# Algorithms for Finding Global Minimizers of Image Segmentation and Denoising Models

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#### **Abstract**

We show how certain non-convex minimization problems that arise in image processing and computer vision can equivalently be restated as convex minimization problems. This allows, in particular, finding global minimizers via standard convex minimization schemes.

#### 1 Introduction

Image denoising and segmentation are two related, fundamental problems of computer vision. The goal of denoising is to remove noise and/or spurious details from a given possibly corrupted digital picture while maintaining essential features such as edges. The goal of segmentation is to divide the given image into regions that belong to distinct objects in the depicted scene.

Variational and partial differential equations based approaches to denoising and segmentation have had great success. An important reason for their success is that these models are particularly well suited to imposing geometric constraints (such as regularity) on the solutions sought. Among the best known and most influential examples are the Rudin-Osher-Fatemi (ROF) total variation based image denoising model, and the Mumford-Shah image segmentation model.

Denoising models such as the ROF model can be easily adapted to different situations. An interesting scenario is denoising of shapes: Here, the given image is

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binary (representing the characteristic function of the given shape), and the noise is a perturbation in the geometry of the shape: Its boundary might be very rough, and the user might be interested in smoothing out its boundary, and perhaps removing small, unnecessary connected components of the shape. This task is a common first step in many object detection and recognition algorithms.

A common difficulty to many variational image processing models is that the energy functional to be minimized has local minima (which are not global minima). This is a much more serious drawback than non-uniqueness of global minimizers (which is also a common phenomenon) because local minima of segmentation and denoising models often have completely wrong levels of detail and scale: whereas global minimizers of a given model are usually all reasonable solutions, the local minima tend to be blatantly false. Many solution techniques for variational models are based on gradient descent, and are therefore prone to getting stuck in such local minima. This makes initial guess for gradient descent based algorithms sometimes critically important for obtaining satisfactory results.

In this paper we propose algorithms which are guaranteed to find global minimizers of certain denoising and segmentation models that are known to have local minima. As a common feature, the models we consider involve minimizing functionals over characteristic functions of sets, which is a non-convex collection; this feature is responsible for the presence of local minima. Our approach, which is based on observations of Strang in [22, 23], is to extend the functionals and their minimization to all functions in such a way that the minimizers of the extended functionals can be subsequently transformed into minimizers for the original models by simple thresholding. This allows, among other things, computing global minimizers for the original non-convex variational models by carrying out standard convex minimization schemes.

Our first example is binary image denoising, where the given noisy image for the ROF model is taken to be binary, and the solution is also sought among binary images. This problem has many applications where smoothing of geometric shapes is relevant. Some examples are the denoising of text documents, and the fairing of surfaces in computer graphics. Because the space of binary functions is nonconvex, the minimization problem involved is actually harder than minimizing the original ROF model. In Section 2, we show how this problem can be written very simply as a convex optimization problem.

In Section 3 we generalize the observation of Section 2 to the two-phase, piecewise constant Mumford-Shah model. The formulation we obtain turns out to be very closely related to the segmentation algorithm of Chan and Vese [7, 8]. Our observations indicate why this algorithm is successful in finding interior contours and other hard-to-get features in images.

#### 2 Previous Work

The results and the approach of this paper follow very closely observations of Strang in [22, 23]. In those papers, optimization problems of the following form, among others, are studied:

$$\inf_{\{u:\int fu\,dx=1\}}\int |\nabla u|\tag{1}$$

where f(x) is a given function. It is shown in particular that the minimizers of (1) turn out to be characteristic functions of sets. The main idea involved is to express the functional to be minimized and the constraint in terms of the level sets of the functions u(x) and f(x). The coarea formula of Felming and Rishel [12] is the primary tool.

In this paper, the idea of expressing the functionals in terms of level sets is applied to some simple image processing models. For instance, in Section 4 where we study the piecewise constant Mumford-Shah energy, the relevant energy can be expressed in a way that is almost the same as (1).

We should point out that our emphasis in this paper is in some sense opposite that of [22, 23]. Indeed, in those works the main point is that some energies of interest that need to be minimized over all functions turn out to have minimizers that take only two values. In our case, we start with a variational problem that is to be minimized over only functions that take two values (i.e. characteristic functions of sets), but show that we may instead minimize over all functions (that are allowed to take intermediate vales), i.e. we may ignore the non-convex constraint. This allows us to end up with a convex formulation of the original non-convex problem.

## **3** The ROF Model for Binary Images

Rudin, Osher, and Fatemi's total variation based image denoising model is one of the best known and successful of PDE based image denoising models. Indeed, being convex it is one of the simplest denoising techniques that has the all important edge preserving property.

Let  $D \subset \mathbf{R}^N$  denote the image domain, which will be taken to be bounded and with Lipschitz boundary. Usually, D is simply a rectangle, modelling the computer screen. For convenience, in this section we will take D to be the entire space  $\mathbf{R}^N$ . Let  $f(x): \mathbf{R}^N \to [0,1]$  denote the given (grayscale) possibly corrupted (noisy) image. The energy to be minimized in the standard ROF model is then given by:

$$E_2(u,\lambda) = \int_{\mathbf{R}^N} |\nabla u| + \lambda \int_{\mathbf{R}^N} \left( u(x) - f(x) \right)^2 dx \tag{2}$$

where  $\lambda > 0$  is a parameter to be chosen by the user (or is estimated from the level of noise, in case this information is available).

An interesting application of the ROF model described above is to binary image denoising. This situation arises when the given image f(x) is binary (i.e.  $f(x) \in \{0,1\}$  for all  $x \in \mathbf{R}^N$ ) and is known to be the corrupted version of another binary image  $u: \mathbf{R}^N \to \{0,1\}$  that needs to be estimated. Naturally, f(x) can then be expressed as

$$f(x) = \mathbf{1}_{\Omega}(x)$$

where  $\Omega$  is an arbitrary bounded measurable subset of  $\mathbf{R}^N$ . In this case, the noise is in the *geometry*; for example, the boundary  $\partial\Omega$  of  $\Omega$  might have spurious oscillations, or  $\Omega$  might have small connected components (due to presence of noise) that need to be eliminated. The ROF model (2) can be specialized to this scenario by restricting the unknown u(x) to have the form  $u(x) = \mathbf{1}_{\Sigma}(x)$ , where  $\Sigma$  is a subset of  $\mathbf{R}^N$ . One then obtains the following optimization problem:

$$\min_{\substack{\Sigma \subset \mathbf{R}^N \\ u(x) = \mathbf{1}_{\Sigma}(x)}} \int_{\mathbf{R}^N} |\nabla u| + \lambda \int_{\mathbf{R}^N} \left( u(x) - \mathbf{1}_{\Omega}(x) \right)^2 dx \tag{3}$$

Problem (3) is non-convex because the minimization is carried out over a non-convex set of functions. Recalling that the total variation of the characteristic function of a set is its perimeter, and noticing that the fidelity term in this case simplifies, we write (3) as the following geometry problem:

$$\min_{\Sigma \subset \mathbf{R}^N} \operatorname{Per}(\Sigma) + \lambda |\Sigma \triangle \Omega| \tag{4}$$

where  $Per(\cdot)$  denotes the perimeter,  $|\cdot|$  is the N-dimensional Lebesgue measure, and  $S_1 \triangle S_2$  denotes the symmetric difference between the two sets  $S_1$  and  $S_2$ .

Usual techniques for approximating the solution. A very successful method of solving problems of the type (4) has bee via some *curve evolution* process, sometimes referred to as *active contours*. Indeed, the unknown set  $\Sigma$  can be described by its boundary  $\partial \Sigma$ . The boundary  $\partial \Sigma$  is then updated iteratively, usually according to gradient flow for the energy involved.

Numerically, there are several ways of representing  $\partial \Sigma$ . For the applications mentioned above, explicit curve representations as in Kaas, Witkin, Terzopoulos [15] are not appropriate, since such methods do not allow changes in curve topology (and have a number of other drawbacks). Instead, the most successful algorithms are those based on either the level set method of Osher and Sethian [20], or on the variational approximation approach known as Gamma convergence theory [9].

In the level set formulation, the unknown boundary  $\partial \Sigma$  is represented as the 0-level set of a (Lipschitz) function  $\phi : \mathbf{R}^N \to \mathbf{R}$ :

$$\Sigma = \left\{ x \in \mathbf{R}^N : \phi(x) > 0 \right\}$$

so that  $\partial \Sigma = \{x \in \mathbf{R}^N : \phi(x) = 0\}$ . The functional to be minimized in (3), which we called  $E_2(\cdot, \lambda)$ , can then be expressed in terms of the function  $\phi(x)$  as follows:

$$\int_{\mathbf{R}^N} |\nabla H(\phi(x))| \, dx + \lambda \int_{\mathbf{R}^N} \left( H(\phi(x)) - \mathbf{1}_{\Omega}(x) \right)^2 dx \tag{5}$$

Here, the function  $H(x): \mathbf{R} \to \mathbf{R}$  is the Heaviside function:

$$H(\xi) = \begin{cases} 0 & \text{if } x < 0, \\ 1 & \text{if } x \ge 0. \end{cases}$$

In practice, one takes a smooth (or at least Lipschitz) approximation to H(x), which we shall call  $H_{\varepsilon}(\xi)$ , where  $H_{\varepsilon}(\xi) \to H(\xi)$  in some manner as  $\varepsilon \to 0$ .

The Euler-Lagrange equation for (5) is easy to obtain. It leads to the following gradient flow:

$$\phi_t(x,t) = H_{\varepsilon}'(\phi) \left\{ \operatorname{div} \left( \frac{\nabla \phi}{|\nabla \phi|} \right) + 2\lambda \left( \mathbf{1}_{\Omega}(x) - H_{\varepsilon}(\phi) \right) \right\}.$$
 (6)

When equation (6) is simulated using reinitialization for the level set function  $\phi(x)$  and a compactly supported approximation  $H_{\varepsilon}(x)$  to H(x), it is observed to define a continuous evolution (with respect to, say, the  $L^1$ -norm) for the unknown function  $u(x) = \mathbf{1}_{\Sigma}(x)$  and decreases the objective energy (3) through binary images. It is analogous to the gradient descent equation in [7], which is natural since (3) is the restriction to binary images of also the energy considered in that work, namely the two phase, piecewise constant Mumford-Shah segmentation energy. In Section 4 we will consider this energy for general (not necessarily binary) images.

Another representation technique for the unknown set  $\Sigma$  in (4) is, as we mentioned, based on the Gamma convergence ideas. Here, the given energy is replaced by a sequence of approximate energies with a small parameter  $\varepsilon>0$  in them. The sequence converges to the original energy as  $\varepsilon\to0$ . The approximations have the form:

$$E_{\varepsilon}(u,\lambda) = \int_{\mathbf{R}^N} \varepsilon |\nabla u|^2 + \frac{1}{\varepsilon} W(u) + \lambda \left\{ u^2 (c_1 - f)^2 + (1 - u)^2 (c_2 - f)^2 \right\} dx$$

In this energy,  $W(\xi)$  is a double-well potential with equidepth wells at 0 and 1; for instance, a simple choice is  $W(\xi) = \xi^2 (1 - \xi)^2$ . The term  $\frac{1}{\varepsilon} W(u)$  can be thought

of as a penalty term that forces the function u to look like the characteristic function of a set: u is forced to be approximately 0 or 1 on most of  $\mathbf{R}^N$ . The term  $\varepsilon |\nabla u|^2$ , on the other hand, puts a penalty on the transitions of u between 0 and 1. Taken together, it turns out that these terms both impose the constraint that u should be a characteristic function, and approximate its total variation. Precise versions of these statements have been proved in [16]. The remaining terms in  $E_\varepsilon$  are simply the fidelity term written in terms of u.

We now argue by the help of a very simple example that these techniques will get stuck in local minima in general, possibly leading to resultant images with the wrong level of detail. This fact is already quite familiar to researchers working with these techniques from practical numerical experience.

**Example:** Consider the two dimensional case, where the observed binary image f(x) to be denoised is the characteristic function of a ball  $B_R(0)$  of radius R, which is centered at the origin. In other words, we take  $\Omega = B_R(0)$ . Implementing the gradient descent algorithm defined by (6) requires the choice of an initial guess for the interface  $\phi(x)$  (or, equivalently, an initial guess for the set  $\Sigma$  that is represented by  $\phi(x)$ ). A common choice in practical applications is to take the observed image itself as the initial guess. In our case, that means we initialize with  $\Sigma = B_R(0)$ .

Now, one can see without much trouble that the evolution defined by (6) will maintain radial symmetry of  $\phi(x)$ . That means, at any given time  $t \geq 0$ , the set (i.e. the candidate for minimization) represented by  $\phi(x)$  is of the form

$$\left\{x \in \mathbf{R}^2 : \phi(x) > 0\right\} = B_r(0)$$

for some choice of the radius  $r \ge 0$ . We can write the energy of  $u(x) = \mathbf{1}_{B_r(0)}(x)$  in terms of r, as follows:

$$E(r) := E_2(\mathbf{1}_{B_r(0)}(x), \lambda) = 2\pi r + \lambda \pi |R^2 - r^2|$$

A simple calculation shows that if  $\lambda < \frac{2}{R}$ , then the minimum of this function is at r = 0. Hence, if we fix  $\lambda > 0$ , then the denoising model prefers to remove disks of radius smaller than the critical value  $\frac{2}{R}$ .

But now, once again an easy calculation shows that if  $R>\frac{1}{\lambda}$ , then E(r) has a local maximum at  $r_{max}(\lambda)=\frac{1}{\lambda}$ . See Figure 1 for the plot of E(r) in such a case. Thus the energy minimization procedure described by (6) cannot shrink disks of radius  $R\in(\frac{1}{\lambda},\frac{2}{\lambda})$  to a point, even though the global minimum of the energy for an original image given by such a disk is at  $u(x)\equiv 0$ .

We can easily say a bit more: There exists  $\delta>0$  such that if  $\Sigma\subset\mathbf{R}^N$  satisfies  $|\Sigma\ \triangle\ B_R(0)|<\delta$  then  $E_2(\mathbf{1}_\Sigma(x),\lambda)>E_2(\mathbf{1}_{B_R(0)}(x),\lambda)$ . In words, all binary images nearby, but not identical with, the observed image  $\mathbf{1}_{B_R(0)}(x)$  have strictly

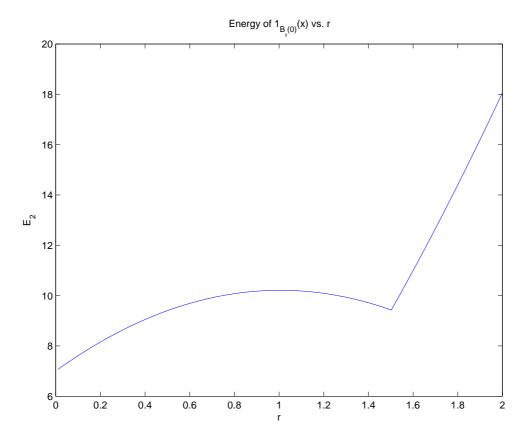


Figure 1: Energy (3) of  $u(x)=\mathbf{1}_{B_r(0)}(x)$  as a function of  $r\in[0,2]$  when the observed image is given by  $f(x)=\mathbf{1}_{B_R(0)}(x)$ . Here,  $R=\frac{3}{2}$  and the parameter  $\lambda$  was chosen to be  $\lambda=1$ . There is clearly a local minimum, corresponding to  $r=R=\frac{3}{2}$ .

higher energy. This can be seen simply by noting that the energy of any region that is *not* a disk is strictly larger than the energy of the disk having the same area as the given region and its center at the origin.

To summarize: If  $f(x) = \mathbf{1}_{B_R(0)}(x)$  with  $R \in (\frac{1}{\lambda}, \frac{2}{\lambda})$ , and if the initial guess for the continuous curve evolution based minimization procedure (6) is taken to be the observed image f(x) itself, then the procedure gets stuck in the local minimizer u(x) = f(x). The unique global minimizer is actually  $u(x) \equiv 0$ .

**Proposed Method for Finding the Global Minimum:** We now turn to an alternative way of carrying out the constrained, non-convex minimization problem (3)

that is guaranteed to yield a global minimum.

The crux of our approach is to consider minimization of the following convex energy, defined for any given observed image  $f(x) \in L^1(\mathbf{R}^N)$  and  $\lambda \geq 0$ :

$$E_1(u(x),\lambda) := \int_{\mathbf{R}^N} |\nabla u| + \lambda \int_{\mathbf{R}^N} |u(x) - f(x)| \, dx. \tag{7}$$

This energy differs from the standard ROF model only in the fidelity term: The  $L^2$ -norm square of the original model is replaced by the  $L^1$ -norm as a measure of fidelity. It was previously introduced and studied in signal and image processing applications in [1, 2, 3, 17, 18, 6]. This variant of the ROF model has many interesting properties and uses; the point we'd like to make in this section is that it also turns out to solve our geometry denoising problem (4).

First, let us state the obvious fact that energies (2) and (7) agree on binary images (i.e. when both u and f are characteristic functions of sets). On the other hand, energy (7) is convex, but unlike energy (7), it is not strictly so. Accordingly, its global minimizers are not unique in general. Nevertheless, being convex, it does not have any local minima that are not global minima, unlike the constrained minimization (3). We therefore adopt the following notation: For any  $\lambda \geq 0$ , we let  $M(\lambda)$  denote the set of all minimizers of  $E_1(\cdot, \lambda)$ . It is easy to show that for each  $\lambda \geq 0$  the set  $M(\lambda)$  is non-empty, closed, and convex.

The relevance of energy (7) for our purposes is established in Theorem 1, which is taken from [6], where additional geometric properties of it are noted. The proof is based on the following proposition that expresses energy (7) in terms of the level sets of u and f. We include its short proof for completeness.

**Proposition 1** The energy  $E_1(u, \lambda)$  can be rewritten as follows:

$$E_{1}(u,\lambda) = \int_{-\infty}^{\infty} Per\left(\{x : u(x) > \mu\}\right) + \left|\{x : u(x) > \mu\} \triangle \{x : f(x) > \mu\}\right| d\mu \quad (8)$$

**Proof:** As it is done in [23], we try to write everything in terms of level sets. Recall the coarea formula for functions of bounded variation [12, 13, 11]):

$$\int_{\mathbf{R}^N} |\nabla u| = \int_{-\infty}^{\infty} \Pr(\{x : u(x) > \mu\}) d\mu \tag{9}$$

Also, there is the following "layer cake" formula:

$$\begin{split} \int_{\mathbf{R}^N} |u - f| \, dx &= \int_{\{u > f\}} |u - f| \, dx + \int_{\{f > u\}} |u - f| \, dx \\ &= \int_{\{u > f\}} \int_{f(x)}^{u(x)} d\mu \, dx + \int_{\{f > u\}} \int_{u(x)}^{f(x)} d\mu \, dx \\ &= \int_{\mathbf{R}^N} \int_{\mathbf{R}} \mathbf{1}_{\{u > f\}}(x) \mathbf{1}_{[f(x), u(x))}(\mu) + \mathbf{1}_{\{f > u\}}(x) \mathbf{1}_{[u(x), f(x))}(\mu) \, d\mu \, dx \\ &= \int_{\mathbf{R}} \int_{\mathbf{R}^N} \mathbf{1}_{\{u > f\}}(x) \mathbf{1}_{[f(x), u(x))}(\mu) + \mathbf{1}_{\{f > u\}}(x) \mathbf{1}_{[u(x), f(x))}(\mu) \, dx \, d\mu \end{split}$$

where we simply changed the order of integration in the last step. But now we have:

$$\mathbf{1}_{\{u>f\}}(x)\mathbf{1}_{(f(x),u(x))}(\mu) = 1 \text{ iff } x \in \{u>f\} \cap \{u>\mu\} \cap \{f>\mu\}^c$$

and 0 otherwise, and

$$\mathbf{1}_{\{f>u\}}(x)\mathbf{1}_{(u(x),f(x))}(\mu) = 1 \text{ iff } x \in \{f>u\} \cap \{u>\mu\}^c \cap \{f>\mu\}$$

and 0 otherwise. That means

$$\mathbf{1}_{\{u>f\}}(x)\mathbf{1}_{(f(x),u(x))}(\mu) + \mathbf{1}_{\{f>u\}}(x)\mathbf{1}_{(u(x),f(x))}(\mu) = \mathbf{1}_{\{u>\mu\}\triangle\{f>\mu\}}(x)$$

Therefore

$$\int_{\mathbf{R}^{N}} |u - f| \, dx = \int_{-\infty}^{\infty} \left| \{x : u(x) > \mu\} \, \Delta \, \{x : f(x) > \mu\} \right| \, d\mu$$

Putting these formulae together gives the one in the statement of the claim.  $\Box$ 

We can now state the theorem that forms the basis of our entire approach. It is also taken directly from [6]:

**Theorem 1** If the observed image f(x) is the characteristic function of a bounded domain  $\Omega \subset \mathbf{R}^N$ , then for any  $\lambda \geq 0$  there is a minimizer of  $E_1(\cdot, \lambda)$  that is also the characteristic function of a (possibly different) domain. In other words, when the observed image is binary, then for each  $\lambda \geq 0$  there is at least one  $u(x) \in M(\lambda)$  which is also binary.

In fact, if  $u_{\lambda}(x) \in M(\lambda)$  is any minimizer of  $E_1(\cdot, \lambda)$ , then for almost every  $\mu \in [0, 1]$  we have that the binary function

$$\mathbf{1}_{\{x:u_{\lambda}(x)>\mu\}}(x)$$

is also a minimizer of  $E_1(\cdot, \lambda)$ .

**Proof:** The proof is based on the observation that once energy (7) is written as (8), it can be minimized pointwise in  $\mu$ . Moreover, this pointwise problem is precisely the geometry problem (4). The conclusion now follows easily. See [6] for details.

Let us state how the theorem above helps us solve (3) in the form of a corollary. It follows immediately from the theorem:

**Corollary 1** To find a solution (i.e. a global minimizer) u(x) of the non-convex variational problem (3), it is sufficient to carry out the following three steps:

- 1. Find any minimizer of the **convex** energy (7); call it v(x).
- 2. Let  $\Sigma = \{x \in \mathbf{R}^N : v(x) > \mu\}$  for some  $\mu \in (0,1)$ .
- 3. Set  $u(x) = \mathbf{1}_{\Sigma}(x)$ .

Then u(x) is a global minimizer of (3) for almost every choice of  $\mu$ .

The most involved step in the solution procedure described in Corollary 1 is finding a minimizer of (7). One can approach this problem in many ways; for instance, one possibility is to simply carry out gradient descent.

Numerical Exmaple: The synthetic image of Figure 2 represents the given binary image f(x), which is a simple geometric shape covered with random (binary) noise. The initial guess was an image composed of all 1's (an all white image). In the computation, the parameter  $\lambda$  was chosen to be quite moderate, so that in particular the small circular holes in the shape should be removed while the larger one should be kept. The result of the minimization is shown in Figure 3; in this case the minimizer is automatically very close to being binary, and hence the thresholding step of the algorithm in Corollary 1 is almost unnecessary.

Figure 4 shows the histograms of intermediate steps during the gradient descent based minimization. As can be seen, the intermediate steps themselves are very far from being binary. The histogram in the lower right hand corner belongs to the final result shown in Figure 3. Thus the gradient flow goes through non-binary images, but in the end reaches another binary one. Although this is not implied by Theorem 1, it seems to hold in practice.

### 4 Piecewise Constant Segmentation

We extend the discussion of Section 3 to the two-phase, piecewise constant Mumford-Shah segmentation model. Unlike in the previous section, this time we let the corrupted image f(x) be non-binary; it is assumed to be some measurable function

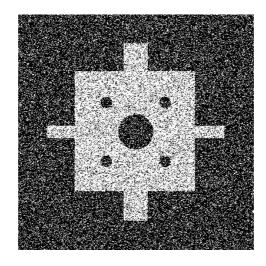
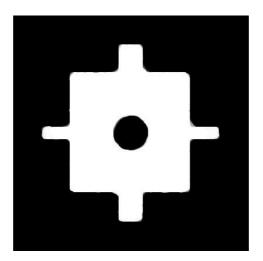


Figure 2: Original binary image.



 $Figure \ 3: \ Final \ result \ found \ (no \ need \ to \ threshold \ in \ this \ case).$ 

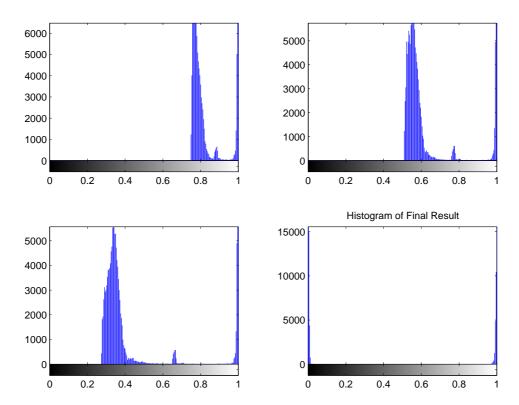


Figure 4: Histograms for intermediate images as the gradient descent proceeds. As can be seen, the intermediate images themselves are not binary; however, by the time the evolution reaches steady state, we are back to a binary image.

that takes its values in the unit interval. Thus, the discussion of this section supersedes that of the previous. Also, from now on we will assume that the image domain D is a bounded subset of  $\mathbf{R}^N$  with Lipschitz boundary. The segmentation energy, which we will call MS, can then be written as:

$$MS(\Sigma, c_1, c_2) := \operatorname{Per}(\Sigma; D) + \lambda \int_{\Sigma} (c_1 - f(x))^2 dx + \lambda \int_{D \setminus \Sigma} (c_2 - f(x))^2 dx.$$
(10)

The model says we should solve:

$$\min_{\substack{c_1, c_2 \in \mathbf{R} \\ \Sigma \subset D}} MS(\Sigma, c_1, c_2).$$
(11)

This optimization problem can be interpreted to be looking for the best approximation in the  $L^2$  sense to the given image f(x) among all functions that take only two values. These values, denoted  $c_1,c_2$ , and where each is taken, namely  $\Sigma$  and  $D\setminus \Sigma$ , are unknowns of the problem. As before, there is a penalty on the geometric complexity of the interface  $\partial \Sigma$  that separates the regions where the two values  $c_1$  and  $c_2$  are taken. Functional (10) is non-convex, and can have more than one minimizer. Existence of at least one minimizer follows easily from standard arguments. Notice that if  $\Sigma$  is fixed, the values of  $c_1$  and  $c_2$  that minimize  $MS(\Sigma,\cdot,\cdot)$  read  $c_1=\int_{\Sigma}f(x)dx/|\Sigma|$  and  $c_2=\int_{D\setminus\Sigma}f(x)dx/|D\setminus\Sigma|$ . A natural way to approximate the solution is a two-step scheme where at the first step one computes  $c_1$  and  $c_2$  according to these formulae, and at the second step updates the shape  $\Sigma$ . Even the minimization of  $MS(\cdot,c_1,c_2)$  is a difficult problem since this functional is non-convex. In what follows we focus on the minimization of  $MS(\cdot,c_1,c_2)$ .

**Chan-Vese Algorithm:** In [7] Chan and Vese (CV) proposed a level set based algorithm for solving the optimization problem (11). The idea is to represent the boundary  $\partial \Sigma$  with the 0-level set of the function  $\phi: D \to \mathbf{R}^N$ . Energy (10) can then be written in terms of the level set function  $\phi$ ; it turns out to be:

$$CV(\phi, c_1, c_2) = \int_D |\nabla H_{\varepsilon}(\phi)|$$

$$+ \lambda \int_D H_{\varepsilon}(\phi)(c_1 - f(x))^2 + (1 - H_{\varepsilon}(\phi))(c_2 - f(x))^2 dx. \quad (12)$$

The function  $H_{\varepsilon}$  is, as before, a regularization of the Heaviside function. The precise choice of the regularization  $H_{\varepsilon}$  of H is a crucial ingredient of the CV algorithm. We will return to this topic.

Variations of energy (12) with respect to the level set function  $\phi$  lead to the following gradient descent scheme:

$$\phi_t = H_{\varepsilon}'(\phi) \left\{ \operatorname{div} \left( \frac{\nabla \phi}{|\nabla \phi|} \right) - \lambda \left( (c_1 - f(x))^2 - (c_2 - f(x))^2 \right) \right\}.$$

The optimal choice for the constants  $c_1, c_2$  is determined in the way described above.

**The Proposed Algorithm:** The CV algorithm chooses a non-compactly supported, smooth approximation  $H_{\varepsilon}$  for H. As a result, the gradient descent equation given above and the following one have the same stationary solutions:

$$\phi_t = \operatorname{div}\left(\frac{\nabla\phi}{|\nabla\phi|}\right) - \lambda\left((c_1 - f(x))^2 - (c_2 - f(x))^2\right)$$

where we simply omitted the approximate Heaviside function altogether. This equation, in turn, is gradient descent for the following energy:

$$\int_{D} |\nabla \phi| + \lambda \int_{D} \left( (c_1 - f(x))^2 - (c_2 - f(x))^2 \right) \phi \, dx. \tag{13}$$

This energy is homogeneous of degree 1 in  $\phi$ . As a result, it does not have a minimizer in general. In other words, the gradient descent written above does not have a stationary state: If the evolution is carried out for a long time, the level set function  $\phi$  would tend to  $+\infty$  wherever it is positive, and to  $-\infty$  wherever it is negative. This issue is related to the nonuniqueness of representation with level sets, and is easy to fix: one can simply restrict minimization to  $\phi$  such that  $0 \le \phi(x) \le 1$  for all  $x \in D$ . With this fix, we arrive at the following statement:

**Theorem 2** For any given fixed  $c_1, c_2 \in \mathbf{R}$ , a global minimizer for  $MS(\cdot, c_1, c_2)$  can be found by carrying out the following convex minimization:

$$\min_{0 \le u \le 1} \underbrace{\int_{D} |\nabla u| + \lambda \int_{D} \left\{ (c_1 - f(x))^2 - (c_2 - f(x))^2 \right\} u(x) \, dx}_{:= \tilde{E}(u, c_1, c_2)}.$$

and then setting  $\Sigma = \{x : u(x) \ge \mu\}$  for a.e.  $\mu \in [0, 1]$ .

**Proof:** We once again rely on the coarea formula; since u takes its values in [0,1], we have:

$$\int_{D} |\nabla u| = \int_{0}^{1} \text{Per}(\{x : u(x) > \mu\}; D) \, d\mu.$$

For the other terms that constitute the fidelity term, we proceed as follows:

$$\int_{D} (c_{1} - f(x))^{2} u(x) dx = \int_{D} (c_{1} - f(x))^{2} \int_{0}^{1} \mathbf{1}_{[0, u(x)]}(\mu) d\mu dx$$

$$= \int_{0}^{1} \int_{D} (c_{1} - f(x))^{2} \mathbf{1}_{[0, u(x)]}(\mu) dx d\mu$$

$$= \int_{0}^{1} \int_{D \cap \{x: u(x) > \mu\}} (c_{1} - f(x))^{2} dx d\mu.$$

Also, we have:

$$\int_{D} (c_2 - f(x))^2 u(x) dx = \int_{0}^{1} \int_{D \cap \{x : u(x) > \mu\}} (c_2 - f(x))^2 dx d\mu$$
$$= C - \int_{0}^{1} \int_{D \cap \{x : u(x) > \mu\}^c} (c_2 - f(x))^2 dx d\mu.$$

where  $C=\int_D (c_2-f)^2\,dx$  is independent of u. Putting it all together, and setting  $\Sigma(\mu):=\{x:u(x)>\mu\}$ , we get the following formula that is valid for any  $u(x)\in L^2(D)$  such that  $0\leq u(x)\leq 1$  for a.e.  $x\in D$ :

$$\tilde{E}(u, c_1, c_2) = \int_0^1 \left\{ \operatorname{Per}(\Sigma(\mu); D) + \lambda \int_{\Sigma(\mu)} (c_1 - f(x))^2 dx + \lambda \int_{D \setminus \Sigma(\mu)} (c_2 - f(x))^2 dx \right\} d\mu - C$$

$$= \int_0^1 MS(\Sigma(\mu), c_1, c_2) d\mu - C.$$

It follows that if u(x) is a minimizer of the convex problem, then for a.e.  $\mu \in [0,1]$  the set  $\Sigma(\mu)$  has to be a minimizer of the original functional  $MS(\cdot,c_1,c_2)$ .  $\square$ 

**Remark:** The optimization problem that forms the content of the Theorem 2 can be interpreted as follows: The level set formulation of the two-phase model depends on the level set function  $\phi$  only through the term  $H(\phi)$ . The term  $H(\phi)$  represents a parametrization of binary functions (since, for any given function  $\phi$ , the function  $H(\phi)$  is binary). So the minimization of (12) is thus a minimization over binary functions. Minimization of (13) on the other hand, corresponds to removing the non-convex constraint of being binary; instead we minimize over functions that are allowed to take intermediate values. The content of the theorem above is that the minimizers (essentially) automatically satisfy the more stringent constraint.

We now turn to the question of how to minimize the convex problem stated in the theorem. In that connection, we have the following claim: **Claim 1** Let  $s(x) \in L^{\infty}(D)$ . Then the following convex, constrained minimization problem

$$\min_{0 \leq u \leq 1} \int_D |\nabla u| + \lambda \int_D s(x) u \, dx$$

has the same set of minimizers as the following convex, unconstrained minimization problem:

$$\min_{u} \int_{D} |\nabla u| + \int_{D} \alpha \nu(u) + \lambda s(x) u \, dx$$

where  $\nu(\xi) := \max\{0, 2|\xi - \frac{1}{2}| - 1\}$ , provided that  $\alpha > \frac{\lambda}{2} ||s(x)||_{L^{\infty}(D)}$ .

**Proof:** The term  $\alpha\nu(u)$  that appears in the second, unconstrained minimization problem given in the claim is an *exact penalty* term [14]; see Figure 5 for a plot of its graph. Indeed, the two energies agree for  $\{u\in L^\infty(D): 0\leq u(x)\leq 1\forall x\}$ . So we only need to show that any minimizer of the unconstrained problem automatically satisfies the constraint  $0\leq u\leq 1$ . This is immediate: If  $\alpha>\frac{\lambda}{2}\|s(x)\|_{L^\infty}$ , then

$$|\lambda s(x)| \max\{|u(x)|, |u(x) - 1|\} < \alpha \nu(u(x)) \text{ whenever } u(x) \in [0, 1]^c,$$

which means that the transformation  $u \to \min\{\max\{0,u\},1\}$  always decreases the energy of the unconstrained problem (strictly if  $u(x) \in [0,1]^c$  on a set of positive measure). That leads to the desired conclusion.  $\square$ 

**Numerical Examples:** Here we detail how we obtained the numerical results pertaining to the two phase piecewise constant segmentation model that are presented in this paper. Given  $c_1, c_2$ , the "exact penalty" formulation of the equivalent minimization problem described above leads to the following Euler-Lagrange equation:

$$\operatorname{div}\left(\frac{\nabla u}{|\nabla u|}\right) - \lambda s(x) - \alpha \nu'(u) = 0$$

where  $s(x) = (c_1 - f(x))^2 - (c_2 - f(x))^2$ . The following explicit gradient descent scheme was used to solve the last equation:

$$\frac{u^{n+1} - u^n}{\delta t} = D_x^- \left( \frac{D_x^+ u^n}{\sqrt{(D_x^+ u^n)^2 + (D_y^+ u^n)^2 + \varepsilon_1}} \right) + D_y^- \left( \frac{D_y^+ u^n}{\sqrt{(D_x^+ u^n)^2 + (D_y^+ u^n)^2 + \varepsilon_1}} \right) - \lambda s(x) - \alpha \nu_{\varepsilon_2}'(u^n) \quad (14)$$

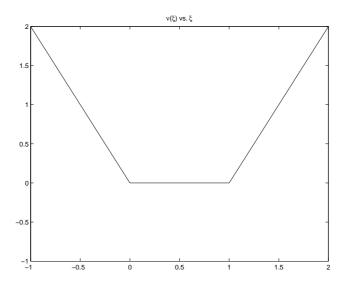


Figure 5: The function  $\nu(\xi)$  is used for exact penalization as a method to impose the constraint  $0 \le u \le 1$  in the minimization of Claim 1.

where  $\varepsilon_1, \varepsilon_2 > 0$  are small constants, and  $\nu_{\varepsilon_2}(\xi)$  is a regularized version of  $\nu(\xi)$  that smooths the latter's kinks at 0 and 1.

The image shown in the Figure 5 is not piecewise constant with two regions; in fact it is not very well approximated by any image that takes only two values. This makes it a challenging test case for the two-phase segmentation problem (images that are already approximately two valued are easily and very quickly segmented by these algorithms, and thus are easier examples).

Figure 6 shows the result found (i.e. the function u) using algorithm (14) to update the unknown function u(x) that represents the two phases, when the given image f(x) is the one shown in Figure 5. The two constants  $c_1$  and  $c_2$  were initially chosen to be 1 and 0, and updated occasionally; they eventually converged to 0.4666 and 0.0605, respectively. Although the considerations above (in particular, Theorem 2) do not imply that the minimizers of the objective functional turn out to be binary themselves (which would make the thresholding step in the algorithm of Theorem 2 unnecessary), in practice they seem to. Indeed, the image of Figure 6 is very close to being binary. Furthermore, it gets even closer to being binary if the computation is repeated using smaller values of the regularization parameters  $\varepsilon_1$  and  $\varepsilon_2$  that appear in scheme (14).

Figure 7 displays the histograms of u(x) at intermediate stages of the gradient descent computation. During the evolution, the function u certainly takes a con-

tiuum of values; however, as steady state approaches, the values that u can take accumulate at the extreme ends of its allowed range. In this case, the extreme values seem to be about 0.04 at the low end and 0.97 at the high end. They are not 0 and 1 because the exact penalty function  $\nu$  that appears in (14) is regularized.

The theory above says that for fixed  $c_1$  and  $c_2$ , all level sets of the function u(x) are minimizers. The table below shows the value of energy (10) computed by taking  $\Sigma = \{x : u(x) = \mu\}$  for several different values of  $\mu$ , where u(x) is the numerical example of Figures 6 and 7, and the constants  $c_1$  and  $c_2$  have the values quoted above.

$\mu$	Energy
0.2	17.7055
0.4	17.6458
0.5	17.6696
0.6	17.6655
0.8	17.6740

For comparison, we note that the energy of a disk centered at the middle of the image with radius a quarter of a side of the image domain has energy of about 112. Thus, even though there is some minor variation among different level sets of the function u(x) (see Figure 8 for a plot of several level contours) and their corresponding energies (due to some of the approximations, such as in the penalty function  $\mu$ , that were made to get a practical numerical algorithm), the difference in energy between them is quite small; they are all almost minimizers.

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Figure 6: The given image f(x) used in the two-phase segmentation numerical results discussed in Section 4.

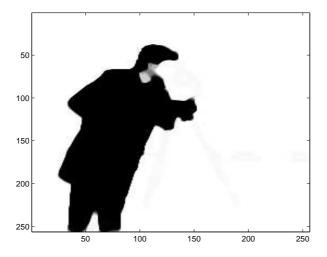
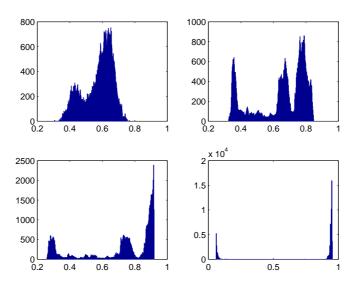


Figure 7: The solution u(x) obtained by the numerical scheme of Section 4. Although our claims do not imply the solution itself turns out to be binary, this seems to be always the case in practice. As can be seen, the computed solution is very close to being binary.



Figure~8:~Histograms~of~intermediate~solutions~of~the~fbw~in~(14).

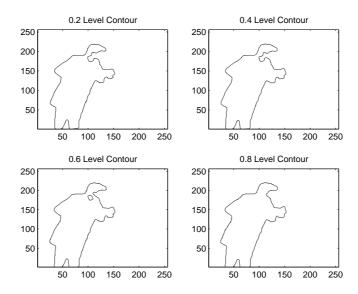


Figure 9: Plot of several level contours of the solution obtained. They are all very close to each other.

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