

# Initialization Framework for Latent Variable Models

Heydar Maboudi Afkham, Carl Henrik Ek and Stefan Carlsson

*Computer Vision and Active Preception Lab., KTH, Stockholm, Sweden*

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**Abstract:** In this paper, we discuss the properties of a class of latent variable models that assumes each labeled sample is associated with set of different features, with no prior knowledge of which feature is the most relevant feature to be used. Deformable-Part Models (DPM) can be seen as good example of such models. While Latent SVM framework (LSVM) has proven to be an efficient tool for solving these models, we will argue that the solution found by this tool is very sensitive to the initialization. To decrease this dependency, we propose a novel clustering procedure, for these problems, to find cluster centers that are shared by several sample sets while ignoring the rest of the cluster centers. As we will show, these cluster centers will provide a robust initialization for the LSVM framework.

## 1 INTRODUCTION

Latent variable models are known for their flexibility in adapting to the variations of the data. In this paper, we focus on a specific class of latent variable models for discriminative learning. In these models, it is assumed that a set of feature is associated with each labeled sample and the role of the latent variable is select a feature from this set to be used in the calculations. In both training and testing stages, these models do not assume that a prior knowledge of which feature to be used is available. Deformable Part Models (DPM) (Felzenszwalb et al., 2010; Felzenszwalb and Huttenlocher, 2005) can be seen as a good example of these models. With the aid of Latent SVM framework (LSVM), DPM provides a level of freedom for samples, in terms of relocatable structures, to adapt to the intra-class variation. As the result of this flexibility, the appearance of the samples becomes more unified and the training framework can learn a more robust classifier over the training samples. A good example of the model discussed in this paper can be found within the original DPM framework. In their work (Felzenszwalb et al., 2010), the method does not assume the ground truth bounding boxes are perfectly aligned and leaves it to the method to relocate the bounding boxes, to find a better alignment between the samples and the location of this alignment is considered as a latent variable. In a more complex example (Yang et al., 2012; Kumar et al., 2010), the task is to train an object detector without having a prior knowledge of the location of

the object in the image and considering it as a latent variable. Here, it is left to the learning framework to both locate the object and train the detector for finding it in the test images. Looking at the solutions provided for these examples, we can see that they are either guided by a high level of supervision, such as considering the alignment to be close to the user annotation (Felzenszwalb and Huttenlocher, 2005; Azizpour and Laptev, 2012), or guided by the bias of the dataset, such as considering the initial location to be in the center of the image, in a dataset in which most of the objects are already located at the center of the images (Yang et al., 2012; Kumar et al., 2010). In general, such weakly supervised learning problems are considered to be among the hardest problems in computer vision and to our knowledge no successful general solutions has been proposed for them. This is because with no prior knowledge of how an object looks like and acknowledging the fact that different image descriptors such as HOG (Dalal and Triggs, 2005) and SIFT (Lowe, 2004) are not accurate enough, finding the perfect correspondence between the samples becomes a very challenging problem.

In this paper, we address the problem of supervision in the mentioned models and ask the questions, “Will the training framework still hold if no cue about the object is given to the model?”, and if the model doesn’t hold, “How can we formulate the desirable solution and automatically push the latent variables toward this solution?”. To answer these questions,

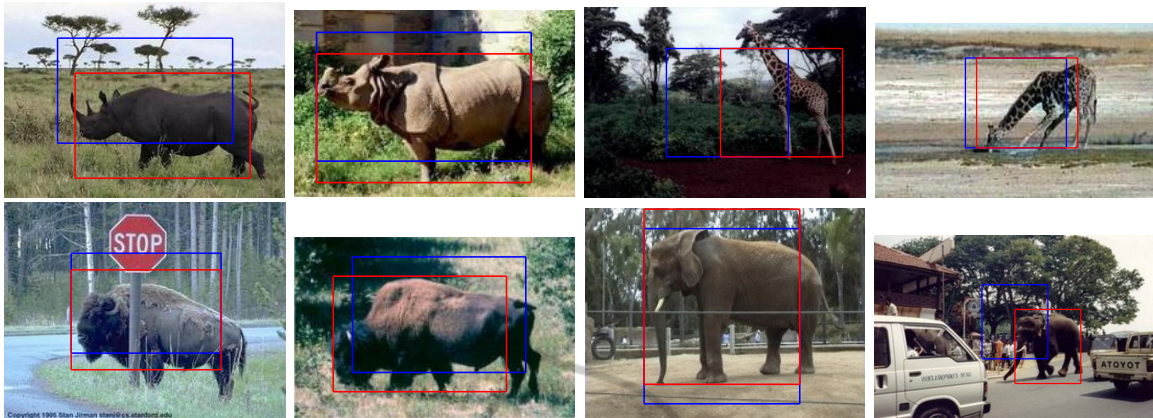


Figure 1: This figure shows, the final output of the object localization. In these examples, the training method was initialized using the blue bounding boxes located at the center of the image. In the training framework, it is left to the training algorithm to correctly converge to the location the objects with only the knowledge that this object has appeared in a number of images and does appear in the others. As it can be seen in these examples, final output (red box) correctly points to the objects.

we formulate this problem as a weakly-supervised clustering problem and show that the cluster centers provide an efficient initialization for the Latent SVM model (LSVM) (Felzenszwalb et al., 2010). To experimentally evaluate our method, we look at the problem of object classification with latent localization. This setup will provide us with an easy to evaluate framework which is very challenging to solve. Fig. 1, shows examples of this problem. In each image the blue box is the location that is initially considered to be the location of the object (In this case top left of the image) and the red box is the location found after the model is trained using the discussions in §3.

We organize this paper as following : In §2, we provide a proper definition of the problem and discuss strategies that can be used for initializing the latent variable models. In §3, we propose an initialization algorithm and in §4, we experimentally evaluate the properties of this initialization and compare it with other strategies. Finally, §5 concludes the paper.

## 2 PROBLEM DEFINITION

To formulate the problem, we assume that a dataset of labeled images  $\mathcal{D} = \{(\mathbf{x}_i, y_i)\}_{i=1}^N$  is provided with  $\mathbf{x}_i$  being the image and  $y_i \in \{-1, 1\}$  being the binary label associated with it. For each image, there is a latent variable  $h_i \in Z(\mathbf{x}_i)$  which localizes a fixed size bounding box. The content of this bounding box is encoded by the feature vector  $\Phi(\mathbf{x}_i, h_i) \in \mathbb{R}^d$ . In this problem, the task of the learning algorithm is to classify the images  $\mathbf{x}_i$  according to the labeling  $y_i$  while correctly localizing the object. If the accurate value of  $h_i$  is known for the training examples then the

problem becomes a standard detector training problem. However, with the assumption that the value of  $h_i$  of the training images is undetermined, the training task becomes significantly more challenging, because for each image we do not know which value of  $h_i$  points to the object and a wrong fixation of this value can lead to training of inefficient models. The Latent SVM model (LSVM) (Felzenszwalb et al., 2010) addresses this problem by minimizing the objective function

$$L_{\mathcal{D}}(\beta) = \frac{1}{2} \|\beta\|^2 + C \sum_{i=1}^N \max(0, 1 - y_i f_{\beta}(\mathbf{x}_i)), \quad (1)$$

where

$$f_{\beta}(\mathbf{x}_i) = \max_{z \in Z(\mathbf{x}_i)} \beta^T \Phi(\mathbf{x}_i, z). \quad (2)$$

This optimization is usually done by iterating between fixing the latent variables based on computed  $\beta$  and optimizing the model parameters  $\beta$  over the fixed problem. These iterations usually start by an initial fixation of the latent variables. In this paper, we will show that the outcome of this method is very sensitive to the initialization and will discuss the effect of different initializations on the solution.

When no prior knowledge of the object locations is available, we can take several strategies to pick the initial bounding box. In strategies such as picking the *center*, *top left* or a *random* bounding box in the image, the selection is done independent of the content of the image. In these cases, the performance of the final classifier depends on how good these initials overlap with the interest object. Unfortunately, guaranteeing this overlap without previous knowledge of the object is not possible. To fix this, it is more desirable for the initialization to be derived from the content

of the images. Since, we already know that the solution is an object that exists in within all the positive images, it is natural to use a clustering procedure on positive images to obtain a proper initialization. An example of such a procedure, is to use the *kmeans* algorithm to cluster the feature vectors coming from the positive images. Each cluster center produces by *kmeans* corresponds to certain feature vectors that repeats across the positive training samples. Using a cross validation process it is possible to pick the most representative center and use it to initially fix the latent variables.

The problem with such a selection is the fact that very few bounding boxes in each image actually correspond to the object and these feature vectors will most likely be ignored by a generic clustering algorithm, due to lack of data. This brings the need for an algorithm that can ignore the large and irrelevant feature vectors and only focus on what is shared between the positive images. This is desired because a feature vector that exists in all the positive images is more likely to represent the object category. Such an algorithm is presented in the following section.

### 3 LATENT OPTIMIZATION

To find feature vectors that are shared between the images, let  $\mathcal{D}^+$  contain only the positive images ( $|\mathcal{D}^+| = M$ ). For a given point  $p \in \mathbb{R}^d$  we define  $\Psi(\mathbf{x}_i, p) = \Phi(\mathbf{x}_i, z_i^*(p))$  where

$$z_i^*(p) = \arg \min_{z \in Z(\mathbf{x}_i)} \|p - \Phi(\mathbf{x}_i, z)\|^2. \quad (3)$$

The aim of this section is to find a point  $p \in \mathbb{R}^d$  such that the objective function

$$C_{\mathcal{D}^+}(p) = \frac{1}{M} \sum_{\mathbf{x}_i \in \mathcal{D}^+} \|p - \Psi(\mathbf{x}_i, p)\|^2, \quad (4)$$

is minimized. In other words, we wish to find a feature vector that each positive image has a feature vector similar to it. To minimize this objective function, we use an iterative method starting with a point  $p^{(0)}$  (or equivalently an initial fixation of the latent variables) and calculate the next point as

$$p^{(k+1)} = \frac{1}{M} \sum_{\mathbf{x}_i \in \mathcal{D}^+} \Psi(\mathbf{x}_i, p^{(k)}). \quad (5)$$

Clearly, if for each  $\mathbf{x}_i$ ,  $|Z(\mathbf{x}_i)| = 1$  then this iteration converges to the mean of the data after one step. However, because of the latent variables, the feature vector obtained from  $\Psi(\mathbf{x}_i, p^{(k)})$  changes with the change of  $p^{(k)}$ . This fact makes the behaviour this algorithm more complex. The following theorem shows that this iterative method minimizes the objective function 4.

**Theorem 1.** *Given an imageset  $\mathcal{D}^+$  and  $p^{(k)}$  and  $p^{(k+1)}$  defined as above, the following statement always holds :*

$$C_{\mathcal{D}^+}(p^{(k+1)}) \leq C_{\mathcal{D}^+}(p^{(k)}) \quad (6)$$

*Proof.* To avoid clutter in the proof, for all  $p, q \in \mathbb{R}^d$  we define

$$\Delta_i(p, q) = \|p - \Psi(\mathbf{x}_i, q)\|^2. \quad (7)$$

Since  $p^{(k+1)}$  is calculated by Eq. 5, we can write

$$\frac{1}{M} \sum_{\mathbf{x}_i \in \mathcal{D}^+} \Delta_i(p^{(k+1)}, p^{(k)}) \leq \frac{1}{M} \sum_{\mathbf{x}_i \in \mathcal{D}^+} \Delta_i(p^{(k)}, p^{(k)}). \quad (8)$$

This inequality assumes that the latent variables are fixed by the point  $p^{(k)}$ , and  $p^{(k+1)}$  is simply the mean of these fixed points. Due to the properties of mean, the value of  $\frac{1}{M} \sum_{\mathbf{x}_i \in \mathcal{D}^+} \Delta_i(\hat{p}, p^{(k)})$  is the lowest when  $\hat{p} = p^{(k+1)}$ . We can also conclude from the definition of  $\Psi(\mathbf{x}_i, p)$  that

$$\Delta_i(p^{(k+1)}, p^{(k+1)}) \leq \Delta_i(p^{(k+1)}, p^{(k)}). \quad (9)$$

Combining the inequalities 8 and 9, gives us

$$C_{\mathcal{D}^+}(p^{(k+1)}) \leq C_{\mathcal{D}^+}(p^{(k)}). \quad (10)$$

□

This theorem shows that the iterative method discussed in this section will converge to a mode in  $\mathbb{R}^d$  with each points of this cluster coming from a different image. Since this method only uses the positive images, there is no guarantee that the found point actually corresponds to the object we are looking for. To address this problem, Alg 1 sequentially finds  $k$  distinct such points.

Fig. 2(Left), shows the behaviour of Alg. 1 on synthetic data. In this data, each image is simulated as a set of points randomly sampled from different distributions (Explained in the appendix). This figure shows, how Alg. 1 converges to the data modes and avoids already found distributions. In this example, every set (image) contains at least one point from each distribution. To highlight the difference between *kmeans* and Alg. 1, we construct a slightly different synthetic data. Here, rather than populating the sets with points coming from all distributions, we divide the sets in to two groups and follow a different strategy for populating each group. As it can be see in Fig. 2(Right), eight distributions are marked by three colors  $\{blue, cyan, magenta\}$ . We sample from *blue* and *cyan* distributions to construct the sets of the first group and from *magenta* and *cyan* distributions to construct the sets of the second group. Clearly, the solution we are interested in should belong to all

**Algorithm 1:** Finding shared representations.**Input:**  $\{(\mathbf{x}_n, \mathcal{L}_{x_n})\}_{n=1}^N, k$ **Output:**  $k$  cluster centers in the joint feature space

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1:  $\mathbf{P} \leftarrow \emptyset$  // Computed Centers
2: for  $i \leftarrow 1$  to  $k$  do
3:    $p_i^{(0)} \leftarrow$  Randomly pick a vector from  $\mathcal{D}^+$ 
4:    $j \leftarrow 1$ 
5:   while not converged do
6:      $\mathbf{A}_{p_j^{(i)}} = \{\psi(\mathbf{x}, \mathbf{p}_j^{(i)}) : \mathbf{x} \in \mathcal{D}^+\}$ 
7:      $\mathbf{A}_j^{(i)} \leftarrow \{a \in \mathbf{A}_{p_j^{(i)}} : \forall p \in \mathbf{P} (\|a - p_j^{(i)}\| < \|a - p\|)\}$ 
8:      $p_j^{(i+1)} \leftarrow (\sum_{a \in \mathbf{A}_j^{(i)}} a) / |\mathbf{A}_j^{(i)}|$ 
9:      $j \leftarrow j + 1$ 
10:  end while
11:   $\mathbf{P} \leftarrow \mathbf{P} \cup \{p_j^*\}$ 
12: end for

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sets and come from a *cyan* distribution. We execute kmeans on all the data points to find two cluster centers and execute Alg. 1 using the initial points demonstrated in this figure. Here, it is expected from kmeans to divide all the data points into two clusters. As it can be seen, neither of the cluster centers found by kmeans is close to the *cyan* distributions. On the other hand, the centers found by 1, are located at the center of both *cyan* distributions. In other words, while kmeans tends to divide the data into several partitions, Alg. 1 focuses on locating modes of the data with the property that a feature vector close to them exists in every sample, a property that is not necessarily true for the centers found by kmeans or other existing clustering algorithms.

## 4 EXPERIMENTS AND RESULTS

To experimentally analyze the effects of the initialization on the outcome of latent variable models, this paper uses the mammals dataset (Heitz et al., 2009) which has been used to benchmark the methods in (Kumar et al., 2010; Yang et al., 2012) and follows their experimental setting. In these experiments, it is assumed that the objects have the same size and the main challenge is considered to be the localization of the object. To describe the image, we have used the HOG descriptor (Dalal and Triggs, 2005; Vedaldi and Fulkerson, 2008). The latent svm implementation used in this paper is based on (Felzenszwalb et al., 2010) and as we can see in table 1, our implementation slightly outperforms the results presented in (Yang et al., 2012) for linear models, using the same assumptions. Each experiment is repeated

10 times on random splits of the dataset into training and testing sets and the mean performance is reported.

Table 1: Comparison between the classification rates obtained using different initialization methods. The large difference between these numbers shows the sensitivity of the local variable models to initialization and how important it to have robust methods for initializing them. In this table, each experiment was repeated 10 times and the average performance is reported. (\*) Result from (Yang et al., 2012).

Init. Type	Acc. %
Center	$80.15 \pm 2.79$
Center (*)	$75.07 \pm 4.18$
Random	$66.93 \pm 3.56$
Top Left	$61.75 \pm 3.06$
Kmeans (10 Centers)	$69.85 \pm 2.15$
Alg. 1 (10 Points)	$78.47 \pm 3.91$

As discussed in §2 and §3, we compare several strategies of initialization  $\{center, random, top\ left, kmeans, Alg. 1\}$  and measure their effect as the performance of the resulting detector on the test set.

- *Center*: In this case the initial location is selected to be the center of each image. As we can see in table 1, this initialization provides us with the best performance despite the fact that this initialization has nothing to do with the content of the image.
- *Random*: In this case the initial location is selected randomly. Ideally on a non-biased dataset the performance of the random initialization should be close to the center localization but as we can see in table 1, there is a significant performance drop when the random initialization is used.
- *Top Left*: This initialization type was chosen to make sure that initial locations has minimal over-

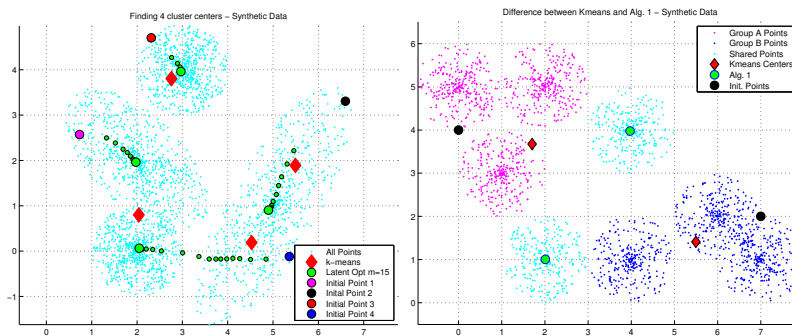


Figure 2: This figure shows the behaviour of Alg. 1 on synthetic data and compares it with the kmeans algorithm. As we can see Alg. 1 converges to the center of the distributions and confirms our analysis in §3. **(LEFT)** To build the synthetic data which simulates the conditions of the problem, 1500 data points were randomly sampled from 4 distributions. These points were randomly distributed in 100 sets each containing 15 points. In this toy problem, each set was considered to be a data sample with 15 different latent locations to choose from. **(RIGHT)** To show the difference between kmeans and Alg. 1, the synthetic data in this figure is produced by sampling from eight distributions are marked by three colors  $\{blue, cyan, magenta\}$ . We have divided the sets into two groups with the first groups sampling from *blue* and *cyan* distributions and the second sampling from *magenta* and *cyan* distributions. The distributions shared by all sets are the *cyan* distributions and as we can see while the centers found by kmeans are not close to these distributions, Alg. 1 converges to the center of these distributions.

lap with the target objects. As it can be seen in table 1, the lowest performance is achieved when this overlap is minimized.

- *Kmeans*: We use kmeans as a baseline for the performance of Alg. 1. In order to use kmeans for the initialization, we first cluster all the feature vectors coming from the positive training set into 10 clusters. To pick which cluster center which is the most representative, we divide the training set in half and cross validate LSVM method while initialized with different centers. As it can be seen in table 1, the performance significantly improves compared to choosing a random initialization. Here, once the most representative center is selected, the LSVM is trained over the whole training set and only this boundary is used for evaluating the test images and no other information is used at the testing stage.
- *Alg. 1*: Similar to the setting for kmeans, 10 modes were produced using Alg. 1 and the most representative was used for initialization of LSVM. Similarly, only the decision boundary trained using the most representative center is used at the testing stage. As it can be seen in table 1 the results significantly outperforms the baselines. A

It should be mentioned that in most cases there should be no difference between the performance of *center*, *random*, *top left* initialization strategies, due to the fact that these initializations have nothing to do with the content of the image. In the case of this problem, the significant performance gained when the initial lo-

cation placed at the center of the image, comes from the fact that most objects in this dataset are located in the center and placing the initial location at the center of the images gives the largest cover of the objects. In other words, by doing so we assume that for most objects, we already know the location of the object. In reality and on larger datasets, the performance of such initialization should be closer to *top left* initialization since the chance of covering the object using random selection or picking the center location decreases.

## 5 CONCLUSIONS

In this paper, we have shown how different initialization strategies can effect the outcome of the LSVM framework. To reduce the effects of the initialization, we have formulated what a desired solution looks like in terms of cluster centers and proposed an algorithm for finding these cluster centers. As our experiments show, LSVM framework trains a reasonably accurate model using the initialization provided by our method, without taking advantage of dataset bias or being guided by user annotation.

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## APPENDIX

### SYNTHETIC DATA

Each image in the dataset  $\mathcal{D}$ , can be seen as a collection of different feature vectors. These feature vectors usually come from many different distributions. In the problem discussed in this paper, among all these distributions we are interested in the distributions that each image in the dataset has a feature vector coming from that distribution. To build this data synthetically, we assume that  $\gamma_1, \dots, \gamma_n$  are given distributions in  $\mathbb{R}^d$  and assume that each image is set containing several points sampled from these distributions. Each set is populated with  $m$  vectors where each is obtained by randomly selecting a distribution and sampling from it. This way each set simulates an image with  $|Z(\mathbf{x})| = m$ . For  $d = 2$ , it is possible to visualize the data points and get better understanding of how the algorithms behave and visually compare them with other algorithms.