



University of Brasilia at Gama – FGA/UnB
Software Engineering

**COMPRESSED SENSING WITH GAUSSIAN PROBABILITISTIC
PRIOR INFORMATION FOR
MAGNETIC RESONANCE IMAGE RECONSTRUCTION**

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RESUMO

A obtenção de imagens de ressonância magnética de alta qualidade é uma tarefa árdua pelo funcionamento das máquinas e pela complexidade dos tecidos analisados, tornando praticamente impossível a obtenção de todas as medidas do objeto de estudo em um exame de breve execução. Para lidar com este problema, diversas técnicas foram aplicadas para gerar imagens de alta qualidade com quantidades menores de medidas, entre elas, Compressed Sensing (CS) tem sido a técnica com melhores resultados e mais aprofundada no contexto de reconstrução de imagens de ressonância magnética nas últimas duas décadas. CS é capaz de produzir imagens de altíssima qualidade com proporções muito menores de amostragens, como 10%, 15% do total de medidas disponíveis utilizando técnicas de *undersampling* por sua boa capacidade de lidar com sinais esparsos.

O desempenho do CS evoluiu com as contribuições de uso de filtros esparsificantes e com o uso de informação a priori. Recentemente, o estado-da-arte veio sendo composto pelo uso de redes neurais profundas, muitas vezes em conjunto com o uso de CS e outras de maneira singular. Desta maneira, esta monografia tem como intuito trazer o uso de CS com pré-filtragem e aplicar o uso de redes adversárias generativas para a extração de informação a priori.

A metodologia é composta pela utilização de formas variadas de CS para a reconstrução de imagens comparando métodos, pré-processamentos e parâmetros, a fim de explicitar os diferentes resultados atingíveis com CS. Experimentos com redes adversárias generativas também foram conduzidos com o intuito de demonstrar a capacidade das mesmas para geração de sinais com alta fidelidade.

Os resultados obtidos demonstram que o uso de pré-filtragem é preferível para a reconstrução de imagens, melhorando as métricas de relação sinal-ruído (SNR), relação sinal-ruído de pico (PSNR) e similaridade estrutural (SSIM). Os resultados das GANs mostram que estas são capazes de gerar sinais com alto teor de realismo quando comparadas as imagens criadas com as imagens de treino. Os próximos passos da pesquisa envolvem aplicar o uso de informação a priori de forma manual e posteriormente usando GANs, comparando todas as técnicas e transformações aplicadas.

Contents

1	Introduction	1
1.1	Context	1
1.2	Scientific Problem Definition and Proposal	2
1.3	Objectives	3
1.3.1	General Objective	3
1.3.2	Specific Objectives	3
2	Machine Learning in MRI Concepts and State-of-Art	4
2.1	Magnetic Resonance Imagery	4
2.1.1	K-space	4
2.1.2	Undersampling	5
2.2	Compressed Sensing	6
2.2.1	Introduction	6
2.2.2	Reconstructing using the L1 Regularization	7
2.3	Prior Information	7
2.3.1	Prior Information Retrieval	7
2.3.2	Context-Dependent Probabilistic Prior Information	8
2.3.3	L1-minimization Adaptation	11

2.3.4	Parameter Fine-tuning	11
2.3.5	Future Contributions	11
2.4	Artificial Neural Networks	12
2.4.1	Biological Inspirations	12
2.4.2	Neuron	12
2.4.3	Multilayer Perceptron	13
2.4.4	Activation Functions	14
2.4.5	Loss Functions	17
2.4.6	Backpropagation	18
2.4.7	Gradient Descent	18
2.5	Generative Adversarial Networks	19
3	Preliminary Reconstructing and Signal Generating Experiments	21
3.1	1-D Direct vs Indirect L1-Minimization	21
3.2	2D Compressed Sensing Reconstruction with Pre Filtered Signal	22
3.2.1	Subsampling	22
3.2.2	Pre-filtering sparsifying transform	23
3.3	Preliminary Tests with Generative Adversarial Networks	24
3.3.1	Data Transformation	25
3.3.2	Generator Network Architecture	25
3.3.3	Discriminator Network Architecture	26
4	Preliminary Reconstruction and Generation Results	27
4.1	1-D Compressed Sensing Reconstruction	27
4.2	Compressed Sensing Reconstruction with Pre Filtered Signal	28

4.2.1	Brain Sagittal Reconstruction	29
4.2.2	Prior Information Cardiac Sagittal Reconstruction	30
4.3	Preliminary Tests with Generative Adversarial Networks	31
5	Conclusion	33

List of Tables

2.1	Activation functions and respective derivatives.	15
2.2	Loss functions and formulas.	17
4.1	Phantom reconstruction metrics.	29
4.2	Sagittal reconstruction metrics.	30
4.3	Sagittal reconstruction metrics.	31

List of Figures

2.1	Single-coil k-space points from fully-sampled knee. Source: [1, 2]	5
2.2	Single-coild fully-sampled knee spatial image. Source: [1, 2]	5
2.3	Under-sampling patterns. (a) Cartesian undersampling , (b) radial undersampling , (c) spiral undersampling , (d) isolated samples in the k-space, according to the realisation of a random process [3].	6
2.4	Fully-sampled cardiac MRI frame.	9
2.5	Manual segmentation: the lighter the colour, the more pixels are dislocated in the region throughout frames.	9
2.6	2D <i>weights</i> matrix visualization.	10
2.7	3D <i>weights</i> matrix visualization.	10
2.8	2D <i>weights</i> matrix visualization.	11
2.9	3D <i>weights</i> matrix visualization.	11
2.10	Schematic of an Artificial Neuron. Source: [4]	13
2.11	Artificial Neural Networks (ANN) Architecture Sample.	14
2.12	Rectified linear units (ReLU) activation function	15
2.13	Leaky ReLU activation function	16
2.14	<i>Tanh</i> activation function. Source: [5]	16
2.15	Sigmoid activation function. Source: [5]	17
2.16	Gradient descent example visualization. Source: [6]	19

2.17 Generative Adversarial Network (GAN) training diagram. Source: [7]	20
3.1 First 10 random 1-D signals	22
3.2 Shepp-Logan phantom reference image.	23
3.3 Spiral undersampling method with 30.95% data points.	23
3.4 2-D High pass horizontal filter.	23
3.5 2-D High pass vertical filter.	23
3.6 2-D High pass diagonal filter.	23
3.7 Sample of digits from MNIST	24
3.8 Generator network architecture of a 3-D image. Source: [8, 9]	26
3.9 Discriminator network architecture of a 3-D image. Source: [8, 9]	26
4.1 Signal-to-Noise Ratio (SNR) x L size for direct and indirect L1-minimization for 200 random 1D signals	28
4.2 zero-filled reconstruction	29
4.3 L1-minimization reconstruction	29
4.4 L1-minimization with pre-filtering reconstruction	29
4.5 Shepp-Logan phantom reference image	29
4.6 Zero-filled reconstruction	30
4.7 L1-minimization reconstruction	30
4.8 L1-minimization with pre-filtering reconstruction	30
4.9 Reference image	30
4.10 Dynamic cardiac zero-filled reconstruction	31
4.11 Dynamic cardiac L1-minimization reconstruction	31
4.12 Dynamic cardiac L1-minimization with pre-filtering reconstruction	31

4.13	Dynamic cardiac reference image	31
4.14	Discriminator fake loss over epochs	32
4.15	Discriminator real loss over epochs	32
4.16	Generator loss over epochs	32
4.17	GAN generated MNIST digits	32
5.1	Activies cronogram	33

NOMENCLATURE AND ABBREVIATIONS

MRI Magnetic Resonance Imaging	1
MR Magnetic Resonance	1
ANN Artificial Neural Networks	x
DL Deep Learning	2
ML Machine Learning	1
GAN Generative Adversarial Network	xi
CNN Convolutional Neural Network	2
GPU Graphics Processing Units	2
CV Computer Vision	2
NLP Natural Language Processing	2
CS Compressed Sensing	1
MLP Multilayer Perceptron	13
DFN Deep Feedforward Networks	13
SNR Signal-to-Noise Ratio	xi

PSNR Peak Signal-to-Noise Ratio	29
DCGAN Deep Convolutional GAN	14
ReLU Rectified linear units	x
SGD Stochastic Gradient Descent	18
IRLS iteratively reweighted least squares	7

1 INTRODUCTION

In this thesis, we propose a GAN approach for prior information extraction to feed a Compressed Sensing (CS) algorithm, aiming to reconstruct images with both reduced signal-to-noise error and less acquisition time compared to conventional CS. Achieving higher quality with a reduced number of samples allows faster exam procedures, making Magnetic Resonance Imaging (MRI) cheaper, faster and more convenient for both patients and clinics.

1.1 Context

MRI is a widely used imaging modality in medical practice because of its great tissue contrast capabilities, it has evolved into the richest and most versatile biomedical imaging technique today [10], making MRI the best option for medical imaging whenever it is possible to use.

However, like everything in life, there is a trade-off to consider when using MRI. Typically, reconstructing an MRI is an ill-posed linear inverse task (a problem that has either none or infinite solutions in the desired class). Problems of this nature impose a trade-off between *accuracy* and *speed* [11]. The information obtained from Magnetic Resonance (MR) is commonly represented by individual samples in the k-space, which translates to the Fourier transform of the image to be reconstructed [12]. This MR sampling sparse nature makes CS a liable technique to use when reconstructing MRI, hence we here propose a novel CS prior information approach for better results.

CS has been for years the state-of-art technique in MRI reconstruction and has been improved later by the use of sparsifying pre-filtering techniques and prior information [13, 14]. CS uses the premise that given a signal with a sparse representation in some known domain, it is possible to reconstruct the signal using limited linear measurements taken from a non-sparse representation.

Machine Learning (ML) methods have been utterly developed and improved recently

with the use of higher computing power derived from the invention of Graphics Processing Units (GPU) and other hardware improvements, allowing ANN to come to practicality. These ANN models, often referenced as Deep Learning (DL), have become the state-of-art in various areas, such as Computer Vision (CV), Natural Language Processing (NLP), Recommendation Systems, amongst other fields [15, 16, 17]. These fast-paced developments led to improvements in medical data processing using DL as well. ML techniques can be used in several different manners to improve medical analysis, here we focus on applying GAN in the process of attaining improved prior information to feed the CS algorithm obtaining higher signal-to-noise ratios and faster computation procedures.

1.2 Scientific Problem Definition and Proposal

MRI is great for high-quality tissue images, but there are some drawbacks: MRI exams are often very long and require the patient to be in a static position throughout the whole process, this makes the exam challenging for patients that have difficulties in keeping a still position for several minutes. Another intrinsic complication in MRI procedures is that it is nearly impossible to get images from moving tissues like a beating heart or flowing blood veins as that would require an enormous amount of samples, which with current technologies used in clinics is not viable. Algorithms that reconstruct MRI try to tackle this sampling issue by producing the best possible quality images for the least amount of samples collected, making the exams faster and less sample-dependent.

CS algorithms have been the state-of-art in MRI reconstruction for the past few years, and now with the advances of DL, new techniques are being produced taking advantages of how ANN are powerful in imaging processing, especially Convolutional Neural Network (CNN) and, more recently, GAN networks are becoming the new state-of-art techniques in several computer vision areas. Most CS contributions without the use of machine learning cite the usage of sparsifying pre-filtering techniques and prior information that have been proven to improve efficiency and achieve better reconstructions [14, 13]. More recently, the contributions in MRI reconstruction has been more focused in deep learning approaches, with some architectures using CS along ANNs [18, 19, 20, 21, 22, 23].

Prior information for CS can go from previous MRI frames and exams to even medical records. Prior information is normally generated by simplistic mathematical approaches like filtering and thresholding on the images. Besides the simpler technique applied, these information extraction procedures oftentimes is restricted to few frames and does not take into account the nature of organs and tissues structures, a feature that DL should be able to identify and use to generate better quality information. This means that there is a lot

of room for improvement towards prior information engineering techniques, as DL models have been proven superior in tasks of this nature.

Within this context, we propose a modern prior information engineering system with the usage of GAN, aiming for higher quality prior information to feed the CS and reducing the number of samples dependability. This will reduce the number of samples needed, making the MRI exams faster and, consequently, cheaper.

1.3 Objectives

1.3.1 General Objective

This thesis' goal is to develop an MR prior information system retriever based on GAN architecture to analyse if the quality of the prior information fed to CS algorithms can be improved, hence improving quality in reconstructed MRI and decreased necessity for larger sampling.

1.3.2 Specific Objectives

In order to achieve the general objective described above, we have set the following specific goals:

- Implement direct and indirect CS MRI reconstruction algorithm using undersampled *k*-space measurements.
- Evaluate CS MRI reconstructions against fully sampled MRI.
- Implement a GAN for ordinary image generation with the well-known MNIST dataset.
- Implement a GAN architecture for optimized prior information retrieval.
- Evaluate the use of GAN architecture for prior information retrieval against state-of-art MRI reconstruction techniques.

2 MACHINE LEARNING IN MRI CONCEPTS AND STATE-OF-ART

2.1 Magnetic Resonance Imagery

MRI is an indirect process that produces cross-sectional images with high spatial resolution from nuclear magnetic resonances, gradient fields and hydrogen atoms of the subject’s anatomy [24]. The acquisition of these measurements is performed by a measuring instrument called *receiver coil* and it can be done by using one receiver coil or in some cases with multiple coils [1, 2]. These receiver coils are placed in proximity to a specific region in the subject to be imaged. During the imaging process, the MRI machine generates a sequence of spatially and temporally-varying magnetic fields which induce the body to emit resonant electromagnetic response fields which are then measured by the receiver coil [1, 2].

2.1.1 K-space

The k-space is the output generated by the MRI machine scan after extracting measurements from a given subject tissue. The k-space is represented in the spatial frequency in two or three dimensions of a subject and may also be referred to as the Fourier space. This k-space representation contains an implicit sparsity that is exploited when performing undersampling [25] and reinforce the usage of algorithms like CS for MRI reconstruction as CS depends on signals that have a sparse representation in an orthonormal basis.

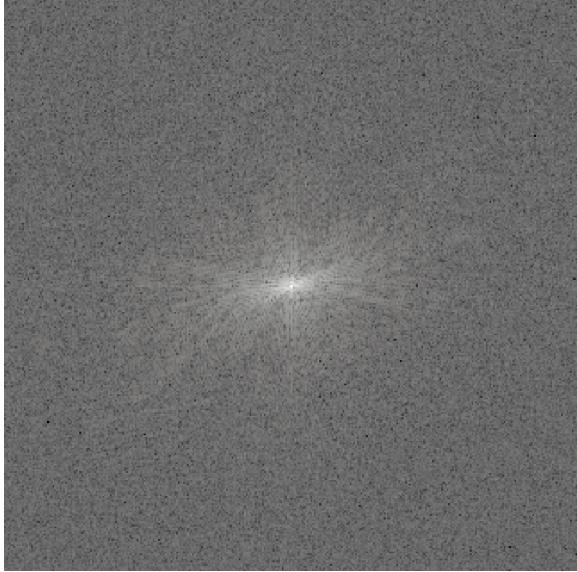


Figure 2.1. Single-coil k-space points from fully-sampled knee.
Source: [1, 2]



Figure 2.2. Single-coil fully-sampled knee spatial image.
Source: [1, 2]

2.1.2 Undersampling

The time required to acquire all the measurements responses from every single atom in a subject would be extremely high and problematic to everyone involved (patients, doctors and clinics). The way machines can do faster MRI is by performing *undersampling*, also referred as subsampling and sampling, when scanning the subject.

Undersampling is performed by giving the machine a known prescribed path in which it will extract measurements from the multidimensional k-space representation. This allows machines to collect only a fraction of data measurements needed for image reconstruction hence speeding up the data acquisition process without critical quality loss.

There are some undersampling patterns to use and each has its benefits depending on several parameters, such as: the subject's region extraction, the algorithm used for reconstruction, acquisition time.

In the figure below we can see some of the most used patterns. In this research, we will focus mostly on the cartesian undersampling method, as that is the one used in the FastMRI dataset [1, 2], which we will use for our experiments.

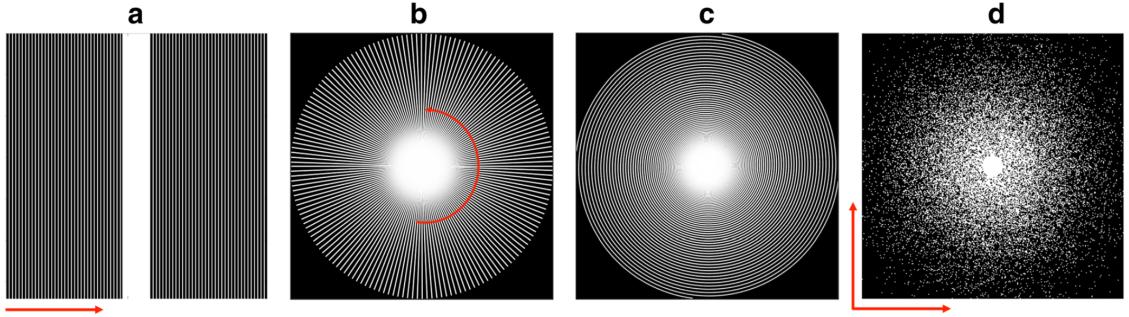


Figure 2.3. Under-sampling patterns. (a) **Cartesian undersampling**, (b) **radial undersampling**, (c) **spiral undersampling**, (d) **isolated samples in the k-space**, according to the realisation of a random process [3].

2.2 Compressed Sensing

2.2.1 Introduction

Compressed sensing, often referred to as Compressive Sensing, represented a major breakthrough in signal processing when it was introduced by Donoho, Candès, Romberg, and Tao in 2004 [26, 27, 28] since it allows a sampling at a rate much lower than the Nyquist-Shannon's theorem: the sampling signal rate must be at least twice the maximum frequency present in the signal (Nyquist rate).

CS is a powerful algorithm that implements a novel technique for the acquisition of signals of sparse or compressible nature. CS theory aims to provide accurate reconstruction of unknown sparse signals from underdetermined linear measurements. CS parts from the principle that if given x , a digital image or signal has a sparse representation in an orthonormal basis (e.g. wavelet, Fourier, cosine, etc), then the N most important coefficients in that expansion allow reconstruction with l_2 error $O(N^{1/2-1/p})$ [26].

The idea was inspired from questioning the necessity of extracting large portions of samples when much of these samples are discarded, exposing the inefficiency of trying to gather all signal.

“Why go to so much effort to acquire all the data when most of what we get will be thrown away? Can we not just directly measure the part that will not end up being thrown away?” [26]

In MRI, the so-called *k-space* is a sparse representation of the image to be reconstructed and when it is undersampled, the Nyquist criterion is violated [25]. These are favorable

conditions for the CS theory, hence all thousands of papers citing the usage of CS for MRI reconstruction and state-of-the-art results involving any application of CS for MRI reconstruction.

2.2.2 Reconstructing using the L1 Regularization

2.3 Prior Information

The application of prior information in MRI reconstruction was first introduced with the general idea of exploiting the common information shared throughout sequential frames acquired from the MR scan [12, 14]. This would not only improve the reconstruction of MR images but would also make dynamic scans more feasible as large portions of the image could be used as prior information for the reconstruction of future and previous frames. The MRI exam is known for requiring the patient to stand still throughout the exam so that there are not grand differences between different frames.

A dynamic cardiac MRI for instance is naturally a situation that will be impossible for the subject to be still throughout the exam, as the heart will keep beating constantly; hence increasing the level of difficulty for the reconstruction. Although a dynamic cardiac exam has this innate motion characteristic, it also contains crucial portions of *support regions* shared between most of the frames that can be exploited to improve the reconstruction. These support regions are essentially structures that will hold – practically – the same position therefore, these are elements in the image that share the same information and prior information application exploits this nature by slightly reducing the number of variables in our underdetermined system in the CS algorithm.

2.3.1 Prior Information Retrieval

To leverage prior information for a frame – as it was first introduced –, one could apply edge detection filters to first generate support position candidates and then extract the position of these edges and apply a τ factor to increase the value in these specific positions when running the CS algorithm, based on an iteratively reweighted least squares (IRLS) method [14]. In practice, this operation tells the l_p -minimization that those are values with a higher magnitude and also, likely a support position. These procedures are extensively explored in the experiments documented in the experiments chapter.

Prior information was then introduced as a deterministic approach – whether the pixel

was a support position or it was not. This approach, of course, raises a few questions: what if the subject moved slightly and the support position has been moderately altered? What if the next frame contains support positions that are getting narrower or wider? E.g. Brain MRIs, where the cranial structure's diameter is increased or decreased frame-by-frame. Prior information theory responds to these question with a high tolerance impact in the reconstruction. It has been documented that using the same amount of mistaken support positions and correct ones will improve SNR in the next frame's reconstruction.

□

Alternatively, a non-deterministic approach for prior information generation was proposed to achieve more robustness based on more accurate information. This technique uses a Gaussian – or normal – distribution with fixed covariance matrix value for each pixel of the support positions in the space domain instead of a deterministic 0 or 1 value. Using probabilistic support positions improves reconstruction metrics as shown in □

2.3.2 Context-Dependent Probabilistic Prior Information

We propose a novel approach to the non-deterministic prior information extraction by using knowledge about regions in the image that are more dislocated than others to apply different variance in the normal distributions procedure. When moving from a frame to another, portions of the image are dislocated in different degrees, that is, some regions will be more likely to contain support positions than others. By taking this into account, we segmented the image into four levels of movement – *high*, *medium*, *low* and *close-to-zero* movement. The regions labelled as *high* movement receive a covariance matrix with a higher variance; thus assigning this region as a low-confidence support position, whilst the regions labelled as *low* or *close-to-zero* receive a Gaussian distribution with low variance, representing it as a highly likely support position prior information. Similarly, the *medium* labeled positions are assigned an intermediate variance for an intermediate confidence.

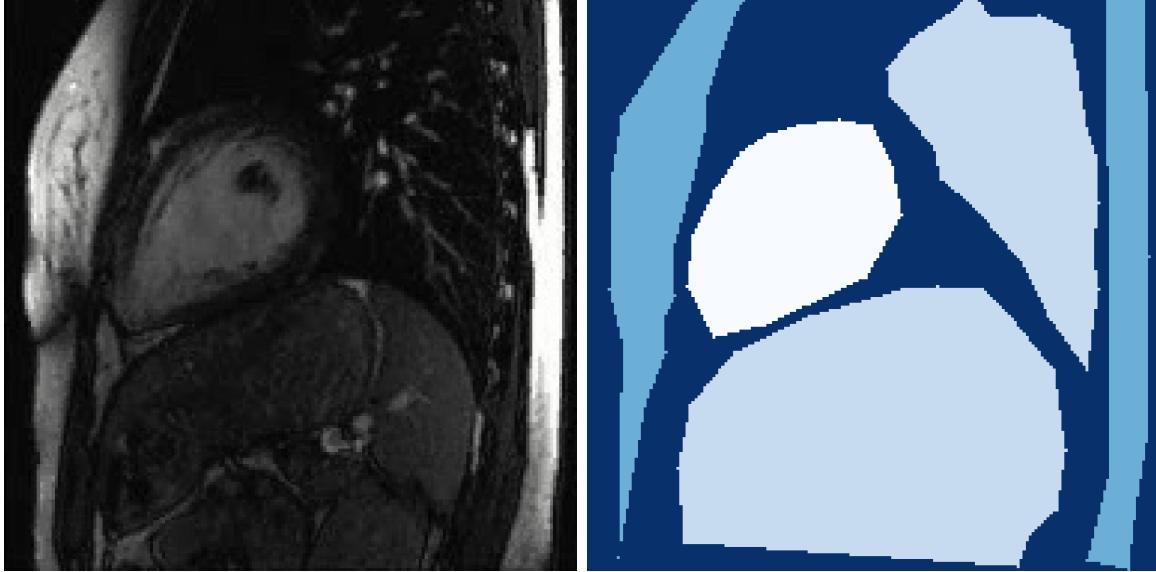


Figure 2.4. Fully-sampled cardiac MRI frame.

Figure 2.5. Manual segmentation: the lighter the colour, the more pixels are dislocated in the region throughout frames.

Identifying these different regions gives us the possibility to apply different weights for each region in the image. The regions labelled as *high* movement receive a covariance matrix with a higher variance; thus assigning this region as a low-confidence support position, whilst the regions labelled as *low* or *close-to-zero* receive a Gaussian distribution with low variance, representing it as a support position prior information. A *weights* matrix is generated by computing an element-wise Gaussian distribution with the according covariance matrix on the segmented mask and then adding these distributions each iteration, forming accumulations on the centre of several support positions. The regions labelled *high* are specifically multiplied by -1 so that the confidence is represented with negative magnitudes. After computing the *weights* matrix, we expect to see distinguishably different magnitude levels, representing different confidences.

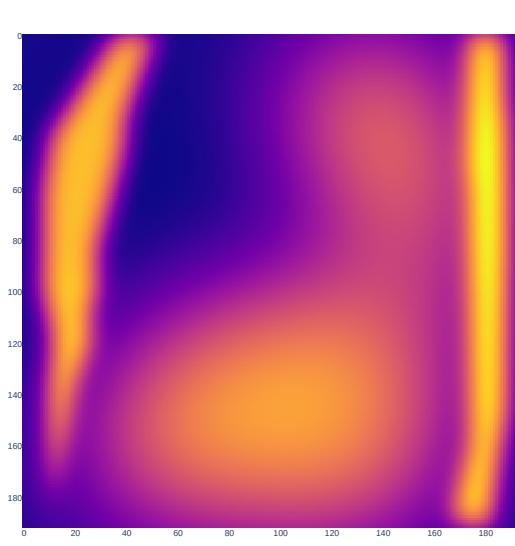


Figure 2.6. 2D *weights* matrix visualization.

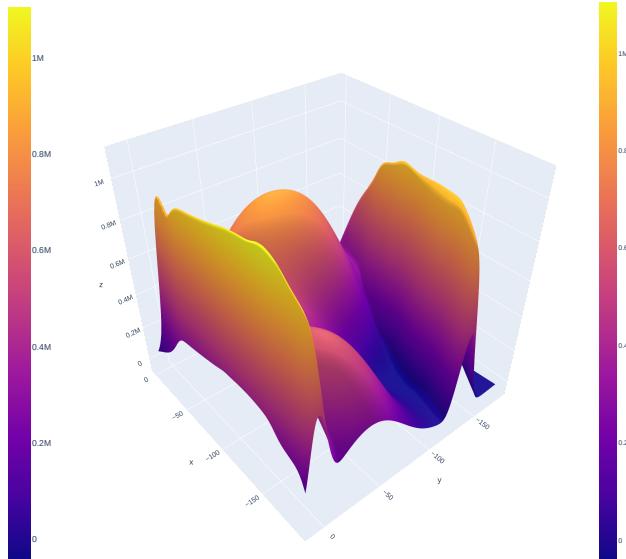


Figure 2.7. 3D *weights* matrix visualization.

Inverting the normal distributions in regions annotated as *high* is key to our implementation due to a *bias* element-wise addition applied after the normal distributions step. The latter step is used to normalize the magnitudes in our *weight* matrix between some constraints. As observed in the previous figures, the generated *weights* matrix contains elements ranging from values below 0 to values over 1×10^6 , we then apply a customized normalization function with a τ factor and a *bias* constant to re-weight the values in this matrix.

$$\omega(x) = (x \oslash \max x_i) \circ \tau + \beta \quad (2.1)$$

Where ω is our re-weighted *weights* matrix, x is the *weights* matrix, τ is the factor by which the values are multiplied and β is the *bias* term that increments all elements in the re-weighted matrix. The bias term defines $\min x_i$ for every x_i in the regions labelled as *medium*, *low* and *close-to-zero* movement.

At this point, the *weights* matrix is computed and re-weighted, so now this matrix is multiplied element-wise with each of the Φ masks created in the same fashion as the deterministic prior information, generating a *weighted phi* mask κ .

$$\kappa = \omega \circ \Phi \quad (2.2)$$

After this operation, our Φ positions no longer holds deterministic values, but rather

dispose of different levels of importance due to the *weights*' values.

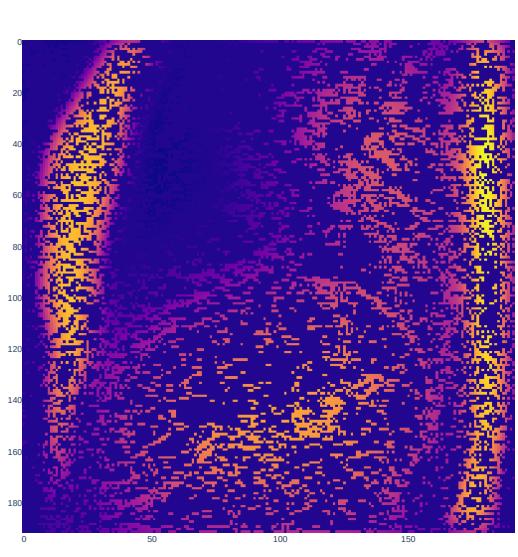


Figure 2.8. 2D *weights* matrix visualization.

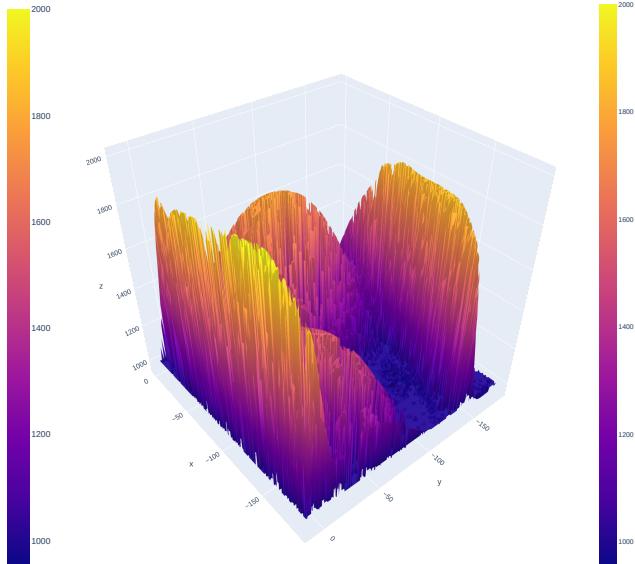


Figure 2.9. 3D *weights* matrix visualization.

2.3.3 L1-minimization Adaptation

After calculating κ , a minor modification is necessary for the *L1*-minimization algorithm. In essence, we perform an element-wise multiplication between the input and κ every time before an iteration so that the prior information is applied in the computation.

$$x_i = x_i \circ \kappa_k \quad (2.3)$$

Where $x, i, k\dots$

2.3.4 Parameter Fine-tuning

After some fine-tuning experiments in the covariance matrices, biases and τ , we have found that these are the optimal configuration for a cardiac MRI:

2.3.5 Future Contributions

Naturally, this implies that this task could gain a lot from a specialized image segmentation system or even a non-movement region detection system, which are more

elaborated in future steps section.

2.4 Artificial Neural Networks

2.4.1 Biological Inspirations

ANNs, as the name suggests, have been (loosely) inspired by biological neural networks (brains) from animals. The concept of using many layers of vector-valued representation is drawn from neuroscience. The choice of the functions $f^{(i)}(x)$ used to compute these representations is also loosely guided by neuroscientific observations about the functions that biological neurons compute [29]. Another trait they share is that just like the human brain can be trained to pass forward only meaningful signals to achieve larger goals of the brain, the neurons on a neural network can be trained to pass along only useful signal [5].

2.4.2 Neuron

The most basic unit in ANNs is the *artificial neuron*. Neurons act as feature detectors and this is one of the advantages of deep learning techniques in contrast to classical machine learning as the ANN is responsible for doing feature engineering and selection, and often outperform humans in this task.

These artificial neurons that are modelled mirroring the behaviour of the biological neuron as both of them are stimulated by inputs and carry some information they receive to other neurons. Artificial neurons take in inputs x_1, x_2, \dots, x_n , each and multiply them by their respective weights w_1, w_2, \dots, w_n . Then these weighted inputs are summed together producing the *logit* of the artificial neuron, $z = \sum_{i=0}^n w_i x_i + b$, with b being a constant number added called *bias*. After this, the logit is passed to a function f in order to generate the value $y = f(z)$.

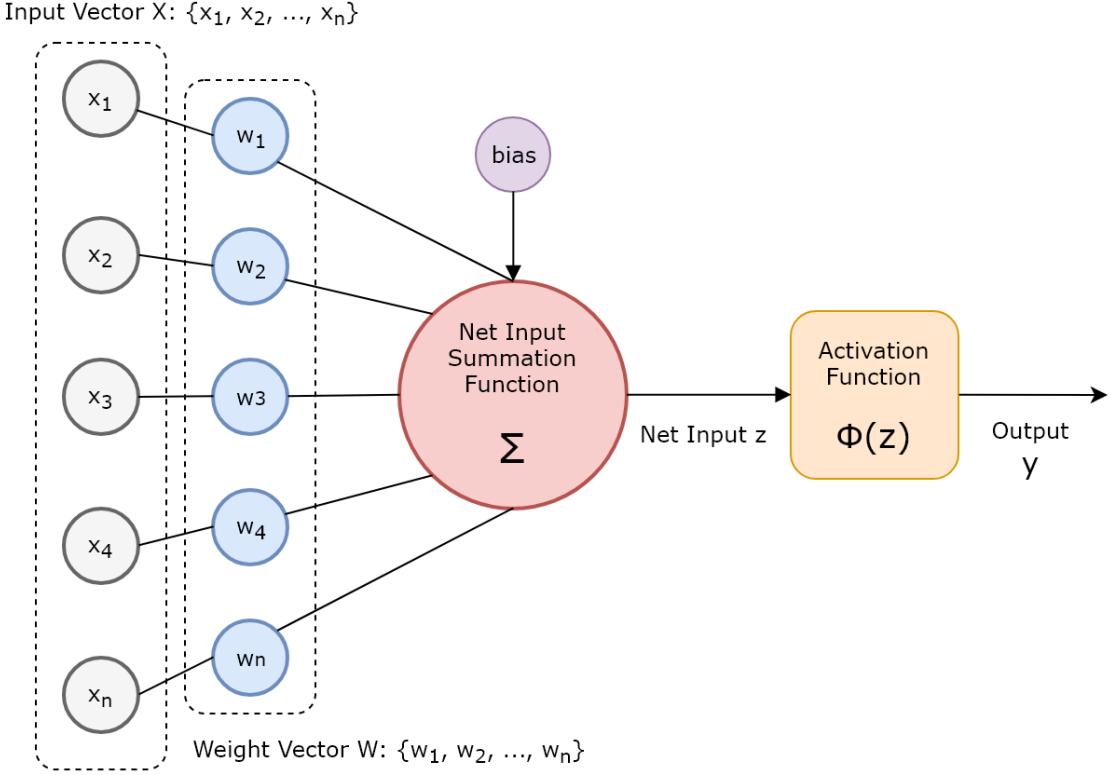


Figure 2.10. Schematic of an Artificial Neuron. Source: [4]

2.4.3 Multilayer Perceptron

Deep Feedforward Networks (DFN) or Multilayer Perceptron (MLP)s are a type of ANN very commonly used. It is the foundation of many famous architectures like CNNs. DFNs have an input layer followed by one or many hidden layers and a single output layer. Each layer is fully connected to the adjacent layer.

MLPs are computational models that flow information through the function f that evaluates x . The goal is to approximate some function f^* . For instance, a classifier $y = f^*(x)$ maps an input x to a category y . The feedforward defines a mapping $y = f(x; \theta)$ and learns the value of the parameters θ that result in the best function approximation [29].

The behaviour of an ANN is shaped by its architecture, which describes the number of units it should have and how these units connect to each other and how complex the model is. Often adding too much complexity to the network will lead to overfitting the training set, which occurs when the model shapes the training data too precisely and cannot generalise new data fed.

Most ANNs are organized into rows of neurons called layers. These layers are arranged

in a chain-like structure, with each layer being a function of the layer before it. These layers' goal is to extract *representations* out of the data fed and generalize what is meaningful towards minimizing the error rate. This architecture scheme is represented by the following equation, where i is the layer index:

$$h^{(i)} = g^{(i)}(W^{(i)T}x + b^{(i)})$$

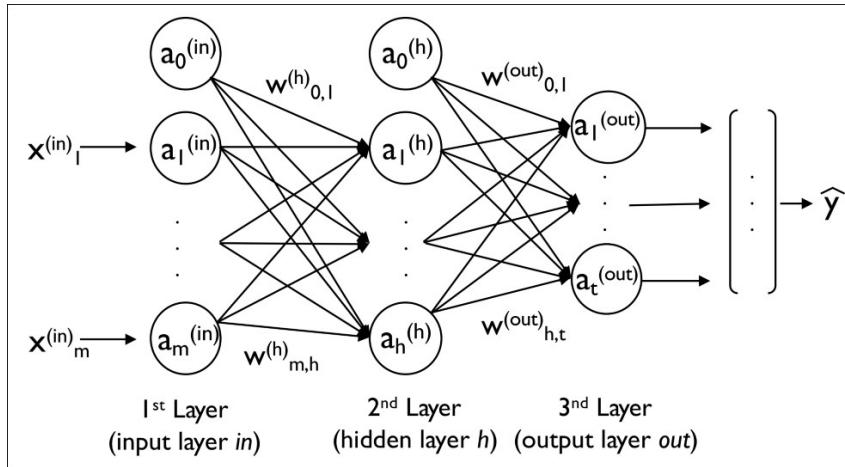


Figure 2.11. ANN Architecture Sample.

2.4.4 Activation Functions

Activation functions are a scalar-to-scalar function used to propagate the output of one layer's neurons forward to the next layer. There are several types of activation functions for different purposes and network architectures.

The Deep Convolutional GAN (DCGAN) architecture used in my MNIST experiment is built using ReLU activations in the generator network and a *tanh* in the output layer. The discriminator network uses LeakyReLU activations for all layers and a *sigmoid* function for the output layer. LeakyReLU activation functions [30, 31] have been proven to work well for higher resolution modelling [8] in contrast to the usage of maxout activation functions that were first proposed in the original GAN paper [32].

Name	Function	Derivative
Sigmoid	$\phi(x) = \frac{1}{1 + e^{-x}}$	$\phi'(x) = \phi(x)(1 - \phi(x))$
TanH	$\phi(x) = \frac{2}{1 + e^{-2x}} - 1$	$\phi'(x) = 1 - \phi(x)^2$
ReLU	$\phi(x) = \begin{cases} 0 & x \leq 0 \\ x & x > 0 \end{cases}$	$\phi'(x) = \begin{cases} 0 & x \leq 0 \\ 1 & x > 0 \end{cases}$
Leaky ReLU	$\phi(x) = \begin{cases} \alpha x & x \leq 0 \\ x & x > 0 \end{cases}$	$\phi'(x) = \begin{cases} \alpha & x \leq 0 \\ 1 & x > 0 \end{cases}$

Table 2.1. Activation functions and respective derivatives.

Some of the most used and also required activation functions in the use of GANs and other widely used neural networks are described below.

2.4.4.1 ReLU

The ReLU [33] transform activates a node only if the input is above a certain threshold having a linear relationship with the dependent variable and outputs zero for every input below zero.

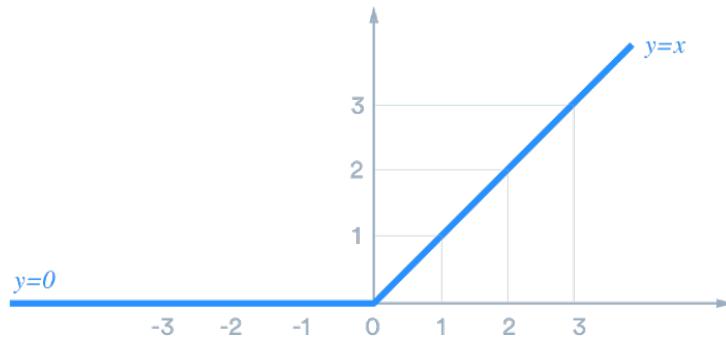


Figure 2.12. ReLU activation function

2.4.4.2 Leaky ReLU

ReLU activation functions have the “dying ReLU” problem, where a ReLU neuron is stuck in the negative side and always outputs 0 [34, 35]. This happens when the slope of ReLU in the negative range is also 0, once a neuron gets negative, it is unlikely for it to recover. These “dead” neurons are not playing any role in discriminating the input and

are essentially useless.

To mitigate this issue within ReLUs, LeakyReLUs are a strategy that opposed to having the function being zero when $x < 0$, it has instead a small negative slope (most times with $\alpha = 0.01$).

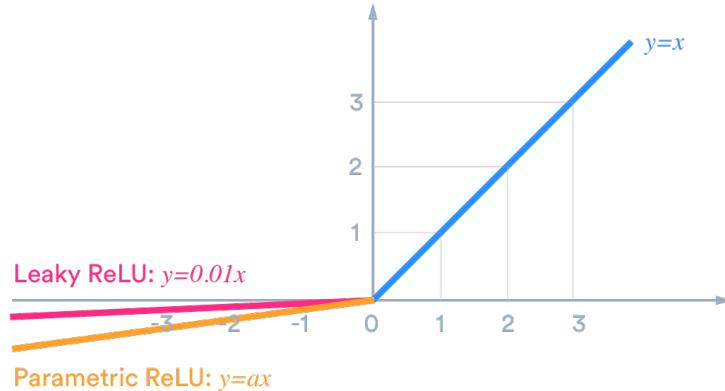


Figure 2.13. Leaky ReLU activation function

2.4.4.3 Tanh

Tanh is a hyperbolic trigonometric function that deals more easily with negative numbers [5]. Unlike the Sigmoid function, tanh ranges from -1 to 1.

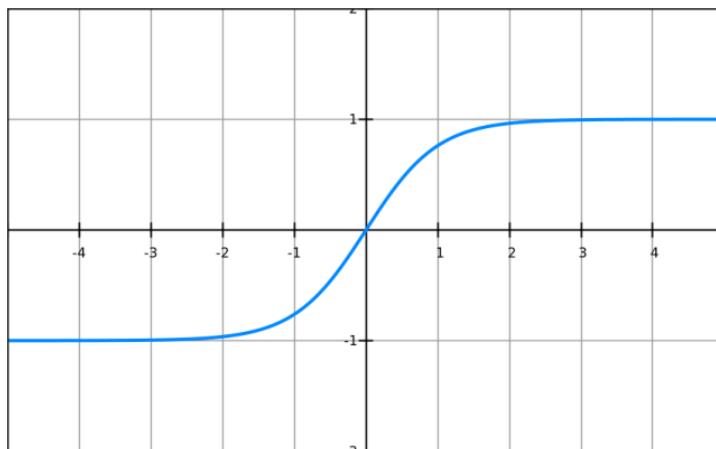


Figure 2.14. Tanh activation function. Source: [5]

2.4.4.4 Sigmoid

Sigmoids can reduce extreme values or outliers in data without removing them, framing the input from 0 to 1 and most outputs will be close to either 0 or 1.

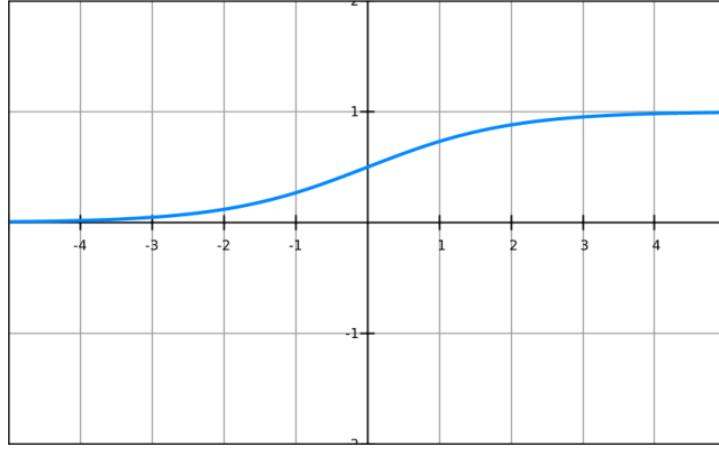


Figure 2.15. Sigmoid activation function. Source: [5]

2.4.5 Loss Functions

Loss functions are used to determine how a neural network is performing on the given data. A metric is calculated based on the error observed in the network's predictions and the model then tries to minimize this error in an optimization problem fashion.

Some of the most commonly used functions are described in the table below.

Mean squared error	MSE	$= \frac{1}{n} \sum_{t=1}^n e_t^2$
Root mean squared error	RMSE	$= \sqrt{\frac{1}{n} \sum_{t=1}^n e_t^2}$
Mean absolute error	MAE	$= \frac{1}{n} \sum_{t=1}^n e_t $
Mean absolute percentage error	MAPE	$= \frac{100\%}{n} \sum_{t=1}^n \left \frac{e_t}{y_t} \right $

Table 2.2. Loss functions and formulas.

In the case of generative networks, the original GAN paper presents a loss function called *minmax* [32], that is described as.

$$\min_G \max_D V(D, G) = E_{x \sim P_{data}(x)}[\log D(x)] + E_{z \sim p_z(z)}[\log(1 - D(G(z)))] \quad (2.4)$$

Where x is the input data representing an image, $D(x)$ is the discriminator network,

$G(x)$ is the generator function and z represents the latent vector that is mapped to data-space by G . Hence, the scalar probability that the output of the generator G is a real image is given by $D(G(z))$ [32].

2.4.6 Backpropagation

Backpropagation is a technique used to implement gradient descent in weight space for an MLP [36, 37]. In essence, backpropagation computes the error partial derivatives of an approximating function $F(w, x)$ computed by the ANN with respect to the weight and input vector for each training example [38].

The development of the backpropagation algorithm is a milestone in neural networks development and research as it made computationally efficient to train MLPs, thus confirming that ANNs research field was filled with potential in the mid-1980s.

In order to evaluate the derivatives of the function $F(w, x)$ with respect to all the elements in the weight vector w for an input vector $x = [x_1, x_2, \dots, x_{m_0}]^T$ for an MLP with layer $l = 2$, we have the following equation where φ is the activation function, w is the ordered weight vector and x is the input vector fed into the MLP [38]:

$$F(\mathbf{w}, \mathbf{x}) = \sum_{j=0}^{m_1} w_{oj} \varphi \left(\sum_{i=0}^{m_0} w_{ji} x_i \right) \quad (2.5)$$

2.4.7 Gradient Descent

Gradient descent is an optimization algorithm frequently used in ANNs to find the values of coefficients of a function that minimizes a cost function. Gradient descent can be very time consuming on large datasets due to the necessity of having a prediction for each instance in the training set. In scenarios where there is a large number of data instances, a variation of gradient descent called Stochastic Gradient Descent (SGD) can be used. SGD updates the coefficients for each training instance or batch instead of at the end after running through all the training set instances.

Most deep learning models are powered by the SGD and it can be visualized as the figure below demonstrates: the function starts in a random point in the loss function and after each iteration, the SGD calculates how it should adjust parameters in order to reach the minimal point in the loss function, hence moving towards the valley as illustrated.

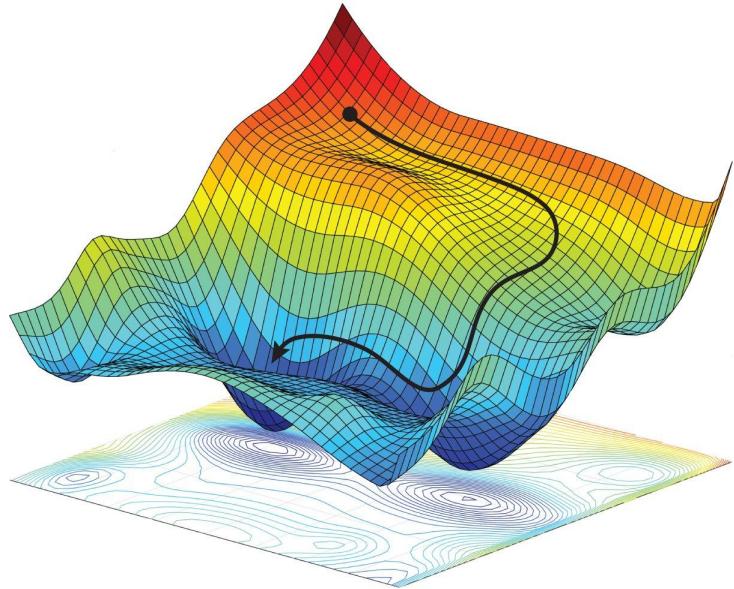


Figure 2.16. Gradient descent example visualization. Source: [6]

2.5 Generative Adversarial Networks

GANs are a machine learning strategy proposed in 2014 by Ian Goodfellow [32] that consists of two simultaneously trained models: the *Generator* $G(x)$ and the *Discriminator* $D(x)$. The generator has the role to generate fake data whilst the discriminator is trained to discern whether the given input is real or fake.

In essence, the generator takes a vector of random numbers (z) as input and outputs a fake example that strives to look as close as possible to the training data pattern. The discriminator takes an image (x) as input from two sources: real examples from the training set and fake examples generated by the generator network, then the discriminator outputs a scalar probability that the image is real [7].

GANs play a minimax two-player game in which D tries to maximize the probability to correctly classify real and fake samples ($\log D(x)$), whilst G tries to minimize the chance that D will correctly predict its generated outputs are fake ($\log(1 - D(G(x)))$).

Ideally, this minimax game would resolve to a solution with $pg = p_{data}$, where the discriminator is incapable of distinguishing real from fake inputs. However, GANs are still a novel neural network approach with its convergence theory is still being highly researched and hardly reaching this point in reality [32].

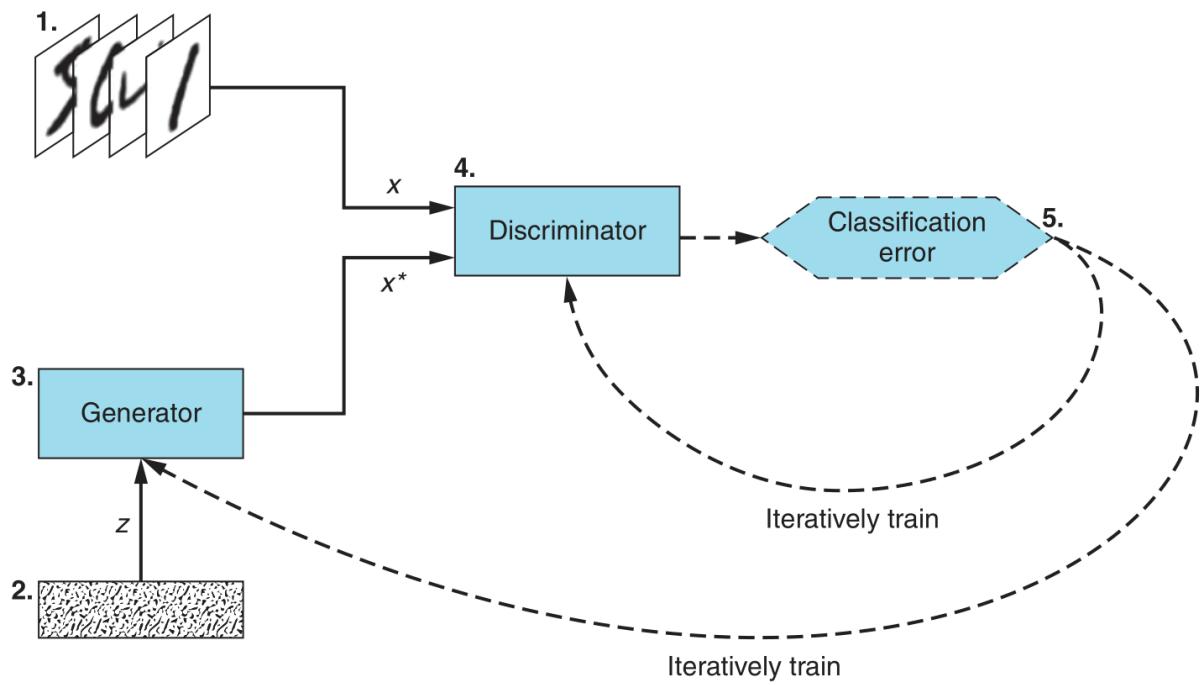


Figure 2.17. GAN training diagram. Source: [7]

3 PRELIMINARY RECONSTRUCTING AND SIGNAL GENERATING EXPERIMENTS

The following tests have had the purpose to elucidate the application of techniques for signal reconstruction with CS using pre-filtering pre-processing step and later with prior-information for better results. The signal generation experiment with GANs is a step before implemeting these neural networks to MRI measurements for prior-information leverage.

3.1 1-D Direct vs Indirect L1-Minimization

L1-minimization admits both direct and indirect approaches, in which there is the accuracy x resources trade-off. The direct method often produces a higher quality reconstruction but is very memory consuming, whilst the indirect method loses a little bit of quality, but requires much less memory to compute the equations system.

To visualize this trade-off, I have created 200 random 1-D arrays ranging values from the standard normal distribution and have taken 10% of data points randomly to reconstruct the whole signal using different L sizes.

L is the number of linear measurements extracted from the 1-D signal. The figure below displays the first 10 signals created in the standard normal distribution range.

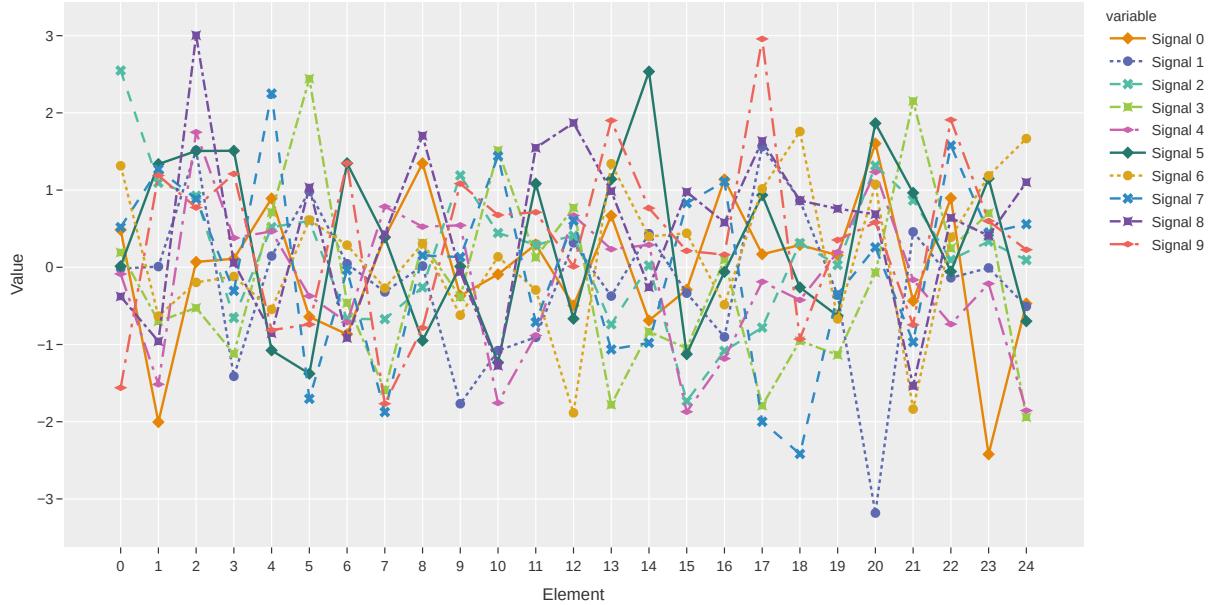


Figure 3.1. First 10 random 1-D signals

3.2 2D Compressed Sensing Reconstruction with Pre Filtered Signal

In order to experiment the sparsifying method of applying pre-filtering processing to the input measurements in the k-space for 2D images, I have conducted experiments some different images. A simple 2D image of the well-known Shepp-Logan phantom [39] for baseline is used for the first experiment and two MR (sagittal head slice and a singlecoil knee obtained from the NYU fastMRI Initiative database (fastmri.med.nyu.edu) [2, 1]) images that are considerably more complex than the phantom for the reconstruction task.

Within these images, *SNR* improvement gained by applying sparsifying pre-filtering [12, 13] can be best visualized and confirmed for these different scenarios.

3.2.1 Subsampling

I then created a phantom image with dimension of 256×256 , hence 65536 data points, using the phantominator python module. Then, I simulated an undersampled phantom image by applying the spiral undersampling pattern achieving approximately 30.95% of data points from the fourier space which resulted in a matrix with 20285 non zero elements.

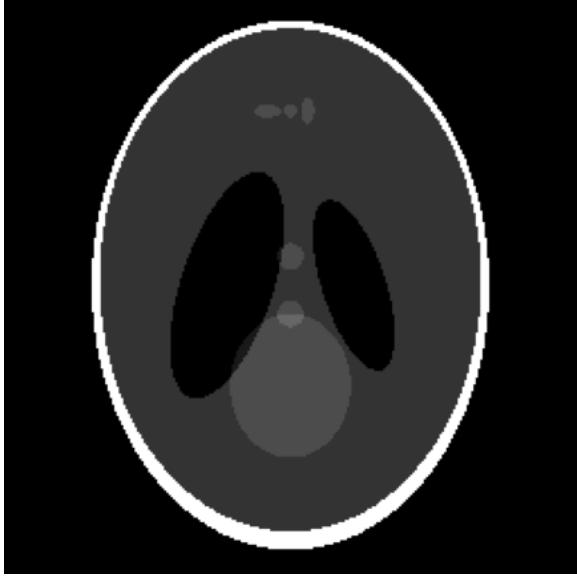


Figure 3.2. Shepp-Logan phantom reference image.

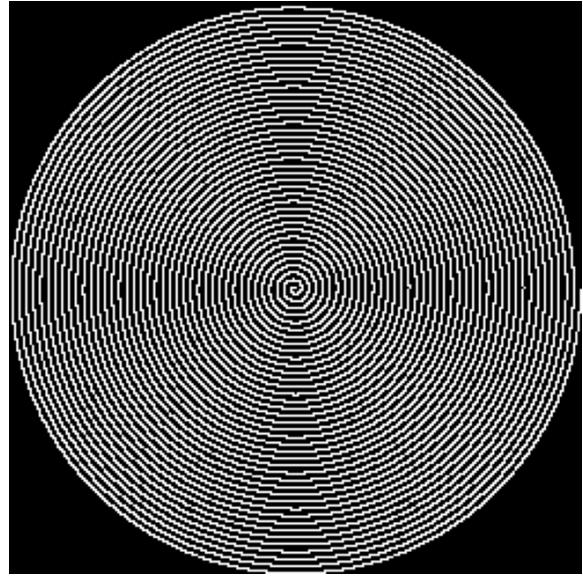


Figure 3.3. Spiral undersampling method with 30.95% data points.

3.2.2 Pre-filtering sparsifying transform

For the pre-filtering step, $f = 3$ is used where f is the number of filters applied to increase sparsity in the signal to be reconstructed. The filters are all 2×2 matrices and increase the sparsity in the signal from different perspectives, using more filters leverages the ability to sparsify the signal. The 3 filtered images are then composed into one single image containing the highest gain each filter could provide given a single pixel in the image [13]. The different filters used can be better seeing in the figure below.

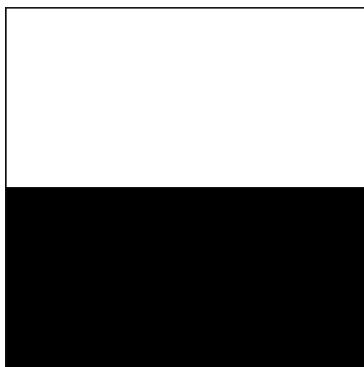


Figure 3.4.
2-D High pass
horizontal filter.

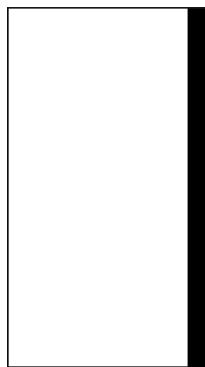


Figure 3.5.
2-D High pass
vertical filter.

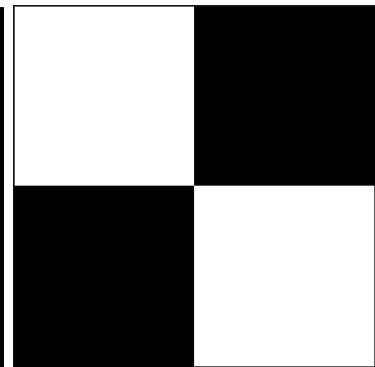


Figure 3.6.
2-D High pass
diagonal filter.

The pre-filtering method is evaluated against the zero-fill reconstruction method (used

as a dummy baseline) and an L1-minimization method without pre-filtering with the very same parameters used in the pre-filtering L1-minimization.

3.3 Preliminary Tests with Generative Adversarial Networks

In order to test the usage of GANs for data generation and in the future use it along with *prior information* for CS systems, I have developed a GAN capable of generating handwritten digits from 0 to 9 using the notable MNIST dataset. The MNIST dataset contains 60,000 examples for training and 10,000 examples for testing. The digits have been size-normalized and centred in a fixed-size image (28×28 pixels) with values from 0 to 9. For simplicity, each image has been flattened and converted into a 1-dimensional numpy array of 784 features (28×28).

The idea is to test if the neural network can output liable digits that look both readable (to the extent in which the MNIST dataset is) and also like it has been made by a human, just like the dataset itself.

Each MNIST image contains a 28×28 black and white image, like the following:



Figure 3.7. Sample of digits from MNIST

A DCGAN was used for the experiment. A DCGAN is an extension of the GAN, except that it explicitly uses convolutional and convolutional-transpose layers in the discriminator and generator networks, respectively [8].

3.3.1 Data Transformation

Each input image used by the *dataloader* went through a computer-vision pre-processing step that includes:

- Grayscale transform: convert the image to greyscale. When loaded, the MNIST digits are in RGB format with three channels. Greyscale reduces these three to one.
- ToTensor: convert the image to a PyTorch Tensor, with dimensions (channels, height, width). This also rescales the pixel values, from integers between 0 and 255 to floats between 0.0 and 1.0.
- Normalize: scale and translate the pixel values from the range 0.0, 1.0 to -1.0, 1.0. The first argument is μ and the second argument is σ , and the function applied to each pixel is:

$$\rho \leftarrow \frac{(\rho - \mu)}{\sigma} \quad (3.1)$$

3.3.2 Generator Network Architecture

The generator network architecture is implemented using PyTorch as:

- A linear *fully-connected* module (or layer) to map the latent space to a $7 \times 7 \times 256 = 12544$ -dimensional space that will later be undersampled several times until we reach $1 \times 28 \times 28$.
- An optional 1-dimensional batch normalization module
- A leaky ReLU module.
- A 2-dimensional convolutional layer with $padding = 2$, $stride = 1$ and 5×5 kernel (or filter).
- Two 2-dimensional transposed convolutional layers with $padding = 1$, $stride = 2$ and 4×4 kernel.
- Two optional 2-dimensional batch normalization modules after each 2-dimensional transposed convolutional layer.
- A *Tanh* activation function, rescaling the images to a $[-1, 1]$ range.

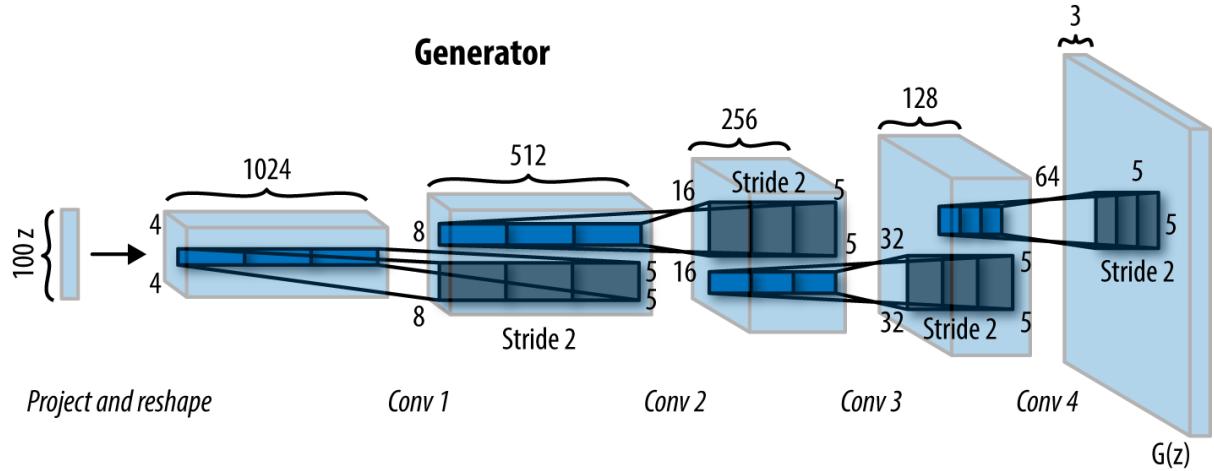


Figure 3.8. Generator network architecture of a 3-D image. Source: [8, 9]

The latent space (random signal) input goes through each layer being upscaled until it reaches the target image dimension 28×28 and then fed into the discriminator network.

3.3.3 Discriminator Network Architecture

The discriminator is a CNN-based image binary classifier network that takes an image as input and outputs a scalar probability that the given image is real or generated. The architecture is quite similar to the Generator network, except backwards. Here, the discriminator takes a $1 \times 28 \times 28$ input image, processes it through a series of convolutions, batch normalizations, and LeakyReLU layers, and outputs the final probability through a Sigmoid activation function.

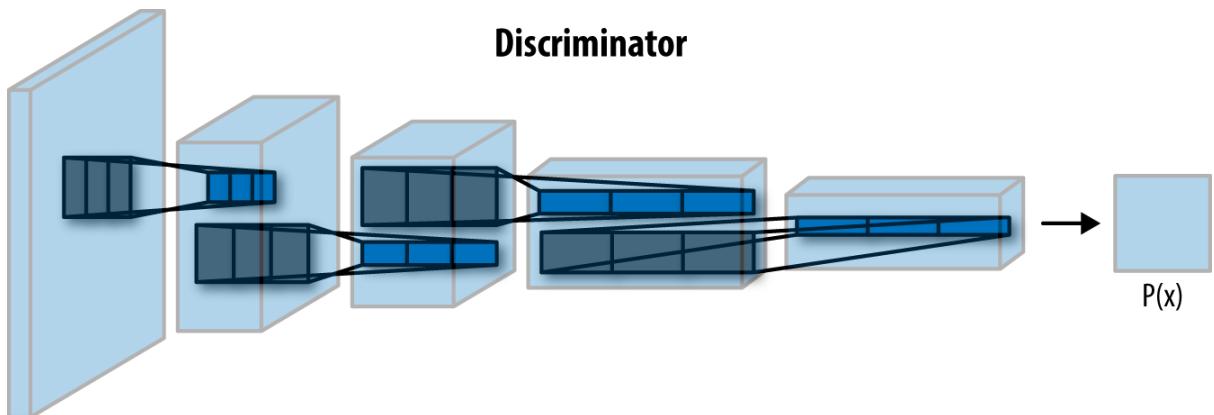


Figure 3.9. Discriminator network architecture of a 3-D image. Source: [8, 9]

4 PRELIMINARY RECONSTRUCTION AND GENERATION RESULTS

The results below serve to confirm the viability of the proposed methods as well as to create baselines with different images and signals that will later be compared to the fully-working GAN + CS proposed method. The following preliminary results are produced with a CS algorithm version that is still under refinement and is not yet producing the highest reconstruction standard intended for the end of the present thesis.

4.1 1-D Compressed Sensing Reconstruction

The reconstruction quality over computational resources demanded trade-off starts making a difference with L size around 140, which is over 50% of the fully sampled signal, a proportion a lot higher than the ones normally used in MRI reconstruction. This demonstrates that there is no big prejudice in using indirect reconstruction method for very undersampled signals (close to 10% of the data) such as the ones used in MRI.

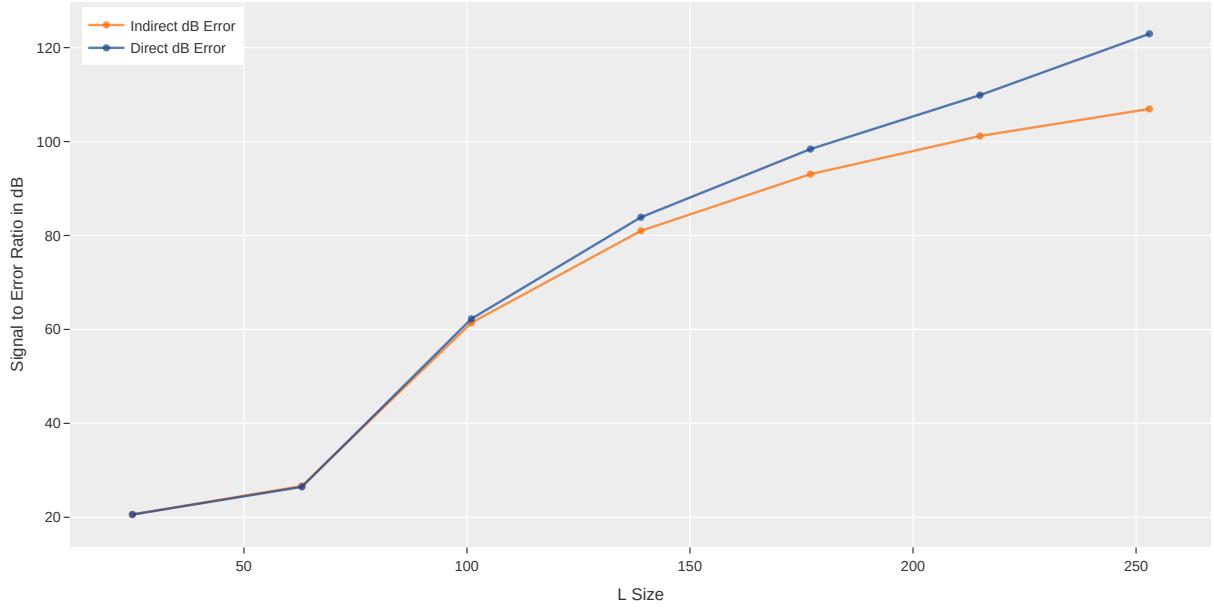


Figure 4.1. SNR x L size for direct and indirect L1-minimization for 200 random 1D signals

4.2 Compressed Sensing Reconstruction with Pre Filtered Signal

The results in the experiment show a huge gain of resolution in the pre-filtering method reconstructed image compared to the L1-minimization alone.

The zero-filled reconstruction (dummy baseline) was unable to reconstruct a high fidelity image and performed very poorly in the PSNR and SNR metrics. The L1-minimization compressed sensing approach reconstructed the image with some noticeable noise artefacts, yet much better than the zero-filling approach. Finally, the L1-minimization along the usage of sparsifying pre-filtering delivered a great looking image without eye-catching artefacts and also increased the metrics hugely.

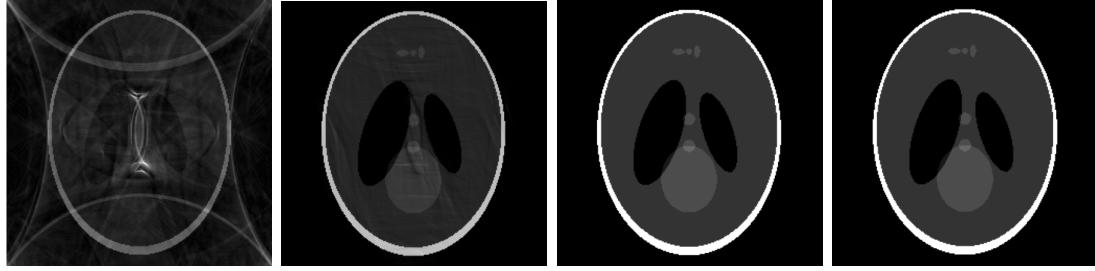


Figure 4.2.
zero-
filled
reconstruction

Figure 4.3.
L1-
minimization
reconstruction

Figure 4.4.
L1-
minimization
with
pre-
filtering
reconstruction

Figure 4.5.
Shepp-
Logan
phantom
reference
image

The usage of L1-minimization certainly improves the MRI reconstruction, but the metrics reinforce how adding the pre-filtering step to preprocess the image achieves incredibly higher scores in Peak Signal-to-Noise Ratio (PSNR) and SNR.

	PSNR	SSIM	SNR	MSE
Zero-fill	22.41	0.36	4.22	0.02
L1-minimization	38.76	0.96	20.57	5.3e-5
Pre-filtering L1-minimization	91.89	0.99	73.70	2.5e-9

Table 4.1. Phantom reconstruction metrics.

4.2.1 Brain Sagittal Reconstruction

Another experiment done was using a reference sagittal head image of shape (256×256) . The same spiral undersampling pattern with 30.95% data points used in the phantom experiment was used here for artificial undersampling.

This image poses a harder reconstruction challenge as it is filled with more details and more complex structures than the phantom. That said, it is clear that the undersampling pattern and amount of data points has not been sufficient to reconstruct a high fidelity image in any scenario, but the L1-minimization with pre-filtering reconstruction looks like the winner again, reinforcing the idea that pre-filtering is a good pre-processing strategy.

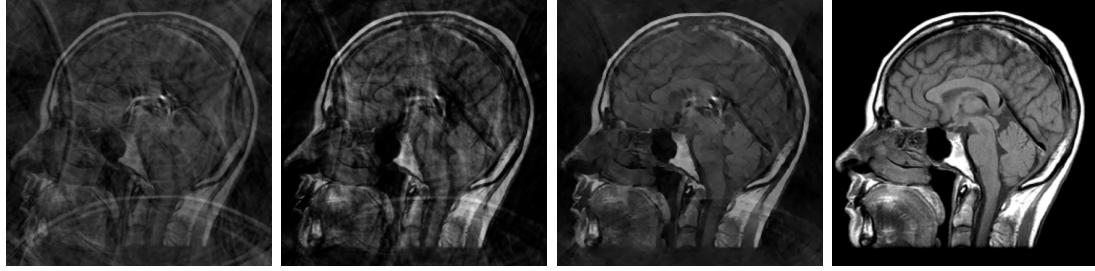


Figure 4.6.
Zero-filled
reconstruction

Figure 4.7.
L1-minimization
reconstruction

Figure 4.8.
L1-minimization
with
pre-filtering
reconstruction

Figure 4.9.
Reference
image

The metrics evaluated were not affected as much as they were for the phantom experiment, but they were mostly improved by pre-filtering usage.

	PSNR	SSIM	SNR	MSE
Zero-fill	17.68	0.32	5.52	2392.57
L1-minimization	21.51	0.43	8.76	1134.88
Pre-filtering L1-minimization	21.03	0.45	8.97	1081.53

Table 4.2. Sagittal reconstruction metrics.

4.2.2 Prior Information Cardiac Sagittal Reconstruction

In order to test if the extracted prior information from one frame is relevant and how it impacts the reconstruction of the next frame, we have tested reconstructing frames from a dynamic cardiac MRI exam. The extracted frames are $(n \times n)$ in shape

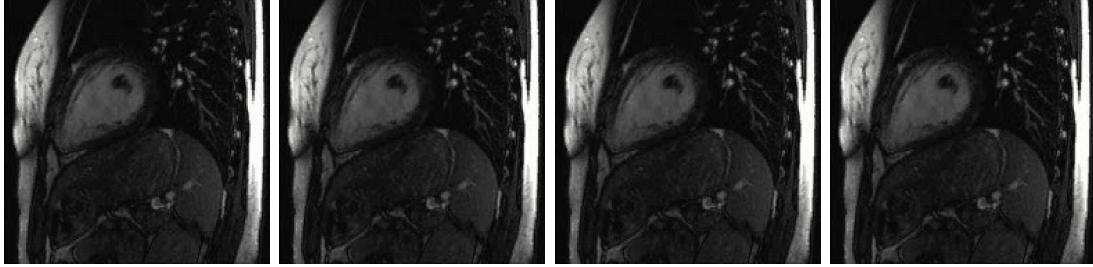


Figure 4.10.
Dynamic
cardiac
zero-
filled
reconstruction

Figure 4.11.
Dynamic
cardiac
L1-
minimization
reconstruction

Figure 4.12.
Dynamic
cardiac
L1-
minimization
with
pre-
filtering
reconstruction

Figure 4.13.
Dynamic
cardiac
reference
image

	PSNR	SSIM	SNR	MSE
Zero-fill	17.68	0.32	5.52	2392.57
L1-minimization	21.51	0.43	8.76	1134.88
Pre-filtering L1-minimization	21.03	0.45	8.97	1081.53

Table 4.3. Sagittal reconstruction metrics.

4.3 Preliminary Tests with Generative Adversarial Networks

Both discriminator losses (fake and real) start very high and quickly decreases as the generator loss curve goes up in an invertedly correlated manner. This happens especially because the generator starts by tricking the discriminator network very easily as it is naïve to determine if an image is real or generated. Quickly the discriminator starts to detect how the data is disposed and manages to interpret the generated images are different from the training examples it is seeing.

This phenomenon exposes how bad the generator is in the first epochs and how easily the discriminator can distinguish between created and real. Then as the epochs go by and both networks get more sophisticated, the generator starts to get better at creating the desired signal style and makes the discriminator's loss get higher again as it is observed in the loss curves below.

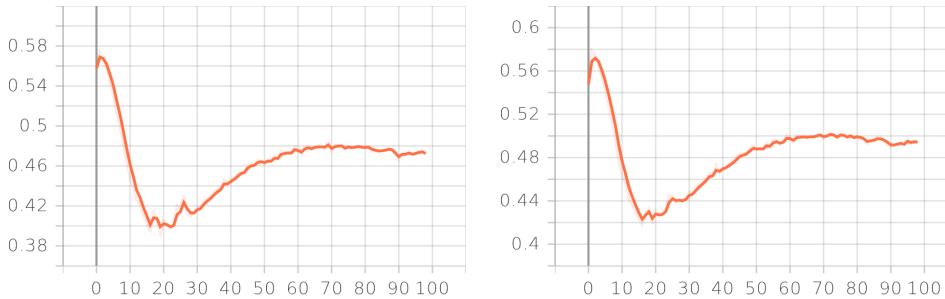


Figure 4.14.
Discriminator
loss over epochs

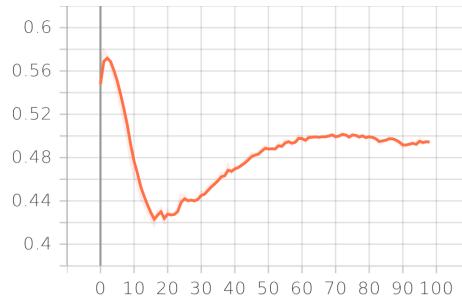


Figure 4.15.
Discriminator
real loss
over epochs

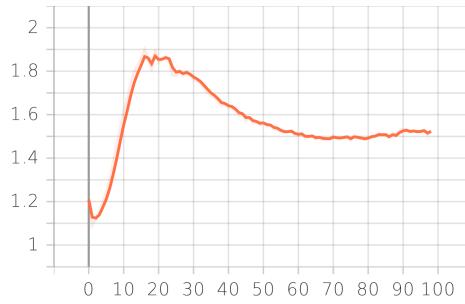


Figure 4.16. Generator
loss over epochs

After 100 epochs, the DCGAN for MNIST number generation had an exceptionally good performance when the generated images are displayed. It is hard to tell if these are generated images or if they are part of the training set. The generated images sometimes have a bit more blur to them, but certainly with more epochs and more training samples this could be minimized. The following results were not cherry-picked in any manner.

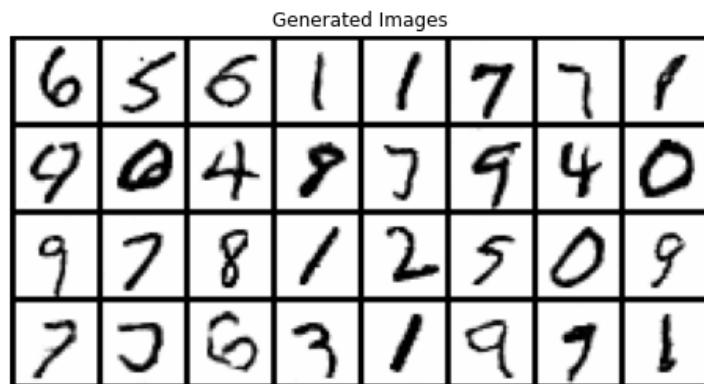


Figure 4.17. GAN generated MNIST digits

5 CONCLUSION

The next steps in this thesis involve using prior-information for best CS results and training a GAN for prior information leverage aiming for better resolution reconstructions and dynamic MRI.

The cronogram has been split in four major activities: prior information with CS, GAN architecture design and implementation, experiment GAN for prior information generation and evaluate GAN results for image reconstructions. These tasks have been organized within the expected for each task to be well done and documented. There is not yet an official date for the conclusion of this thesis, but the tasks have more than enough time to be well developed and evaluated.

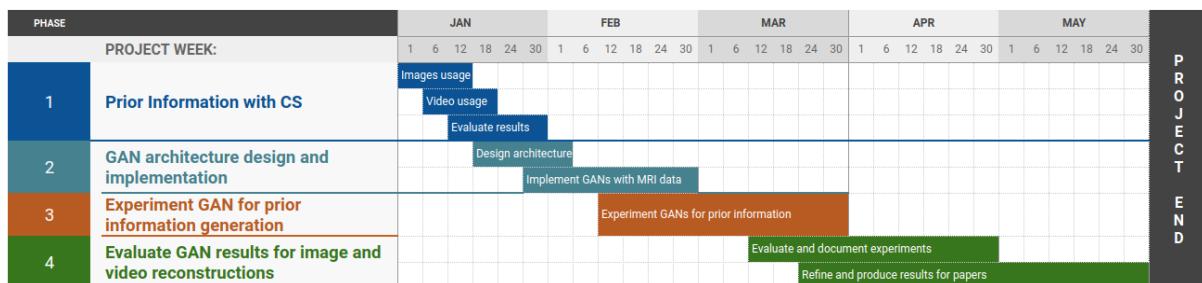


Figure 5.1. Activities cronogram

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