# An Iterative Method for Multi-class Cost-sensitive Learning

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## ABSTRACT

Cost-sensitive learning addresses the issue of classification in the presence of varying costs associated with different types of misclassification. In this paper, we present a method for solving multi-class cost-sensitive learning problems using any binary classification algorithm. This algorithm is derived using three key ideas: 1) iterative weighting; 2) expanding data space; and 3) gradient boosting with stochastic ensembles. We establish some theoretical guarantees concerning the performance of this method. In particular, we show that a certain variant possesses the *boosting property*, given a form of weak learning assumption on the component binary classifier. We also empirically evaluate the performance of the proposed method using benchmark data sets and verify that our method generally achieves better results than representative methods for cost-sensitive learning, in terms of predictive performance (cost minimization) and, in many cases, computational efficiency.

## Categories and Subject Descriptors

I.2.6 [**Artificial Intelligence**]: Learning

## General Terms

Algorithms

## Keywords

cost-sensitive learning, multi-class classification, boosting

## 1. INTRODUCTION

Classification in the presence of varying costs associated with different types of misclassification is important for practical applications, including many data mining applications, such as targeted marketing, fraud and intrusion detection among others. A body of work on this subject has become known as *cost-sensitive learning*, in the areas of machine learning and data mining.



Research in cost-sensitive learning falls into three main categories. The first category is concerned with making particular classifier learners cost-sensitive, including methods specific for decision trees [14, 4], neural networks [13] and support vector machines [12]. The second category uses Bayes risk theory to assign each example to its lowest expected cost class [8, 18]. This requires classifiers to output class membership probabilities and sometimes requires estimating costs [18] (when the costs are unknown at classification time). The third category concerns methods that modify the distribution of training examples before applying the classifier learning method, so that the classifier learned from the modified distribution is cost-sensitive. We call this approach *cost-sensitive learning by example weighting*. Work in this area includes stratification methods [7, 6] and the *costing* algorithm [19]. This approach is very general since it reuses arbitrary classifier learners and does not require accurate class probability estimates from the classifier. Empirically this approach attains similar or better cost-minimization performance.

Unfortunately, current methods in this category suffer from a major limitation: they are well-understood only for two-class problems. In the two-class case, it is easy to show that each example should be weighted proportionally to the difference in cost between predicting correctly or incorrectly [19]. However, in the multi-class case there is more than one way in which a classifier can make a mistake, breaking the application of this simple formula. Heuristics, such as weighting examples by the average misclassification cost, have been proposed [6, 15], but they are not well-motivated theoretically and do not seem to work very well in practice when compared to methods that use Bayes risk minimization [8].

In this paper, we propose a method for multi-class cost-sensitive learning based on an iterative scheme for example weighting. There are a number of key techniques we employ in this method: 1) an iterative process which empirically adjusts the example weighting according to the performance of the learning algorithm; 2) data space expansion for multi-class labels; 3) gradient boosting [16] with stochastic ensembles. The first two ideas are combined in a unifying framework given by the third.

We establish theoretical performance guarantee for a variant of our algorithm. In particular, we show that this variant possesses the so-called *boosting property*, given that the component classification algorithm satisfies a certain *weak learning assumption*. We test our method on several multi-class data sets and discover excellent predictive performance (i.e.

cost minimization) as compared to existing cost-sensitive algorithms. Moreover, our results show that when the distribution of costs is skewed (as is common in many data mining applications) our method has the added advantage that it uses drastically smaller sample sizes and hence requires much less computational resources.

## 2. PRELIMINARIES

We begin by introducing some general concepts and notation we use in the rest of the paper.

### 2.1 Cost-sensitive learning and related problems

A popular formulation of the cost-sensitive learning problem is via the use of a cost matrix. A cost matrix, $C(y_1, y_2)$, specifies how much cost is incurred when an example is predicted to belong to class $y_1$ when its correct label is $y_2$, and the goal of a cost-sensitive learning method is to minimize the expected cost. Zadrozny and Elkan [18] noted that this formulation is not applicable in situations in which misclassification costs depend on particular instances, and proposed a more general form of cost function, $C(x, y_1, y_2)$, that allows dependence on the instance $x$. Here we adopt a formulation based on this (although slightly more general).

Once we allow the costs to depend on each example, it is natural to assume that the costs are generated according to some distribution, along with the examples, which leads to the following formulation. In (multiclass) cost sensitive classification, examples of the form $(x, \vec{C})$ are drawn from a distribution $D$ over a domain $X \times R^{+^k}$. (Throughout the paper, we will let $k$ denote $|Y|$.) Here, for each label $y \in Y$, $C_y$ equals the cost of misclassifying instance $x$ as $y$, i.e. $C_y = C(x, y, y^*)$ where $y^*$ is the minimum cost label for $x$.

Given a set of examples, $S = (x, \vec{C})^m$, the goal is to find a classifier $h : X \to \{1, ..., k\}$ which minimizes the expected cost of the classifier:

$$\arg\min_h \mathop{\mathrm{E}}_{(x,\vec{C})\sim D}[C_{h(x)}] \tag{1}$$

We can assume without loss of generality that the costs are normalized so that

$$\forall x \in X \ \min_{y\in Y} C_y = 0.$$

Note that with this normalization, the above formulation is equivalent to the common formulation in terms of misclassification cost, i.e.,

$$\arg\min_h \mathop{\mathrm{E}}_{(x,\vec{C})\sim D}[C_{h(x)} I(h(x) \neq \arg\min_y C_y)]$$

where we used $I(\cdot)$ to denote the indicator function which takes on the value 1 whenever the statement is true, and the value 0 otherwise.

Normally a learning method attempts to do this by minimizing the empirical cost in the given training data, given some hypothesis class $\mathcal{H}$:

$$\arg\min_{h\in\mathcal{H}} \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[C_{h(x)}] = \frac{1}{|S|}\sum_{(x,\vec{C})\in S} C_{h(x)} \tag{2}$$

Here the empirical expectation notation, $\hat{\mathrm{E}}$, refers to the averaged empirical cost.

As a building block of our method, we make use of methods for solving importance weighted classification problems, which we define below. In importance weighted classification, examples of the form $(x, y, c)$ are drawn from a distribution $D$ over a domain $X \times Y \times R^+$. Given a set of examples $S = (x, y, c)^m$, the goal is to find a classifier $h : X \to Y$ having minimum importance-weighted misclassification error:

$$\arg\min_h \mathop{\mathrm{E}}_{(x,y,c)\sim D}[c \cdot I(h(x) \neq y)]$$

Again, usually, a learning method attempts to meet this goal by minimizing the empirical weighted error in some hypothesis class $\mathcal{H}$:

$$\arg\min_{h\in\mathcal{H}} \hat{\mathrm{E}}_{(x,y,c)\sim S}[c \cdot I(h(x) \neq y)] \tag{3}$$

We emphasize that the importance weighted formulation critically differs from the per example formulation of multiclass cost-sensitive learning in that there is a single weight associated with each instance $x$, whereas in multi-class cost-sensitive learning there is a weight (misclassification cost) associated with each label $y$. We note that importance weighted classification can be solved very well with a classifier learning method, by use of weighted rejection sampling techniques [19].

### 2.2 Hypothesis representations and other notation

In the above, we assumed that the hypotheses output by a cost-sensitive learner is a functional hypothesis $h$, i.e. $h : X \to Y$. It is also possible to allow hypotheses that are *stochastic*, namely

$$h : X \times Y \to [0, 1]$$

subject to the stochastic condition:

$$\forall x \in X \ \sum_{y\in Y} h(y|x) = 1.$$

With stochastic hypotheses, *stochastic* cost-sensitive learning is defined as the process of finding a hypothesis minimizing the following expected cost:

$$\arg\min_h \mathop{\mathrm{E}}_{(x,\vec{C})\sim D} \mathop{\mathrm{E}}_{y\sim h(y|x)}[C_y]$$

In general, we sometimes use the following short-hand notation for the expected cost of a stochastic hypothesis $h$ on an instance $x$.

$$C_{h(x)} \equiv \mathop{\mathrm{E}}_{y\sim h(y|x)}[C_y]$$

Note that in the special case that $h$ is deterministic, the above formulation is equivalent to the definition given in Eq. 1. Also, this is a *convexification* of the standard objective function that we usually expect a stochastic cost-sensitive learner to minimize, i.e.

$$\mathop{\mathrm{E}}_{(x,\vec{C})\sim D}[C_{\arg\max_y h(y|x)}]$$

We also consider a variant of cost-sensitive learning in which *relational* hypotheses are allowed. Here relational hypotheses $h$ are relations over $X \times Y$, i.e. $h : X \times Y \to \{0, 1\}$. In general $h$ is neither functional nor stochastic, and in particular it may violate the stochasticity: $\sum_{y\in Y} h(x, y) = 1$.

- An input sample $S = (x, \vec{C})^m$.
- A component learner $A$ that takes an importance-weighted sample $S' = (x, y, c)^m$ and outputs a functional hypothesis $h$, possibly by weighted sampling.
- An integer $T$ specifying the number of iterations to be performed.

1. Let $S' =$
$$\{(x, \arg\min_y C_y, \max_y C_y) | (x, \vec{C}) \in S\}$$
2. **For** $t := 1$ to $T$ **do**
   (a) Let $h_t := A(S')$.
   (b) Choose $\alpha_t \in [0, 1)$, for example $\alpha_t := 1/t$.
   (c) For each example $(x, y, c) \in S'$ with an error $h_t(x) \neq y$ update the importance weight by
   $$c := \alpha_t C_{h(x)} + (1 - \alpha_t)c$$
3. Return $h_T$.

**Figure 1: Method *IW (Iterative Weighting)***

# 3. THE METHODOLOGY

Our methodology can be interpreted as a *reduction*, which translates a multi-class cost-sensitive learning problem to a classifier learning problem.

This methodology is derived using three key ideas: 1) iterative weighting; 2) expanding data space; and 3) gradient boosting with stochastic ensembles. The first two ideas are combined in a unifying framework given by the third.

Below we will explain the first two key ideas by exhibiting a prototypical method based on each, and then derive our main learning method that makes use of them in a gradient boosting framework.

## 3.1 Iterative cost weighting

We note that the weighting scheme proposed in [19], called *costing*, exploits the following observation: For the binary class case, the above formulation in terms of per example cost for each class can be further reduced to a formulation in terms of a single importance number per example. This is possible by associating a number indicating the importance of an example $(x, y)$, given by $|C_0 - C_1|$. This conversion allows us to reduce the cost-sensitive learning problem to a weighted classifier learning problem, but it is not immediately obvious how that would be done for the multi-class scenario. It is therefore natural to consider iterative weighting schemes, in which example weights are iteratively modified in search for the optimal weighting. The technique, which we call *IW (Iterative Weighting)*, presented in Figure 1, is an example of such a scheme. We can show that the final weights of IW are the optimal weights in some sense, *provided* the algorithm converges.

THEOREM 1. *Assume IW converges and the final hypothesis is $h$. Then, the following holds:*

$$\hat{\mathrm{E}}_{(x,y,c)\sim S'}[c \cdot I(h(x) \neq y)] = \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[C_{h(x)}].$$

- An input sample $S = (x, \vec{C})^m$.
- A component learner $A$ that takes an importance-weighted sample $S' = (x, y, c)^m$ and outputs a functional hypothesis $h$, possibly by weighted sampling.

1. Set $S' = \{(x, y, \max_{y'} C_{y'} - C_y) | (x, \vec{C}) \in S, y \in Y\}$
2. Return $h := A(S')$

**Figure 2: Method *DSE (Data Space Expansion)***

This theorem says that, if the iterative algorithm converges, then the loss of the component learner with respect to the final weights (left hand side) is equal to the expected cost for the problem that we wish to solve (right hand side).

**Proof**

At convergence, for every example $(x, y, c)$ that the $h$ errs on, we must have:

$$c = \alpha_t C_{h(x)} + (1 - \alpha_t)c$$

and thus

$$c = C_{h(x)}$$

Noting that when $h$ does not err on $x$, we have $I(h(x) \neq y) = 0$, Thus, it follows that

$$\hat{\mathrm{E}}_{(x,y,c)\sim S'}[c \cdot I(h(x) \neq y)] = \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[C_{h_t(x)}]$$

**Q.E.D**

## 3.2 Data space expansion

One drawback to iterative weighting is an inability to directly take into account the different costs associated with multiple ways of misclassifying examples. This translates to non-convergence of the method in practice. We address this issue by the technique of expanding data space, as we describe below.

Given a labeled sample $S$ consisting of $(x, \vec{C})$ of size $m$, we define an *expanded* sample $S'$ of size $mk$ for weighted classification, where $k$ is the size of the label set, i.e. $k = |Y|$, as follows.

$$S' = \{(x, y, \max_{y'} C_{y'} - C_y) | (x, \vec{C}) \in S, y \in Y\}$$

Note here that the newly defined weights are more like *benefits* than costs, since larger costs are mapped to smaller weights.

It turns out that minimizing the importance weighted loss,

$$\hat{\mathrm{E}}_{(x,y,c)\sim S'} c \cdot I(h(x) \neq y)$$

on this new data set also minimizes the cost on our original sample. The algorithm DSE (Data Space Expansion), shown in Figure 2, is based on this observation, which is summarized as theorem below.

THEOREM 2. *With the definitions given in Figure 2, a hypothesis $h$ minimizing the weighted classification error on the expanded weighted sample $S'$,*

$$\hat{\mathrm{E}}_{(x,y,c)\sim S'}[c \cdot I(h(x) \neq y)]$$

*also minimizes the cost on the original sample $S$,*

$$\hat{\mathrm{E}}_{(x,\vec{C})\sim S}[C_{h(x)}]$$

**Proof**

$$\begin{aligned}
&\arg\min_h \hat{\mathrm{E}}_{(x,y,c)\sim S'}[c \cdot I(h(x) \neq y)] \\
=\ &\arg\min_h \hat{\mathrm{E}}_{(x,\vec{C})\sim S} \sum_{y\in Y}[(\max_{y'\in Y} C_{y'} - C_y) \cdot I(h(x) \neq y)] \\
=\ &\arg\max_h \hat{\mathrm{E}}_{(x,\vec{C})\sim S} \sum_{y\in Y}[C_y \cdot I(h(x) \neq y)] \\
=\ &\arg\max_h \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[(\sum_{y\in Y} C_y) - C_{h(x)}] \\
=\ &\arg\min_h \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[C_{h(x)}]
\end{aligned}$$

**Q.E.D.**

## 3.3 Gradient Boosting with Stochastic Ensembles

Having described two key ideas, namely iterative weighting and data space expansion, we now apply them together to arrive at our main method. We do so by casting the *stochastic* multiclass cost-sensitive learning in the framework of gradient boosting [16], with the objective function defined as the expected cost of the *stochastic ensemble*, obtained as a mixture of individual hypotheses, on the *expanded* data set. As we stated in Section 2, a functional hypothesis of the form $h : X \to Y$ can be viewed as a special case of a stochastic hypothesis. We then define a stochastic ensemble hypothesis $H$, given multiple functional hypotheses, $h_t, t = 1, ..., T$, as the conditional distribution defined as the mixture of the component hypotheses, namely,

$$\forall x \in X, \forall y \in Y, H(y|x) = \frac{1}{T}\sum_{t=1}^{T} I(h_t(x) = y)$$

Let $H_t$ denote the mixture hypothesis of the learning procedure at round $t$. The procedure is to update its current combined hypothesis by the mixture of the previous combined hypothesis and a new hypothesis, i.e. by setting

$$H_t(y|x) = (1-\beta)H_{t-1}(y|x) + \beta I(h(x) = y)$$

Thus, the expected cost of $H_t$ on $x$ is

$$\mathrm{E}_{y\sim H_t(y|x)} C_y = (1-\beta)\,\mathrm{E}_{y\sim H_{t-1}(y|x)} C_y + \beta C_{h_t(x)}$$

If we now take a derivative of this function with respect to $\beta$, we get:

$$\frac{\partial\, \mathrm{E}_{y\sim H_t(y|x)} C_y}{\partial\beta} = C_{h_t(x)} - \mathrm{E}_{y\sim H_{t-1}(y|x)} C_y$$

Note that this is the difference between the average cost of the current *ensemble* hypothesis and the new weak hypothesis assigning probability one to the specified label.

We then take the expectation of this derivative with respect to all data points $(x, y)$ in the expanded data set $S'$, and thus the gradient is $mk$-dimensional. The weak learner is to find a hypothesis $h$ whose inner-product with the negative gradient is large. That is, the output $h$ of the weak learner seeks to maximize the following sum.

$$-\langle h, \nabla C\rangle = \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[\,\mathrm{E}_{y'\sim H_{t-1}(y|x)} C_{y'} - C_{h(x)}] \qquad (4)$$

This leads to the following example weighting on the expanded sample:

$$w_{x,y} = C_{H_{t-1}(x)} - C_y$$

where we recall that $C_{H_{t-1}(x)}$ denotes $\mathrm{E}_{y'\sim H_{t-1}(y|x)} C_{y'}$.

Note that these weight updates are similar to those used in IW and DSE. In fact, the IW weight update rule is essentially equivalent to the GBSE rule, except IW has, for each instance $x$, the weight for only the best (least cost) label, and hence $C_y = 0$ holds. This is because the IW update rule mixes the weights from earlier iterations, which is equivalent to taking the average over the stochastic ensemble as is done in GBSE. The DSE update rule differs from the GBSE rule in that $\max_{y'} C_{y'}$ is used in place of $C_{H_{t-1}(x)}$, which ensures that the weight is always non-negative, even though DSE has weights for all labels. Thus, the GBSE weights can be viewed as IW weights, applied on the expanded data set, as in DSE. As a consequence, the GBSE weights $w_{x,y} := C_{H_{t-1}(x)} - C_y$ can be negative, since $y$ is not necessarily the best label. This means that the weak learner now receives both positive and negative weights. While the minimization of weighted misclassification with positive and negative weights makes perfect sense as an optimization problem, its interpretation as a classification problem is not immediately clear. In particular, it prohibits the use of weighted sampling as a means of realizing the weighted classification problem.

We deal with this problem by converting a relational version of the weighted multi-class classification problem (i.e. of finding $h$ to maximize Eq. 4) in each iteration to a weighted binary classification problem. Specifically we convert each example pair $(x, y)$ to $((x, y), l)$, and set $l = 1$ if the weight on $(x, y)$ is positive, and $l = 0$ if the weight is negative. The output hypothesis of the binary classifier is in general relational, so it is converted to a stochastic hypothesis by the procedure **Stochastic** shown in Figure 4. (The particular way this procedure is defined is motivated by the theoretical guarantee, which will be shown in the next subsection.) The overall process, consisting of multiple iterations of such a reduction, constitutes a reduction of the *stochastic* multi-class cost-sensitive classification to binary weighted classification.

With the foregoing definitions, we can now state our main method, *GBSE* (Gradient Boosting with Stochastic Ensembles), which is shown in Figure 3.

## 3.4 Theoretical Performance Guarantee on a Variant

It turns out that a strong theoretical performance guarantee can be proved on a variant of this method. The variant is obtained by simply replacing the weight updating rule of GBSE by the following:

$$w_{x,y} = \frac{C_{H_{t-1}(x)}}{k} - C_y$$

The resulting variant, which we call GBSE-T (Gradient Boosting with Stochastic Ensembles - Theoretical version), is summarized in Figure 5.

We can show that GBSE-T has a boosting property given a version of weak learning condition on the component classifier. This weak learning condition, which we make precise below, is one that is sensitive to class imbalance.

DEFINITION 1. *We say that an algorithm $A$ for the binary importance weighted classification problem, as defined*

- An input sample $S = (x, \vec{C})^m$.
- A component learner $A$ for (importance weighted) binary classification that takes a sample of the form $((x, y), l, w)^*$, and outputs a relational hypothesis $h$.
- A subprocedure **Stochastic**, as specified in Figure 4.
- An integer $T$ specifying the number of iterations to be performed.

1. Set $S' = \{(x, y) | (x, \vec{C}) \in S, y \in Y\}$.
2. Initialize $H_0$ by $\forall x \in X, y \in Y\ H_0(y|x) = 1/k$.
3. **For** $t := 1$ to $T$ **Do**
   (a) $w_{x,y} := \mathrm{E}_{y' \sim H_{t-1}(x)}[C_{y'}] - C_y$ for all $(x, y)$ in $S'$.
   (b) $S_t = \{((x, y), I(w_{x,y} \geq 0), |w_{x,y}|) | (x, y) \in S'\}$.
   (c) Let $h_t := A(S_t)$
   (d) $f_t :=$ **Stochastic**$(h_t, H_{t-1})$.
   (e) Choose $\alpha_t \in [0, 1)$, for example $\alpha_t = \frac{1}{t}$.
   (f) Set $H_t := (1 - \alpha_t)H_{t-1} + \alpha_t f_t$.
4. **End For**
5. Return $H_T$.

**Figure 3: Method *GBSE (Gradient Boosting with Stochastic Ensembles)***

**Stochastic**($h$: a relational hypothesis, $H$: a stochastic hypothesis)

1. Define $f$ by setting for each $x \in X$:
   - (default) if $|\{y \in Y | h(x, y) = 1\}| = 0$ then define for all $y \in Y$, $f(y|x) = H(y|x)$.
   - else $f(y|x) = \frac{I(h(x,y)=1)}{|\{y \in Y | h(x,y)=1\}|}$.
2. Output $f$.

**Figure 4: Sub-procedure Stochastic**

*in Section 2, satisfies the* weak learning condition *for a given classification sample $S = (x, y)^m$, if for all weighted samples $S' = (x, y, c)^m$ for it, its output $h$ satisfies the following, for some fixed $\gamma > 0$:*

$$\hat{\mathrm{E}}_{(x,y,c) \sim S'} cI(h(x) = y) \geq \hat{\mathrm{E}}_{(x,y,c) \sim S'} cI(y = 0) + \gamma \hat{\mathrm{E}}_{(x,y,c) \sim S'} cI(y = 1) \quad (5)$$

Intuitively, this weak learning condition requires that the weak learner achieve better weighted accuracy than that attainable trivially by assigning all examples to the negative class.

THEOREM 3. *Suppose that the component learner $A$ satisfies the weak learning condition for sample $S'$ as defined by GBSE-T. If we set $\alpha_t = \alpha$ for all $t$, the output of GBSE-T*

Identical to Method *GBSE* of Figure 3, except for the following change:
3.(a) Set $w_{x,y} := \frac{C_{H_{t-1}(x)}}{k} - C_y$, for all $(x, y)$ in $S'$.

**Figure 5: Method *GBSE-T (Gradient Boosting with Stochastic Ensembles - Theoretical variant)***

*satisfies:*

$$\hat{\mathrm{E}}_{(x,\vec{C}) \sim S} C_{H_T(x)} \leq \exp\{-\frac{\gamma\alpha}{k} T\} \hat{\mathrm{E}}_{(x,\vec{C}) \sim S} C_{H_0(x)}$$

This theorem shows that the empirical cost of the output hypothesis of GBSE-T converges exponentially fast, given the weak learning assumption.

**Proof**

We first establish the following simple correspondence between the *weak learning conditions* on the relational multi-class classification problem that we wish to solve in each iteration, and the weighted binary classification problem that is given to the component algorithm to solve it.

DEFINITION 2. *Let $S$ be a weighted sample of the form $S = (x, y, c)^m$, where weights $c$ can be both positive and negative. Then define a transformed sample $S'$ for weighted classification as $S' = ((x, y), I(c > 0), |c|)^m$.*

1. *The relational weighted multi-class classification problem for $S$ is to find a relational hypothesis $h : X \times Y \to \{0, 1\}$ that maximizes the following sum:*
$$\hat{\mathrm{E}}_{(x,y,c) \sim S} c \cdot h(x, y)$$
2. *The weighted binary classification problem for the transformed sample $S'$ is to find a hypothesis $h : X \times Y \to \{0, 1\}$ that maximizes the following weighted classification accuracy:*
$$\hat{\mathrm{E}}_{((x,y),I(c>0),|c|) \sim S'} |c| \cdot I(h(x, y) = I(c > 0))$$

Note that, in a relational weighted classification problem as defined in Definition 2, the goal of a learner is to try to assign 1 to pairs with positive weights and assign 0 to those with negative weights as much as possible.

LEMMA 1. *For all $h$:*

$$\hat{\mathrm{E}}_{(x,y,c) \sim S} c \cdot h(x, y) = \hat{\mathrm{E}}_{((x,y),I(c>0),|c|) \sim S'} [|c| I(h(x, y) = I(c > 0))] - \hat{\mathrm{E}}_{((x,y),I(c>0),|c|) \sim S'} [|c| I(c < 0)]$$

**Proof of Lemma 1**

$$\begin{aligned}
& \hat{\mathrm{E}}_{((x,y),I(c>0),|c|) \in S'} |c| \cdot I(h(x, y) = I(c > 0)) \\
= & \hat{\mathrm{E}}_{(x,y,c) \in S} |c| \cdot I(h(x, y) = I(c > 0)) \\
= & \hat{\mathrm{E}}_{(x,y,c) \sim S} c \cdot I(h(x, y) = 1 \text{ and } c \geq 0) \\
& + \hat{\mathrm{E}}_{(x,y,c) \sim S} -c \cdot I(h(x, y) = 0 \text{ and } c < 0) \\
= & \hat{\mathrm{E}}_{(x,y,c) \sim S} c \cdot h(x, y) I(c \geq 0) + c(h(x, y) - 1) I(c < 0) \\
= & \hat{\mathrm{E}}_{(x,y,c) \sim S} c \cdot h(x, y) + \hat{\mathrm{E}}_{(x,y,c) \sim S} |c| I(c < 0)
\end{aligned}$$

Hence the lemma follows. **Q.E.D.**

This lemma establishes that getting positive weighted accuracy on the original relational weighted multi-class classification problem is equivalent to the weak learning condition on the transformed weighted binary classification problem.

**Proof of Theorem 3**

First, note that applying **Stochastic** to $h_t$ can increase the expected cost only for $x$'s such that $|\{y|h_t(x,y)=1\}|=0$, and for such $x$'s the cost of the output function $f$ equals that of $H_{t-1}$ by the definition of **Stochastic**. Hence, the average empirical cost of $f$ on the original sample $S$, satisfies the following:

$$\begin{aligned} & \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[C_{f(x)} - \sum_y h_t(x,y)C_y] \\ \leq & \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[C_{H_{t-1}(x)} I(\forall y\, h(x,y)=0)] \end{aligned} \quad (6)$$

Now recall that the expected empirical cost of $H_t$ equals the following, where we drop the subscript $t$ from $\alpha_t$.

$$\hat{\mathrm{E}}_{(x,\vec{C})\sim S} C_{H_t(x)} = \hat{\mathrm{E}}_{(x,\vec{C})\sim S}(1-\alpha)C_{H_{t-1}(x)} + \alpha C_{f(x)} \quad (7)$$

Hence, by combining Eq. 6 and Eq. 7, we can show the following bound on the decrease in empirical cost in each iteration. Here, we also drop the subscript $t$ on $h$.

$$\begin{aligned} & \hat{\mathrm{E}}_{(x,\vec{C})\sim S} C_{H_{t-1}(x)} - \hat{\mathrm{E}}_{(x,\vec{C})\sim S} C_{H_t(x)} \\ = & \hat{\mathrm{E}}_{(x,\vec{C})\sim S} \alpha(C_{H_{t-1}(x)} - C_{f(x)}) \\ = & \alpha \hat{\mathrm{E}}_{(x,\vec{C})\sim S} C_{H_{t-1}(x)} - \sum_y h(x,y)C_y \\ & + \sum_y h(x,y)C_y - C_{f(x)} \\ \geq & \alpha \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[C_{H_{t-1}(x)} \\ & - \sum_y h(x,y)C_y - C_{H_{t-1}(x)} I(\forall y\, h(x,y)=0)] \\ \geq & \alpha \hat{\mathrm{E}}_{(x,\vec{C})\sim S}[\sum_{y:h(x,y)=1} h(x,y)\left(\frac{C_{H_{t-1}(x)}}{k} - C_y\right) \\ & + \sum_{y:h(x,y)=0} \frac{C_{H_{t-1}(x)}}{k} - C_{H_{t-1}(x)} I(\forall y\, h(x,y)=0)] \\ = & \alpha \hat{\mathrm{E}}_{(x,\vec{C})\sim S} \sum_y h(x,y)\left(\frac{C_{H_{t-1}(x)}}{k} - C_y\right) \\ & + C_{H_{t-1}}(x)\left(\frac{|\{y: h(x,y)=0\}|}{k} - I(\forall y\, h(x,y)=0)\right) \\ \geq & \alpha \hat{\mathrm{E}}_{(x,\vec{C})\sim S} \sum_y h(x,y)\left(\frac{C_{H_{t-1}(x)}}{k} - C_y\right) \\ = & \alpha \hat{\mathrm{E}}_{(x,\vec{C})\sim S} \sum_y \left|\frac{C_{H_{t-1}(x)}}{k} - C_y\right| \\ & (I(h(x,y) = I(\frac{C_{H_{t-1}(x)}}{k} > C_y)) - I(\frac{C_{H_{t-1}(x)}}{k} < C_y)) \end{aligned} \quad (8)$$

Here Lemma 1 was applied to get the last equality. Next apply the weak learning assumption on the induced measure over $(x', y', c')$ defined by: $x' = (x,y)$, $y' = I\left(\frac{C_{H_{t-1}(x)}}{k} > C_y\right)$, and $c' = \left|\frac{C_{H_{t-1}(x)}}{k} - C_y\right|$ to get:

$$\begin{aligned} & \geq \alpha\gamma \hat{\mathrm{E}}_{(x,\vec{C})\sim S} \sum_y \left|\frac{C_{H_{t-1}(x)}}{k} - C_y\right| I\left(\frac{C_{H_{t-1}(x)}}{k} > C_y\right) \\ & \geq \frac{\alpha\gamma}{k} \hat{\mathrm{E}}_{(x,\vec{C})\sim S} C_{H_{t-1}(x)} \end{aligned}$$

The last inequality follows because for all $y$ $C_y \geq 0$, there exists $y$ such that $C_y = 0$, and the sum is bounded below by its largest term.

Since the expected cost is convex (in fact linear), this implies convergence to the global optimum. Noting that in each iteration, the empirical cost is reduced at least by a factor of $1 - \frac{\gamma\alpha}{k}$, the theorem follows. **Q.E.D.**

Note that at earlier iterations, the binary classifier used as the component learner is likely to be given a weighted sample with balanced positive and negative examples. As the number of iterations increases and progress is made, however, it will receive samples that are increasingly more negative. (This is because the positive examples correspond to labels that can further improve the current performance.) It therefore becomes easier to attain high weighted accuracy by simply classifying all examples to be negative. The weak learning condition of Eq. 5 appropriately deals with this issue, as it requires that the weak learner achieve better weighted accuracy than that attainable by assigning all examples to the negative class, as we mentioned earlier.

# 4. EXPERIMENTAL EVALUATION

We use the C4.5 decision tree learner [17] as the base classifier learning method, because it is a standard for empirical comparisons and it was used as the base learner by Domingos for the *MetaCost* method [8].

We compare our methods against three representative methods: Bagging [5], Averaging cost [7, 8] and MetaCost. The Averaging cost method was also used for comparison in [8]. Note that Bagging is a cost-insensitive learning method. Here we give a brief description of these methods, and refer the reader to [5, 8] for the details.

- *Bagging* obtains multiple sub-samples from the original training set by sampling with replacement, feeds them to the base learner (C4.5), and takes the average over the ensemble of output hypotheses
- *Averaging Cost* (AvgCost) obtains a sub-sample by weighted sampling with weights defined as the average cost for each $x$, and then feeds it to the base learner (C4.5).
- *MetaCost* uses bagging to obtain an ensemble of hypotheses, then uses the ensemble to estimate class probabilities, relabels the examples with the labels that minimize the expected risk according to the probability estimates and finally runs the base learner (C4.5) to obtain a single classifier that predicts the new labels.

There are some deviations from these methods in our implementation, which we clarify below. The main deviation is that we use *rejection* sampling for all methods (including bagging), while other sampling schemes such as resampling with replacement are used in the original methods.[1] We do

[1] In weighted rejection sampling, the original data are

this for two reasons: (1) inadequacy of resampling with replacement (or over-sampling), especially for C4.5, has been noted by various authors [19, ?]; (2) since our proposed methods use rejection sampling, we do the same for the other methods for fairness of comparison. We stress that this deviation should only improve their performance. Another deviation is that we use a variant of MetaCost that skips the last step of learning a classifier on the relabeled training data set, but directly minimizes the expected risk on the test data. It has been observed that this variant performs at least as well as MetaCost, in terms of cost minimization. (This variant has been called *BagCost* by Margineantu [15].) Also, in our implementation of AvgCost, we perform weighted sampling multiple times to obtain an ensemble of hypotheses, then output their average as the final hypothesis. (In the original implementation, only a single iteration of weighted sampling was performed.) We note that, due to our normalization assumption that the minimum cost for each instance $x$ is always zero, our version of AvgCost is identical to a more sophisticated variant in which the difference between the average cost and the minimum cost is used for sampling weights. Our experience shows that this variant of AvgCost performs better than the original method.

The methods were applied to five benchmark data sets available from the UCI machine learning repository [3] and one data set from the UCI KDD archive [2]. These data sets were selected by the criterion of having approximately 1000 examples or more, besides being multi-class problems. A summary of these data sets is given in Table 1. Here *class ratio* is defined as the class frequency of the least frequent class divided by that of the most frequent one. We note that the KDD-99 data set is actually a larger data set. We used the so-called 10% training data set, which consists roughly of 500 thousand instances, and further sampled down by random sampling 40% of them, to get the data set of size 197,710 which we used for our experimentation. Emphatically, we only used data from the original training set, and not data from the test set. We do this because of the idiosyncratic property of this data set that the test data are generated from a considerably different data distribution. While this property is both realistic and interesting for empirical evaluation of a method for intrusion detection, we judged it not to be desirable for the current purpose of evaluating general purpose cost-sensitive classification algorithms.

Except for the KDD-99 data set, these data sets do not have standard misclassification costs associated with them. For this reason, we follow Domingos and generate cost matrices according to a model that gives higher costs for misclassifying a rare class as a frequent one, and the lower costs for the reverse. (Note therefore that our experiments do not exploit the full generality of the instance-dependent cost formulation presented in Section 2.) This reflects a situation that is found in many practical data mining applications, including direct marketing and fraud detection, where the rare classes are the most valuable to identify correctly.

Our cost model is as follows: Let $\hat{P}(y_1)$ and $\hat{P}(y_2)$ be the empirical probabilities of occurrence of classes $y_1$ and $y_2$ in the training data. We choose the non-diagonal entries of the cost matrix $C(y_1, y_2)$, $y_1 \neq y_2$ with uniform probability from the interval $[0, 2000\hat{P}(y_1)/\hat{P}(y_2)]$. In [8], the diagonal entries were then chosen from the interval $[0, 1000]$, which often leads to cost matrices in which the correct label is not the least costly one. Besides being unreasonable (see Elkan [9]), these cost matrices can give an unfair advantage to cost-sensitive methods over cost-insensitive ones. We therefore set the diagonal entries to be identically zero, which is consistent with our normalization assumption.

scanned once (without replacement), and each example is accepted with probability equal to (or proportional to) its weight.

| Data Set | # of examples | # of classes | Class ratio |
|---|---|---|---|
| Annealing | 898 | 5 | 0.01316 |
| KDD-99 | 197710 | 5 | 0.0001278 |
| Letter | 20000 | 26 | 0.9028 |
| Satellite | 6435 | 6 | 0.4083 |
| Solar flare | 1389 | 7 | 0.002562 |
| Splice | 3190 | 3 | 0.4634 |

**Table 1: Data set characteristics: data size, number of classes, and the ratio between the frequency of the most common class to the least common.**

In all experiments, we randomly select two thirds of the examples in the data set for training and use the remaining one third for testing. Also, for each training/test split we generate a different cost matrix according to the rules above. Thus, the standard deviations that we report reflect both variations in the data and in the misclassification costs.

We remark on certain implementation details of the proposed learning methods in our experimentation. First, we note that in all of the methods used for comparison, except IW, C4.5 was used as the component algorithm with weighted rejection sampling, and the final hypothesis is expressed essentially as an ensemble of output decision tress of C4.5. IW, as a meta-method, does not use ensembles; instead we used an ensemble method of *costing* [19] on C4.5, as the component algorithm. Its output hypothesis is therefore also an ensemble of decision trees. DSE, as stated in its definition in Figure 2, is not an ensemble method, but analogously to AvgCost, we performed multiple iterations of weighted sampling according to the weighting scheme of DSE and averaged the resulting hypotheses to define the final hypothesis. Finally, the choice of the mixture weight $\alpha_t$ was set at $1/t$ for all methods.

The results of these experiments are summarized in Table 2 and Table 3. Table 2 lists the average costs attained by each of these methods on the 6 data sets, and their standard errors. These results were obtained by averaging over 20 runs, each run consisting of 30 iterations of the respective learning method. These results appear quite convincing: GBSE outperforms all comparison methods on all data sets, except on Splice, for which it ranks second after MetaCost. Also, GBSE is the best performing among the proposed methods in the paper, confirming our claim that the combination of various techniques involved is indeed necessary to attain this level of performance.

Table 3 lists the average total data size used by each of the methods in 30 iterations. The data size for IW is not listed, since it consists of 30 iterations of 10 rounds of costing and the direct comparison of total data size does not seem to make as much sense for this method. Examining these results in conjunction with the data characteristics in Table 1 reveals a definite trend. First, note that the data sets are divided into two groups: those having very large skews, or very low class ratios (Annealing, KDD-99 and Solar flare), and

| Data Set | Bagging | AvgCost | MetaCost | IW | DSE | GBSE |
|---|---|---|---|---|---|---|
| Annealing | $1059 \pm 174$ | $127.4 \pm 12.2$ | $206.8 \pm 42.8$ | $67.38 \pm 9.22$ | $127.1 \pm 14.9$ | $\mathbf{33.72 \pm 4.29}$ |
| Solar | $5403 \pm 397$ | $237.8 \pm 37.5$ | $5317 \pm 390$ | $174.2 \pm 32.7$ | $110.9 \pm 28.7$ | $\mathbf{48.17 \pm 9.52}$ |
| KDD-99 | $319.4 \pm 42.2$ | $42.43 \pm 7.95$ | $49.39 \pm 9.34$ | $50.43 \pm 10.0$ | $46.68 \pm 10.16$ | $\mathbf{1.69 \pm 0.78}$ |
| letter | $151.0 \pm 2.58$ | $91.90 \pm 1.36$ | $129.6 \pm 2.44$ | $247.7 \pm 4.15$ | $114.0 \pm 1.43$ | $\mathbf{84.63 \pm 2.44}$ |
| Splice | $64.19 \pm 5.23$ | $60.78 \pm 3.65$ | $\mathbf{49.95 \pm 3.05}$ | $67.26 \pm 4.18$ | $135.5 \pm 14$ | $57.50 \pm 4.38$ |
| Satellite | $189.9 \pm 9.57$ | $107.8 \pm 5.95$ | $104.4 \pm 6.43$ | $140.1 \pm 18.2$ | $116.8 \pm 6.28$ | $\mathbf{93.05 \pm 5.57}$ |

**Table 2: Experimental results: the average cost and standard error.**

| Data Set | Bagging | AvgCost | MetaCost | IW | DSE | GBSE |
|---|---|---|---|---|---|---|
| Annealing | $11991 \pm 13.1$ | $1002.8 \pm 183$ | $11987 \pm 9.84$ | – | $3795.5 \pm 688$ | $1260.2 \pm 224$ |
| Solar | $18499 \pm 20.4$ | $334.80 \pm 37.5$ | $18510 \pm 14.4$ | – | $2112.8 \pm 276$ | $486.45 \pm 53.3$ |
| KDD-99 | $395310 \pm 143$ | $2551.9 \pm 428.6$ | $395580 \pm 143$ | – | $12512 \pm 2450$ | $4181 \pm 783.6$ |
| letter | $40037 \pm 44.3$ | $159720 \pm 2028$ | $40052 \pm 41$ | – | $479130 \pm 2710$ | $363001 \pm 5557$ |
| Splice | $42515 \pm 26.6$ | $33658 \pm 1697$ | $42501 \pm 21$ | – | $52123 \pm 592$ | $50284 \pm 3659$ |
| Satellite | $86136 \pm 123$ | $60876 \pm 1641$ | $85984 \pm 127$ | – | $218870 \pm 6516$ | $140810 \pm 3335$ |

**Table 3: Experimental results: the average data size used by each method in 30 iterations, and standard error.**

those having moderate skews (Satellite, Splice and Letter). It is evident that the methods based on example weighting (AvgCost, GBSE, DSE) use magnitudes smaller data sizes for the 3 data sets in the first group (i.e. with large skews), as compared to the other methods (Bagging and MetaCost). The performance of GBSE is especially impressive on this group, achieving much lower cost while requiring very small data sizes. It is worth mentioning that it is these data sets in the first group with large skews, that require *cost-sensitive* learning the most.

## 5. DISCUSSION

It is not the first time that the issue of incorporating cost-sensitivity to boosting has been addressed. For example, AdaCost [10] suggested a way of modifying AdaBoost's exponential loss using a function (called *cost adjustment function*) of the cost and confidence. The rational choice of this cost adjustment function, however, appears not to be well-understood. The stochastic ensemble that we employ in the present paper provides a straightforward but reasonable way of incorporating cost and confidence, i.e. in terms of *expected cost*. An interesting future direction is to investigate the relationship between these alternative approaches to cost-sensitive boosting. Also note that AdaCost, being a modification of AdaBoost, is restricted to two-class problems. Comparing and studying possible relationships between GBSE and other (both cost-sensitive and insensitive) multi-class extensions of boosting, such as AdaBoost.M2 [11], is another interesting topic. Finally, GBSE can also be viewed as a reduction from multi-class classification to binary classification. Comparison with existing methods for such reductions (e.g. [1]) is another important research issue.

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