ALZHEIMER'S DISEASE DIAGNOSTICS BY ADAPTATION OF CONVOLUTIONAL NETWORK

By Tara Rasti, Fall 2020

Introduction

Alzheimer's disease (AD) is a progressive brain disorder and the most common case of dementia in the late life. AD leads to the death of nerve cells and tissue loss throughout the brain, thus reducing the brain volume in size dramatically through time and affecting most of its functions [1]. The estimated number of affected people will double for the next two decades, so that one out of 85 persons will have the AD by 2050 [2]. An estimated 5.5 million people aged 65 and older are living with AD, and AD is the sixth-leading cause of death in the United States. The global cost of managing AD, including medical, social welfare, and salary loss to the patients' families, was \$277 billion in 2018 in the United States, heavily impacting the overall economy and stressing the U.S. health care system [3]. Because the cost of caring the AD patients is expected to rise dramatically, the necessity of having a computer-aided system for early and accurate AD diagnosis becomes critical [4]. Several popular non-invasive neuroimaging tools, such as structural MRI (sMRI), functional MRI (fMRI), and positron emission tomography (PET), have been investigated for developing such a system [5, 6]. The sMRI has been recognized as a promising indicator of the AD progression [4, 7]. Early diagnosis, playing an important role in preventing progress and treating the Alzheimer's disease (AD), is based on classification from brain images [8].

Methodology

- Dataset

For this purpose we use a dataset which is available through this link: https://www.kaggle.com/legendahmed/alzheimermridataset. The dataset consists of 6400 MRI axial slices. The dataset is categorized into four groups: non demented, very mild demented, mild demented and moderate demented categories. The dimensions of all photos are 176 * 208 and all of them are jpgs. That resulted in a total of 6400 training images, 3200 of which were AD and the other 3200 were Healthy Control (HC).

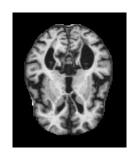
Non Demented



Very mild Demented



Mild Demented



Moderate Demented



To train deep learning algorithms, the following ingredients are required:

- Training data (inputs and targets)
- The model (also called the network)

- The loss function (also called the objective function or criterion)
- The optimizer

Training

The training process for deep learning algorithms is an iterative process. In each iteration, we select a batch of training data. Then, we feed the data to the model to get the model output. After that, we calculate the loss value. Next, we compute the gradients of the loss function with respect to the model parameters (also known as the weights). Finally, the optimizer updates the parameters based on the gradients. This loop continues. We also use a validation dataset to track the model's performance during training. We stop the training process when the performance plateaus [9].

Model

- Convolutional Neural Networks

Convolutional neural network, which has shown remarkable performance in the field of image recognition, has also been used for the diagnostic classification of AD with multimodal neuroimaging data. This project proposed to predict the AD with a deep convolutional neural network, which can learn generic features capturing AD biomarkers and adapt to different domain datasets.

Convolutional networks (also known as convolutional neural networks or CNNs) are a specialized kind of neural network for processing data that has a known, grid-like topology. Examples include image data, which can be thought of as a 2D grid of pixels. Convolutional networks have been tremendously successful in practical applications. Convolutional networks are simply neural networks that use convolution in place of general matrix multiplication in at least one of their layers. Convolution leverages three important ideas that can help improve a machine learning system: sparse interactions, parameter sharing and equivariant representations. Moreover, convolution provides a means for working with inputs of variable size.

Convolutional networks, typically have sparse interactions (also referred to as sparse connectivity or sparse weights). This is accomplished by making the kernel smaller than the input. For example, when processing an image, the input image might have thousands or millions of pixels, but we can detect small, meaningful features such as edges with kernels that occupy only tens or hundreds of pixels. Parameter sharing refers to using the same parameter for more than one function in a model. In a convolutional neural net, each member of the kernel is used at every position of the input. The parameter sharing used by the convolution operation means that rather than learning a separate set of parameters for every location, we learn only one set. In the case of convolution, the particular form of parameter sharing causes the layer to have a property called equivariance to translation. To say a function is equivariant means that if the input changes, the output changes in the same way.

A typical layer of a convolutional network consists of three stages. In the first stage, the layer performs several convolutions in parallel to produce a set of linear activations. In the second stage, each linear activation is run through a nonlinear activation function, such as the rectified linear activation function. This stage is sometimes called the detector stage. In the third stage, we use a pooling function to modify the output of the layer further. A pooling function replaces the output of the net at a certain location with a summary statistic of the nearby outputs. For example, the max pooling operation reports the maximum output within a rectangular neighborhood [10].

- Transfer Learning

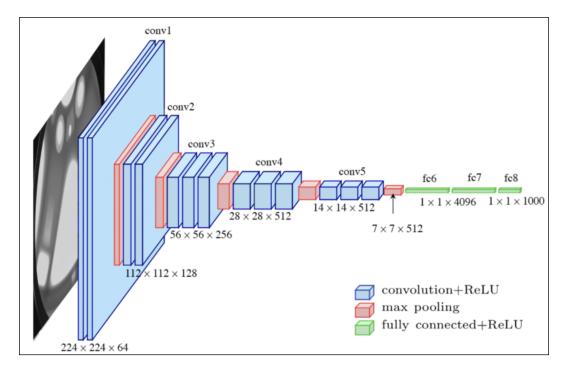
Transfer learning is a research problem in machine learning that focuses on storing knowledge gained while solving one problem and applying it to a different but related problem. For example, knowledge gained while learning to recognize cars could apply when trying to recognize trucks. In transfer learning approach knowledge is transferred from one model to another. Transfer learning is an optimization that allows rapid progress or improved performance when modeling the second task. Transfer learning is the improvement of learning in a new task through the transfer of knowledge from a related task that has already been learned. To deal with the lack of data, we make use of a technique called transfer learning. Briefly, transfer learning is a method that allows us to use knowledge gained from other tasks in order tackle new but similar problems quickly and effectively. This reduces the need for data related to the specific task we are dealing with [10].

VGG Neural Network

VGG is an innovative object-recognition model that supports up to 19 layers. Built as a deep CNN, VGG also outperforms baselines on many tasks and datasets outside of ImageNet. VGG is now still one of the most used image-recognition architectures [11].

• VGG16

VGG16 is a convolutional neural network model proposed by K. Simonyan and A. Zisserman from the University of Oxford in the paper "Very Deep Convolutional Networks for Large-Scale Image Recognition". The model achieves 92.7% top-5 test accuracy in ImageNet, which is a dataset of over 14 million images belonging to 1000 classes.



Architecture of VGG 16

• VGG19

VGG-19 is a convolutional neural network that is 19 layers deep. You can load a pre-trained version of the network trained on more than a million images from the ImageNet database. The pre-trained network can classify images into 1000 object categories, such as keyboard, mouse, pencil, and many animals.

The loss function

As with other machine learning models, to apply gradient-based learning we must choose a cost function, and we must choose how to represent the output of the model. An important aspect of the design of a deep neural network is the choice of the cost function. In most cases, our parametric model defines a distribution $p(y \mid x; \theta)$ and we simply use the principle of maximum likelihood. This means we use the cross-entropy between the training data and the model's predictions as the cost function. Most modern neural networks are trained using maximum likelihood. This means that the cost function is simply the negative log-likelihood, equivalently described as the cross-entropy between the training data and the model distribution [10].

The optimizer

Optimization used as a training algorithm for a machine learning task differs from pure optimization. Sometimes, the loss function we actually care about (say classification error) is not one that can be optimized efficiently. For example, exactly minimizing expected 0-1 loss is typically intractable (exponential in the input dimension), even for a linear classifier. In such situations, one typically optimizes a surrogate loss function instead, which acts as a proxy but has advantages. For example, the negative log-likelihood of the correct class is typically used as a surrogate for the 0-1 loss. The negative log-likelihood allows the model to estimate the conditional probability of the classes, given the input, and if the model can do that well, then it can pick the classes that yield the least classification error in expectation. A very important difference between optimization in general and optimization as we use it for training algorithms is that training algorithms do not usually halt at a local minimum. Instead, a machine learning algorithm usually minimizes a surrogate loss function but halts when a convergence criterion is satisfied [10].

Dealing with Overfitting

What is Overfitting?

Overfitting occurs when the gap between the training error and test error is too large. A central problem in machine learning is how to make an algorithm that will perform well not just on the training data, but also on new inputs. Many strategies used in machine learning are explicitly designed to reduce the test error, possibly at the expense of increased training error. These strategies are known collectively as regularization. There are a great many forms of regularization available to the deep learning practitioner [10].

• Data Augmentation

The best way to make a deep learning model generalize better is to train it on more data. Of course, in practice, the amount of data we have is limited. One way to get around this problem is to create fake data and add it to the training set. The performance of deep learning neural networks often improves with the amount of data available. Data augmentation is a technique to artificially create new training data from existing training data. Dataset augmentation applies transformations to your training examples: they can be as simple as flipping an image, or as complicated as applying neural

style transfer. The idea is that by changing the makeup of your data, you can improve your performance and increase your training set size [10].

• Dropout

Dropout provides a computationally inexpensive but powerful method of regularizing a broad family of models. The term dropout refers to dropping out units (hidden and visible) in a neural network. Dropout is a simple way to prevent neural networks from overfitting. Dropout may be implemented on any or all hidden layers in the network as well as the visible or input layer. It is not used on the output layer [10].

• Weight Decay

Weight decay is a regularization technique by adding a small penalty, usually the L2 norm of the weights (all the weights of the model), to the loss function. Because we are subtracting a constant times the weight from the original weight. This is why it is called weight decay. We use weight decay to prevent overfitting, to keep the weights small and avoid exploding gradient [10].

Why PyTorch

There has been significant progress in computer vision because of deep learning in recent years. This helped to improve the performance of various tasks such as image recognition, object detection, image segmentation, and image generation. Deep learning frameworks and libraries have played a major role in this process. PyTorch, as a deep learning library, has emerged since 2016 and gained great attention among deep learning practitioners due to its flexibility and ease of use. PyTorch is a deep learning framework developed by Facebook's artificial intelligence research group [9].

Accuracy Results

In the first five trials we utilized the dataset with a ratio of 20% for the validation and 80% for the train set. We implemented a CNN model with 5 convolutional layers, 3 max pooling layers and three linear layers. For activation functions we used ReLU, Leaky ReLU and Softmax functions. For regularization we used dropout. After testing several figures we understood that a 0.3 dropout performs better for the model. For the loss function we used cross entropy and for the optimizer, Adam with 1e-4 learning rate and 1e-5 weight decay. After 100 epochs the result was 65% on validation set and 87% for train, set which shows a great gap between training and validation set. In the second trial we did the same except that we trained the model for 200 epochs. The results show that the model includes overfitting. In the 3rd and 4th trials we used the VGG16 model and trained it for 20 epochs with two optimizers; first SGD with an accuracy of 75% on validation set, second Adam with an accuracy of 76% on validation set. After comparing these two models we found that Adam optimizer performs better on the data, hence we used this optimizer for the other proceeding models. In the 5th trial we applied VGG19 with 50 epochs and a performance of 79% on the validation set. However, the model still shows overfitting.

After those five first trials we decided to change the ratio of the validation and train sets. Therefore, instead of 20% for the validation set we changed it to 30%. Incredibly, the results changed and the overfitting was gone. In the next four trials we used CNN model with 100 and 200 epochs with accuracies of 85% and 92% on validation set respectively. Then I used the VGG16 again and this time we got a result of 93% on validation and 97% on training set. In the final trial we exerted the VGG19 model and got the best results within all previous models. For the training set 98% accuracy and for the validation set 96.3%. As you can see below, a table is illustrated which shows accuracy and other information about the models:

Trial	Model	Validation Ratio	Optimizer	Epochs	Accuracy(val)	Accuracy(train)
1st	CNN	0.2	Adam	100	65%	87%
2nd	CNN	0.2	Adam	200	68%	96%
3rd	VGG16	0.2	SGD	20	75%	95%
4th	VGG16	0.2	Adam	20	76%	96%
5th	VGG19	0.2	Adam	50	79%	99%
6th	CNN	0.3	Adam	100	85%	90%
7th	CNN	0.3	Adam	200	92%	96%
8th	VGG16	0.3	Adam	30	93%	97%
9th	VGG19	0.3	Adam	30	96%	98%

Table Showing Accuracy for all Implemented Models

The codes for this project can be found here:

https://github.com/TaraRasti/Deep_Learning_F20_Project

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