

Mistake-Driven Learning in Text Categorization

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Abstract

Learning problems in the text processing domain often map the text to a space whose dimensions are the measured features of the text, e.g., its words. Three characteristic properties of this domain are (a) very high dimensionality, (b) both the learned concepts and the instances reside very sparsely in the feature space, and (c) a high variation in the number of active features in an instance. In this work we study three mistake-driven learning algorithms for a typical task of this nature – text categorization.

We argue that these algorithms – which categorize documents by learning a linear separator in the feature space – have a few properties that make them ideal for this domain. We then show that a quantum leap in performance is achieved when we further modify the algorithms to better address some of the specific characteristics of the domain. In particular, we demonstrate (1) how variation in document length can be tolerated by either normalizing feature weights or by using negative weights, (2) the positive effect of applying a threshold range in training, (3) alternatives in considering feature frequency, and (4) the benefits of discarding features while training.

Overall, we present an algorithm, a variation of Littlestone’s Winnow, which performs significantly better than any other algorithm tested on this task using a similar feature set.

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1 Introduction

Learning problems in the natural language and text processing domains are often studied by mapping the text to a space whose dimensions are the measured features of the text, e.g., the words appearing in a document. Three characteristic properties of this domain are (a) very high dimensionality, (b) both the learned concepts and the instances reside very sparsely in the feature space and, consequently, (c) there is a high variation in the number of active features in an instance.

Multiplicative weight-updating algorithms such as Winnow (Littlestone, 1988) have been studied extensively in the theoretical learning literature. Theoretical analysis has shown that they have exceptionally good behavior in domains with these characteristics, and in particular in the presence of irrelevant attributes, noise, and even a target function changing in time (Littlestone, 1988; Littlestone and Warmuth, 1994; Herbster and Warmuth, 1995), but only recently have people started to use them in applications (Golding and Roth, 1996; Lewis et al., 1996; Cohen and Singer, 1996). We address these claims empirically in an important application domain for machine learning – text categorization. In particular, we study mistake-driven learning algorithms that are based on the Winnow family, and investigate ways to apply them in domains with the above characteristics.

The learning algorithms studied here offer a large space of choices to be made and, correspondingly, may vary widely in performance when applied in specific domains. We concentrate here on the text processing domain, with the characteristics mentioned above, and explore this space of choices in it.

In particular, we investigate three variations of on-line prediction algorithms and evaluate them experimentally on large text categorization problems. The algorithms we study are all learning algorithms for linear functions. They are used to categorize documents by learning, for each category, a linear separator in the feature space. The algorithms differ by whether they allow the use of negative or only

positive weights and by the way they update their weights during the training phase.

We find that while a vanilla version of these algorithms performs rather well, a quantum leap in performance is achieved when we modify the algorithms to better address some of the specific characteristics we identify in textual domains. In particular, we address problems such as wide variations in document sizes, word repetitions and the need to rank documents rather than just decide whether they belong to a category or not. In some cases we adopt solutions that are well known in the IR literature to the class of algorithms we use; in others we modify known algorithms to better suit the characteristics of the domain. We motivate the modifications to the basic algorithms and justify them experimentally by exhibiting their contribution to improvement in performance. Overall, the best variation we investigate, performs significantly better than any known algorithm tested on this task, using a similar set of features.

The rest of the paper is organized as follows: The next section describes the task of text categorization, how we model it as a classification task, and some related work. The family of algorithms we use is introduced in Section 3 and the extensions to the basic algorithms, along with their experimental evaluations, is presented in Section 4. In Section 5 we present our final experimental results and compare them to previous works in the literature.

2 Text Categorization

In text categorization, given a text document and a collection of potential classes, the algorithm decides which classes it belongs to, or how strongly it belongs to each class. For example, possible classes (categories) may be $\{bond\}$, $\{loan\}$, $\{interest\}$, $\{acquisition\}$. Documents that have been categorized by humans are usually used as training data for a text categorization system; later on, the trained system is used to categorize new documents. Algorithms used to train text categorization systems in information retrieval (IR) are often ad-hoc and poorly understood. In particular, very little is known about their generalization performance, that is, their behavior on documents outside the training data. Only recently, some machine learning techniques for training linear classifiers have been used and shown to be effective in this domain (Lewis et al., 1996; Cohen and Singer, 1996). These techniques have the advantage that they are better understood from a theoretical standpoint, leading to performance guarantees and guidance in parameter settings. Continuing this line of research we present different algorithms and focus on adjusting them to the unique characteristics of the domain, yielding good performance on the categorization task.

2.1 Training Text Classifiers

Text classifiers represent a document as a set of features $d = \{f_1, f_2, \dots, f_m\}$, where m is the number of *active* features in the document, that is, features that occur in the document. A feature f_i may typically represent a word w , a set w_1, \dots, w_k of words (Cohen and Singer, 1996) or a phrasal structure (Lewis, 1992; Tzeras and Hartmann, 1993). The *strength* of the feature f in the document d is denoted by $s(f, d)$. The strength is usually a function of the number of times f appears in d (denoted by $n(f, d)$). The strength may be used only to indicate the presence or absence of f in the document, in which case it takes on only the values 0 or 1, it may be equal to $n(f, d)$, or it can take other values to reflect also the size of the document.

In order to rank documents, for each category, a text categorization system keeps a function F_c which, when evaluated on d , produces a score $F_c(d)$. A decision is then made by assigning to the category c only those documents that exceed some threshold, or just by placing at the top of the ranking documents with the highest such score.

A *linear* text classifier represents a category as a weight vector $w_c = (w(f_1, c), w(f_2, c), \dots, w(f_n, c)) \equiv (w_1, w_2, \dots, w_n)$, where n is the total number of features in the domain and $w(f, c)$ is the weight of the feature f for this category. It evaluates the score of the document by computing the dot product:

$$F_c(d) = \sum_{f \in d} s(f, d) \cdot w(f, c).$$

The problem is modeled as a supervised learning problem. The algorithms use the training data, where each document is labeled by zero or more categories, to learn a classifier which classifies new texts. A document is considered as a positive example for all categories with which it is labeled, and as a negative example to all others.

The task of a training algorithm for a linear text classifier is to find a weight vector which best classifies new text documents. While a linear text classifier is a linear separator in the space defined by the features, it may not be linear with respect to the document, if one chooses to use complex features such as conjunctions of simple features. In addition, a training algorithm may give also advice on the issue of feature selection, by reducing the weight of non-important features and thus effectively discarding them.

2.2 Related Work

Many of the techniques previously used in text categorization make use of linear classifiers, mainly for reasons of efficiency. The classical vector space model, which ranks documents using a nonlinear similarity measure (the “cosine correlation”) (Salton and Buckley, 1983) can also be recast as a linear classification by incorporating length normalization into

the weight vector and the document vector features values. State of the art IR systems determine the strength of a term based on three values: (1) the frequency of the feature in the document (*tf*), (2) an inverse measure of the frequency of the feature throughout the data set (*idf*), and (3) a normalization factor that takes into account the length of the document. In Sections 4.1 and 4.3 we discuss how we incorporate those ideas in our setting.

Most relevant to our work are non-parametric methods, which seem to yield better results than parametric techniques. Rocchio's algorithm (Rocchio, 1971), one of the most commonly used techniques, is a batch method that works in a relevance feedback context. Typically, classifiers produced by the Rocchio algorithm are restricted to having non-negative weights. An important distinction between most of the classical non-parametric methods and the learning techniques we study here is that in the former case, there was no theoretical work that addressed the generalization ability of the learned classifier, that is, how it behaves on new data.

The methods that are most similar to our techniques are the on-line algorithms used in (Lewis et al., 1996) and (Cohen and Singer, 1996). In the first, two algorithms, a multiplicative update and additive update algorithms suggested in (Kivinen and Warmuth, 1995a) are evaluated in the text categorization domain, and are shown to perform somewhat better than Rocchio's algorithm. While both these works make use of multiplicative update algorithms, as we do, there are two major differences between those studies and the current one. First, there are some important technical differences between the algorithms used. Second, the algorithms we study here are mistake-driven; they update the weight vector only when a mistake is made, and not after every example seen. The Experts algorithm studied in (Cohen and Singer, 1996) is very similar to a basic version of the BalancedWinnow algorithm which we study here. The way we treat the negative weights is different, though, and significantly more efficient, especially in sparse domains (see Section 3.1). Cohen and Singer experiment also, using the same algorithm, with more complex features (sparse n-grams) and show that, as expected, it yields better results.

Our additive update algorithm, Perceptron, is somewhat similar to what is used in (Wiener, Pedersen, and Weigend, 1995). They use a more complex representation, a multi-layer network, but this additional expressiveness seems to make training more complicated, without contributing to better results.

2.3 Methodology

We evaluate our algorithms on the the Reuters-22173 text collection (Lewis, 1992), one of the most commonly used benchmarks in the literature.

For the experiments reported In Sections 3.2 we explore and compare different variations of the al-

gorithms; we evaluate those on two disjoint pairs of a training set and a test set, both subsets of the Reuters collection. Each pair consists of 2000 training documents and 1000 test documents, and was used to train and test the classifier on a sample of 10 topical categories. The figures reported are the average results on the two test sets.

In addition, we have tested our final version of the classifier on two common partitions of the complete Reuters collection, and compare the results with those of other works. The two partitions used are those of Lewis (Lewis, 1992) (14704 documents for training, 6746 for testing) and Apte (Apte, Damerau, and Weiss, 1994) (10645 training, 3672 testing, omitting documents with no topical category).

To evaluate performance, the usual measures of recall and precision were used. Specifically, we measured the effectiveness of the classification by keeping track of the following four numbers:

- p_1 = number of correctly classified class members
- p_2 = number of mis-classified class members
- n_1 = number of correctly classified non-class members
- n_2 = number of mis-classified non-class members

In those terms, the *recall* measure is defines as p_1/p_1+p_2 , and the *precision* is defined as p_1/p_1+n_1 . Performance was further summarized by a *break-even point* – a hypothetical point, obtained by interpolation, in which precision equals recall.

3 On-Line learning algorithms

In this section we present the basic versions of the learning algorithms we use. The algorithms are used to learn a classifier F_c for each category c . These algorithms use the training data, where each document is labeled by zero or more categories, to learn a weight vector which is used later on, in the test phase, to classify new text documents. A document is considered as a positive example for all categories with which it is labeled, and as a negative example to all others. The algorithms are on-line and mistake-driven. In the on-line learning model, learning takes place in a sequence of trials. On each trial, the learner first makes a prediction and then receives feedback which may be used to update the current hypothesis (the vector of weights). A mistake-driven algorithm updates its hypothesis only when a mistake is made. In the training phase, given a collection of examples, we may repeat this process a few times, by iterating on the data. In the testing phase, the same process is repeated on the test collection, only that the hypothesis is not updated.

Let n be the number of features of the current category. For the remainder of this section we denote a training document with m active features

by $d = (s_{i_1}, s_{i_2}, \dots, s_{i_m})$, where s_{i_j} stands for the strength of the i_j feature. The label of the document is denoted by y ; y takes the value 1 if the document is relevant to the category and 0 otherwise. Notice, that we care only about the active features in the domain, following (Blum, 1992). The algorithms have three parameters: a threshold θ , and two update parameters, a *promotion* parameter α and a *demotion* parameter β .

Positive Winnow (Littlestone, 1988):

The algorithm keeps an n -dimensional weight vector $w = (w_1, w_2, \dots, w_n)$, w_i being the weight of the i th feature, which it updates whenever a mistake is made. Initially, the weight vector is typically set to assign equal positive weight to all features. (We use the value θ/d , where d is the average number of active features in a document; in this way initial scores are close to θ .) The promotion parameter is $\alpha > 1$ and the demotion is $0 < \beta < 1$.

For a given instance $(s_{i_1}, s_{i_2}, \dots, s_{i_m})$ the algorithm predicts 1 iff

$$\sum_{j=1}^m w_{i_j} s_{i_j} > \theta,$$

where w_{i_j} is the weight corresponding to the active feature indexed by i_j . The algorithm updates its hypothesis only when a mistake is made, as follows: (1) If the algorithm predicts 0 and the label is 1 (positive example) then the weights of all the active features are promoted — the weight w_{i_j} is multiplied by α . (2) If the algorithm predicts 1 and the received label is 0 (negative example) then the weights of all the active features are demoted — the weight w_{i_j} is multiplied by β . In both cases, weights of inactive features maintain the same value.

Perceptron (Rosenblatt, 1958)

As in PositiveWinnow, in Perceptron we also keep an n -dimensional weight vector $w = (w_1, w_2, \dots, w_n)$ whose entries correspond to the set of potential features, which is updated whenever a mistake is made. As above, the initial weight vector is typically set to assign equal weight to all features. The only difference between the algorithms is that in this case the weights are updated in an *additive* fashion. A single update parameter $\alpha > 0$ is used, and a weight is promoted by *adding* α to its previous value, and is demoted by *subtracting* α from it. In both cases, all other weights maintain the same value.

Balanced Winnow (Littlestone, 1988):

In this case, the algorithm keeps two weights, w^+, w^- , for each feature. The overall weight of a feature is the difference between these two weights, thus allowing for negative weights. For a given instance $(s_{i_1}, s_{i_2}, \dots, s_{i_m})$ the algorithm predicts 1 iff

$$\sum_{j=1}^m (w_{i_j}^+ - w_{i_j}^-) s_{i_j} > \theta, \quad (1)$$

where $w_{i_j}^+, w_{i_j}^-$ correspond to the active feature indexed by i_j . In our implementation, the weights w^+ are initialized to $2\theta/d$ and the weights w^- are set to θ/d , where d is the average number of active features in a document in the collection.

The algorithm updates the weights of active features only when a mistake is made, as follows: (1) In the promotion step, following a mistake on a positive example, the positive part of the weight is promoted, $w_{i_j}^+ \leftarrow \alpha \cdot w_{i_j}^+$ while the negative part of the weight is demoted, $w_{i_j}^- \leftarrow \beta \cdot w_{i_j}^-$. Overall, the coefficient of s_{i_j} in Eq. 1 increases after a promotion. (2) In the demotion step, following a mistake on a negative example, the coefficient of s_{i_j} in Eq. 1 is decreased: the positive part of the weight is demoted, $w_{i_j}^+ \leftarrow \beta \cdot w_{i_j}^+$ while the negative part of the weight is promoted, $w_{i_j}^- \leftarrow \alpha \cdot w_{i_j}^-$. In both cases, all other weights maintain the same value.

In this algorithm (see in Eq. 1) the coefficient of the i th feature can take negative values, unlike the representation used in PositiveWinnow. There are other versions of the Winnow algorithm that allow the use of negative features: (1) Littlestone, when introducing the Balanced version, introduced also a simpler version — a version of PositiveWinnow with a duplication of the number of features. (2) A version of the Winnow algorithm with negative features is used in (Cohen and Singer, 1996). In both cases, however, whenever there is a need to update the weights, *all* the weights are being updated (actually, n out of the $2n$). In the version we use, only weights of *active* features are being updated; this gives a significant computational advantage when working in a sparse high dimensional space.

3.1 Properties of the Algorithms

Winnow and its variations were introduced in Littlestone’s seminal paper (Littlestone, 1988); the theoretical behavior of multiplicative weight-updating algorithms for learning linear functions has been studied since then extensively. In particular, Winnow has been shown to learn efficiently any linear threshold function (Littlestone, 1988). These are functions $F : \{0, 1\}^n \rightarrow \{0, 1\}$ for which there exist real weights w_1, \dots, w_n and a real threshold θ such that $F(s_1, \dots, s_n) = 1$ iff $\sum_{i=1}^n w_i s_i \geq \theta$. In particular, these functions include Boolean disjunctions and conjunctions on $k \leq n$ variables and r -of- k threshold functions ($1 \leq r \leq k \leq n$). While Winnow is guaranteed to find a perfect separator if one exists, it also appears to be fairly successful when there is no perfect separator. The algorithm makes no independence or any other assumptions on the features, in contrast to other parametric estimation techniques (typically, Bayesian predictors) which are commonly used in statistical NLP.

Theoretical analysis has shown that the algorithm has exceptionally good behavior in the presence of

irrelevant features, noise, and even a target function changing in time (Littlestone, 1988; Littlestone, 1991; Littlestone and Warmuth, 1994; Herbster and Warmuth, 1995), and there is already some empirical support for these claims (Littlestone, 1995; Golding and Roth, 1996; Blum, 1995). The key feature of Winnow is that its mistake bound grows linearly with the number of *relevant* features and only logarithmically with the total number of features. A second important property is being mistake driven. Intuitively, this makes the algorithm more sensitive to the relationships among the features — relationships that may go unnoticed by an algorithm that is based on counts accumulated separately for each attribute. This is crucial in the analysis of the algorithm as well as empirically (Littlestone, 1995; Golding and Roth, 1996).

The discussion above holds for both versions of Winnow studied here, PositiveWinnow and BalancedWinnow. The theoretical results differ only slightly in the mistake bounds, but have the same flavor. However, the major difference between the two algorithms, one using only positive weights and the other allowing also negative weights, plays a significant role when applied in the current domain, as discussed in Section 4.

Winnow is closely related, and has served as the motivation for a collection of recent works on combining the “advice” of different “experts” (Littlestone and Warmuth, 1994; Cesa-Bianchi et al., 1995; Cesa-Bianchi et al., 1994). The features used are the “experts” and the learning algorithm can be viewed as an algorithm that learns how to combine the classifications of the different experts in an optimal way.

The additive-update algorithm that we evaluate here, the Perceptron, goes back to (Rosenblatt, 1958). While this algorithm is also known to learn the target linear function when it exists, the bounds given by the Perceptron convergence theorem (Duda and Hart, 1973) may be exponential in the optimal mistake bound, even for fairly simple functions (Kivinen and Warmuth, 1995b). We refer to (Kivinen and Warmuth, 1995a) for a thorough analysis of multiplicative update algorithms versus additive update algorithms. In particular, it is shown that the number of mistakes the additive and multiplicative update algorithms make, depend differently on the domain characteristics. Informally speaking, it is shown that the multiplicative update algorithms have advantages in high dimensional problems (i.e., when the number of features is large) and when the target weight vector is sparse (i.e., contain many weights that are close to 0). This explains the recent success in using these methods on high dimensional problems (Golding and Roth, 1996) and suggests that multiplicative-update algorithms might do well on IR applications, provided that a good set of features is selected. On the other hand, it is

shown that additive-update algorithms have advantages when the examples are sparse in the feature space, another typical characteristics of the IR domain, which motivates us to study experimentally an additive-update algorithm as well.

3.2 Evaluating the Basic Versions

We started by evaluating the basic versions of the three algorithms. The features we use throughout the experiments are single words, at the lemma level, for nouns and verbs only, with minimal frequency of 3 occurrences in the corpus. In the basic versions the strength of the feature is taken to indicate only the presence or absence of f in the document, that is, it is either 1 or 0. The training algorithm was run iteratively on the training set, until no mistakes were made on the training collection or until some upper bound (50) on the number of iterations was reached. The results for the basic versions are shown in the first column of Table 1.

4 Extensions to the Basic algorithms

4.1 Length Variation and Negative features

Text documents vary widely in their length and a text classifier needs to tolerate this variation. This issue is a potential problem for a linear classifier which scores a document by summing the weights of all its active features: a long document may have a better chance of exceeding the threshold merely by its length.

This problem has been identified earlier on and attracted a lot of work in the classical work on IR (Salton and Buckley, 1983), as we have indicated in Section 2.2. The treatment described there addresses at the same time at least two different concerns: length variation of documents and feature repetition. In this section we consider the first of those, and discuss how it applies to the algorithms we investigate. The second concern is discussed in Section 4.3.

Algorithms that allow the use of negative features, such as BalancedWinnow and Perceptron, tolerate variation in the documents length naturally, and thus have a significant advantage in this respect. In these cases, it can be expected that the cumulative contribution of the weights and, in particular, those that are not indicative to the current category, does not count towards exceeding the threshold, but rather averages out to 0. Indeed, as we found out, no special normalization is required when using these algorithms. Their significant advantage over the unnormalized version of PositiveWinnow is readily seen in Table 1.

In addition, using negative weights gives the text classifier more flexibility in capturing “truly negative” features, where the presence of a feature is indicative for the *irrelevance* of the document to the category. However, we found that this phenomenon

Algorithm	Version					
	Basic	Norm	θ -range	Linear Freq.	Sqrt. Freq.	Discard
BalancedWinnow	64.87	NA	69.66	72.11	71.56	73.2
PositiveWinnow	55.56	63.56	65.80	67.20	69.67	70.0
Perceptron	65.91	NA	63.05	66.72	68.29	70.8

Table 1: Recall/precision break-even point (in percentages) for different versions of the algorithm. Each figure is an average result for two pairs of training and testing sets, each containing 2000 training documents and 1000 test documents.

only rarely occurs in text categorization and thus the main use of the negative features is to tolerate the length variation of the documents.

When using PositiveWinnow, which uses only positive weights, we no longer have this advantage and we seek a modification that tolerates the variation in length. As in the standard IR solution, we suggest to modify $s(f, d)$, the strength of the feature f in d , by using a quantity that is normalized with respect to the document size.

Formally, we replace the strength $s(f, d)$ (which may be determined in several ways according to feature frequency, as explained below) by a *normalized strength*,

$$s^n(f, d) = \frac{s(f, d)}{\sum_{f \in d} s(f, d)}.$$

In this case (which applies, as discussed above, only for PositiveWinnow), we also change the initial weight vector and initialize all the weights to θ .

Using normalization gives an effect that is similar to the use of negative weights, but to a lesser degree. The reason is that it is used uniformly; in long documents, the number of indicative features does not increase significantly, but their strength, nevertheless, is reduced proportionally to the total number of features in the document. In the long version of the paper we present a more thorough analysis of this issue.

The results presented in Table 1 (second column) show the significant improvements achieved in PositiveWinnow performance, when normalization is used. In all the results presented from this point on, positive winnow is normalized.

4.2 Using Threshold range

Training a linear text classifier is a search for a weight vector in the feature space. The search is for a linear separator that best separates documents that are relevant to the category from those that are not. In general, there is no guarantee that a weight vector of this sort exists, even in the training data, but a good selection of features make this more likely. While the basic versions of our algorithms search for linear separators, we have modified those so that our search for a linear classifier is biased to look for “thick” classifiers. To understand this, consider, for

the moment, the case in which all the data is perfectly linearly separable. Then there will generally be many linear classifiers that separate the training data we actually see. Among these, it seems plausible that we have a better chance of doing well on the unseen test data if we choose a linear separator that separates the positive and negative training examples as “widely” as possible. The idea of having a wide separation is less clear when there is no perfect separator, but we can still appeal to the basic intuition.

Using a “thick” separator is even more important when documents are ranked rather than simply classified; that is, when the actual score produced by the classifier is used in the decision process. The reason is that if $F_c(d)$ is the score produced by the classifier F_c when evaluated on the document d then, under some assumptions on the dependencies among the features, the probability that the document d is relevant to the category c is given by $Prob(d \in c) = \frac{1}{1+e^{-F_c(d)}}$. This function, known as the *sigmoid* function, “flattens” the decision region in a way that only scores that are far apart from the threshold value indicate that the decision is made with significant probability.

Formally, among those weight vectors we would like to choose the hyper-plane with the largest “separating parameter”, where the separating parameter τ is defined as the largest value for which there exists a classifier F_c (defined by a weight vector w) such that for all positive examples d , $F_c(d) > \theta + \tau/2$ and for all negative d , $F_c(d) < \theta - \tau/2$.

In this implementation we do not try to find the optimal τ (as is done in (Cortes and Vapnik, 1995)), but rather determine it heuristically. In order to find a “thick” separator, we modify, in all three algorithms, the update rule used during the training phase as follows: Rather than using a single threshold we use two separate thresholds, θ^+ and θ^- , such that $\theta^+ - \theta^- = \tau$. During training, we say that the algorithm predicts 0 (and makes a mistake, if the example is labeled positive) when the score it assigns an example is below θ^- . Similarly, we say that the algorithm predicts 1 when the score exceeds θ^+ . All examples with scores in the range $[\theta^-, \theta^+]$ are considered mistakes. (Parameters used: $\theta^- = 0.9$, $\theta^+ = 1.1$, $\theta = 1$).

The results presented in the third column of Table 1 show the improvements obtained when the threshold range is used. In all the results presented from this point on, all the algorithms use the threshold range modification.

4.3 Feature Repetition

Due to the bursty nature of term occurrence in documents, as well as the variation in document length, a feature may occur in a document more than once. It is therefore important to consider the frequency of a feature when determining its strength. On one hand, there are cases where a feature is more indicative to the relevance of the document to a category when it appears several times in a document. On the other hand, in any long document, there may be some random feature that is not significantly indicative to the current category although it repeats many times. While the weight of f in the weight vector of the category, $w(f, c)$, may be fairly small, its cumulative contribution might be too large if we increase its strength, $s(f, d)$, in proportion to its frequency in the document.

As mentioned in Section 2.2, the classical IR literature has addressed this problem using the *tf* and *idf* factors. We note that the standard treatment in IR suggests a solution to this problem that suits batch algorithms – algorithms that determine the weight of a feature after seeing all the examples. We, on the other hand, seek a solution that can be used in an on-line algorithm. Thus, the frequency of a feature throughout the data set, for example, cannot be taken into account and we take into account only the *tf* term. We have experimented with three alternative ways of adjusting the value of $s(f, d)$ according to the frequency of the feature in the document: (1) Our default is to let the strength indicate only the activity of the feature. That is, $s(f, d) = 1$, if the feature is present in the document (active feature) and $s(f, d) = 0$ otherwise. (2) $s(f, d) = n(f, d)$, where $n(f, d)$ is the number of occurrences of f in d ; and (3) $s(f, d) = \sqrt{n(f, d)}$ (as in (Wiener, Pedersen, and Weigend, 1995)). These three alternatives examine the tradeoff between the positive and negative impacts of assigning a strength in proportion to feature frequency. In most of our experiments, on different data sets, the choice of using $\sqrt{n(f, d)}$ performed best. The results of the comparative evaluation appear in columns 3, 4, and 5 of Table 1, corresponding to the three alternatives above.

4.4 Discarding features

Multiplicative update algorithm are known to tolerate a very large number of features. However, it seems plausible that most categories depend only on fairly small subsets of indicative features and not on all the features that occur in documents that belong to this class. Efficiency reasons, as well as the occasional need to generate comprehensible explanations

to the classifications, suggest that discarding irrelevant features is a desirable goal in IR applications. If done correctly, discarding irrelevant features may also improve the accuracy of the classifier, since irrelevant features contribute noise to the classification score.

An important property of the algorithms investigated here is that they do not require a feature selection pre-processing stage. Instead, they can run in the presence of a large number of features, and allow for discarding features “on the fly”, based on their contribution to an accurate classification. This property is especially important if one is considering enriching the set of features, as is done in (Golding and Roth, 1996; Cohen and Singer, 1996); in these cases it is important to allow the algorithm to decide for itself which of the features contribute to the accuracy of the classification.

We filter features that are irrelevant for the category based on the weights they were assigned in the first few training rounds.

The algorithm is given as input a range of weight value which we call the *filtering range*. First, the training algorithm is run for several iterations, until the number of mistakes on the training data drops below a certain threshold. After this initial training, we filter out all the features whose weight lie in this filtering range. Training then continues as usual.

There are various ways to determine the filtering range. The obvious one may be to filter out all features whose weight is very close to 0, but there are a few subtle issues involved due to the normalization done in the PositiveWinnow algorithm. In the results presented here we have used, instead, a different filtering range: Our filtering range is centered around the initial value assigned to the weights (as specified earlier for each algorithm), and is bounded above and below by the values obtained after one promotion or demotion step relative to the initial value. Thus, with high likelihood, we discard features which have not contributed to many mistakes – those that were promoted or demoted at most once (possibly, with additional promotions and demotions which canceled each other, though).

The results of classification with feature filtering appear in the last column of Table 1. We hypothesize that the improved results are due to reduction in the noise introduced by irrelevant features. Further investigation of this issue will be presented in the long version of this paper. Typically, about two thirds of the features were filtered for each category, significantly reducing the output representation size.

5 Summary of Experimental Results

The study described in Section 3.2 was used to determine the version that performs best, out of those we have experimented with. Eventually, we have selected the version of the BalancedWin-

Algorithm	Apte's split	Lewis's split
BalancedWinnow⁺	83.3	74.7
Experts unigram (Cohen and Singer, 1996)	64.7	65.6
Neural Network (Wiener, Pedersen, and Weigend, 1995)	77.5	NA
Rocchio (Rocchio, 1971)	74.5	66.0
Ripper (Cohen and Singer, 1996)	79.6	71.9
Decision trees (Lewis and Ringuette, 1994)	NA	67.0
Bayes (Lewis and Ringuette, 1994)	NA	65.0
SWAP (Apte, Damerau, and Weiss, 1994)	78.9	NA

Table 2: Break-even points comparison. The data is split into training set and test set based on Lewis's split - (Lewis, 1992), 14704 documents for training, 6746 for testing, and Apte's split - (Apte, Damerau, and Weiss, 1994), 10645 training, 3672 testing, omitting documents with no topical category.

now algorithm, which incorporates the θ -range modification, a square-root of occurrences as the feature strength and the discard features modification (BalancedWinnow⁺ in Table 2).

We have compared this version with a few other algorithms which have appeared in the literature on the complete Reuters corpus. Table 2 presents break-even points for BalancedWinnow⁺ and the other algorithms, as defined in Section 2.3.

The results are reported for two splits of the complete Reuters corpus as explained in Section 2.3. The algorithm was run with iterations, threshold range, feature filtering, and frequency-square-root feature strength.

The first two rows in Table 2 compare the performance of BalancedWinnow⁺ with the two algorithms that most resemble our approach, the Experts algorithm from (Cohen and Singer, 1996) and a neural network approach presented in (Wiener, Pedersen, and Weigend, 1995). (see Section 2.2).

Rocchio's algorithm is one of the classical algorithms for this tasks, and it still performs very good compared to newly developed techniques (e.g, (Lewis et al., 1996)). We also compared with the Ripper algorithm presented in (Cohen and Singer, 1996) (we present the best results for this task, with negative tests), a simple decision tree learning system and a Bayesian classifier. The last two figure are taken from (Lewis and Ringuette, 1994) where they were evaluated only on Lewis's split. The last comparison is with the learning system used by (Apte, Damerau, and Weiss, 1994), SWAP, which was evaluated only on Apte's split.

Our results significantly outperform (by at least 2-4%) all results which appear in that table and use the same set of features (based on single words). Of the results we know of in the literature, only a version of the Experts algorithm of (Cohen and Singer, 1996) which uses a richer feature set - sparse word trigrams - outperforms our result on the Lewis split, with a break-even point of 75.3%, compared with 74.6% for the unigram-based BalancedWinnow⁺. However,

this version achieves only 75.9% on the Apte split (compared with 83.3% of BalancedWinnow⁺). In the long version of this paper we plan to present the results of our algorithm on a richer feature set as well.

6 Conclusions

Theoretical analyses of the Winnow family of algorithms have predicted an exceptional ability to deal with large numbers of features and to adapt to new trends not seen during training. Until recently, these properties have remained largely undemonstrated.

We have shown that while these algorithms have many advantages there is still a lot of room to explore when applying them to a real-world problem. In particular, we have demonstrated (1) how variation in document length can be tolerated through either normalization or negative weights, (2) the positive effect of applying a threshold range in training, (3) alternatives in considering feature frequency, and (4) the benefits of discarding irrelevant features as part of the training algorithm. The main contribution of this work, however, is that we have presented an algorithm, BalancedWinnow⁺, which performs significantly better than any other algorithm tested on these tasks using unigram features.

We have exhibited that, as expected, multiplicative-update algorithms have exceptionally good behavior in high dimensional feature spaces, even in the presence of irrelevant features. One advantage this important property has is that it allows one to decompose the learning problem from the feature selection problem. Using this family of algorithms frees the designer from the need to choose the appropriate set of features ahead of time: A large set of features can be used and the algorithm will eventually discard those that do not contribute to the accuracy of the classifier. While we have chosen in this study to use a fairly simple set of features, it is straight forward to plug in instead a richer set of features. We expect that this will further improve the results of the algorithm, although further research is

needed on policies of discarding features and avoidance of over-fitting. In conclusion, we suggest that the demonstrated advantages of the Winnow-family of algorithms make it an appealing candidate for further use in this domain.

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