

# How to train your dragon: best practices in pedestrian classifier training

Remi Trichet, *remi.trichet@inria.fr* Francois Bremond, *Francois.Bremond@inria.fr*

**Abstract**—The present year witnesses another milestone in Pedestrian detection’s journey: it has achieved remarkable progresses in the course of the past 15 years, and experts foresee an everyday use of numerous stemming applications within the next 15. Standing on the tipping point between yesterday and tomorrow pushes to the field’s retrospect. Features, context, the combination of approaches, and feature data are responsible for most of the breakthroughs in pedestrian detection. If the first three elements cover the bulk of the literature, much less efforts have been dedicated to feature data. In many aspects, the construction of a training set remains similar to what it was at the birth of the domain, some related problems are not well studied, and sometimes still tackled empirically. This paper gets down to the study of pedestrian classifier training conditions. More than a survey of existing training classifiers or features, our goal is to highlight impactful parameters, potential new research directions, and combination dilemmas. Our findings are experimentally verified on two major datasets: the INRIA and Caltech-USA datasets.

**Index Terms**—pedestrian detection, real-time, data selection, parameter estimation, big data, dataset training

## I. INTRODUCTION

Accurately detecting human beings in an image is a task that an infant can perform at the age of one, but a computer algorithm still can’t reliably achieve. Nevertheless, a wide range of applications such as congestion analysis, automotive, abnormal event detection, retail data mining, robotics, human gait characterization have boosted the field’s progresses which is now at the core of many computer vision advances.

In a recent study [5], features, context, feature data, and the combination of approaches have been identified as the key performance elements for pedestrian detection. Roughly 30% of approaches focus on developing, combining or adapting features, and this direction has lead to most of the breakthroughs over the past years, such as channel features [18], or convolutional network [47]. Context adaptation [28], harnessing frequent geometry [32][50] or environment patterns [48][64], also successfully received attention from the community, but this algorithm adaptation is often application or even dataset dependent.

However, less progress has been stated concerning other areas such as feature data, threshold estimation, or non-maximum suppression. Training sets are still generated by applying the same techniques than a decade ago: Increasing the data variability is done by blindly mixing datasets [5], and Bootstrapping [16][25] remains among the preferred choices for training set generation. To the best of our knowledge, alternatives to greedy non-maximum suppression are never considered, and the majority of training parameters are overlooked in short-length papers.

Consequently, lots of question have never been clearly answered and the related approaches are often empirical. How to augment the number of positive examples in my training set ? Is a soft-cascade always working ? What are the alternatives to non-maximum suppression ? Should we always consider data cleaning before training ? What is the best strategy to set up the classifier cascade thresholds ? These details are often understated in short-length papers. Hence, by conducting thorough parameter evaluation and comprehensive method variation tests, this detail oriented paper strives to determine what are the best practices to train a pedestrian classifier. To the best of our knowledge, these experiments are new to the community. They bring a new light on some overlooked parameters, and precise or disprove some of the community empirical findings that were never experimentally validated before.

More specifically, our contributions are manifold. We provide extensive experiments and analysis in the domains of model size, instance selection, data cleaning, data generation, as well as near real-time classifier choice and thresholding, soft-cascading, and non-maximum suppression.

In order to study underexplored state-of-the-art areas and assess parameters impact independently, we first define a baseline that performs on par with common pedestrian detectors while remaining generic enough to test on the most common competitors. To validate our experiments, we provide tests with two types of features: HoG [16] and LBP [46]; and on two major datasets, INRIA [16] and Caltech USA [20].

The rest of this paper is organized as follows. Section 2 reviews the related state-of-the-art. Section 3 describes the reference classifier training methodology along with the variants that we will utilize throughout this study. Section 4 covers the large set of experiments we have undertaken and Section 5 sums up our findings and discusses possible future directions.

## II. RELATED WORK

Classifier optimization always has been a major concern in computer vision. Techniques ranging from active learning [39] to feature augmentation [11][40] have been employed to fine-tune the classifiers on tasks as various as person re-identification [42][62][68] or event recognition [12][41].

More specifically, data selection from imbalanced datasets has been a concern for pedestrian detection since the birth of the field. [34] proved that disproportioned datasets degrade SVMs prediction accuracy, especially for non-linearly separable data. Subsequent research on these experiments [65] showed that best performance was obtained for approximately

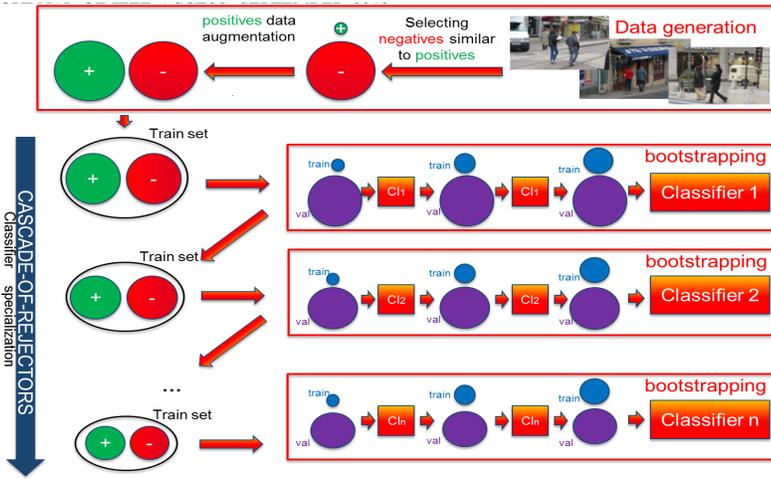


Fig. 1. Training pipeline. The initial training set generation selects data while balancing negative and positive sample cardinalities (Top). A cascade classifiers is then trained on it (Left), each independent classifier being learnt through bootstrapping (right). balanced positive and negative sets is sought all along the cascade. *cl*:classifier

comparable class cardinalities when over-sampling the minority set.

Limited work carries out these issues. Bootstrapping [16][25] is probably the most common adaptation to the problem. This data training mechanism improves accuracy thanks to successive training set updates focusing on samples difficult to classify. It is first initialized on uniformly drawn negative samples. At each iteration, the negative training set is augmented with the false positives of the previous run. The process keeps repeating until the performance converges or a memory threshold is reached. [57] further showed that 2 iterations lead to optimal performance. [27] adopted a different practice for random forest. They first train  $t$  trees with uniformly-drawn samples, use them to obtain harder positive and negative sample to train the next  $t$  trees, and iteratively repeat this process. Each tree group bears its own bias, therefore improving the overall forest performance. [24] incorporated latent values to optimise hard negative mining.

Generative techniques consider over-sampling the minority class through artificial data generation. This type of approach suffers from both, over-generalization and the risk to create erroneous data. The SMOTE algorithm [58] is probably the most renowned one. It creates new data as the linear combination of a randomly selected data point and one of its (same class)  $k$  nearest neighbors. Borderline-SMOTE [29] subsequently proposed to improve SMOTE through clever selection and strengthening of weak pairs. ADASYN [30] considered limiting the over-generalization through density-based data generation. In the same vein, [35] introduced cluster-based oversampling, simultaneously tackling the intra- and inter-class imbalance issue. Finally, [6] proposed curriculum learning to progressively incorporate into the training set data ordered according to their generative power (i.e. from easy samples to hard ones).

Cost sensitive learning methods [22][53] are an alternative to generative techniques that learn the cost of misclassifying each data sample. They optimize the classification performance by up-weighting important data and provide a natural way to enhance the minority class.

Finally, [33] highlighted and tackled the imbalance issue with a two classifier cascade: The first one aims for a high recall whereas the second one enhances precision.

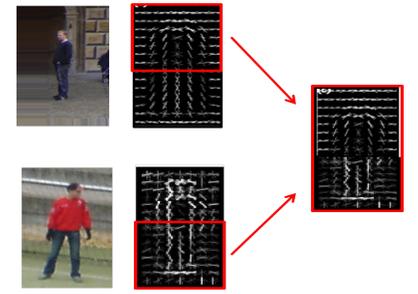


Fig. 2. data generation example with horizontal cut. the Feature values corresponding to upper and lower body parts are concatenated to form a new data sample.

An alternate strategy to deal with the excess of negative data is to successively prune out easy-to-classify instances via the use of several consecutive classifiers. This cascade-of-rejectors [69][7] can also be considered to speed up the detection. The technique, inspired by Viola & Jones face detector [56], builds up a classifier cascade that consists of successive rejection stages that get progressively more complex, therefore rejecting more difficult candidates as the classifiers get more specialized. [52] further upgraded the principle from features to data by performing hierarchical clustering on the dataset, leading to improved classification results. However, it requires, for each new datum, to find the corresponding set of clusters it belongs to. See [31] for more details on learning from imbalanced datasets.

### III. REFERENCE TRAINING METHODOLOGY

The definition of a reference method is a difficult task. On one hand, it needs to be generic enough to allow the evaluation of a wide spectrum of common techniques and its reproducibility under various application environments. But on the other hand, results competitive with the state-of-the-art are required for the conclusions to be usable with up-to-date algorithms. Bearing this compromise in mind, we chose the following reference method. For the sake of conciseness, we restricted it to real-time or near real-time techniques, but in most cases, the extension to more compute-intensive strategies remain possible.

Our training methodology, illustrated in figure 1, unfolds as follows. It decomposes in two distinct parts: The initial training set generation and the classifier training. The initial training set generation carefully selects data from a set of images while balancing negative and positive sample cardinalities. We then train a cascade of 1 to  $n$  classifiers. This cascade could include a cascade-of-rejectors [15][69][7][52][56], a soft cascade [8], or both. In addition, each independent classifier is learnt through bootstrapping [16][25] to improve performance. One key aspect is to seek balanced positive and negative sets at all time. Hence, all along the cascade, the minority class is oversampled to create balanced positive and negative sets.

The body of this section describes each of these components independently while putting the emphasis on important

aspects or parameters that will be analysed in the next section dedicated to experiments.

### A. Minority class oversampling

As the studies on the topic prove that class imbalance leads to deteriorated performance, one may seek class balance before any training. Two options exist. The first one consists in undersampling the majority class. This solution doesn't satisfy our needs as a large training set is preferable to fully represent the data variability and avoid overfitting during the training process. The remaining solution oversamples the minority class. This process implies the creation of extra data that will not match any existing instance in the dataset images. Nevertheless, these newly created instances have to credibly represented a hypothetical data. The risk of introducing spurious data in the dataset should be minimized, especially in our case, where we may generate over 90% of one class instances. Consequently, the large majority of data generation algorithms create new data by fusing the values of existing samples. In order to respect the credibility constraint, the parent data often are feature space neighbors, leading to the new data to be created in their feature space local vicinity. This is the over-generalization issue that cripples most generative techniques, like SMOTE [58]: Instead of exploring the emptiest areas of the feature space, these algorithms keep populating the denser areas, thus reinforcing the dataset biases.

In order to free the process from this issue, [55] takes advantage of the spatial localization of most pedestrian feature values. Indeed, due to their block or region decomposition, geometrically structured descriptors, such as HoG [16], LBP [1], DPM [24], or Haar-inspired strategies [18], each of their feature values have a precise localization within the proposal window. The new data generation function  $f(\cdot)$  unfolds as follows. We randomly select 2 instances  $x_1 = [x_{11}, \dots, x_{1n}]$  and  $x_2 = [x_{21}, \dots, x_{2n}]$  from the same class  $S_i$  and create a new sample with their lower and upper body features, while avoiding doublons. Uniform sampling with replacement is employed. More formally:

$$\forall x_1, x_2 \in S_i \quad f(x_1, x_2) = [x_{11}, \dots, x_{1\frac{n}{2}}, x_{2\frac{n}{2}+1}, \dots, x_{2n}] \quad (1)$$

So, a minority class of  $n$  samples can spawn a maximum of  $n \times (n - 1)$  new instances. Not normalizing the final vector gives better results than a re-normalization.

In their approach, [55] solely used the horizontal dichotomy in order to guarantee the creation of valid data. In this paper, we choose to explore the limits of this strategy by relaxing this constraint in the experiments section.

### B. Training Set Generation

This first step aims to form the initial training set by selecting insightful instances from the swarm of features extracted from the training images  $I$ . Let  $D$  be the desired training set cardinality, with training set positive and negative instances cardinality  $P$  and  $N$  so that  $D = P + N$ . To get a representative negative subset of the training set, the common approach is to randomly extract  $n = D/|I|$  data from each image. This typically leads to a large imbalanced training set

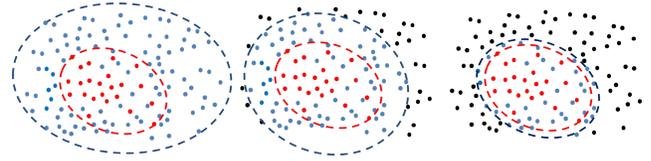


Fig. 3. Representation of the parameter  $n_0$  for instances selection. Positive samples are in red, selected negative ones are in blue, unselected ones in black. The rough class boundaries are depicted with dashed lines. From left to right,  $n_0=100\%$ ,  $n_0=50\%$ ,  $n_0=10\%$

with few positives  $P$  and a large subset  $N$  of negative instances. Data generation techniques [58][35] can then be employed to limit the class bias. Positive instances horizontal flipping is also frequently used.

To get a better control over the selected instances, we borrowed the strategy from [55]. In this approach, ground truth positive examples are augmented with *near positives*, defined as negative data having a spatial overlap with the ground truth positive samples higher than 90%. Let  $P_i$  be the set of already selected positive data after the processing of  $i$  images. When processing image  $i + 1$ , negative samples are selected as follows:

- 1) Data negative samples are sorted according to their similarity with positive samples  $P_i$ .
- 2) Select the subset of the  $n_0$  most similar negative samples (based on euclidean distance), with  $n_0 \geq n$ .
- 3)  $n = D/|I|$  negative instances are randomly selected among the remaining subset  $n_0$ .

Finally, the positive class set is oversampled to reach a balance with the negative instances.

$n_0$  represents the overall similarity between the selected negative  $N$  and positive data  $P$ . Setting  $n_0$  low restricts the selected negative data to those similar to positive ones, therefore building a classifier that focusses on the often misclassified data that lie on the border between the two classes. Inversely,  $n_0=100\%$  comes back to the common approach for instance selection where the full training set is represented. Figure 3 depicts the idea. Only ground truth positive and the same amount of original negative samples are fed to the training set, the rest of the original data and the augmented data are directed to the validation set.

### C. Data cleaning

The purpose of data cleaning is to spruce the dataset off data that could potentially mislead the classifier training. In this paper, we experimented with two data cleaning methods with distinct spirits: Tomek links [54] focusses on mixed classes neighborhoods while DBSCAN cleans overpopulated areas [23].

**Tomek links** are defined by pairs of opposite classes data that are respective nearest neighbors. More formally, given an instance pair  $(x_i, x_j)$ , where  $x_i$  and  $x_j$  are respectively positive and negative instances, and  $d(x_i, x_j)$  is the distance between  $x_i$  and  $x_j$ , the  $d(x_i, x_j)$  pair is called a Tomek link if there is no instance  $x_k$ , such that  $d(x_i, x_k) < d(x_i, x_j)$  or  $d(x_j, x_k) < d(x_i, x_j)$ . Therefore, Tomek links are drawn either from noise, or on classes border. So they are a powerful tool, for data clean-up, especially after synthetic sampling. Tomek links can then be removed until all nearest neighbor

pairs belong to the same class. In our case, we employ euclidian distance and restrict the deletion to negative data to spare the scarce minority class.

**DBSCAN cleaning** makes use of the renowned density-based clustering algorithm [23]. It basically performs a density correction, removing data from the mostly populated areas of the feature space. The algorithm unfolds as follows. With number of neighbors present within distance  $d$  associated to each point, the data points are ranked according to their neighbor cardinality. The algorithm then iteratively removes the sample in the most populated neighborhood, updates the cardinalities accordingly, and stops when all data samples have less than  $Nmin$  points in their vicinity. We used the euclidian distance,  $d=0.8$  and  $Nmin=5$  for all our experiments.

#### D. Classifier

The two employed classifiers in this work are typical Adaboost [63] and random forest [9]. After specifying our random forest setting, we will remind bootstrapping, the cascade-of-rejectors and soft cascade methods in the next subsections.

In our paper, the tree construction is grounded on typical entropy optimization  $S$ :

$$S = \sum_{C_i=0}^1 -p_{C_i} \log(p_{C_i}) \quad (2)$$

And the final confidence score for instance  $x$  is obtained by voting:

$$P(C_i(x)) = \frac{1}{F} \sum_{t=0}^F c_t(C_i) \quad (3)$$

where  $F$  is the number of trees,  $c_t(C_i)$  is the count for category  $C_i$  at the leaf node  $L_t$  of the  $t^{th}$  decision tree.

#### E. Classifier bootstrapping

We use bootstrapping [16][25] to augment the training set. At each iteration, we incorporate the most badly classified data from the validation set into the training set and retrain the classifier. We aim to repeat this process until the training set reaches its desired initial size  $T$ .

However, a maximum of 2 iterations is advised to avoid overfitting [57]. Therefore, we employ [55] method to deal with this issue, the idea being to introduce new data to prevent overfitting. Instead of simply selecting data from the validation set, we generate it the same way as detailed in the minority class oversampling section. Each new datum is generated based on 2 data samples selected among the misclassified data. Weighted sampling is utilized in this case, the probability of an instance to be drawn being set in relation to the extent of the error. The training set reduction algorithm, illustrated in figure 4, iterates until the desired training set size is reached.

#### F. Cascade-of-Rejectors

The cascade-of-rejectors is an efficient strategy for sieving out a large proportion of false positives, and has been applied several times in the context of pedestrian detection [15][69][7][52][56]. However, tuning this cascade of classifiers remains problematic. First, the classifiers are prone to

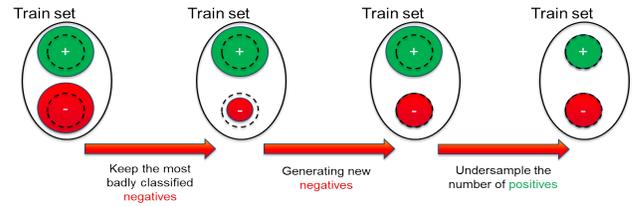


Fig. 4. Set cardinality balancing pipeline. '+' and '-' signs represent the positive and negative classes. Best viewed in color.

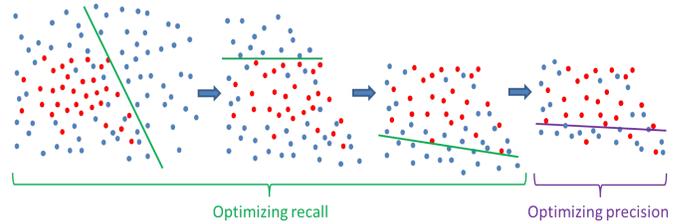


Fig. 5. Toy example of a cascade-of-rejectors training for binary classification. Blue and red colors represent the classes. The line depicts a (hypothetical) separating hyperplane. The first iterations optimize on recall, the final one optimizes for precision. Minority and majority classes are respectively under- and over-sampled between each iteration to learn on balanced classes. Best viewed in color.

errors that propagate along the cascade. Second, most authors employing this technique solely reject negative data after each classification step to avoid the creation of classifiers likely to reject positive data during the subsequent stages. However, this leads to progressively more imbalanced datasets. Fine tuning of the original positive versus negative instances ratio is usually necessary to obtain satisfactory results (a 2 to 4 times bigger negative set is the common practice). But this strategy is still under-optimal as balanced datasets are proved [65] to be the best performing ones.

To deal with these issues, we propose a new cascade-of-rejectors. Our approach borrows [33]'s main idea with a two-stage classification: The first one aims for a high recall whereas the second one focuses on precision.

The first stage embodies a cascade of  $n - 1$  rejectors. To avoid increasing the false negative rate, these  $n - 1$  initial classifiers enforce a high recall, moderately optimizing the performance on the precision. The metric formalizes as:

$$M = (TN \times W_{TN} + TP) / (P + N) \quad (4)$$

with  $P, N, TP, TN$ , and  $W_{TN}$  being respectively positive, negative, true positive, true negative instances, and the weight associated to  $TN$ .  $W_{TN}$  enforces the recall. Note that  $W_{TN} = 0$  comes back to optimizing for recall. When  $W_{TN} \leq 0.5$ , its parametrization has little impact on the results. We used  $W_{TN} = 0.5$  for all our experiments. Negative data, with a confidence below 0.5, are rejected at each stage. Negative data over-sampling is then undertaken to avoid emptying the validation set. When insufficient to match the positive instances cardinality, random positive data under-sampling is then considered. We employ Adaboost as classifiers for this part. These decision trees have demonstrated competitive performance whilst retaining computational efficiency [61][44]. During testing, typically over 95% of the data are safely discarded after 4-5 iterations. This cascade-of-rejectors guaranties a high recall, increased performance induced by



Fig. 6. Difficulties that non-maximum suppression tackles. Detections are in green. Best viewed in color.

balancing the datasets at each iteration, and few parameters to tune.

The second stage performs a finer classification, aiming for high precision results. Since the bulk of the data have been removed, more demanding computation can be performed at this stage without slowing down the detector. We employ a dense forest classifier, that has shown [5] giving slightly better performance than its counterparts on pedestrian detection. We optimize for the log-average miss rate [49]. Figure 5 illustrates the cascade effect on a toy dataset.

The only parameters that require careful tuning are the rejection thresholds  $t(C_i)$ , with  $C_i$  the  $i^{\text{th}}$  classifier of the cascade. During  $i^{\text{th}}$  stage of the testing cascade, any data  $d_j$  with confidence  $C_i(d_j) < t(C_i)$  will be rejected. Since the method optimizes for the recall, the large majority of data around the separating border between the two classes are negative ones. It is then sensible to set  $t(C_i) > 0.5$ . See the experiment section for a thorough evaluation of this parameter impact on performance.

### G. Non-Maximum Suppression

Objects of interest generating lots of detections in their neighborhood, non-maximum suppression is a method that keeps the maxima while deleting other proposals. Its purpose is two-fold. First, aggregating the information of nearby detections. It also works as an overlap control mechanism, balancing overlapping false detections and detections over self-occluding objects. Figure 6 illustrates these difficulties.

In this study, we use 3 different implementations of the threshold non-maximum suppression. First, Greedy non-maximum suppression is the systematically employed technique that compares the bounding boxes that overlap over a pre-defined threshold and keeps the best scoring one. We set the overlap threshold to 0.6 for all our experiments. We also question its supremacy by testing out to other methods. Threshold non-maximum suppression (t-NMS) [10] groups the detection according to the bounding boxes overlap with the group best candidate, keeps the  $d$  candidates with highest confidence for each group, and builds the final candidate position with the group mean border positions. Despite its simplicity, t-NMS doesn't require heavy parametrization and combines well with the cascade algorithms. Even though the complexity is  $O(n^2)$ , it is quite fast in practice since most of the negatives have already been suppressed by the cascade-of-rejectors. We set the overlap threshold to 0.4 and  $d = 2$  for all our experiments. We also considered Scale non-maximum suppression (s-NMS) [10] that differs from t-NMS by treating each scale separately.

In the occurrence of nested detections, our parametrization always favors the biggest one. Threshold non-maximum suppression is utilized by default.

## IV. EXPERIMENTS

This section, dedicated to our experimental validation, breaks down into several sub-parts. After presenting the datasets and the experimental setup details, we run individual experiments on each part main parameters or overlooked variations. The last part compares our work to the state-of-the-art.

### A. Datasets

We experimented on the INRIA and Caltech-USA datasets. The INRIA [16] features 1832 training and 741 testing high resolution pictures. This is among the most widely used dataset for person detection. Despite its small size compared to more recent benchmarks, the INRIA dataset boasts high quality annotations and a large variety of angles, scenes, and backgrounds.

The Caltech-USA [20] dataset totals 350000 images. Despite some annotation errors [67], its large size along with crowded environments, tiny pedestrians, and numerous occlusions probably make the Caltech-USA dataset the most widely used one. We experimented on the "reasonable set" that restricts algorithms to pedestrians over 50 pixels in size and a maximum of 35% occlusion.

With respectively a large variability and tiny occluded detections, everyday pictures and automotive application, these two datasets offer complementary settings for our experiments.

### B. Experimental setup

We experimented on the INRIA [16] and Caltech-USA (reasonable set) [20] datasets with HoG [16] and Haar-LBP [14] features. HoGs are configured with  $12 \times 6$  cells,  $2 \times 2$  blocks and 12 angle orientations, for a total of 2640 values. Block as well as full histogram normalisation are performed. The LBP descriptor follows the same structure, features a value per channel for each cell, and non-uniform patterns are pruned out [45]. Candidates are selected using a multi-scale sliding window approach [69] with a stride of 4 pixels. Approximately 60% are loosely filtered according to "edgeness" and symmetry. No background subtraction or motion features are employed on the Caltech dataset. The training set size is up-bounded at 12K samples. Increasing this threshold doesn't improve performance.

No generative resampling is performed. The model height is set to 100 pixels for the INRIA dataset, 50 pixels for Caltech.

We also implemented a cascade-of-rejectors baseline with the commonly used settings: Rejector optimisation is done according to the log-average miss rate metric. We tested this baseline with a typical initial validation set containing 4 times more negative samples than positive ones. To compare with the common multi-dataset training technique, we also used an augmented version with PETS2009 [26] positive samples. This set gathers 50K positive instances from both datasets and 110K

Component	default method	default parameter
Model height pixels		INRIA=100 Caltech=80
# data selected	FairTrain [55]	INRIA=2000 Caltech=1600
Data oversampling	FairTrain [55]	horizontal cut
Data Cleaning	Tomek links	
Resampling	Bootstrapping	5 iterations
Cascade	cascade-of-rejectors	6 iterations
Cascade classifier	Adaboost	512 weak classifiers
Cascade final classifier	Adaboost	512 weak classifiers
NMS	t-NMS	$t = 0.4$ and $d = 2$

TABLE I  
DEFAULT METHODS AND PARAMETERS USED FOR THE TRAINING

negatives from the sole INRIA benchmark. Table I summarizes the default methods and parameters that will be utilized for the different components alternatives.

We utilised the openCV implementation for the Adaboost and random forest classifiers. Random forest parameter set includes the number of decision trees, the number of sampled feature dimensions and the max tree depth. They were selected by measuring out-of-bag errors (OOB) [9]. It was computed as the average of prediction errors for each decision tree, using the non-selected training data. Adaboost is initialised with a tree depth of 2 and 256 weak classifiers. Each extra run adds 64 weak classifiers. This is a low number of weak classifiers compared to typical settings (i.e. [1024, 2048]). However, in practice, increasing this value leads us to lower performance and speed.

We use a variant of the threshold non-maximum suppression (t-NMS) [10] that groups the detections according to the bounding boxes overlap with the group  $d$  best candidate, keeps the  $d$  candidates with highest confidence for each group, and builds the final candidate position with the group mean border positions. We set the overlap threshold to 0.6 and  $d = 3$  for all our experiments. Log-Average Miss Rate (LAMR) [49] is employed as metric for all runs.

### C. Model size

state-of-the-art detectors empirically set the model size [49]. In this subsection, we conducted thorough experimentation to determine what should be the optimal model size given the initial conditions and what is its actual impact on performance. We varied the model height size in relation to the pedestrian minimal size observed in the dataset. We identified the smallest detections to be 50 and 60 pixels in height for respectively the Caltech (reasonable set) and INRIA datasets. And, as shown in table II, the best corresponding model sizes for these 2 benchmarks are 70 and 100 pixels in height.

The overall rule should be "the bigger, the better" as implicitly stated in [49]. Indeed, a bigger model will harness more information. However, the quality of the detection deteriorates when the model height size exceeds 150% of the smallest dataset pedestrian. The performance gain is then a factor of small detection frequency. As shown in [20], in Caltech's reasonable set, small detections represents the bulk of occurring pedestrians whereas they remain pretty rare in INRIA, which explains the discrepancy between the two datasets.

**lesson learned:** Model size matters. Optimal setting is in the  $[m \times 140\%, m \times 160\%]$  range, depending on the frequency of

Model height size	INRIA	Caltech
$m$	24.61%	55.78%
$m \times 110\%$	24.54%	52.82%
$m \times 120\%$	23.92%	49.58%
$m \times 130\%$	22.95%	47.51%
$m \times 140\%$	21.44%	<b>46.25%</b>
$m \times 150\%$	20.7%	47.73%
$m \times 160\%$	<b>20.31%</b>	51.32%
$m \times 170\%$	20.83%	57.28%
$m \times 180\%$	21.34%	61.23%

TABLE II  
LAMR PERFORMANCE ACCORDING TO THE MODEL HEIGHT SIZE THAT IS SET IN RELATION TO THE MINIMAL PEDESTRIAN SIZE  $m$ . BEST RESULTS IN BOLD.

the size  $m$  within the testset. When the frequency is unknown, we advise setting it up to  $m \times 150\%$ .

### D. Instances selection

Experiments on the training set  $D=P+N$  size have been conducted in [55], with the performance reaching a plateau around 200K data instances. Therefore we will restrict our tests to the parameter  $n_0$  tuning the similarity of the selected negative instances to the positive ones. We implemented 3 strategies for this purpose.

- 1) **Percentage:** Select the  $n_0$  percent most similar data.
- 2) **Threshold:** Select the data with similarity inferior to a threshold  $n_0$ .
- 3) **Amount:** Select the  $n_0$  most similar data.

In every case, a minimum of  $n = D/m$  data are selected to insure the same training set cardinality. Figure 7 compares these variants performance.

Despite the rather simple strategies employed here, data selection leads to a significant classifier performance improvement. The best strategy clearly consists in selecting the  $p$  data most similar to positives. The other two strategies, exploring parts of the feature space further away from positive instances lead to less discriminative classifiers. We explain the lower amount of data preselected for the Caltech dataset by the lower data variability of this dataset. Indeed, the angle and commonly encountered background objects are similar over the frames.

**lesson learned:** data selection and adaptation is an important area that deserves more attention from the community. As shown in these experiments, careful initial instances selection is crucial, and can lead to major performance boost. The selected negatives similarity with positive data must be set in accordance to the dataset variability.

### E. Data cleaning

Data cleaning is a rarely used pre-processing step. Is it overlooked or just dispensable? To answer this question, we evaluated its true importance for pedestrian training through the 2 proposed data cleaning methods. Since our training set generation method is not deterministic, each result is averaged over a total of 6 runs on the two datasets. Approximately 0.01 to 0.2% of the negative data are removed. No significant dataset discrepancy has been observed. Results are detailed in table III.

Dataset size	Tomek links [54]	DBSCAN [23]
80K	-0.71 ± 0.05%	-0.7 ± 0.13%
100K	-0.85 ± 0.04%	-0.54 ± 0.11%
120K	-0.7 ± 0.1%	<b>-0.62 ± 0.16%</b>
140K	<b>-0.59 ± 0.11%</b>	-0.34 ± 0.09%
160K	-0.34 ± 0.05%	-0.16 ± 0.07%
180K	-0.19 ± 0.03%	-0.12 ± 0.07%
200K	-0.17 ± 0.05%	-0.05 ± 0.05%

TABLE III

AVERAGE LAMR PERFORMANCE IMPROVEMENT INDUCED BY DATA CLEANING TECHNIQUES OVER 6 RUNS ON TWO DATASETS. BEST RESULTS IN BOLD.

The use of data cleaning shows a small but steady mean improvement. Tomek links outperform DBSCAN cleaning on this experiment. We assume that the DBSCAN cleaning may remove important data on the fuzzy class borders that fine tuned training sets tend to densely populate in order to improve the classifier discriminative power. Also, its use typically counterbalances the over-generalization problem which does not impair the training set construction presented in this study (see section III-A). We also observe that the data cleaning efficiency is inversely correlated to the training set size. We assume that the repercussions of the few spurious data instances weaken in accordance to the population density in their neighbourhood.

**lesson learned:** If cleaning the data could be dispensable, especially for big fine tuned training sets, it never harms the classification. Counter-intuitively, the popular DBSCAN does not perform best for pedestrian detection.

#### F. Additional data generation

Let's define a *credible* datum as a generated datum that visually looks like the object it represents. Note that the credibility of newly generated data can only be checked visually, so, the overall credibility of a generated set can hardly be quantified. In their original work, [55] restricted their data generation methodology to horizontal cuts, arguing that the newly generated instance credibility was paramount. In this experiment we test the limits of this hypothesis through the implementation of the following variations:

- 1) **Horizontal cuts** in 2 parts of equal areas. The original strategy proposed in [55]. See figure 2.
- 2) **Vertical cuts** in 2 parts of equal areas. Note that cutting along a vertical axis doesn't equal to an instance horizontal flip since the two parent instances are likely to be different data.
- 3) **Random cuts** in 2 parts of equal areas. This implies the cut passing through the centre of the bounding box. Each new data will be generated according to a different cut.
- 4) **Random walks**. The first half of the blocks are selected with the random walk algorithm [21] from the first sample. The other half of the blocks are selected from the second sample. Each new data will be generated according to a different walk. Using random walks guarantees that, at least, one of the 2 sets will be fully connected.
- 5) **Random walks and horizontal cut combination**. This strategy generates negative data utilizing the random walks and positive data with horizontal cuts.

Model height size	Feature	INRIA	Caltech
Horizontal cut	HoG	20.31%	<b>46.25%</b>
Vertical cut	HoG	21.28%	46.51%
Random cut	HoG	22.68%	48.64%
Random walks	HoG	25.32%	50.11%
1+4 combination	HoG	22.38%	48.14%
Horizontal cut	LBP	<b>18.42%</b>	51.26%
Vertical cut	LBP	19.12%	52.02%
Random cut	LBP	20.43%	56.91%
Random walks	LBP	22.87%	60.89%
1+4 combination	LBP	19.88%	53.28%

TABLE IV

DATA GENERATION ALGORITHM ACCORDING TO 4 DIFFERENT STRATEGIES. SEE TEXT FOR THE DESCRIPTION OF THE METHODS. BEST RESULTS IN BOLD.

The methods are ranked according to their generative power and the loss of credibility of the newly generated data. For instance, with the same initial set, random walks will generate a much greater variety of new data than horizontal cuts, leading to a better exploration of the feature space. However, its likelihood to create spurious data values, that doesn't correspond to a legitimate pedestrian image is also much higher. Results are provided in table IV.

With performance inversely decreasing with the loss of newly produced data credibility, this experiment confirms that the credibility is paramount. But counter-intuitively, the same rule also applies for negative example generation that do not have this restriction. Indeed, the combination of strategies 1 and 4 doesn't lead to any improvement. This implies that the negative samples structural geometry (i.e. a window frame, a car cabin, or tree branches) also cannot be tampered with.

**lesson learned:** Data generation credibility is essential and should not be jeopardized for both positive and negative instances. Horizontal cuts is the best performing method.

#### G. Classifier type

The general predominance of Random forest classifiers over Adaboost for pedestrian recognition have already been observed in the past [5]. In this study, We aim to push the experiment further by evaluating its actual impact within a cascade-of-rejectors. For this purpose, we tune the number of Adaboost and Random forest classifiers, keeping to 6 their total amount. Adaboost classifiers always come first in the cascade in order to benefit from their higher speed-up. This experiment is only presented for HoG features and the INRIA dataset as a similar behaviour is observed with other datasets and features.

Results are reported in table V. The only noticeable improvement is observed when the 4 or 5 last classifiers are random forests. This is sensible as over 90% of the data have been removed after these stages and the remaining ones are mostly grouped around the pedestrians. Using random forest as first classifier doesn't increase the results. We assume that a more discriminative classifier is less important during this first stage that focusses on negative data broadly different from the positive ones.

We also state that random forest only brings in a marginal improvement of 0.42% compared to Adaboost classifiers, contrary to what is usually reported [5][27]. We assume that

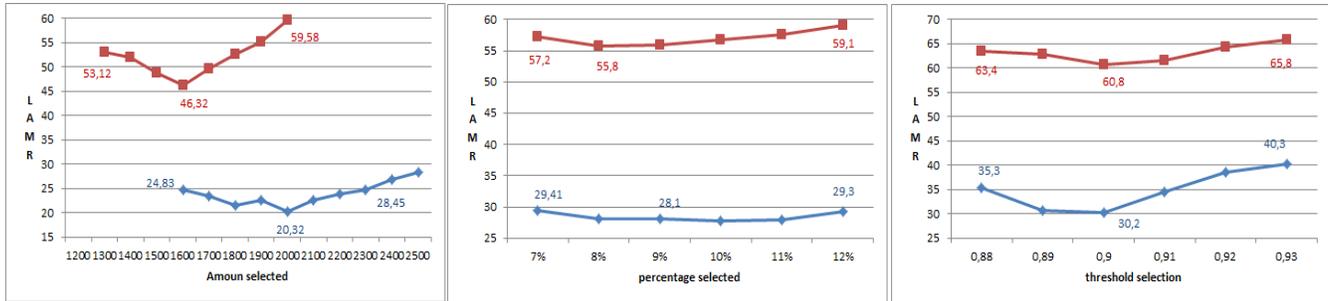


Fig. 7. Comparison of the instance selection strategy variants, setting the log-average miss rate performance in relation to the parametrization. The INRIA dataset is in red, Caltech in blue.

# Random forest classifiers	INRIA	Speed
0/6	20.31%	2.11fps
1/6	20.31%	1.09fps
2/6	20.29%	1.98fps
3/6	20.42%	1.75fps
4/6	20.13%	1.47fps
5/6	<b>19.89%</b>	1.24fps
6/6	19.92%	0.84fps

TABLE V

LAMR PERFORMANCE AND SPEED OF A HOG-BASED CASCADE ON THE INRIA DATASET, WHILE VARYING THE CASCADE CLASSIFIER TYPES (ADABOOST OR RANDOM FOREST). BEST RESULTS IN BOLD.

the Adaboost and random forest performance converge with the overall result improvement.

**lesson learned:** Powerful classifiers perform best when used in all but first classifiers of the cascade. general progresses in the field also call back into question the random forest classifier predominance [5].

### H. Soft cascade

We tested our pipeline with several variations of the famous soft-cascade [8], ranked according to their speed-up and their loss of discriminative power:

- 1) **soft-cascade:** The standard soft cascade [8] that adds one feature value at each step of the cascade.
- 2) **block-cascade:** A more discriminative but slower version of the soft cascade that adds one feature block at each step of the cascade.
- 3) **Reverse engineering :** This method Learns first the classifier based on the full length descriptor and recursively removes the less discriminative half until only 2 blocks are left. It can be considered as a preprocessing step that quickly prunes out most of the negatives before running the standard cascade.

Results are presented in table VII.

Standard soft cascade provides results on par with the state-of-the-art [8], which are significantly worse than our baseline. This makes sense as classifiers trained on low feature dimensionality are far less discriminative than full-length descriptors. The preprocessing classifier based on reverse engineering doesn't degrade the results but provides only a 2-fold speedup.

**lesson learned:** These results may sound the death knell of soft-cascade. If a soft-cascade provides a huge speed boost, its classifiers are independently not discriminative enough to compete with a the most recent classifiers trained on finely selected instances.

### I. Classifiers threshold setting

Training various classifiers on different data, with varying features of a classifier cascade invariably leads to a set of disparate classifiers. In addition to that, variations between train and test sets may induce more discrepancies. The use of any classifier cascade is pointless without a mechanism to cope with these variations. To deal with it, the actual strategy estimates and sets a different threshold  $t(C_i)$  for each classifier  $C_i$ . Setting up a set of cascade thresholds is not an easy task and the methods that tackle this particular pedestrian detection sub-problem are, at best, empirical. During testing, the optimal threshold setting is supposed to replicate the training phase dataset reduction across the cascade. In other words, each step should progressively eliminate harder negative examples while the recall should remain close or equal to 1.

In order to compare the efficacy of various commonly employed thresholding technique, we first needed to find the optimal setting. We manually looked for the threshold set leading to the best performance, exhaustively trying out all possibilities with a 0.01 step. Then, we tested the following methods:

- 1) **Constant:** Using the same threshold  $t(C_i) = T$  for each cascade classifier  $t(C_i); i = 0...1$
- 2) **Increasing:** Monotonously increasing the classifiers threshold.  $t(C_i) = T + c \times 0.01$ .
- 3) **Recall:** Estimating thresholds during training such that recall equals to  $r$ . This is the method most commonly used in the literature.
- 4) **Regular:** Using the same purposely low threshold  $t(C_i) = T$  for each cascade classifier  $t(C_i)$  and insuring that  $d\%$  of the remaining data are removed at each step of the cascade.

the parameter  $T$ , and  $d$  are manually set to the best performing value. However, the same value is employed for all datasets.  $T=0.58$  is utilized for experiment 1,  $T=0.55$  for experiment 2, and  $d=40\%$  for experiment 4. Results are displayed in table in table VIII.

This experiment revealed several interesting points. First of all, while looking for the optimal thresholds, we stated that they are not data but feature dependent. In other words, the best threshold varies a lot when the feature dimensionality changes but remain quite similar otherwise.

Second, the most common threshold estimation strategy (3), along with the removal of a predefined amount of data at

Method	INRIA	Speed(CPU/GPU)	Method	CALETCH	Speed(CPU/GPU)
HoG [16]	46%	0.5fps	HoG [16]	69%	0.5fps
HoG-LBP [59]	39%	Not provided	DPM [24]	63.26%	< 1fps
MultiFeatures [60]	36%	< 1fps	FeatSynth [2]	60.16%	< 1fps
FeatSynth [2]	31%	< 1fps	MultiFeatures+CSS [57]	60.89%	No
MultiFeatures+CSS [57]	25%	No	FPDW [17]	57.4%	2-5fps
Channel Features [18]	21%	0.5fps	Channel Features [18]	56.34%	0.5fps
FPDW [17]	21%	2-5fps	Roerei [4]	48.35%	1 fps
DPM [24]	20%	< 1fps	MOCO [13]	45.5%	< 1fps
RF local experts [43]	15.4%	3fps	JointDeep [47]	39.32%	< 1fps
PCA-CNN [36]	14.24%	< 0.1fps	InformedHaar [66]	34.6%	< 0.63fps
<b>Ours - HoG</b>	<b>19.01%</b>	<b>4/60fps</b>	katamari-v1 [5]	22.49%	< 1fps
VeryFast [3]	18%	8/135fps	<b>FRCNN [51]</b>	56%	7fps
WordChannels [15]	17%	0.5/8fps	CrossTalk cascades [19]	53.88%	30-60fps
crossTalk cascades [19]	17%	30-60fps	<b>Ours - LBP</b>	<b>51.12%</b>	<b>3.6/54fps</b>
<b>Ours - LBP</b>	<b>17.28%</b>	<b>3.7/54fps</b>	<b>Ours - HoG</b>	<b>45.4%</b>	<b>3.9/58fps</b>
<b>Ours - HoG+LBP</b>	<b>17.08%</b>	<b>2.6/39fps</b>	<b>Ours - HoG+LBP</b>	<b>45.2%</b>	<b>1.8/27fps</b>
SSD [38]	15%	56fps	WordChannels [15]	42.3%	0.5/8fps
FRCNN [51]	13%	7fps	SSD [38]	34%	56fps
RPN+PF [37]	7%	6fps	RPN+PF [37]	10%	6fps

TABLE VI

COMPARISON WITH THE STATE-OF-THE-ART. NEAR REAL-TIME METHODS ARE SEPARATED FROM OTHERS. *Ours* IS IN BOLD. DEEP LEARNING TECHNIQUES ARE IN RED. COMPUTATION TIMES ARE CALCULATED ACCORDING TO 640×480 RESOLUTION FRAMES. THE USED METRIC IS THE LOG-AVERAGE MISS RATE (THE LOWER THE BETTER).

Soft cascade type	Feature	INRIA	Caltech
No soft-cascade	HoG	20.31% / 2.1fps	46.25% / 1.9fps
soft-cascade	HoG	32.21% / 18.1fps	58.26% / 17.8fps
block soft-cascade	HoG	27.86% / 9.9fps	53.97% / 10.1fps
reverse engineering	HoG	20.32% / 4 fps	<b>46.23%</b> / 3.9fps
No soft-cascade	LBP	18.42% / 1.9fps	51.36% / 1.7fps
soft-cascade	LBP	30.87% / 17.8fps	61.32% / 17.1fps
block soft-cascade	LBP	28.33% / 9.8fps	55.45% / 9.5fps
reverse engineering	LBP	<b>18.41%</b> / 3.7fps	51.33% / 3.6fps

TABLE VII

LAMR PERFORMANCE AND CPU SPEED IN FRAME-PER-SECOND OF DIFFERENT SOFT-CASCADE STRATEGIES. SEE TEXT FOR THE DESCRIPTION OF THE METHODS. BEST RESULTS IN BOLD.

each step (4) perform poorly when used with these classifiers settings. The best performing thresholding technique uses a constant threshold (1). This technique leads to an average deletion of respectively 80% and 35% of the data during the first 2 steps of the cascade, and 5-10% during the remaining stages. Hence, we assume that the thresholding technique (3) and (4) might also be dependent to classifier sets displaying a rather regular deletion behaviour across the cascade steps.

Finally, all the tested techniques, are, on average performing 4.48% worse than the optimal. This elects threshold estimation as a potentially rewarding research direction.

**lesson learned:** Counter-intuitively, the variability between the threshold values is not much caused by the data, but mostly by the features. None of existing approaches are close to the optimal and there is still space for improvement in that particular direction.

J. Non maximum suppression

Table IX covers our evaluation the 3 existing non-maximum suppression algorithms previously described in section III-G. The best results for both HoG and LBP features are always obtained by the threshold non-maximum suppression with an average 2% boost compared to the common greedy non-maximum suppression.

**lesson learned:** Surprisingly, the choice of the non maximum suppression algorithm significantly impacts the over-

Thresholding strategy	Feature	INRIA	Caltech
Optimal	HoG	17.89%	42.67%
Constant	HoG	20.23%	46.25%
Increasing	HoG	21.44%	47.25%
Recall	HoG	25.01%	50.88%
Regular	HoG	27.11%	51.84%
Optimal	LBP	<b>15.44%</b>	<b>42.45%</b>
Constant	LBP	18.63%	51.26%
Increasing	LBP	19.96%	51.78%
Recall	LBP	22.34%	54.88%
Regular	LBP	22.36%	62.53%

TABLE VIII

THRESHOLD ESTIMATION TECHNIQUES EVALUATION ON INRIA AND CALTECH BENCHMARKS. SEE TEXT FOR THE DESCRIPTION OF THE METHODS. BEST RESULTS IN BOLD.

NMS type	Feature	INRIA	Caltech
Greedy-NMS	HoG	22.91%	49.62%
t-NMS	HoG	20.23%	46.25%
s-NMS	HoG	21.47%	<b>46.98%</b>
Greedy-NMS	LBP	20.21%	52.32%
t-NMS	LBP	<b>18.63%</b>	51.26%
s-NMS	LBP	19.87%	52.65%

TABLE IX

NON-MAXIMUM SUPPRESSION ALGORITHM EVALUATION ON INRIA AND CALTECH BENCHMARKS. BEST RESULTS IN BOLD.

all detector performance. The over-employed greedy non-maximum suppression, that underperforms in every case, shouldn't be blindly chosen.

K. Comparison with the state-of-the-art

Table VI shows our method's ranking compared to the state-of-the-art, in terms of performance and speed. Table I parameters and random forest were employed for this experiment.  $T=0.58$  is used for these runs. This work compares favorably to the state-of-the-art. We used using Adaboost classifiers and soft coding reverse engineering for this experiment. While not being the best detector on the market, with respectively 17% and 46% log-average miss rate on the INRIA and Caltech datasets, our approach still provides a very competitive performance/speed ratio. For instance, it scores better than

the famous DPM [24] and similarly to the integral channel features [17] while running 7 times faster. VeryFast [3] and the crosstalk cascades provide a viable alternative in terms of speed while displaying lower performance. LBP is the best descriptor on the INRIA dataset, while HoGs show the best performance on Caltech. We assume that the numerous little pedestrians on the latter impact the texture descriptor more strongly than the gradients. The HoG and LBP fusion yields little improvement. Packed crowds, teeny pedestrians and awkward poses remain the main failure cases. Finally, Deep learning technique [36], [51] perform worse and are slower, [37] performs much better but is slower, while [38] is slightly more efficient and faster. It is worth stating that careful machine learning choices and parametrisation can bring standard features close to state-of-the-art deep learning performance. The end-to-end system is trained in 4-14hours on one core for 10-200K data and processes 3-4 fps on an Intel Xeon 2.1GHz CPU, calculated over frames of  $640 \times 480$  pixels in size. Further speed up is possible when using a GPU, leading our detector to perform in real-time. This makes this detector one of the best performing real-time pedestrian detector to date.

## V. CONCLUSION

This paper studied through extensive experimentation underexplored mechanisms at various stages of a pedestrian classifier training, with an emphasis on data selection, dataset generation at the initial stage and during the classifier cascade. Our study highlighted data selection and non-maximum suppression as underexploited components for high performance, and soft-cascade thresholding assessment as potential research direction. We also disproved a few empirically received ideas, namely the predominance of the random forest classifier (with high performing detectors) or the use of the DBSCAN algorithm for data cleaning.

Besides extra experiments in domains such as active learning or part-based models, the application of the optimized *Fair-Train* pipeline to deep learning is an obvious possible next step for this work.

## VI. ACKNOWLEDGMENT

To appear in camera ready version.

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