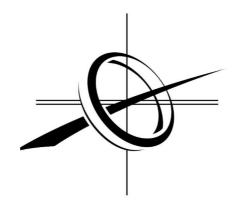
UNIVERSIDAD POLITÉCNICA DE MADRID

ESCUELA TÉCNICA SUPERIOR DE INGENIERÍA DE SISTEMAS INFORMÁTICOS

GRADO EN INGENIERÍA DE SOFTWARE

TRABAJO DE FIN DE GRADO

A NEURAL NETWORKS BENCHMARK FOR IMAGE CLASSIFICATION



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A través de este documento, el lector puede hacerse hacerse una idea de como ha sido la historia de los métodos usados para clasificación de imagen, tanto con métodos clásicos como con redes neuronales artificiales; todo ello en el contexto de visión para robots. Primero revisaremos los métodos clásicos, pasando por sus restricciones y limitaciones, escogeremos uno y sacaremos diferentes medidas sobre cómo se comporta. Después, exploraremos, brevemente, los tipos de redes neuronales que se utilizan para esta tarea, pasando por el estado del arte y su aportación; también escogeremos una red y mediremos su eficacia. Con todo ello, explicaremos el método empleado para medir y los resultados experimentales obtenidos. Por último discutiremos estos resultados y expondremos nuestras conclusiones, así como posibles lineas futuras de investigación.

Through this document, the reader can get an idea of the history of modern and classic methods for the task of image classification, we present a simple image classification task, in the context of robotic vision, and how different neural networks reach to stable solutions. First, we'll review different classic methods, evaluating their constraints and limitations, only to pick one up and benchmark it. Then, briefly, we will explore more modern methods, choose one, and benchmark it. Then, both benchmarks will be compared, and experimental results will be analyzed and explained. We'll conclude with a discussion of the results, pointing out future lines of research.

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1 Introduction

In this work we present a simple image classification task, in the context of robotic vision, and how different neural networks reach to stable solutions. First, we'll review different classic methods, evaluating their constraints and limitations, only to pick one up and benchmark it. Then, briefly, we will explore more modern methods, choose one, and benchmark it. Then, both benchmarks will be compared, and experimental results will be analyzed and explained. We'll conclude with a discussion of the results, pointing out future lines of research.

1.1 Context and Motivation

Computer Vision (CV) is a continuously developing field, with applications in a broad range of scientific and everyday life problems, from face recognition in social networks to homeland security. Recent years have seen vertiginous progress; nowadays, the most advanced techniques are able to classify multiple objects in real time streaming video, adjusting perspective, color and brightness enhacement, and virtual rotation. CV, then, deals with the process of perceiving image and video. Perception is not only a matter of reception of images, but comprises multiple and interrelated subroutines. All those subroutines that are involved in the act of perception must be addressed when programming CV, and, more often than not, will be inspired by biology [1].

The story of CV is short, as it started in the 60's, with the invention of the perceptron (Rosenblatt, 19622]). A perceptron is, roughly speaking, a simple linear classifier. Similarly to a transistor, a perceptron gives, for a particular input, a binary output. Thus, it can be used to generate simple silogisms, as a logic gate, or, more abstractly, to classify one dimensional problems, which is a kind of binary feature extraction. Feature extraction is a major issue in CV. The invention was a revolution until the 70's, in which computer science served to other, more "complicated" problems, not solved yet by neural networks. At that time, only the perceptron was available, with manifest pitfalls. Logic problems of the form NOR (relatively simple as an abstract problem) are out of scope of a one-layer perceptron. It is not linearly divisible, thus, at first sight, it couldn't be solved with the neural networks of the time. In the 80's, the field was reborn with the appearance of more intrincate and sophisticated architectures (multilayer, convolutional networks, tree networks, etc) and learning rules. Emergent memory processes and feature extraction and match were studied [3], until now, where CV modules are able to identify objects in different, changing environments and perspectives, in lack of good visibility conditions, with astonishing accuracy, and in real time.

1.2 Goals

The motivation behind this work is to find a way of dealing with the vast amount of methods relevant to the task of Robotic Vision and finding the ones that work best. In the process we'll explore the common, shared basis of all these methods trying to get some insights for future elections of one method or

another, based on the task at hand. This will be useful for researchers as it alleviates the decision process needed when researching higher levels of abstraction that depends on Robotic Vision, making them able to forget about the technical issues and focusing on higher level subjects.

All in all, the benchmark proposed in this study wants to be plausible and realistic with nowaday's robotics' constraints. Thus, some conditions must be taken into account when assessing the different methodologies studied. For example, as a CV module installed in a robot must work "on the fly", real-time processing will be an important issue to be evaluated. In the section dedicated to the method, all the criteria will be explained in detail.

Therefore, the primary goal of this report is to compare classic and modern approaches to solve the task of image classification, in an attempt to grasp some intuition over.

To address this goal, we'll present a benchmark comparing classic methods with modern networks methods, both solving the problem of image classification in the context of robot vision. As benchmarking all the existing methods could be nearly impossible due to the number of procedures in the field, here we'll review some of the most representative ones, from classical, well-established methods to more modern approaches.

Aside from the benchmark itself, this report aims to outline the evolution of the techniques utilized to solve the problem of image classification, trying to evaluate pros and cons of every method.

1.3 Contents of this Report

In this report we state a clue on which method to choose for the task of image classification. In chapter 2 we expose the state of the art. In chapters 3 and 4 we briefly explore the history of image classification tasks and some of the most used methods, reviewing both classic, raw statistical methods and more modern artificial-network-oriented methods, going through it's contributions, constraints and pitfalls. Two experiments are proposed to test each method in chapter 5. In chapter 6 we explain both methods used and it's peculiarities, we continue explaining how the tests were conducted and the results obtained in chapter 7 to end up with the obtained conclusions and further lines of work in the last chapter. Appendixes 1 and 2 contains the gory details about software, hardware and source code.

2 State of the Art

This chapter will explore the state of the art on the subject of image classification with neural networks. We will explore the most modern and stable methods using neural networks and finish with the latest approaches being developed nowadays.

2.1 Neural Networks

A neural network is a computing system inspired by the biological neural networks that constitute animal brains. Such systems learn over time to do concrete and general-purpose tasks. Artificial Neural Networks (ANN) are composed of neurons and links. Neurons are simple imitations of biological neurons, with axes and dendrites, the building block or ANN. Links act as synapses between neurons, connecting them, and transportating the signal from one neuron to a given number of other neurons. Neurons receive a fixed number of weighted inputs from other neurons. An activation function operates with the input as argument, producing an output to be propagated. This is the most general case, but things can be much more complicated. Weights can be binary or discrete (generally in a range between 0 and 1), and act as a way to ponderate the importance of every neuron in the other's behavior. The activation function can be of any mathematical nature, and, depending on the task, some are preferred over other ones. In classical formulations, the activation function could be a binary clasificator, a threshold. If the weighted sum of inputs reach a fixed value, the neuron would "fire", being this firing 0 or 1 (step function). More modern formulations include logarithmic compressions, non-linear functions, and discrete (or continuous) inputs and outputs, that opens the possibility of much more complex tasks to be resolved.

Neurons are connected in layers, and activation propagates (or not) from one neuron to another, through the entire network (or the point where activation stops). Almost every modern ANN is composed of an input layer receiving the information to be processed, a number of hidden layers, and an ouput layer. Figure Fig. 1.3.1 represents schematically this notion. The number of hidden layers is also decisive in the complexity of task accesible to the ANN, and major advances in the field come from research in the number and nature of hidden layers and its effects on the output of the system [4], [5], [6]. It is remarkable that, as the number of hidden layer increase, our understanding of the final output and our capability to predict it decrease dramatically. Input, hidden and output layers can be connected in many fashions, that will, also, determine the performance of the network on different tasks. The final result of any input propagated through the network will depend on the pattern of connections in every layer, the functions governing neuronal activation and the weights fixed for each link between neurons. The structure (pattern of connections) and the activation function play a decisive role in the final result offered by the network; basically determine the global model of the network, that has to be in agreement with the nature of the task to be accomplished. The model refers to the general operations carried out over the input data to generate the output. Weights determine the parameters of the global model implemented by means of structure and activation functions. From this can be followed that, modifying the weights, the output of the model will vary. This is the key to learning rules, as the classical and still prevailling Hebb's Rule. Hebb's rule and its variations are based on the principle that little adjustments in the network's weights will lead to a stable, global solution to the problem. Its application is iterative: in each iteration, ANN's output is

compared to the desired output (target output), and the difference between them calculated. Then, weights are slightly modified, and, if the difference is reduced in the next iteration, the change is kept and repeated, until the real output fits the desired output to a fixed threshold (the maximum accepted error). The ratio of modification per iteration, or *learning rate* (η) can be set to any value, and determines the speed of learning. Its classic formulation is: $\Delta w^i = \eta x^i y$. The change in the *i*th synaptic weight w^i equals the learning rate η times the *i*th input x^i times the postsynaptic response y [3].

The input could be, as in the present work, a stream of video or a set of images. Nevertheless, almost every piece of information can be represented in a suitable form to fit ANN requirements. Thus, ANN could perform statistical analysis over any given data; for example, linear regressions over healthy habits questionnaires, or Principal Component Analysis over cosmic microwave radiation data.

In our case, the input will be a set of pictures, and the general model must perform the task of image classification. The output is precisely the classification. Various models will be tested to inform about their performance.

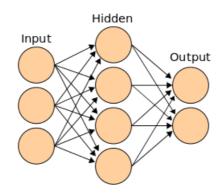


Figure 1.3.1. - An Artificial Neural Network with one hidden layer

To sum up, a neural network can be though of as a conjunction of mathematical operations, usually in the fashion of linear algebra, composed or neurons arranged in layers, and synapses connecting them. It is usually implemented as matrix multiplication operations. This gives us the ability to model many problems, from linear, simple equations to complex, differential equations with many parameters.

2.2 Why Neural Networks?

Neural Networks are task-nonspecific, meaning that, depending on how the problem is modelled, a wide variety of problems can be solved. After categorizing the problem's solutions we can choose between many types of networks, and start testing its performance. Iteratively, weights are adjusted until the solution is stable and accurate enough, and the network can be applied not only to the problem for which it was trained, but also to other analogous tasks. The key feature to keep in mind is that many of the problems we can think of can be expressed in a mathematical way. This is precisely what neural network do best.

2.3 Deep Neural Networks

Deep Neural Networks are one of the many types of neural networks. It's main difference with other networks strives in the amount of hidden layers. Every architecture with more than one layer is considered as deep. It first appeared in [7], where a deep, multi layer perceptron was explored to overcome the problems of the original Rosenblatt's multilayer perceptron. It supposed a revolution in the field, and open the door to real-world complex problems never addressed before by neural networks.

At that time, training networks iteratively was a problem, because computation resources were very limited and no mathematical method was known to work fast and efficiently. In 1989, LeCun et al. successfully applied the backpropagation algorithm (made popular by [8]) to the problem of recognizing handwritten ZIP codes. It took over 3 days to fully train the network [9].

With the develop of better methods (backprop), what was left was the computational cost of the learning process. Nowadays, with modern processors the problem is rapidly fading out. Some authors like Coates et al. [10] explored the use of GPUs over standard processors to train networks, resulting in much more powerful computation capabilities, that conform the standard in modern days.

2.4 Convolutional Neural Networks

A convolution is a mathematical operation that combines two functions to form a third one, it does so by "sliding" one function over the other and multiplying the points that overlap, it is similar to *cross-correlation*. Common applications are *Digital Signal Processors*, which are the core of audio systems or *Image Processing*. Applied to our field of image classification we can think of image as two dimensional functions (we represent images as two dimensional matrices) and we can convolve the image with a small function that we call "kernel". To convolve, we slide the kernel to every position of the image computing a new pixel as the weighted sum of the pixels the kernel is floating over. A visual representation can be found in Fig. 1.3.2

This kind of procedure makes a lot of sense when thinking about images features, for example, we have explained before that adjacent pixels keep some correlation. With convolutions this correlation is captured as the kernel slides over the pixels, the amount of adjacency can be regulated with the *size* parameter of the kernel.

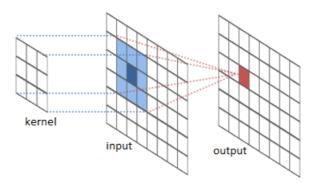


Figure 1.3.2. - A kernel convolution over an image.

One great side effect of this setup is that neurons are only connected locally, conforming regions or groups (often called patches). This saves us from a very big weight matrix, which in turn alleviates the computation process and helps scaling the architecture.

2.5 Residual Neural Networks

It has been stated that depth plays a big role in deep networks[5], [6], but as depth grows difficult in training does so too, as well as overfitting. This happens because of the training techniques used, usually SGD (Stochastic Gradient Descent) and the random initialization of the network [4].

To overcome this problem many techniques have been explored. [11] and [12] tried to initialize the network normalized while [12] introduced intermediate normalization layers. Both methods seem to work well for very deep networks but it comes at a cost: as the network starts converging to a stable solution, accuracy saturates and then degrades very fast [13], [14].

In the last decade of the past century,[15] and [16] proposed a technique for connecting layers called *shortcut connections*, illustrated in Fig. 1.3.3. It works by connecting layers skipping one or two other layers in between, giving a sense of fast forwarding knowledge.

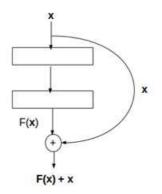


Figure 1.3.3. - A shortcut connection between two layers

In [4] they bring the concept back again and use it to solve this degradation problem with amazing results,

so much that it served to win the *ILSVRC 2015 & COCO 2015 competitions*, which are the most important contests of image classification and segmentation.

Networks setup with *shortcut connections* are what we call *Residual Networks*, the key of this mechanism is it's ability to recover previously stored knowledge that is fading out (degrading accuracy) by bypassing the units causing that fade and forwarding that knowledge, then it is restored with backpropagation to the conflicting units.

3 Neural Networks with Classic Methods

In the following chapter we'll review the most common, classic methods utilized in the past years to solve image classification tasks. Feature extraction and classification will be presented, as well as a standard optimization procedure. At last, a classical method to benchmark will be proposed.

3.1 Classic Methods for Image Classification

The problem of classification has been a tough nut to crack since the beginnings of computer science. The first approaches were based on raw statistical methods, suitable for different kinds of classifications. *Linear Classifiers* appeared first, comprising techniques like LDA (*Linear Discriminant Analysis*) [17], *Logistic Regression, Naive Bayes Classifiers* or the well known *Perceptron* proposed by Rosenblatt back in 1962 [2]). Later, in 1995 the SVM (*Support Vector Machines*) appeared [18], which can be though of as a small network (see below).

Feature Extraction

Image classification has been a subject of study for a while. As soon as 1973, Haralick et al.[19] explored the task by extracting basic features from images with statistical methods. Images are worked out as matrices of pixels to operate upon, just like nowadays. In their method they state what an important characteristic must have to be considered classifiable and then propose various characteristics that can be found in any image and the methods to extract them. Amongst other features, they propose measures of correlation between pixels and entropy.

Along these lines of study, different methods started to appear: points of interest tracking[20] in 1981, improvements over corner detection [21] in 1988, motion recovery (1992)[22] and features detection with a moving correlation window [23] in 1995. All these methods share the same working principle: find characteristic, important features in the most invariant and robust way and work on those features. All of them work with correlations between pixels and gradients made out of these pixels, in the same picture, or along a succession of pictures. In this sense, statistical measures like gradients analysis and correlation have played a central role.

Following these statistical measures we reach to the *Scale Invariant Transform* method (SIFT) [24], a method to extract features from images that can be invariant to image scale and rotation, this method is powerful, but at the time it was slow to compute. Ke and Sukthankar used PCA to normalize gradient patch instead of histograms [25] and showed that PCA-based descriptors can also be distinctive and robust to image deformations. Later on, in 2006, the *Speed Up Robust Features* method was proposed [26], being speed one if its selling point against SIFT and PCA-based descriptors.

Feature Classification

In [27], a decision technique called *K*-Nearest Neighbour (K-NN) is explored. Basically, K-NN assigns to an unclassified item the *class* most heavily represented among its *k* nearest neighbours, to do this, during training we expose the classifier to different items with *n* features, it then creates an n - dimensional feature space and projects the data onto it. With this feature space constructed, we can input an unknown item with similar features and extract the distance of every neighbor to it. Although this method is slow (it has to check the distance of every item with every other item), it is still in use in image classification as it is easy to implement and understand.

A distinctive approach is that of *Support Vector Machines* (SVM), based on Fisher's pattern recognition work [28]. Cortes et al. (1995) in [18] explored the concept creating a *Support Vector Network* able to recognize handwritten digits. The core idea is to non-linearly map input vectors to a high dimensional feature space and construct a decision linear n - dimensional divider inside this feature space. SVMs have been widely used for more than two decades with great results in classification tasks.

3.2 Optimization

All the aforementioned methods share a common problem: the bigger the input dimension, the slower the algorithm becomes. This is due to the fact that the features are extracted operating on the whole pixel matrix, SIFT and SURF pre-process the image best guessing points of interest, but they are slow to converge though. To overcome this problem, the preferred approach is input dimensionality reduction. We'll review basic techniques to achieve this.

Speed up by Dimensionality Reduction

A first step into dimensionality reduction could be acknowledging that, in many cases we can perform the classification over a well distinctive feature descriptor rather than over the original image itself. In this sense we can note that in many situations, color histogram is a great descriptor, in [29] they demonstrate it's effectiveness with an SVM.

We can reduce dimensionality even further by applying some of the revised techniques so far, being PCA (*Principal Components Analysis*) or LDA (*Linear Discriminant Analysis*) based methods the most common. From a classifier perspective, the details are not so important and in fact, they can represent noise. That is why summarization techniques like these work so well.

Another commonly seen technique is to stack processing units in such a way that minimizes computation. A great example are the Support Vector Machines, which are made of feature classifiers. Each classifier is specialized in a certain feature, then we stack all of them ranging from more abstract to more concrete features, this way the top classifiers (abstract) filter out non-interesting points or regions, saving time and computation to the lower classifiers.

3.3 Used Method for the Benchmark

For our benchmark we are going to use the histogram of the images to characterize them and classify with a K-NN. We use this technique because is a well established method to classify features over a data set.

4 Neural Networks with Modern Methods

So far we have reviewed some classic methods for image classification, in this chapter we will do so for the approach of modern neural networks.

In the last two decades, computers computation capabilities have been enormously increased, this fact has made possible for us to create bast neural networks and test them. In the past, nobody could theorize about very big networks because they were not able of test it nor computers able to run it, nowadays, we are breaking that barrier with the usage of GPUs optimized for the task.

4.1 Modern Networks for Image Classification

In 1998 LeCun et al. [30] proposed an architecture, *LeNet5* (Fig. 1.5.1) one of many previously successful iterations of a convolutional architecture for digits recognition.

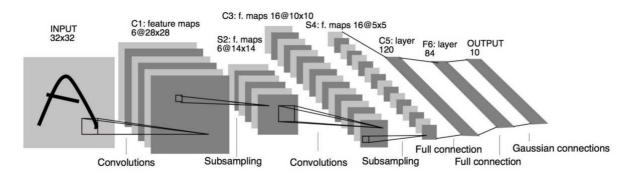


Figure 1.5.1. - LeNet5: Convolutional Neural Network for digits recognition

LeNet5 showed that as images are spatially correlated, using a multi-layer network for the input in a perpixel basis would not take advantage of this correlation, therefore LeNet5 uses a convolutional layer with learnable parameters as input, which is an effective way to extract features distributed across the image with fewer parameters.

Until the arrive of big GPUs, computation has always been a limitation factor for neural networks, but many research has been done. Ciresan et al. proposed it's usage for MLPs (*Multi Layer Perceptrons*) and managed to outperform previous results [31], but it was with[10] that GPUs established as the first choice for neural networks computations.

With powerful performance, Alex Krizhevsky et al. managed to train a very deep network (AlexNet,Fig. 1.5.2) for the ImageNet competition, winning by far to the other competitors[32], it was based on LeNet5 but with some more insights: it uses ReLU (Rectified Linear Unit) as non-linearity (instead of tanh) and introduced dropout and data augmentation, which are techniques to avoid overfitting.

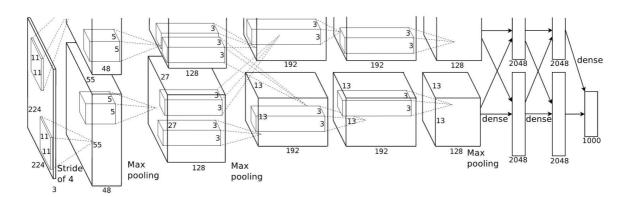


Figure 1.5.2. - AlexNet, winner of the ImageNet 2012

There are many more important deep architectures, we will review some of them in the following sections.

4.2 Why Convolutional Networks?

There are many types of deep networks, but why use convolutions? It has been thoroughly demonstrated that convolutions work well for the task, they are able to capture the correlation between pixels and regions in the image as well to remember lots of different classes.

4.3 Common Convolutional Network Architectures

The most common layers in ConvNets are:

- Convolution: This is the most important layer, it convolves it's input and has some learnable parameters: kernel size, stride, zero padding and number of filters.
- Non-Linearity (Usually ReLU): This layer applies a non linear function to it's input (non saturating also, in the case of ReLU), without this layer the hole network would be just a linear classifier.
- Pool (Usually Max Pooling): This layer is used to downsample the input, this is useful to reduce the conv layers processing overhead.
- FC (Fully connected layer): This layer is a full network by itself, it is normally used as the last layer of the ConvNet, as a classifier of the convolutional part.

A commonly seen configuration is INPUT -> [[CONV -> RELU]*N -> POOL?]*M -> [FC -> RELU]*K -> FC , where N an M are multiplicative factors and ? implies that the layer is optional.

4.4 Used Method for the Benchmark

Training a deep network is hard, requires lots of data and a powerful computer for a long period of time, hence instead of training our own network we are going to use a technique called *transfer learning* (more on this in Comparison section) with a pretrained network. We will use Inception v3 (Google's image classification network) with *tensorflow* (google's proprietary library for artificial intelligence) and Python to retrain the last layers so that it recognizes our data.

5 Design of the Experiment

The main goal of image classification is to categorize digital images under a given criteria or a set of possible choices. To do so, many different algorithms are available. Techniques focus on analyzing the image as a whole, or a portion of it; then, try to assign one or various tags, based on the analysis over the pixels. The classification step is usually a low level tool used as a background for other tasks such as segmentation or edge detection. It is therefore intended to be a semi automated task which runs on a given image or set of images and outputs a classification tag which might be used by upper levels of abstraction.

A simple task of recognition could be approached as follows: first, select objects from the image by segmenting it. To do so, scan pixel by pixel changes in brightness. When an abrupt change is detected, select it as a contour, thus dividing pixels at both sides of the gradient change as different objects. Also, when contours are closed, the inside of that contour can be considered an object. Then, compare the extracted objects to a database of known objects. If the configuration of pixels (in coordinates and brightness) of the objects extracted are under a certain, fixed threshold of variability in comparison to those stored in memory (or a linear transormation of them), then considered both objects the same.

Many other ways exist to solve the problem of image recognition, resembling in variable degree the human vision system. The only way to asses their efficacy is by facing them to real image classification problems and measuring their performance.

This benchmark aims to measure different methods for image classification applied to a real world use case: A robot moving in a environment where it knows the map and tries to go from one (start) point to another (end) recognizing each landmark from a camera. To do it we will be using the images classified as landmarks (split from videos) and the following methods:

- Classic method: Color histogram with a K-NN classifier and opencv.
- Modern method: Transfer learning with Inception v3 and tensorflow.

K-NN works by composing a search tree with the possible pre-specified classes and then matching the input against the nearest class. The search tree is constructed during the training step, it will take as input the color histogram of every image in the training set along with the class the image corresponds to. With the convolutional network we will retrain the last layers of the pretrained network with our train set.

For each method we will: train, test and measure it's response time and accuracy.

5.1 Environment

To ensure the benchmark is as fair as possible with both methods we will apply the following requirements:

- **Isolation**: The benchmark of every method is going to run in an isolated docker container with no other application running, this should minimize resource competition between processes.
- Similar technology: Both benchmarks will be run in a custom python script. I chose python over

other programming languages because it is easy to read and both tensorflow and opencv libraries have bindings for it. If we were to program with different languages time-efficiency issues might be encountered.

5.2 Landmarks

We have established various landmarks (points of the map) from the Computer Science Systems Department (*Departamento de Sistemas Informáticos*) and it's insides. We have captured some videos, split them into images and composed a map with them, we detail every one in the following chart:

#	Name	Description	Sample
1	Floor	The floor in the just before the departments' doors. First floor, fourth block	
2	Doors	The doors to the Computer Science Systems Department and the Economics Department, just before the corridor. First floor, fourth block.	
3	Corridor	The corridor to the Computer Science Systems Department.	
4	Mailboxes	The mailboxes in the corridor, just before the departments.	
5	Fountain	The fountain at the end of the corridor, near the offices.	

#	Name	Description	Sample
6	Bath doors	The bath doors at the right of the fountain.	
7	Office	The Bachelor's director office.	
8	Corridor way back	The corridor from the other side, the way back to the floor.	

5.3 Datasets

We will divide all our images by landmark, then, for every landmark we are going to take 80% for training and 20% for testing (these are typical numbers, not just randomly chosen), we do this to avoid testing with the same images that we train as it would be too easy for any classifier, after all it would have yet seen the image.

The procedure for composing the samples is not complex:

- 1. Take various videos of all the landmarks
- 2. Split the video in frames
- 3. Classify each frame (bunch of frames) by landmark
- 4. Augment samples (more on this in Overfitting section)
- 5. Separate frames from each landmark in trainset (80%) and testset (20%)
- 6. Reduce dimensionality (reduce images from 1920x1080 to 250x141 and 1080x1920 to 141x250 and grayscale) to avoid large computation

Before the augmentation step I was able to obtain a good amount of samples, about 1000 samples average, being 518 samples the minimal by landmark. It is worth noting that for the K-NN algorithm to

work well the dataset must be well-balanced (although much implementations work this out), if the amount of samples in a class is much bigger than in other classes, the probability that a neighbor of that class is found get's increased. Also, as we have images in portrait and in landscape, width and height will differ, which in turn will affect the histogram, for this reason we do not take the histogram itself but it's normalized version.

Although for K-NN we will resize every sample to 250x141 or 141x250 in order to avoid possible histogram pitfalls and process only in graysclae, in the case of the ConvNet this is not necessary as it automatically rescales everything to 299x299x3, being the last x3 the number of channels, so full color images are to be used.

Landmark	No Augmentation	Hflip Augmentation	Balanced
Office	1464	2928	1036
Fountain	518	1036	1036
Mailbox	706	1412	1036
Doors	1560	3120	1036
Bathdoors	662	1324	1036
Back	1443	2886	1036
Corridor	798	1596	1036
Floor	969	1938	1036

To sum up we will have a dataset with the following characteristics:

5.5 Overfitting

A common problem when training is overfitting; that is basically that your model learns too well your data and so it does not generalize correctly. We definitely do not want this to happen, to avoid this effect we are going to *augment* our data, which consists in taking our data and do some transformations on it so that the dataset grows with samples that are slightly different.

There are many techniques applicable here, we could crop the image, change the color map, zoom, rotate or flip it. In our case we are going to augment our data by a factor of two

Also, we must estimate the best values for our models. In the case of the K-NN we have a very important one, the number of neighbors used for predictions \mathbf{k} , to estimate it we are going to use cross-validation with our trainset, this will help us to prevent overfitting too as it vanishes the importance of initialization in our trainset.

5.4 Measures

Accuracy

This is the precision of each method in test. For each class we are going to take all the test images and pass them through the classifier, obtaining a predicted class for each image, then we average that class predictions to obtain an average accuracy for the class, we will call this average it's score. Finally we will average the score of all classes obtaining a global score for the classifier.

Time

This is time spent by each classifier while predicting the class of each image, this will be used as a constraint for real-time. We measure it by getting the delta time before and after the predict function call.

5.7 Algorithm

The algorithm in pseudo-python code can be seen down below. This procedure is done for both methods, we will see the comparison between the scores and global score of each method as well as the times and global time.

Whenever a prediction is done correctly (the predicted class for the image is the actual correct one) we add one point to the score, if it fails no penalization is applied.

```
# our landmarks
landmarks = load_landmarks()
global_scores = []
global_times = []
for landmark in landmarks:
    score = 0.0
    time = 0.0
    # array of tuples like (image, landmark)
    landmark_images = load_images(landmark)
    amount = float(len(landmark_images))
    for (image, correct_class) in landmark_images:
        # time before the call
        dt = current_time_millis()
        predicted_class = predict(image)
        # time elapsed: delta time
        dt = current_time_millis() - dt
        time += dt
        if correct_class == predicted_class:
        score += 1.0
    # average landmark score
    score /= amount
    # average landmark processing time
    time /= amount
    global_scores.append(score)
    global_times.append(time)
global_score = average(global_scores)
global_time = average(global_times)
```

5.8 Features

One downside of the statistical methods like SIFT or SURF is it's slowness, it is has to execute the whole algorithm for each input in runtime, this means that, for every input image the algorithm has to complete entirely, so if it is slow it will delay the results for a while, therefore some of this algorithms may not be suitable for vision as it would violate the real-time constraint.

In the case of K-NN, classification could be slow if we had a dense feature space; this is due to the fact that the training data is stored in memory, as it has to be accessible during classification. Also, as the **k** parameter grows, the number of neighbors to take into consideration does too, which in turn may incur

into computation delay. Our data is not dense in features nor in number, so this will probably not be an issue.

On the other hand, neural nets are slow to train but faster in runtime, this is due to the way networks work. When training, backpropagation is used to transmit the gradients back to the input for the parameters to be learned, that is slow but as it is only used during training, not in normal use, it does not delay so much. A forward pass could be slow too, depending on the number of layers and operations per layer but this is usually been taken care of by design. However large architectures like the one we'll employ here will always have a considerable processing overhead, only surpassable with more computation capabilities, namely a GPU card. This is because tensorflow was designed to scale seamlessly across cpu and gpu available in

runtime.

¹. Docker is a containerization technology that is able to take full cpu and memory, we will use it to avoid installing all necessary software and instead used prepared containers that take full advantage of our resources.

6 Design of the Classifiers

After reviewing the whole setup for the designed experiment, it is time to present each classifier and how we are going to use each one.

6.1 K-NN

We train our classifier with supervised learning, passing through the histogram of each picture along with it's corresponding class. This will setup the whole feature space the classifier is going to memorize . To do so, the classifier understands each histogram as a multidimensional feature vector, to which a class label is to be assigned; therefore the process of training consists in storing each feature vector along it's class label. During the classification phase we ask the classifier to predict the class of each item (histogram of every image). In order to do it, the classifier assigns the label which is most frequent among the **k** nearest neighbors to the query sample. The euclidean distance is widely employed to calculate the distance of the neighbors to the sample although other measures have been used too (Hamming distance, Pearson and Spearman).

OpenCV's k nearest neighbor implementation uses the euclidean distance when employed for classification (Bayes is used for regression), which is a safe choice, but not the best. The distance metric method used can push the classification accuracy significantly, Goldberger et al. explore solutions to learn how to reduce dimensionality before the training step and suitable distance according to the nature of the data [33] and [34] tries to establish a reliable method for the case where the distances between neighbors is big.

Besides the method employed for metric distance, there is another important parameter, the *k* neighbors to use when classifying, in our case we will use cross validation to estimate it, the full process is reviewed hereafter.

K Neighbors Estimation

The k neighbors parameter is of the most importance in the K-NN algorithm, it is the amount of neighbors used for matching. We will use a type of cross validation called *k-folds* to estimate the best *k* neighbors for our data, to do so, during training we will randomly partition the dataset into k equal sized subdatasets, then we will train on k-1 subdatasets and validate on the remaining one, this process is repeated k-1 times more so that every subsample is used for validation only once. By measuring the error on each pass and averaging it we obtain a more accurate error rate for the chosen *k* neighbor. We will repeat this procedure for all the values of *k* we want to test and choose the one that minimizes the error. We pick k - folds = 10 as it is a multiplier of the amount of samples we have on each class. The pseudo-python algorithm showing the whole process can be seen below:

```
k_folds = 10
k_values = [1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14]
error_rates = []
for k_value in k_values:
   folds = get_folds(k_folds)
   error = 0.0
    for fold in folds:
       knn.train(fold.train_set)
       fold error = 0.0
       for (sample, correct_class) in fold.validation_set:
            predicted_class = kn.predict(sample)
            if predicted_class != correct_class:
                fold_error += 1.0
       fold_error /= float(len(fold.validation_set))
       error += fold_error
   error /= float(len(folds))
   error_rates.append(error)
# best k will be the k value at the position where the error is minimal
best_k = k_values[ error_rates.index(min(error_rates)) ]
```

After running the cross validation I found that the setup that works best is for k = 1 (depicted in this chart), so this is the value we will use for the benchmark.

К	Error	К	Error
1	0.03188	8	0.08773
2	0.05156	9	0.09156
3	0.06250	10	0.09414
4	0.07102	11	0.09891
5	0.07492	12	0.10016
6	0.07992	13	0.10312
7	0.08234	14	0.10719

With the chosen value we can start the benchmark for the classic method.

6.2 Convolutional Network

As a convolutional network we will use Google's Inception v3, depicted inFig. 1.7.1 using the *transfer learning* technique in *tensorflow*. The great success of this architecture is featured by one the greatest insights of the authors: *The Inception Module*, which gives the network one of it's names (the other name is *GoogLeNet* in homage to Yann Lecun's LeNet).

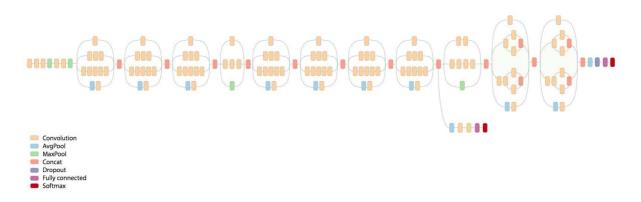


Figure 1.7.1. - Inception v3. Google's classification network

The Inception Module

Being one the greatest contributions of this network to the field development, the Inception module is used as a way to enable the network choose how to combine different filters on type and shape in runtime, as well as to reduce computation overhead in larger convolution filters [6]. The Inception module captures the intuitive idea that different spatial features can be captured at different scales - 5 * 5 and 3 * 3 filters - which can be aggregated later - filter concatenation - so that the next stage can abstract features from different scales simultaneously [6]. To do so it stacks convolutional layers in parallel, splitting the input in various processing pipes and joining them later. To extract different scales, different filters must be used, but it comes at a cost; with large filters computation blows significantly, to over come this issue, various 1 * 1 layers are used, specially before each large convolution filter - 3 * 3 and 5 * 5 -, therefore these 1 * 1 serve mainly for two purposes:

- Reduce filter space dimensionality: By applying 1* convolutions with smaller filter depth than the input we obtain a reduction in filter space (eg: We have a convolution layer that outputs an (N, F, H, W) shaped tensor, where N is the batch size, F the number of convolutional filters and H, W are the spatial dimensions. If we pipe this output through a layer with $F^1 1 * 1$ filters, then the output of this 1 * 1 conv layer will have shape (N, F^1, H, W) . When $F^1 < F$ we reduce filter dimensionality and conversely, if $F^1 > F$ we augment it).
- Augment nonlinearity: Using a ReLU (Rectified Linear Unit) just after each1 * 1 layer gives us more nonlinearity which is specially useful to create richer feature maps.

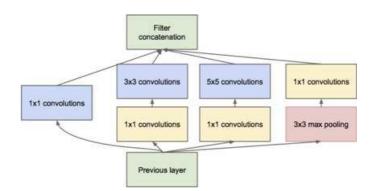


Figure 1.7.2. - GoogLeNet's Inception module

When designing an architecture, one must balance the processing power of the network and the computation requirements of it, this usually implies choosing between different filter sizes. With this parallel approach and the 1 * 1 filter dimensionality reduction technique we are able to use deeper filters and, instead of choose between them, let the model pick the best for the task.

Transfer Learning

Transfer learning is a technique that enables us to use the all or part of the previously stored knowledge of a network and use it as the basement for a new network that can learn more things, this way we can use the more generic part of Google's network and replace the final layers (more concrete) with the knowledge from our data.

there is a handy script that comes with the tensorflow library just for this task. It does basically the same cross validation procedure showed above to estimate the final accuracy of network after the training process. We utilize this procedure instead of training a network from scratch because it's much less exigent in terms of dataset size and setup, computation capabilities and time spent to train.

In the Common Architectures section we explored the common setups for this kind of networks, a common principle can be seen: the network ensembles many layers to extract features from more general to more concrete, until the end where a full connected layer acts as a classifier for the rest of the network; what we derive from it is that the network is a feature extractor and compactor, encoding the information extracted from the image along it's pass through the network, therefore, as the input to the last layer we can observe an encoded and compact summary of the image under classification. In reaction to this input, the last layer must output a label that thinks to be representative of the picture, which is exactly what a classifier does, in fact, as a final layer we could plug in an SVM or a simple softmax. In Inception v3 this layer is trained to output one of the 1000 labels used for ImageNet.

It is this last layer that we're actually going to train, instead of the ImageNet labels, it will output ours. As can be easily seen at first sight, we need much less time to train one layer than the full architecture, that's why this transfer learning procedure is so fast. Additionally, as for the classifier to interpret the network data there is a penultimate layer, informally called "*bottleneck*" that is trained to output meaningful information for the last layer. This layer must be setup for the dataset to retrain with and that is the first step when training. The bottleneck process consists in feeding the network with the samples of the trainset while persisting the output in this penultimate layer to a cache; this cache will be use as the input to the classifier to speed up computation.

During the training phase, the testset provided is traversed extracting the bottleneck file for each image and forwarding it to the classifier, which outputs a label that is tested against the correct label, the last layer's weights are updated accordingly through the back-propagation algorithm.

7 Experimental Results

In this chapter we present the experimental results of the benchmark, explaining how the experiment was conducted. Also, the methodology of each probe will be detailed.

7.1 Results

We do the whole process of training and testing with both methods measuring the time spent to classify and the accuracy, in the following charts we will present our results, first exposing the results of the K-NN method, then the ConvNet results and at the end a comparison about both results will be discussed:

K-NN

Efficacy

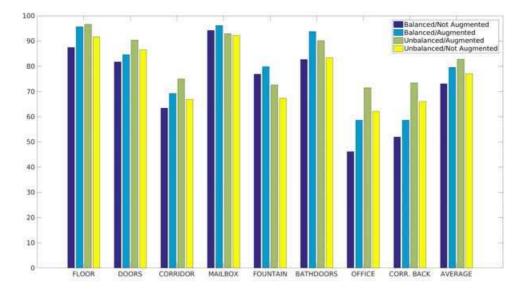


Figure 1.8.1. - K-NN. Average efficacy by landmark with dataset conditions: balanced and unbalanced data, with or without hflip augmentation.

	L1	L2	L3	L4	L5	L6	L7	L8	AVG	MAX
B/NA	87.5	81.73	63.46	94.23	76.92	82.69	46.15	51.92	73.07	94.23
B/A	95.67	84.61	69.23	96.15	79.81	93.75	58.65	58.65	79.56	96.15
U/A	96.65	90.39	75.00	92.93	72.60	90.19	71.50	73.36	82.82	96.65
U/NA	91.75	86.54	66.88	92.25	67.31	83.46	62.12	66.09	77.00	92.25
AVG	92.89	85.81	68.64	93.89	74.15	87.52	59.60	62.50	78.12	
MAX	96.65	90.39	75.00	96.15	79.81	93.75	71.50	73.36	82.82	

At sight of these results, it can be seen that there are some landmarks more difficult than others : for example the officce and the corridor back, and both cases work best with balanced and augmented data, which seems to work better in general. Also, it's worth noting that the other augmented dataset, besides being unbalanced, scores high too. The fact that both augmented datasets are clearly superior might led to the following reasoning: either the OpenCV K-NN algorithm implementations fixes the unbalanced proportions in the trainset or it denotes that the amount of samples radically prevails over the proportions inside the datasets, furthermore, it contributes to the model's resilience to image alterations.

Time

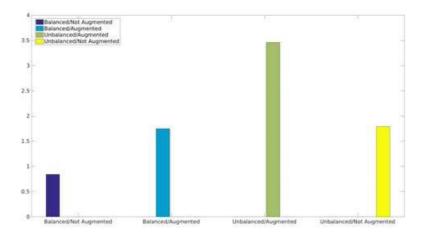


Figure 1.8.2. - K-NN average time to classify by experiment \(in ms\)

	Average time (in ms) to classify
B/NA	0.84
B/A	1.75
U/A	3.46
U/NA	1.79

The chart depicted above shows exactly what we expected: the more samples in the trainset, the more

time spent to classify. This happens due to the way classification works in the K-NN algorithm; when an item is to be classified it is compared to all the **k** nearest neighbors, so the higher the amount of near neighbors, the higher the number of comparisons, hence as the unbalanced, augmented dataset is the biggest, it also annotates the highest time to classify.

ConvNet

Efficacy

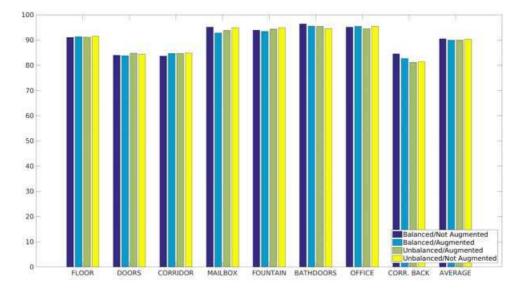


Figure 1.8.3. - Conv Net. Average efficacy by landmark with dataset conditions: balanced and unbalanced data, with or without hflip augmentation.

	L1	L2	L3	L4	L5	L6	L7	L8	AVG	MAX
B/NA	91.08	83.95	83.69	95.12	93.93	96.42	95.13	84.50	90.48	96.42
B/A	91.27	83.84	84.66	92.83	93.47	95.53	95.49	82.70	89.97	95.53
U/A	91.16	84.75	84.69	93.87	94.38	95.47	94.50	81.16	90.00	95.47
U/NA	91.52	84.41	84.86	94.91	94.91	94.56	95.43	81.42	90.25	95.43
AVG	91.25	84.23	84.47	94.18	94.17	95.49	95.13	82.44	90.17	
MAX	91.52	84.75	84.86	95.12	94.91	96.42	95.49	84.50	90.48	

From Fig. 1.8.3 we can clearly advise that the network cares very little about the dataset, they are almost tie. There are a few variations from one dataset to another, but the most significant fact is the landmarks poorly classified, being the corridor back among the most difficult.

Time

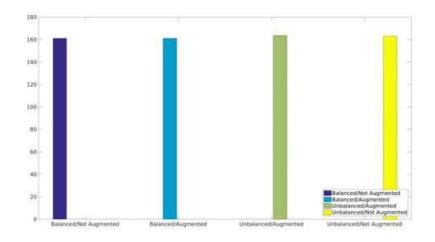


Figure 1.8.4. - Convolutional Network average time to classify by experiment \(in ms\)

	Average time (in ms) to classify
B/NA	161.21
B/A	161.18
U/A	163.71
U/NA	163.12

The convolutional network has a stable classification time, none of the datasets used variate much in the time spent to classify.

7.2 Comparison

Before comparing the results, a side note must be made: The network and the K-NN classifiers give different types of outputs, in the case of K-NN, the output is the class of the most representative neighbor, so it's either correct or incorrect; but in the case of the network, the confidence scores for each label is given, that means that the output is a range, not a boolean. If we were to compare making a boolean out of the confidence (ie: keep the highest score), we would have obtained a 100% score on every dataset and every landmark. As this is not desirable, we took the value of the label with the highest confidence for the benchmark and average over that scores.

After thoroughly examining the results there are some clear assertions to be made: K-NN outperforms the network in classify time but looses abruptly when it comes to classify. While K-NN scores, in average, **78.12**, the network's final average score is **90.17**, that is more than 10 points higher, which in this contexts means a lot. On the other hand, K-NN classifies almost 200x faster in the best case.

It's worth mentioning that, besides the method used, the corridor back landmark always gets the worst

scores. My best guess is that in the case of K-NN, this happens because of lighting conditions, meaning that the histogram is not really representative: as can be seen in the dataset, there is a lot of light, the same principle applies to the *office* which is basically a window of light. As for the network, I would say this happens because of the forms inside the picture which in fact, are pretty similar to the *corridor*, that is basically the same but in the opposite direction and also scores low.

8 Conclusions and Further Work

At sight of the experimental results it is obvious that both the classic and modern methods can be highly improved with further study in the lines of preparation of the dataset, parameter initialization and estimation and data pre-processing, nevertheless the results clearly showed that both methods are a bit narrow-minded, especially the histogram based K-NN which is easily tricked.

After this experiments we can easily extract a rule of thumb: If time is really crucial (in means of preparation, training and classification time), go for the K-NN, for everything else go for the Convolutional Network. Although the process of setting up and training the network is much more intricate and tedious I'd say it is much more reliable in classification and stable in processing time, which means it can escalate to a bigger number of classes and dataset sizes with almost constant classification time.

For the K-NN I would like to explore it's limits, as well as it's possible improvements. This algorithm works surprisingly well, but I feel it could be highly beneficial to use a different metric measure, specially the *Neighborhood Components Analysis* [35] which shows very promising results.

As for the ConvNet, I would be delighted to use other setups as the final classification layer, such as SVMs, a sparsely connected networks or even another little convolutional layer as explained here and here. The idea is that Fully Connected layers can be really slow, but luckily there is a direct transformation between the two types without loosing any information, which could only improve the processing time.

Also, it could be great to explore other methods like SVMs, SURF with PCA or Residual Networks and benchmark them, maybe taking different measures like generalization capabilities or fault-tolerance and test their comfort boundaries. Taking more classic and modern methods and benchmark them could throw more light on which methods to choose over the others.

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1	By en:User:Cburnett [GFDL (http://www.gnu.org/copyleft/fdl.html] (http://creativecommons.org/licenses/by-sa/3.0/)))], via Wikimedia Commons.
6	A kernel convolution over an image. Taken from [River Trail documentation] (http://intellabs.github.io/RiverTrail/tutorial/).
5	He, Kaiming et al Deep Residual Learning for Image Recognition.
4	Yann LeCun et al LeNet5, one of the very first successful convolutional networks aimed to digits recognition.
9	Alex Krizhevsky et al AlexNet.
2	Inception v3 - Google Classification Network.
3	ConvNet. Average efficacy by landmark with dataset conditions: balanced and unbalanced data, with or without horizontal flip augmentation. Plotted with matlab.

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Appendix 1: Technological Environment (Software & Hardware)

- To program the classic methods we use the well known OpenCV library for Python.
- To program the modern, deep neural networks we use transfer learning to Inception v3 with tensorflow.
- Both programs will run in an isolated docker container.
- All the work will be developed in a multi-core processor computer (no gpu) with the following specs:

Component	Specification
CPU	Intel Core i7 - 4790. 4 cores, 3.60GHz, 8MB Smart Cache
RAM	16 GB. DDR 3
Hard Drive	SSD 225 GB
SO	Arch Linux, kernel 4.4.63-1
Docker	17.04.0-ce, build 4845c567eb
OpenCV	OpenCV 3.1.0, Python 2.7
TensorFlow	TensorFlow 1.1.0, Python 2.7

Appendix 2: Resources and Code

Following a brief description of the project's structure and a comprehensible description of each component:

- The project is organized spread across the scripts , src and files directories.
 - Scripts: Contains all the scripts needed to do the whole preparation: split videos into images, resize images, augment them, separate datasets into trainset and testset and manage both docker containers.
 - Src: This is the source code directory, contains the necessary programs to train and test each model as well as measure the results and extract the benchmark measures.
 - Files: Here are contained all the assets for the project: train and test sets of images and dumps of the data and setup of each model.

The source code directory of each method contains one or various scripts to train and test each model and in the case of K-NN, the cross validation k - neighbor estimation is provided too.