analysis

Consider a single filter parameterized in the DFT and spatial domains presented with the exact same data and with the same ℓ_2 regularization λ , and learning rate η . Let the spatial filter at time t be \mathbf{w}_t and the Fourier-parameterized filter be $\hat{\mathbf{w}}_t$, and let

$$\hat{\mathbf{w}}_1 = \text{DFT}\{\mathbf{w}_1\} \quad (\text{C.5})$$

We need to prove that presenting the data to each system in the same order results in the same outputs for all time t .

Proof. Assume that $\hat{\mathbf{w}}_t = \text{DFT}\{\mathbf{w}_t\}$ for a given timestep t .

After presenting both systems with the same minibatch of samples \mathcal{D} and calculating the gradient $\frac{\partial L}{\partial \mathbf{w}_t}$ we update both parameters:

$$\mathbf{w}_{t+1} = \mathbf{w}_t - \eta \left(\frac{\partial L}{\partial \mathbf{w}_t} + \lambda \mathbf{w}_t \right) \quad (\text{C.6})$$

$$= (1 - \eta \lambda) \mathbf{w}_t - \eta \frac{\partial L}{\partial \mathbf{w}_t} \quad (\text{C.7})$$

$$\hat{\mathbf{w}}_{t+1} = \hat{\mathbf{w}}_t - \eta \left(\frac{\partial L}{\partial \hat{\mathbf{w}}_t} + \lambda \hat{\mathbf{w}}_t \right) \quad (\text{C.8})$$

$$= (1 - \eta \lambda) \hat{\mathbf{w}}_t - \eta \frac{\partial L}{\partial \hat{\mathbf{w}}_t} \quad (\text{C.9})$$

We can then compare the effect the new parameters would have on the next minibatch by calculating $\text{DFT}^{-1}\{\hat{\mathbf{w}}_{t+1}\}$. We then get:

$$\text{DFT}^{-1}\{\hat{\mathbf{w}}_{t+1}\} = \text{DFT}^{-1} \left\{ (1 - \eta \lambda) \hat{\mathbf{w}}_t - \eta \frac{\partial L}{\partial \hat{\mathbf{w}}_t} \right\} \quad (\text{C.10})$$

$$= (1 - \eta \lambda) \mathbf{w}_t - \eta \text{DFT}^{-1} \left\{ \frac{\partial L}{\partial \hat{\mathbf{w}}_t} \right\} \quad (\text{C.11})$$

$$= (1 - \eta \lambda) \mathbf{w}_t - \eta \frac{\partial L}{\partial \mathbf{w}_t} \quad (\text{C.12})$$

$$= \mathbf{w}_{t+1} \quad (\text{C.13})$$

where we have used (C.4) for the second last line. Since the IFT of $\hat{\mathbf{w}}$ is the same as \mathbf{w} at timestep $t+1$ given they are the same at timestep t , and we have set them equal to each other at timestep 1, by induction they will be the same for all time. \square

We have proved the simpler case for simple SGD, but the same result holds when momentum terms are added. This does not however hold for the Adam [50] or Adagrad [51] optimizers, which automatically rescale the learning rates for each parameter based on estimates of the parameter's variance.

Appendix D

DTCWT Single Subband Gains

This appendix proves that the DTCWT *gain layer* proposed in chapter 6 maintains the shift invariant properties of the DTCWT.

Recall that with multirate systems, upsampling by M takes $X(z)$ to $X(z^M)$ and down-sampling by M takes $X(z)$ to $\frac{1}{M} \sum_{k=0}^{M-1} X(W_M^k z^{1/M})$ where $W_M^k = e^{\frac{j2\pi k}{M}}$. We will drop the M subscript below unless it is unclear of the sample rate change, simply using W^k .

D.1 Revisiting the Shift Invariance of the DT \mathbb{C} WT

It is easiest to prove the shift invariance of the gain layer by expanding on the shift invariance of the DT \mathbb{C} WT proofs done in [19].

Let us consider one subband of the DT \mathbb{C} WT. This includes the coefficients from both tree A and tree B. For simplicity in this analysis we will consider the 1-D DT \mathbb{C} WT without the channel parameter c .

If we only keep coefficients from a given subband and set all the others to zero, then we have a reduced tree as shown in Figure D.1. The output $Y(z)$ is:

$$Y(z) = \frac{1}{M} \sum_{k=0}^{M-1} X(W^k z) [A(W^k z)C(z) + B(W^k z)D(z)] \quad (\text{D.1})$$

where the aliasing terms are formed from the addition of the rotated z transforms, i.e. when $k \neq 0$.

Theorem D.1. *The aliasing terms for (D.1) cancel out if the impulse responses of B and D are Hilbert transforms of the impulse responses of A and C respectively, and the system in Figure D.1 is nearly shift invariant.*

Proof. See section 4 of [19] for the full proof of this, and section 7 for the bounds on what ‘nearly’ shift invariant means. \square

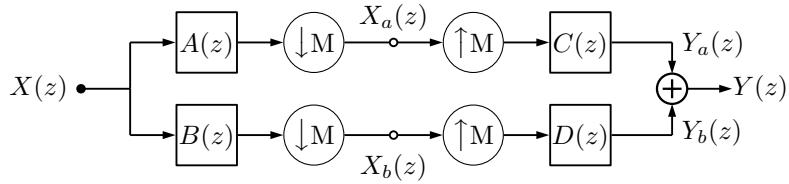


Figure D.1: **Block Diagram of 1-D DTCTWT.** Note the top and bottom paths are through the wavelet or scaling functions from just level m ($M = 2^m$). Figure based on Figure 4 in [19].

Consider the complex filters defined as:

$$P(z) = \frac{1}{2}(A(z) + jB(z)) \quad (\text{D.2})$$

$$Q(z) = \frac{1}{2}(C(z) - jD(z)) \quad (\text{D.3})$$

and define $P^*(z) = \sum_n p^*[n]z^{-n}$ as the Z -transform of p after taking the complex conjugate of the filter taps.

From this, we can rewrite the filters A, B, C and D as:

$$A(z) = P(z) + P^*(z) \quad (\text{D.4})$$

$$B(z) = -j(P(z) - P^*(z)) \quad (\text{D.5})$$

$$C(z) = Q(z) + Q^*(z) \quad (\text{D.6})$$

$$D(z) = j(Q(z) - Q^*(z)) \quad (\text{D.7})$$

Substituting these into (D.1) gives:

$$A(W^k z)C(z) + B(W^k z)D(z) = 2P(W^k z)Q(z) + 2P^*(W^k z)Q^*(z) \quad (\text{D.8})$$

Using (D.8), Kingsbury show that if B is the Hilbert pair of A then P has support only on the right hand side of the frequency plane. Similarly if D is the Hilbert pair of C then Q also has support only on the right hand side of the frequency plane. If P and Q are single side band, then so are P^* and Q^* , but they now have support only on the left hand side of the frequency plane.

Given these properties, the shifted versions of $P(W^k z)$ have negligible overlap with $Q(z)$ except for $k = 0$ (the wanted term) and $k = \pm 1$ where the transition bands overlap. Similarly $P^*(W^k z)$ only overlaps with $Q^*(z)$ when $k = 0$ and a small amount for $k = \pm 1$. [19] quantify the amount of transition band overlap and show it is negligible.

This means $A(W^k z)C(z) + B(W^k z)D(z) = 0$ when $k \neq 0$ and (D.1) reduces to:

$$Y(z) = \frac{1}{M}X(z)[A(z)C(z) + B(z)D(z)] \quad (\text{D.9})$$

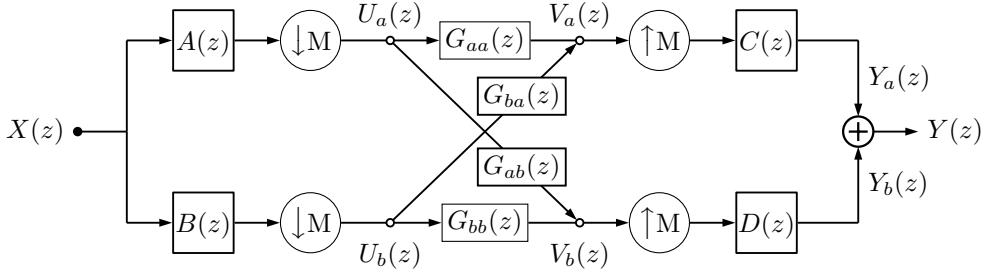


Figure D.2: Block Diagram of 1-D DTWT with subband gains.

D.2 Gains in the Subbands

Figure D.2 shows a block diagram of the extension of the above to general gains. This is a two port network with four individual transfer functions. Let the transfer function from U_i to V_j be G_{ij} for $i, j \in \{a, b\}$. Then V_a and V_b are:

$$V_a(z) = U_a(z)G_{aa}(z) + U_b(z)G_{ba}(z) \quad (\text{D.10})$$

$$= \frac{1}{M} \sum_k X(W^k z^{1/M}) [A(W^k z^{1/M})G_{aa}(z) + B(W^k z^{1/M})G_{ba}(z)] \quad (\text{D.11})$$

$$V_b(z) = U_a(z)G_{ab}(z) + U_b(z)G_{bb}(z) \quad (\text{D.12})$$

$$= \frac{1}{M} \sum_k X(W^k z^{1/M}) [A(W^k z^{1/M})G_{ab}(z) + B(W^k z^{1/M})G_{bb}(z)] \quad (\text{D.13})$$

Further, Y_a and Y_b are:

$$Y_a(z) = C(z)V_a(z^M) \quad (\text{D.14})$$

$$Y_b(z) = D(z)V_b(z^M) \quad (\text{D.15})$$

Then the output is:

$$Y(z) = Y_a(z) + Y_b(z) \quad (\text{D.16})$$

$$\begin{aligned} &= \frac{1}{M} \sum_{k=0}^{M-1} X(W^k z) [A(W^k z)C(z)G_{aa}(z^M) + B(W^k z)D(z)G_{bb}(z) + \\ &\quad B(W^k z)C(z)G_{ba}(z^M) + A(W^k z)D(z)G_{ab}(z)] \end{aligned} \quad (\text{D.17})$$

Theorem D.2. If we let $G_{aa}(z) = G_{bb}(z) = G_r(z)$ and $G_{ab}(z) = -G_{ba}(z) = G_i(z)$ then the end to end transfer function is shift invariant.

Proof. Using the above substitutions, the terms in the square brackets of (D.17) become:

$$G_r(z^M) [A(W^k z)C(z) + B(W^k z)D(z)] + G_i(z^M) [A(W^k z)D(z) - B(W^k z)C(z)] \quad (\text{D.18})$$

[Theorem D.1](#) already showed that the G_r terms are shift invariant and reduce to $A(z)C(z) + B(z)D(z)$. To prove the same for the G_i terms, we follow the same procedure. Using our definitions of A, B, C, D from [Theorem D.1](#) we note that:

$$A(W^k z)D(z) - B(W^k z)C(z) = j \left[P(W^k z) + P^*(W^k z) \right] [Q(z) - Q^*(z)] + \quad (\text{D.19})$$

$$j \left[P(W^k z) - P^*(W^k z) \right] [Q(z) + Q^*(z)] \quad (\text{D.20})$$

$$= 2j \left[P(W^k z)Q(z) - P^*(W^k z)Q^*(z) \right] \quad (\text{D.21})$$

We note that the difference between the G_r and G_i terms is just in the sign of the negative frequency parts, i.e. $AD - BC$ is the Hilbert pair of $AC + BD$. To prove shift invariance for the G_r terms in [Theorem D.1](#), we ensured that $P(W^k z)Q(z) \approx 0$ and $P^*(W^k z)Q^*(z) \approx 0$ for $k \neq 0$. We can use this again here to prove the shift invariance of the G_i terms in (D.18). This completes our proof. \square

Using [Theorem D.2](#), the output is now

$$Y(z) = \frac{2}{M} X(z) [G_r(z^M)(AC + BD) + G_i(z^M)(AD - BC)] \quad (\text{D.22})$$

$$= \frac{2}{M} X(z) [G_r(z^M)(PQ + P^*Q^*) + G_i(z^M)(PQ - P^*Q^*)] \quad (\text{D.23})$$

where we have dropped the z terms on A, B, C, D, P, Q for brevity.

Theorem D.3. *If we treat the two subband coefficients as a complex value $U(z) = U_a(z) + jU_b(z)$ then doing a complex multiply by a gain $G(z) = G_r(z) + jG_i(z)$ maintains shift invariance.*

Proof. This follows from the conditions in [Theorem D.2](#). There we saw that we maintained shift invariance if $G_{aa}(z) = G_{bb}(z) = G_r(z)$ and $G_{ab}(z) = -G_{ba}(z) = G_i(z)$. If we consider V as a complex signal given by $V(z) = V_a(z) + jV_b(z)$, we can see from [Figure D.2](#) that the real and imaginary parts of V are:

$$V_a(z) = G_r(z)U_a(z) - G_i(z)U_b(z) \quad (\text{D.24})$$

$$V_b(z) = G_r(z)U_b(z) + G_i(z)U_a(z) \quad (\text{D.25})$$

which follows the form of a complex multiply. \square

Now we know can assume that our DT \mathbb{C} WT is well designed and extracts frequency bands at local areas, then our complex filter $G(z) = G_r(z) + jG_i(z)$ allows us to modify these passbands (e.g. by simply scaling if $G(z) = C$, or by more complex functions. [Theorem D.3](#) and (D.23) give us an intuition for the real and imaginary parts of a complex gain G . The real part G_r affects how much of the bandpass gain $PQ + P^*Q^*$ propagates through, and the imaginary part G_i affects how much its Hilbert pair $PQ - P^*Q^*$ propagates.

Appendix E

Complex CNN Operations

This appendix lists some of the forward and backward equations for the complex operations we use in the *gain layer* and *wave layer*.

E.1 Convolution

Let us represent the complex input with $\mathbf{x} = \mathbf{x}_r + j\mathbf{x}_i$, which is of shape $\mathbb{C}^{C \times n_1 \times n_2}$. We call $\mathbf{y} = \mathbf{y}_r + j\mathbf{y}_i$ the result we get from convolving \mathbf{x} with $\mathbf{h} = \mathbf{h}_r + j\mathbf{h}_i$, so $\mathbf{y} \in \mathbb{C}^{1 \times (n_1+m-1) \times (n_2+m-1)}$. With appropriate zero or symmetric padding, we can make \mathbf{y} have the same spatial shape as \mathbf{x} . Now, consider the full complex convolution to get \mathbf{y} :

$$y[\mathbf{n}] = \sum_{c=0}^{C-1} \sum_{\mathbf{k}} h[c, \mathbf{k}] x[c, \mathbf{n} - \mathbf{k}] \quad (\text{E.1})$$

We can expand this with real and imaginary terms:

$$\begin{aligned} y[\mathbf{n}] &= \sum_{c=0}^{C-1} \sum_{\mathbf{k}} (h_r[c, \mathbf{k}] + jh_i[c, \mathbf{k}]) (x_r[c, \mathbf{n} - \mathbf{k}] + jx_i[c, \mathbf{n} - \mathbf{k}]) \\ &= \sum_{c=0}^{C-1} \sum_{\mathbf{k}} (x_r[c, \mathbf{n} - \mathbf{k}] h_r[c, \mathbf{k}] - x_i[c, \mathbf{n} - \mathbf{k}] h_i[c, \mathbf{k}]) \\ &\quad + j \sum_{c=0}^{C-1} \sum_{\mathbf{k}} (x_r[c, \mathbf{n} - \mathbf{k}] h_i[c, \mathbf{k}] + x_i[c, \mathbf{n} - \mathbf{k}] h_r[c, \mathbf{k}]) \\ y_r[\mathbf{n}] + jy_i[\mathbf{n}] &= \sum_{c=0}^{C-1} ((x_r * h_r) - (x_i * h_i)) [c, \mathbf{n}] + ((x_r * h_i) + j(x_i * h_r)) [c, \mathbf{n}] \end{aligned} \quad (\text{E.2})$$

Unsurprisingly, complex convolution is the sum and difference of 4 real convolutions.

We can use this fact to find the update and passthrough gradients for complex convolution.

Update Gradients: We need to find $\frac{\partial L}{\partial h_r}$ and $\frac{\partial L}{\partial h_i}$. From (E.2) we can apply the chain rule and the properties of real convolutions to write:

$$\frac{\partial L}{\partial h_r} = \frac{\partial L}{\partial y_r} \frac{\partial y_r}{\partial h_r} + \frac{\partial L}{\partial y_i} \frac{\partial y_i}{\partial h_r} \quad (\text{E.3})$$

$$= \Delta y_r \star x_r + \Delta y_i \star x_i \quad (\text{E.4})$$

where \star is the correlation operation (see [subsubsection 2.4.1.2](#)). Similarly

$$\frac{\partial L}{\partial h_i} = -\Delta y_r \star x_i + \Delta y_i \star x_r \quad (\text{E.5})$$

Passthrough Gradients: Again with application of the chain rule and the properties of real convolution, we have:

$$\frac{\partial L}{\partial x_r} = \frac{\partial L}{\partial y_r} \frac{\partial y_r}{\partial x_r} + \frac{\partial L}{\partial y_i} \frac{\partial y_i}{\partial x_r} \quad (\text{E.6})$$

$$= \Delta y_r \star h_r + \Delta y_i \star h_i \quad (\text{E.7})$$

$$\frac{\partial L}{\partial x_i} = -\Delta y_r \star h_i + \Delta y_i \star h_r \quad (\text{E.8})$$

E.2 Regularization

We must be careful with regularizing complex weights. We want to promote the magnitude of the weights to be small but allow the phase to change unrestricted.

To do ℓ_2 regularization we can apply ℓ_2 to the real and imaginary parts independently:

$$\|r\|_2^2 = \left\| \sqrt{x_r^2 + x_i^2} \right\|_2^2 = \sum_{\mathbf{n}} x_r[\mathbf{n}]^2 + x_i[\mathbf{n}]^2 = \sum_{\mathbf{n}} x_r[\mathbf{n}]^2 + \sum_{\mathbf{n}} x_i[\mathbf{n}]^2 = \|x_r\|_2^2 + \|x_i\|_2^2 \quad (\text{E.9})$$

But this does not hold for ℓ_1 regularization as:

$$\|r\|_1 = \left\| \sqrt{x_r^2 + x_i^2} \right\|_1 = \sum_{\mathbf{n}} \sqrt{x_r[\mathbf{n}]^2 + x_i[\mathbf{n}]^2} \neq \|x_r\|_1 + \|x_i\|_1 \quad (\text{E.10})$$

Also, for ℓ_1 regularization the derivatives have a singularity at the complex origin as:

$$\frac{\partial \ell_1}{\partial x_r[\mathbf{n}]} = \frac{x_r[\mathbf{n}]}{\sqrt{x_r[\mathbf{n}]^2 + x_i[\mathbf{n}]^2}} \quad (\text{E.11})$$

is not defined when $x_r = x_i = 0$. A similar problem was mentioned in [subsection 3.6.1](#) where we wanted to pass gradients through the magnitude operation of a ScatterNet. Since we do not explicitly care if weights are zero or near zero, we choose to handle this by setting the

gradient at the origin to be 0. It is unlikely our weights will ever be zero with this method, but if they are, we cover the case of dividing by zero.

E.3 ReLU applied to the real and imaginary parts independently

If we define a nonlinearity to be $y = \sigma(x)$ where:

$$y = y_r + jy_i = \max(0, x_r) + j\max(0, x_i) \quad (\text{E.12})$$

then the passthrough gradients are:

$$\frac{\partial L}{\partial x_r} = \Delta y_r \mathbb{I}(x_r > 0) \quad (\text{E.13})$$

$$\frac{\partial L}{\partial x_i} = \Delta y_i \mathbb{I}(x_i > 0) \quad (\text{E.14})$$

E.4 Soft Shrinkage

Let $z = x + jy$ and $w = u + jv = \mathcal{S}(z, t)$ where we define the soft shrinkage on a complex number $z = re^{j\theta}$ by a real threshold t as:

$$\mathcal{S}(z, t) = \begin{cases} 0 & r < t \\ (r - t)e^{j\theta} & r \geq t \end{cases} \quad (\text{E.15})$$

This can alternatively be written as:

$$\mathcal{S}(z, t) = \frac{\max(r - t, 0)}{r} z \quad (\text{E.16})$$

$$= gz \quad (\text{E.17})$$

To find the pass through gradients $\frac{\partial L}{\partial x}$, $\frac{\partial L}{\partial y}$ and update equations $\frac{\partial L}{\partial t}$ we need to find all the real and imaginary partial derivatives. We can apply the product rule once we find $\frac{\partial g}{\partial x}$, $\frac{\partial g}{\partial y}$, $\frac{\partial g}{\partial t}$:

$$\frac{\partial g}{\partial x} = \begin{cases} 0 & r < t \\ \frac{r \frac{\partial r}{\partial x} - (r-t) \frac{\partial r}{\partial x}}{r^2} & r \geq t \end{cases} \quad (\text{E.18})$$

$$= \frac{xt\mathbb{I}(g > 0)}{r^3} \quad (\text{E.19})$$

$$\frac{\partial g}{\partial y} = \frac{yt\mathbb{I}(g > 0)}{r^3} \quad (\text{E.20})$$

$$(\text{E.21})$$

$$\frac{\partial g}{\partial t} = \frac{-\mathbb{I}(g > 0)}{r} \quad (\text{E.22})$$

Then from the definition of $w = u + jv$ we have $u = gx$ and $v = gy$, giving us:

$$\frac{\partial u}{\partial x} = g + \frac{x^2 t \mathbb{I}(g > 0)}{r^3} \quad (\text{E.23}) \quad \frac{\partial v}{\partial x} = \frac{xyt \mathbb{I}(g > 0)}{r^3} \quad (\text{E.26})$$

$$\frac{\partial u}{\partial y} = \frac{xyt \mathbb{I}(g > 0)}{r^3} \quad (\text{E.24}) \quad \frac{\partial v}{\partial y} = g + \frac{y^2 t \mathbb{I}(g > 0)}{r^3} \quad (\text{E.27})$$

$$\frac{\partial u}{\partial t} = \frac{-x \mathbb{I}(g > 0)}{r} \quad (\text{E.25}) \quad \frac{\partial v}{\partial t} = \frac{-y \mathbb{I}(g > 0)}{r} \quad (\text{E.28})$$

Putting it all together, our update and passthrough gradients are:

$$\frac{\partial L}{\partial x} = \frac{xt\mathbb{I}(g > 0)}{r^3} (x\Delta u + y\Delta v) + g\Delta u \quad (\text{E.29})$$

$$\frac{\partial L}{\partial y} = \frac{yt\mathbb{I}(g > 0)}{r^3} (x\Delta u + y\Delta v) + g\Delta v \quad (\text{E.30})$$

$$\frac{\partial L}{\partial t} = \frac{-\mathbb{I}(g > 0)}{r} (x\Delta u + y\Delta v) \quad (\text{E.31})$$

These equations are for point-wise application of soft-thresholding. When applied to an entire image, then we sum $\frac{\partial L}{\partial t}$ over all locations.

E.5 Batch Normalization and ReLU applied to the Complex Magnitude

Again let $z = x + jy = re^{j\theta}$ and $w = u + jv = \text{ReLU}(\text{BN}(r))e^{j\theta}$. In (6.5.11) we showed that this nonlinearity is equivalent to soft-thresholding with threshold $t = \mu_r - \frac{\sigma_r \beta}{\gamma}$ and multiplied by a learned gain γ divided by the tracked standard deviation of r : $\frac{\gamma}{\sigma_r}$.

Let us call the action of this nonlinearity \mathcal{B} , defined by:

$$\mathcal{B}(z, \gamma, \beta) = \begin{cases} 0 & r < t \\ \frac{\gamma}{\sigma_r}(r-t)e^{j\theta} & r \geq t \end{cases} \quad (\text{E.32})$$

$$= \mathbb{I}(r > t) \frac{\gamma(r-t)}{\sigma r} z \quad (\text{E.33})$$

$$= g'z \quad (\text{E.34})$$

It is clear from our new definition of g' that the equations (E.29) - (E.31) are now scaled by $\frac{\gamma}{\sigma}$. This immediately gives us the passthrough gradients. For the update equations, we must find some additional information:

$$\frac{\partial t}{\partial \beta} = -\frac{\sigma}{\gamma} \quad (\text{E.35})$$

$$\frac{\partial t}{\partial \gamma} = \frac{\sigma \beta}{\gamma^2} \quad (\text{E.36})$$

$$\frac{\partial g'}{\partial \beta} = \frac{-\gamma \mathbb{I}(g' > 0)}{\sigma r} \frac{\partial t}{\partial \beta} = \frac{\mathbb{I}(g' > 0)}{r} \quad (\text{E.37})$$

$$\frac{\partial g'}{\partial \gamma} = \frac{\mathbb{I}(g' > 0)}{\sigma r} \left(r - t - \frac{\partial t}{\partial \gamma} \right) = \frac{\mathbb{I}(g' > 0)}{\sigma r} \left(r - t - \frac{\sigma \beta}{\gamma^2} \right) \quad (\text{E.38})$$

Therefore, combinining these with (E.31) we get:

$$\frac{\partial L}{\partial \beta} = \frac{\mathbb{I}(g' > 0)}{r} (x \Delta u + y \Delta v) \quad (\text{E.39})$$

$$\frac{\partial L}{\partial \gamma} = \frac{\mathbb{I}(g' > 0)}{\sigma r} \left(r - t - \frac{\sigma \beta}{\gamma^2} \right) (x \Delta u + y \Delta v) \quad (\text{E.40})$$

Appendix F

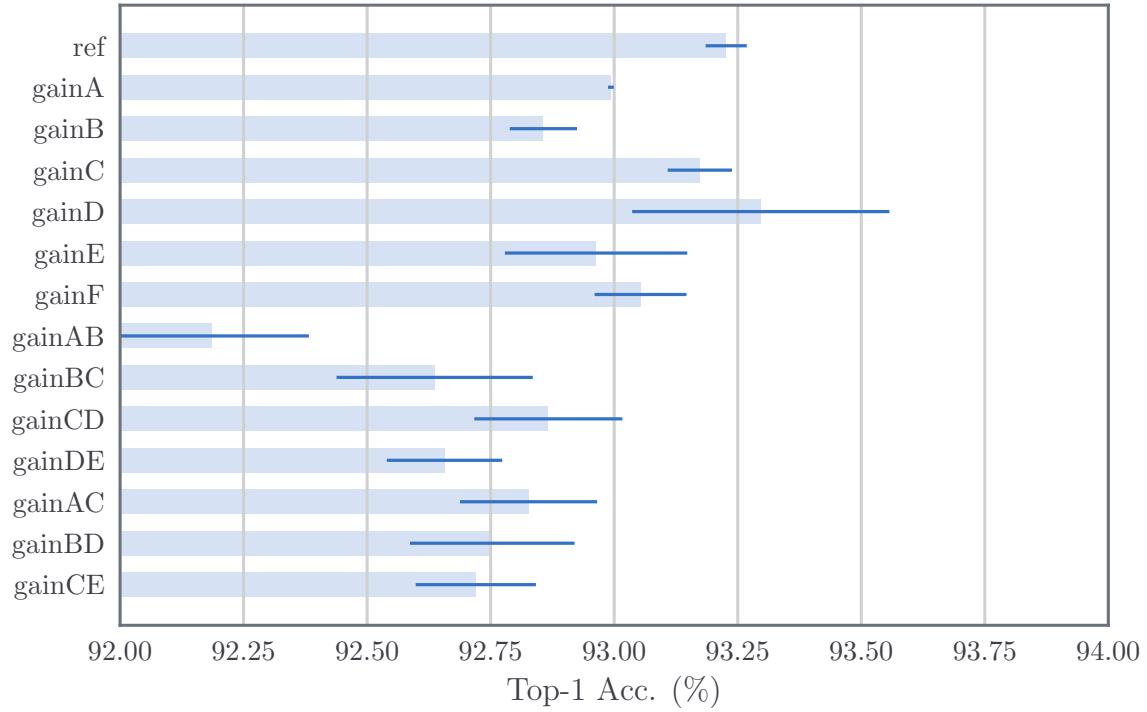
GainLayer Additional Results

This appendix presents some additional results for the gain layer experiments from [subsection 6.4.2](#). Here, we do ablation experiments similar to those done on the invariant layer in [section 5.7](#). In particular, we use the architecture described in [Table 5.5](#) and progressively swap out convolutional layers with gain layers. Again, we run this on CIFAR-10, CIFAR-100 and Tiny ImageNet.

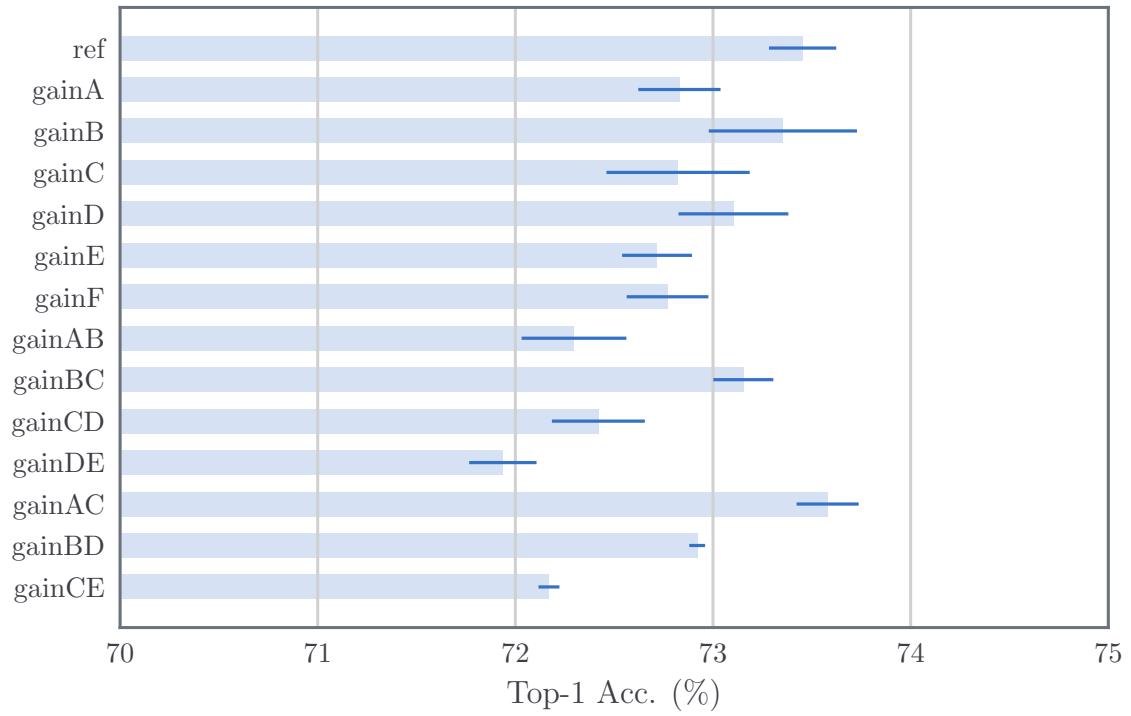
We use the same naming technique as in the previous chapter, calling a network ‘gainX’ means that the ‘convX’ layer was replaced with a wavelet gain layer, but otherwise keeping the rest of the architecture the same. Recall that the CIFAR architectures in [Table 5.5](#) have 6 convolutional layers called ‘convA’ to ‘convF’, and the Tiny ImageNet architectures have 8 convolutional layers called ‘convA’ to ‘convH’.

We train all our networks for with stochastic gradient descent with momentum. The initial learning rate is 0.5, momentum is 0.85, batch size $N = 128$ and weight decay is 10^{-4} . For CIFAR-10/CIFAR-100 we scale the learning rate by a factor of 0.2 after 60, 80 and 100 epochs, training for 120 epochs in total. For Tiny ImageNet, the rate change is at 18, 30 and 40 epochs (training for 45 in total).

[Figure F.1](#) and [Figure F.2](#) show the results from these experiments for CIFAR and Tiny ImageNet. Just as with the large kernel ablation study in [subsection 6.4.2](#), adding a gainlayer typically degrades performance by a small amount.



(a) CIFAR-10



(b) CIFAR-100

Figure F.1: Small kernel ablation results for CIFAR. These graphs show the average and ± 1 standard deviation results for 3 runs. The names on the left represent where the convolutional layer is swapped for a gain layer, with ‘gainAB’ and below indicating two convolutional layers were swapped for gain layers.

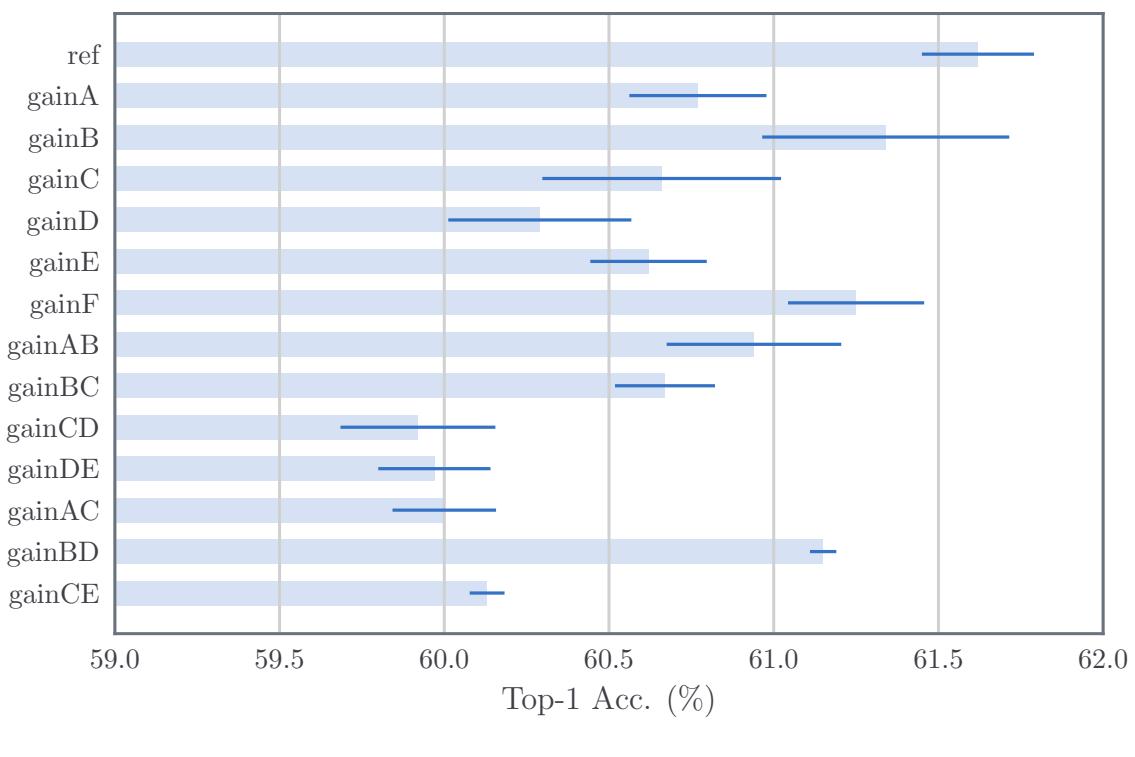


Figure F.2: **Small kernel ablation results for Tiny ImageNet.**

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