

# **Machine Learning methods for Thin Vessel Segmentation**

in Fundus Images



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This thesis is submitted in requirements of the course BITS C421T

May 2015



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# Chapter 1

## Introduction

With the advent of medical imaging, computer aided diagnostic systems have become an integral part of today's medical diagnosis[3]. CAD systems have become a part of routine clinical work and are being used extensively for disease diagnosis. A myriad of different medical imaging systems, like X-Ray, Magnetic Resonance Imaging (MRI), Computed Tomography Scans etc, are used for diagnosis. The output of such systems are multi dimensional digital images, interpretation of which require sophisticated digital image processing methods. Automated medical diagnosis systems can aid in easy and faster interpretation of these images.

Digital fundus imaging in ophthalmology is a vital component in diagnosis of various pathologies. Retinal vessel segmentation forms an important part of diagnosis of such pathologies. Changes in the retinal vasculature is precursor to many diseases such as diabetes, hypertension and stroke. Morphological properties such as diameter, length, branching angle, of the retinal vessel forms an important component in diagnosis and evaluation of ophthalmologic diseases such as diabetic retinopathy [24] and hypertension. For example, vessel diameter measurement can be an aid in diagnosis of hypertension[1].

Vessel segmentation is itself a challenging and a tedious task which may take a couple of hours one done manually. Low contrast between vessel and background, noise in the image and variability in the width, brightness and shape along with the presence of exudates, lesions, hemorrhage spots and other pathological effects make the task much more difficult. Figure 1.1 shows a fundus image of a diseased eye.

Developing an automated retinal vessel segmentation is a first step towards developing a full fledged CAD system for diagnosis of ophthalmology pathologies. There have been a lot



Fig. 1.1 A diseased fundus image

of work in literature on automatic retinal segmentation, including based on matched filtering, tracking methods, morphological methods and learning based methods.

# Chapter 2

## Background and Literature Review

In this chapter we introduce the basic concepts and methods involved in our thesis. We start by introducing machine learning and the various machine learning algorithms. We introduce the idea of clustering and dictionary learning. The sparse coding and dictionary learning formulations are discussed in this section. Later we discuss the relevant previous work relating to our thesis. In the last section we briefly describe some of the latest work which resembles our work and represents the state of art methods.

### 2.1 Machine Learning

Machine learning is the study dealing with the process to learn characteristic information from data. Given some data,  $X$ , a machine learning algorithm learns a function  $f(X)$  which maps the input  $X$  to an output variable  $y$ . The learnt model then can be used to make predictions on previously unseen data  $X'$ . A machine learning algorithm learns the parameters of an adaptive model from a training set, typically optimizing a function. The learnt model is then used to make predictions on previously unseen new data. Broadly, machine learning methods can be divided into two subfields, supervised learning and unsupervised learning. In a supervised setting the target labels or values are known for the test set and we focus on predicting the target value for previously unseen data. In an unsupervised setting no such target information is present for the training data and the focus is to learn structure and compact descriptions from the data.

In a supervised setting, given a set of labelled data points known as the training data:

$$T = \{(x_1, y_1), \dots, (x_n, y_n)\}$$

, where  $x_i \in X$  and  $y_i \in Y$ , we find a function  $f$ , which maps any point in the domain of  $X$  to its corresponding label in  $Y$ . If  $Y$  is a set of discrete values then it is known as a classification problem and if  $Y$  is in a continuous range then it is a problem of regression.

In an unsupervised setting, the task is to find group relations between instances of the unlabeled training dataset with a subsequent aim of categorizing or clustering the data. The algorithm finds the previous unknown structure in the data.

The data point  $x_i$  is typically represented as a vector comprising of feature values and is known as a feature vector. For example, given a dataset  $X = \{x_1, x_2, \dots, x_m\}$  where  $x_i$  is an image patch of size  $\sqrt{n} \times \sqrt{n}$ , each patch can be represented as a vector of length  $n$  formed by concatenating the gray scale values. The entire dataset then can be represented as a matrix:

$$X = \begin{pmatrix} x_{1,1} & x_{1,2} & \cdots & x_{1,n} \\ x_{2,1} & x_{2,2} & \cdots & x_{2,n} \\ \vdots & \vdots & \ddots & \vdots \\ x_{m,1} & x_{m,2} & \cdots & x_{m,n} \end{pmatrix}$$

Here each row of the matrix denotes individual data points and  $x_{i,j}$  is the gray scale intensities of patch  $x_i$  at pixel location  $j$ .

## 2.2 Clustering

A cluster is collection of data points grouped together on basis of some common properties. The data objects or points within a cluster are similar to each other, whereas the points in different groups are dissimilar.

The process of partitioning the data points into smaller groups (called as clusters) with an aim to minimize the intra cluster variance and maximize the inter cluster variance, is known as clustering. The grouping of data points is based on the similarity or dissimilarity of the objects as described by their properties or features.

Similarity or dissimilarity between two objects can be very subjective and hence various measures are used to describe them quantitatively with distance measure being the most common. The distance measure is used to specify the distance between two objects and can be used to create a distance matrix called as similarity/dissimilarity matrix. The most commonly

used distance metric is the Euclidean Distance.

The euclidean distance  $D$  between two points  $p = \{x_1, x_2, \dots, x_n\}$  and  $q = \{y_1, y_2, \dots, y_n\}$  can be defined as:

$$\begin{aligned} dist(p, q) &= \sqrt{(x_1 - y_1)^2 + (x_2 - y_2)^2 + \dots + (x_n - y_n)^2} \\ &= \sqrt{\sum_{i=1}^n (p_i - q_i)^2} \\ &= \|\mathbf{p} - \mathbf{q}\| \end{aligned} \quad (2.1)$$

As the data labels are unknown, cluster analysis is known as an unsupervised method of data partitioning. This is in contrast to classification where the data can be partitioned on the basis of their class labels. Thus, clustering segments the data based on properties of the objects within the dataset and finds previously unknown grouping within the data.

### 2.2.1 K-Means Clustering

In this section we look into the K-Means clustering algorithm (MacQueen et al. [14]). We assume that the number of clusters 'K' is given and we use it to initiate our clustering algorithm. The idea is to partition the dataset into 'k' clusters and assign 'k' centroids, one for each cluster

Given a dataset  $D = [x_1, x_2, \dots, x_n]$  consisting of  $n$  objects, where each object is a feature vector of length 'm'. The clustering algorithm partitions the dataset into 'K' clusters  $C_1, C_2, \dots, C_k$  such that  $C_i \subset D$  and  $C_i \cap C_j = \emptyset$  and with an aim to have a low intra cluster variance and a high inter cluster variance. This means that the points within the same cluster must be highly similar and dissimilar to the points belonging to other clusters.

The partitioning is performed in an iterative manner and follows what is known as Lloyd's algorithm [11]. We initialize the 'k' centroid by some initialization methods such as k-means++ or random initialization. The cluster quality depends on a good initialization. We start by assigning each point in the dataset to the nearest centroid and we get 'k' clusters. We then calculate the mean points of the new clusters and assign them as the new centroids. The process then continues till a stopping criterion is reached. The partitioning is made with an aim to minimize the objective function which is the within cluster sum of squared error.

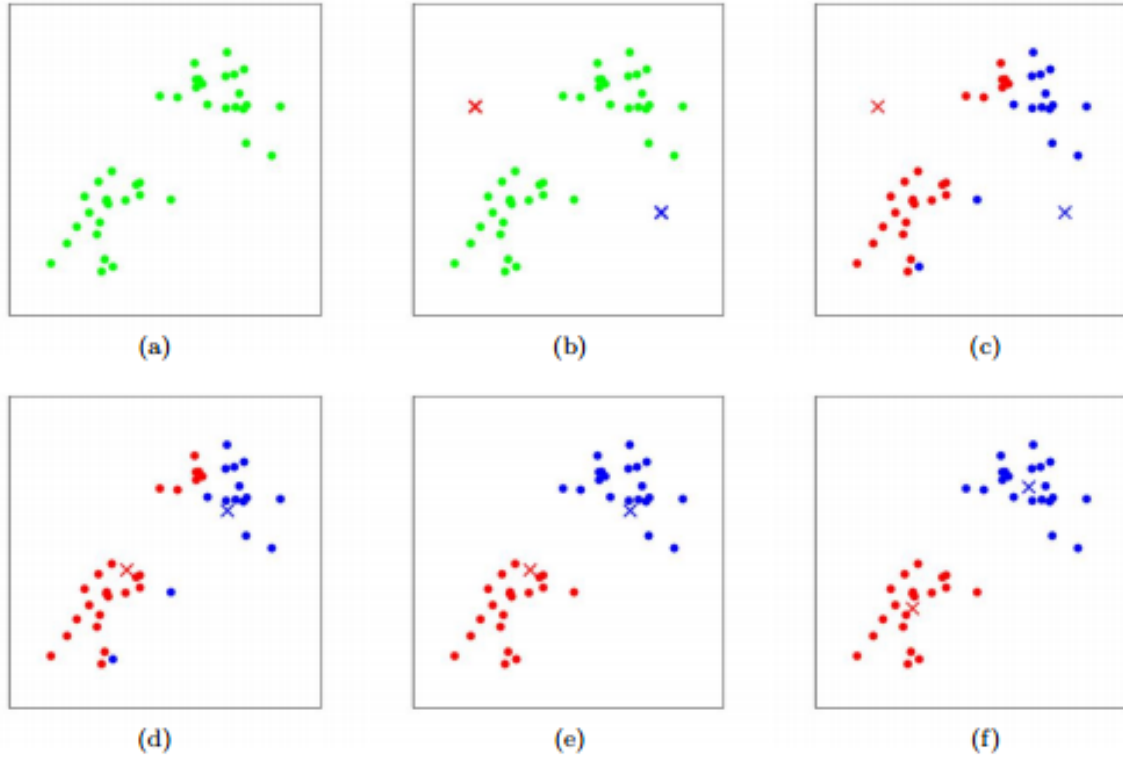


Fig. 2.1 K-means algorithm. Training examples are shown as dots, and cluster centroids are shown as crosses. (a) Original dataset. (b) Random initial cluster centroids. (c-f) Illustration of running two iterations of k-means. In each iteration, we assign each training example to the closest cluster centroid (shown by "painting" the training examples the same color as the cluster centroid to which is assigned); then we move each cluster centroid to the mean of the points assigned to it. (Images courtesy of Michael Jordan. Retrieved from : <http://stanford.edu/~cpiech/cs221/handouts/kmeans.html>)

Each of the clusters  $C_i$  consists of  $p$  points and is represented by its centroid  $c_i$ . The distance between point  $p$  and the cluster centre  $c_i$  is defined as  $\text{dist}(p, c_i)$ . If the data points are in  $m$  dimensional, then the distance can be computed as in Eq 2.1. We can then write our objective function  $J$  as :

$$J = \sum_{j=1}^K \sum_{i=1}^p \|x_i^{(j)} - c_j\|^2 \quad (2.2)$$

K-means clustering is the simplest unsupervised learning algorithm and can be applied to many different problems. An example of K-Means clustering on a toy dataset with 2 clusters is shown in Figure 2.1

## 2.3 Spare Coding and Dictionary Learning

The process of learning a finite set of basis elements ,called atoms, from a given training set of signals  $X = [x_1, x_2, \dots, x_n]$ , such that each signal in  $X$  can be approximately reconstructed by a linear combination of atoms is called Dictionary Learning. This finite sets of atoms is called  $D$ , is the learnt dictionary. Normally we want to find a sparse decomposition of signals in  $D$  i.e, we aim to create a dictionary of sparse elements. This kind of decomposition is different from PCA where the basis elements have to be orthogonal to each other and these dictionaries are in general overcomplete.

By overcomplete here we mean that if any element from the dictionary is removed, we can still approximately construct the signal from the atoms. Also the number of atoms in dictionary is of a greater dimension than the data. Lewicki and Sejnowski [13] describes the process of learning overcomplete set of representations. Some recent work [17] has shown that sparsity captures high-order correlation in data and it captures higher order statistics in data.

Huang and Aviyente [10] talks about how to represent a given signal as a sparse decomposition of a dictionary  $D$ . In a sparse coding task, our goal is to represent a signal  $x \in \mathbb{R}^m$ , as a linear combination of an overcomplete dictionary  $D = [d_1, d_2, \dots, d_k] \in \mathbb{R}^{n \times k}$ . Here 'k' is the no of atoms  $d_k \in \mathbb{R}^n$ . To ensure that the dictionary is overcomplete, we have  $k > n$ . The task is to find a representation in form a sparse code vector  $\alpha$  of  $k \times 1$ , such that:

$$\alpha = \min_{\alpha'} \|\alpha\|_0, \text{ s.t } x = D\alpha, \quad (2.3)$$

where  $\|x\|_0$  is l0 norm, which represents the number of non-zero components in  $x$ .

It has been found that this problem is hard to solve and thus we replace l0-norm with l1-norm, so then as others (Lee et al. [12] , Donoho [5]), the problem is reformulated to find  $\alpha$  by minimizing the objective function  $J$ :

$$J(\alpha; \lambda_1) = \min_{\alpha \in \mathbb{R}^k} \|x - D\alpha\|_2^2 + \lambda_1 \|\alpha\| \quad (2.4)$$

where the parameter  $\lambda_1$  is the regularization parameter which ensures a trade off in sparsity and reconstruction error. It also ensures that the sparse codes are not very large.

To further ensure that the dictionary values  $D$ , do not become arbitrarily large,[15] constrain the columns of  $D$  to have an  $l_2$ -norm less than or equal to one. They call the convex set of matrices verifying this constrain as :

For our purpose we use the Online dictionary learning implementation of [15]. This is a very efficient methods and works brilliantly for millions of training examples. Sparse coding is done using the orthogonal matching pursuit (OMP) method [20, 27] implemented in the SPAMS library provided by Mairal et al. [16]

## 2.4 Literature Review

In this chapter we aim to present a general overview of vessel segmentation methods. We particularly focus on approaches based on machine learning methods both supervised and unsupervised techniques. There has been a considerable work in the domain of automatic retinal vessel segmentation. Fraz et al. [7] give a complete review of retinal vessel segmentation methods. We briefly outline some of the methods before we move on to explain the most recent supervised learning based algorithms.

The earliest work on retinal segmentations are based on tracking methods to trace the blood vessels[2, 26]. In these methods the vessel is traced out from some starting seed points according to some relevant criterion. Recently, Vlachos and Dermatas [29] proposed a methods based on multi scale line tracking. They approach the problem by tracking the retinal vessel at multi scales and combining the individual image maps at different scales. Farnell et al. [6] employ multi scale line operators at various orientations to enhance the blood vessels.

There has been a recent surge in utilizing supervised machine learning based methods for vessel segmentation. In such methods, generally, each image pixel is represented as a feature vector computed using local or global information. A classifier is then trained to label every pixel to a vessel or background. Such classifiers rely on the training dataset, which is not available easily. Ricci and Perfetti [21] proposed a supervised classifier using line operators and using support vector classification. They obtain unsupervised pixel classification by thresholding the response of basic line detector. Additionally, they use orthogonal line detectors along with grey level of target pixel to construct feature vector for support vector machines based supervised classification. Staal et al. [25] proposed a ridge based vessel segmentation method used together with a supervised learning technique. Nguyen et al. [19] proposed a method utilizing gabor filter responses at multiple scales as features



for pixel classification.[9, 28] proposed neural network bases approach to vessel segmentation

We now look in detail some of the recent state-of-art work on retinal vessel segmentation.

Dollár and Zitnick [4] recently proposed a structured learning framework for predicting local edges utilizing random forests. The work exploits the presence of local structure forms like straight lines or T-junctions in image patches to learn an accurate edge detector. They train random forests in structured output spaces. Local image labels are highly interdependent and utilizing the knowledge on such local structures tend to improve the classifiers accuracy. The problem of edge detection is formulated in a way to predict local segmentation maps for input image patches. The structure labels are mapped to proxy discrete labels, in a way that similar structure labels are assigned to same discrete labels. Using this proxy mapping, existing random forest training approaches are used to learn structured random forests. Finally these forests are used to label each pixel to denote the presence or absence of edges.

Rigamonti and Lepetit [22] proposed a method to learn ad hoc features with learned filters. Many hand crafted features have been proposed to solve the problem of extracting linear structures like blood vessels, but are not always effective. In contrast, learned features are better at the task but are computationally expensive. The proposed algorithm complements handcrafted features with learned filters. Linear filters are computed from training images in a convolutional approach. The convolutional filters are constrained to be dissimilar. Further, feature maps are computed by convolving the learned filters with the images. Several such feature maps are computed both for handcrafted features like (OOF, EF) and for learned filters. These feature maps are then used to calculate descriptors at each image location. A random forest classifier is then trained to classify each image pixel as lying on a linear structure or background.

Ganin and Lempitsky [8] proposed a very elegant approach to natural edge detection or thin object segmentation. They employ a combination of convolutional neural networks with the nearest neighbor search. The problem is approached in a patch based framework, where the predictions are made patch by patch.



# Chapter 3

## Framework

In this chapter we present the basic framework of our vessel segmentation method. We start by setting up the various notation used throughout the chapter and give an overview about how the framework is applied. Next we discuss the two different models applied in the framework. This forms the core of this thesis and represents the major work done. At last we mention the various datasets utilized to test our models. In the next chapter we begin by running few experiments on the datasets to test the feasibility of models.

### 3.1 Vessel segmentation Framework

We start by defining the basic notations and methods. The input to our vessel segmentation framework is a training set composed of fundus images and their corresponding segmentation or label maps. The collection of fundus images is represented by  $\bar{I} = (I^{(1)}, I^{(2)}, \dots, I^{(n)})$  and their corresponding ground truth segmentation maps as  $\bar{S} = (S^{(1)}, S^{(2)}, \dots, S^{(n)})$ . Each given image is of size  $M \times N$ . For simplicity, in the rest of the section a single image would be referred as  $I$  and  $S$  and the image  $I$  is a single channel image.

An image can be represented as a collection of overlapping patches computed around every pixel at center. A patch is of size  $\sqrt{n} \times \sqrt{n}$  unless otherwise noted. We represent a patch centered at location  $(x, y)$  as  $p_{x,y}$ , and patches for  $I$  and  $S$  can be represented as  $Ip_{x,y}$  and  $Sp_{x,y}$  respectively. For simplicity we can refer the pixels of patch as  $\mathbb{P}_i$  for  $i = [1 \dots n]$ . A patch is represented as a vector, obtained by concatenating the image patch values.

Now we can represent our datasets as

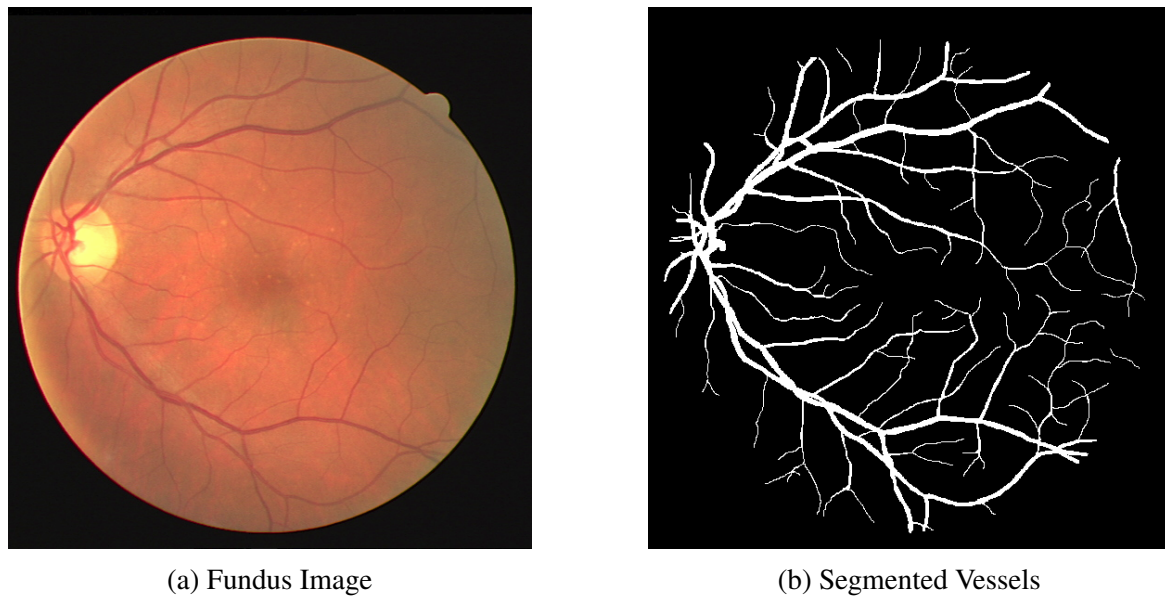


Fig. 3.1 Fundus Image and its segmentation

Fig 3.1 shows an example of fundus image on the left and the corresponding segmentation image on the right.

The aim of our thesis is to learn a function  $F$  which maps image  $I$  to its segmentation  $S$ , i.e, we want to learn a function to segment the vessel in a given fundus image. Inspired by works of [ref][ref] we approach the problem of prediction of segmentation maps in a patch based framework. This means that to predict the segmentation map for an image  $I$ , we decompose the image into overlapping patches and predict the segmentation maps for individual patches. The image and the segmentation map after decomposition can be represented as in Equation [].

$$F : I \rightarrow S$$

Here 'I' denotes the fundus image and 'S' its segmentation map.

At the test time, for a given image , we compute the patches and the mapping fucntion  $F$  is applied on all the test patches to obtain the segmentation. The output segmentation patches are combined by averaging ,thus resulting in an output segmented image. As the computed patches are overlapping, to each pixel several segmentation patches are added. More precisely, each pixel is composed of  $N$  pixel values.

In the next section we discuss how the mapping function  $F$  can be learnt. We present two models to predict the ground truth segmentation of a patch.

## 3.2 Learning

The major part of our work is to learn an effective mapping function  $F$ , to map the fundus image to its segmentation map. In our work, we learn a dictionary of representative image patches and a dictionary of segmentation patches. These dictionaries are then used to map the input image patch to its segmentation map. We explore two different methods to learn these dictionaries. Each image patch in the dictionary is denoted as an atom and our aim is to find the best set of atoms so as to approximately infer closest segmentation. We notice that these dictionaries resemble the local edge structure like junctions, straight lines, crossover etc.

Before we explore the methods we list the common operation performed on our dataset.

### Preprocessing

For our task we have explored some of the preprocessing steps including patch normalization, contrast stretching and local contrast normalization. For all the experiments, we normalize each input patch by subtracting the patch mean and dividing by the standard deviation per patch. Also, all the input patches are vectorized before being fed to the predictor.

### Rotations

To learn better image representations, we augmented the training set by adding the rotated pair of image and segmentation maps at an angle of 30, 60, 90, 120 and 150 degrees. Adding rotations in general would help to learn better cases where the vessels might appear in different orientations.

## 3.3 Cluster Learning

In our cluster learning framework, we learn an intermediate patch and segmentation maps. We then match our new patches to this dictionary and get the segmentation map from these images.

### 3.3.1 Learning

The training procedure for our method requires learning clusters within the input patches. Training is performed on the patches drawn from the input images  $I_1, I_2, \dots, I_n$  and corresponding segmentation images  $S_1, S_2, \dots, S_n$ .

From the training images, we compute a random subset of  $m$  patches of fixed size  $n \times n$  and their corresponding annotation as computed from the segmentation images. This subset of patches is used to further learn a dictionary of patches and their annotations. Let us denote a patch as  $x_i$  and its corresponding segmentation annotation patch as  $y_i$ . Then the new training set can be represented as:

$$X = \{x_1, x_2, \dots, x_n\}$$

$$Y = \{y_1, y_2, \dots, y_n\}$$

We then find 'k' clusters within the patches in  $X$  using K-Means clustering. Let the 'k' learnt cluster centers be represented as  $C = \{c_1, c_2, \dots, c_k\}$  and the  $k$  clusters can be represented as  $C_1, C_2, \dots, C_k$ . A cluster with 'm' patches then can be represented as

$$C_k = \{x_1^k, x_2^k, \dots, x_m^k\}$$

Now, since we know the annotation  $y_i$  for each training patch  $x_i$ , the image patches in each cluster can be replaced with their corresponding annotation patches to obtain the segmentation cluster as:

$$SC_k = \{y_1^k, y_2^k, \dots, y_m^k\}$$

We then find the representative point for the segmentation cluster by averaging all the patches belonging to it.

$$sc_k = \frac{1}{m} \sum_{i=1}^m y_m^k$$

Finally we create a dictionary mapping the cluster to their ground truth representatives as.

$$D = \{(C_k : gc_k) | k = 1..m\}$$

An example of dictionary learnt over image patches on DRIVE dataset is shown in figure[].

### 3.3.2 Prediction

At test time, the patches of an image are assigned to a cluster based on the learnt model. The desired annotation is then obtained by a lookup in the dictionary  $D$ .

### 3.3.3 Algorithm

The entire learning algorithm than can be summarized in following steps:

1. Extract a set of random Image-Segmentation patches of size  $m \times m$  from the training images.
2. Train a K-Means clustering model on the input image patches and learn  $k$  clusters on the input.
3. Compute the ground truth patch for each cluster by averaging the ground truth annotations for each patch belonging to a cluster.
4. Construct a dictionary of Cluster and annotations.

## 3.4 Dictionary Learning

Our method is based on learning dictionary of patches from training data. In our methods we learn two dictionaries each for representing the image patches and their segmentations. The methods in a way is very similar to the above method where we associate a learnt patch to its ground truth segmentation and use the dictionary for predicting new segmentations. The difference is in the way dictionary of patches is learnt and the way in which a patch is matched to the dictionary.

Our proposed method is based on learning a sparse dictionary of patches, of which a linear combination can be used to represent an image patch. For each such image patches, denoted as dictionary atoms, we find a corresponding segmentation label patch.

For example, given an overcomplete dictionary of atoms  $D \in R^{m,k}$ , with  $k$  columns as the learnt image atoms, we can reconstruct a given image patch  $x \in R^m$  as a linear combination of atoms in  $D$ . More formally, we learn a sparse code  $\alpha$  over  $D$  to represent  $x$  as

$$x \approx D\alpha$$

### 3.4.1 Learning

The learning phase starts by extracting image and label patches from the training data. Let us represent our training set composed of random patches as,

$$X = \{x_1, x_2, \dots, x_n\}$$

$$Y = \{y_1, y_2, \dots, y_n\}$$

where  $X$  is composed of image patches and  $Y$  the corresponding label patches. Each image patch can be represented as a vector by concatenating each of the pixel values. So give, 'm' patches of size  $\sqrt{n} \times \sqrt{n}$ , can be represented in form of matrices as  $X, Y \in R^{m,n}$ . Each row in the matrix denotes a patch or image atom.

We learn an overcomplete dictionary  $D$  of basis atoms from the image patches using the dictionary learning methods as described in[1]. The aim is to learn a dictionary  $D$  with  $k$  atoms which can best represent  $x_i$  using linear combination of a few atoms  $d_k$  of the learnt dictionary.

In the next step we learn a sparse code  $\alpha$  to represent our input  $X$  in terms of dictionary atoms.

$$X = \alpha D$$

Since we already know the ground truth annotation  $Y$  for  $X$ , we can infer that  $Y$  can be approximately represented as a linear combination of atoms from a dictionary of labels  $L$ , corresponding to  $D$ .

As the sparse code remains same for both the representations, we can compute the dictionary atoms from  $L$  by taking a weighted average of all the ground truth patches in  $Y$ , weighed by sparse codes for all the patches.

So we learn two dictionaries,  $D$  of image patches and  $L$  of the respective label patches for atoms in  $D$ .

### 3.4.2 Prediction

At the test time, given an image  $I$ , dense patches can be computed and the image can be represented as a matrix  $I \in R$ , where each row represents a patch of size  $n \times n$ . We then compute a sparse decomposition of  $I$  over dictionary  $D$  represented by a sparse code  $\alpha$  such that:

$$I = D\alpha$$



We can then compute the ground truth segmentation  $S$ , for the image patches in  $I$ , as :

$$S = L\alpha$$

where  $\alpha$  is the sparse code computed over dictionary  $D$  and  $L$  is the label dictionary previously computed during the learning phase.

## 3.5 Datasets

For testing the performance of our algorithm , we train and test our system on the following publically available datasets. In this section we describe the characteristics of these datasets

### 3.5.1 DRIVE

The Digital Retinal Images for Vessel Extraction (DRIVE) dataset consists of 40 color retinal images randomly selected from diabetic retinopathy screening program for 400 diabetic patients. Each of these images in dataset are JPEG compressed and have a dimensions of 768 x 584 pixels captured at a resolution on bits per pixel. The images are captures with a 45degree field of view (FOV). Of the 40 images in the dataset, 7 show sign of diabetic retinopathy , while the remaining 33 do not consist of any pathology. Each image is provided with a corresponding mask delineating the FOV.

The dataset is divided is provided with divisions in terms of training and testing set, with each set consisting of 20 images. Each of the 40 images have been manually segmented by human observers trained by an experienced optamologist. For the training set, single ground truth segmentation of the vessels is provided. The test set is provided with two ground truth segmentations, of which the first one is used as gold standard and the other is used to compare the performance with an independent human observer.

### 3.5.2 STARE

The STARE dataset consists of 20 images with blood vessel segmentations, out of which 10 show signs of pathology. The images have been capture with a FOV of 35degrees at 8 bit per pixel resolution, with dimesnsions of each image as 605 x 700 pixels. The dataset consists of segmentation provided by two human observers. In our experiements, we consider the

Table 3.1 Summary of Datasets

Datasets	FOV	# of Images
DRIVE	45	20+20
STARE	30	20
HRF	60	45
ARIA	50	212

segmentations provided by the first observer as ground truth.

For the experiments, the dataset is randomly divided into training and test sets each consisting of 10 images.

### 3.5.3 HRF

The High Resolution Fundus (HRF) image database consists of 45 images of which 15 images come from healthy patients, 15 from patients with diabetic retinopathy and 15 of glaucomatous patients. The images were captured with a FOV of 60 degrees, at a high resolution of 24 bits per pixel. The size of each image is 3504 x 2336 pixels and stored with JPEG compression.

Each image is provided a manual segmentation of vessels as segmented by three independent human observers trained by experienced ophthalmologists. The dataset also provides a corresponding mask image of each image delineating the FOV.

### 3.5.4 ARIA

The ARIA dataset consists of three groups of images. One of the groups consists of 92 images with age-related macular degeneration, the other with 59 images from diabetic patients and the last group with 61 images from a control group.

The images are captured with a 50-degree FOV, stored in uncompressed TIFF format, with a resolution of 8 bits per pixel. Each image has dimensions of 768 x 576 pixels. The dataset provides with blood vessel segmentation images as manually segmented by experts and a corresponding mask delineating the FOV region.

## Chapter 4

# Experimental Evaluation and Results

In this chapter we present the quantitative and qualitative evaluations of our learning algorithms as described in the last chapter. All the experiments are performed on the green channel of the image as it has the maximum contrast between the vessel and background as can be seen in Fig 4.1.

The performance of the proposed vessel segmentation algorithm is evaluated using the segmented vaculature considered as the gold standard and the manually marked segmentations by the human observer. All prior work with which we compare our method, have done a segmentation performance analysis with the manual segmentations provided by the first human observer. Additionally for some of the datasets we have a manual segmentation provided by the second observer which can be used to compare the automated methods to that of manual segmentations.

To assess the performance of segmentation methods, various metrics have been defined in literature [18, 23]. As reported in prior works, we also compute the performance of our vessel segmentation model defined as, true positives (TP): number of correctly classified vessel pixels; false positives (FP): number of pixels falsely classified as vessels; true negatives (TN): number of correctly classified non-vessel pixels; false negatives (FN): number of pixels falsely classified as non-vessels.

Using these metrics we can compute the accuracy (number of TP+TN / total pixels), pixel classification sensitivity and specificity. We also calculate the area under curve for receiver operation characteristics and also the AUC under the precision recall curve obtained by varying the threshold value for the segmented images.

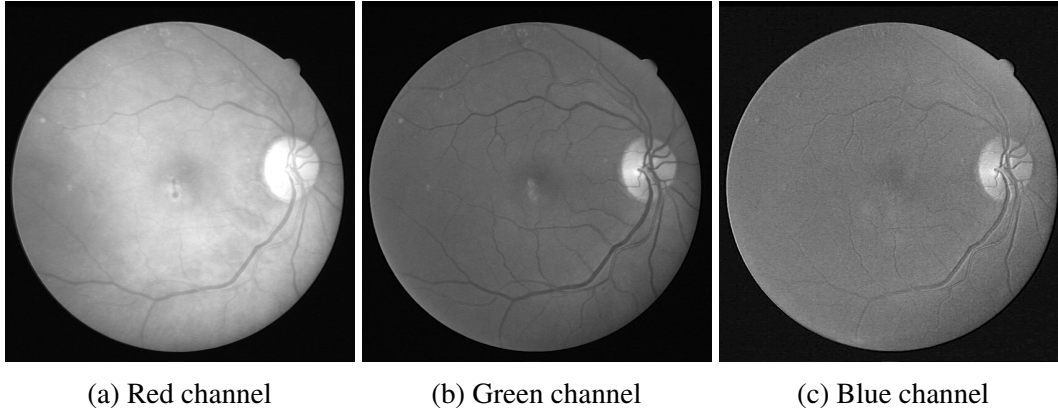


Fig. 4.1 Different channels of a fundus image showing variation in contrast of blood vessels against background. Green channel shows the maximum contrast in blood vessels and background.

For a complete assesment of our segmentation model, we perform various experiments. We start by performing the experiment on all the 4 datasets and compute the performance metrics as described above. Next, to test the generalization of our system we do cross training, i.e, training on images from one dataset and prediction on other.

We test the performance of both our models.

## 4.1 Vessel Segmenetation Assessment

We start by training our model on the DRIVE training dataset. The classifier is trained with a patch size of (10,10) and 1000 clusters. We also run the algorithm with the various preprocessing settings as in sections. Contrast enhancement of the images didn't have a major effect on our performance.

We applied our model with varying patch sizes ranging from (10,10) to (21,21). We note that with increase in patch size, we lose details on the thin vessels, but our confidence on thick vessels increases. We combine the results on various patch sizes by simple averaging of all the cases. This had a slight imporvement over our results.

Results for our model applied on DRIVE test dataset are shown in the table. As we obtain an average value at each pixel by merging the ground truth annotation coming from different overlapping patches, we can threshold the intensities to obtain the binar segmentation. ROC and PRC plots are made by varying the thresholds in steps of 0.02.

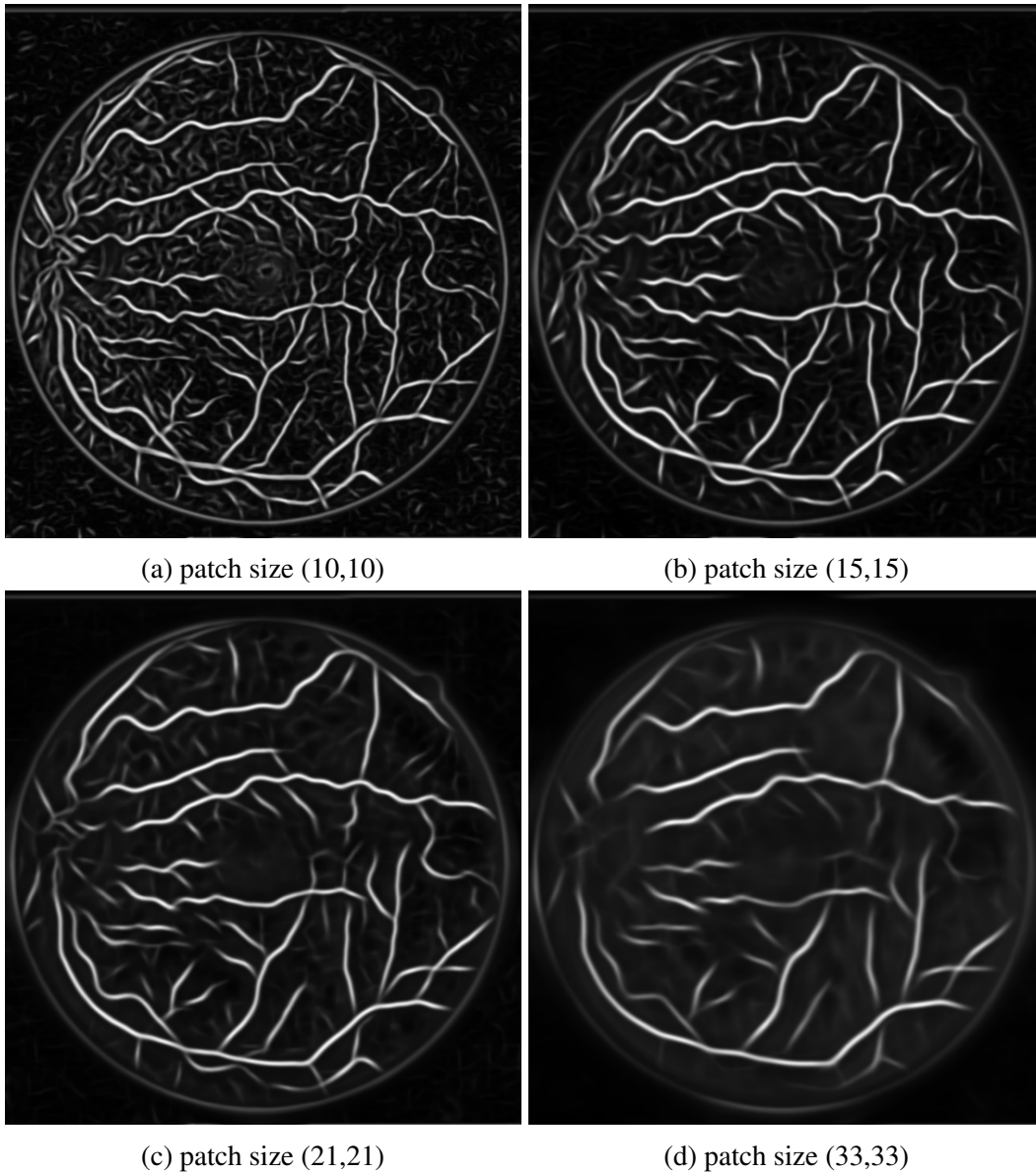


Fig. 4.2 Here we show the effect of varying patch size in our patch based framework. As we increase the patch size, we lose details on thin vessels, but our confidence on thick vessels increases.



## **Chapter 5**

### **Conclusion and Future Work**





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