

Imbalanced Dataset Techniques to Improve Classification of Uneven Class Distributions

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Abstract—This document analyzes the impact of imbalanced datasets on machine learning classifiers. It discusses several techniques to deal with the imbalance. The results are applied to train a Twitter fake account detection algorithm where the performance of these techniques is compared.

Index Terms—imbalance

I. INTRODUCTION

A. Motivation

Many canonical machine learning algorithms used for classification assume that the number of objects in the respective classes is roughly the same. However, in reality, classes are rarely represented equally. In fact, class distribution skews are not only common, but many times expected [1], especially in decision systems aiming to detect rare but important cases. For instance, Covid-19 testing at Georgia Institute of Technology showed that less than 1% of the samples contained the virus. This means that a naive classifier could achieve a 99% accuracy just by labeling all samples as negative for Covid-19.

Imbalanced datasets significantly compromise the performance of many traditional learning algorithms. The disparity of classes in the training dataset may lead the algorithm to bias the classification towards the class with more instances, or even to ignore the minority class altogether. Therefore, it is vital to find efficient ways of dealing with data imbalances.

The overall goal of our project is to provide an overview of the state-of-the-art approaches to solve the issues introduced by imbalanced datasets. Including, a performance comparison of the various techniques. We also aim to implement an efficient scheme that is able to deal with highly complex and imbalanced datasets.

TO-DO: INCLUDE HERE THAT WE WILL DO RESAMPLING, COST-SENSITIVE, ENSEMBLED METHODS AND HYBRID ALGORITHMS, AND EXPLAIN WHY ENSEMBLED METHODS MAY WORK BETTER THAN TRADITIONAL ALGORITHMS.

B. Methodology

Firstly, we study a synthetic dataset characterized by its simplicity. It is made up of two classes following $N(\mu_0, \Sigma_0)$ and $N(\mu_1, \Sigma_1)$, where

$$\mu_0 = \begin{pmatrix} -1 \\ -0.5 \end{pmatrix}, \mu_1 = \begin{pmatrix} 0 \\ 1 \end{pmatrix}, \Sigma_0 = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}, \Sigma_1 = \begin{pmatrix} 4 & 0 \\ 0 & 2 \end{pmatrix}$$

and where the minority class only accounts for 15% of the samples. This simple dataset is especially useful to analyze the imbalance-compensating techniques from a mathematical perspective. Not only do we study the concepts learnt in class at a theoretical level, but we also use plugin machine learning models to illustrate how they affect density distributions.

Secondly, we target a more complex dataset, which gather data of more than thirty-seven thousand Twitter accounts to determine whether they are a bot or human accounts. The number of bot accounts represents a 30% of the total accounts. The dataset contains thirteen features, such as the existence of a profile image, location or favorites and followers count. Some variables are categorical, while others are binary or integer variables.

The performance of the classification will be evaluated using the F_1 score $\in [0, 1]$, where the best possible score is 1. This metric is computed as

$$F_1 = 2 \frac{\text{precision} \times \text{recall}}{\text{precision} + \text{recall}},$$

where the precision is the ratio between correctly identified minority samples and the total number of minority samples, while the recall is given by the fraction of correctly identified minority samples over all samples.

C. Accomplishments

II. OVERVIEW OF THE TECHNIQUES

A. Undersampling

Undersampling is frequently employed to balance datasets before any machine learning algorithm is applied. Undersampling involves randomly removing entries from the majority class. Figure 1 shows the effects of undersampling on the Gaussian training dataset. The class imbalance is somewhat countered. However, the algorithm learnt from this undersampled dataset will be affected. Namely, its ability to generalize and its posterior distribution.

1) *Generalization Ability*: Induction algorithms require a sufficient amount of data to learn a model that generalizes well. If the training set is not large, a classifier may just memorize the characteristics of the training data. Moreover, undersampling has the potential of eliminating valuable samples from consideration of the classifier entirely [2], so it may exacerbate this problem of lack of data. The obtained training set may

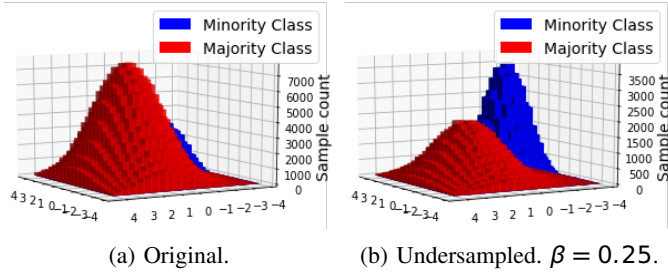


Fig. 1: Gaussian dataset.

vary greatly from one undersampling to another, this leads to a high variance of the learned model. Hence, the achievable complexity of the hypothesis set must be reduced to ensure a good generalization.

2) *Posterior Bias*: One goal of undersampling is to change the priori probabilities of the classes to make them more balanced. The classifier assumes that the features it encounters at testing follow the same distribution as the training set. This mismatch introduced by design is known as sampling selection bias [3] on posterior distribution.

Let $(\mathcal{X}, \mathcal{Y})$ denote the pairs of feature vectors, $\mathbf{x} \in \mathbb{R}^n$, and binary labels, $y \in \{0, 1\}$, contained in our original dataset. We assume that the number of samples labeled as zero is small compared with the number of samples in class one. Undersampling randomly removes points from the majority class, we describe this sampling with the binary random variable S , which takes the value 1 if a sample is selected.

It is reasonable to assume that the selection is independent of the features given the class. Then, applying Bayes rule, the law of total probability and noting that the samples from the minority class are always selected we obtain

$$\begin{aligned} p' &= P(y = 0 | \mathbf{x}, s = 1) = \frac{P(s = 1 | y = 0, \mathbf{x})P(y = 0 | \mathbf{x})}{P(s = 1 | \mathbf{x})} = \\ &= \frac{P(y = 0 | \mathbf{x})}{P(y = 0 | \mathbf{x}) + P(s = 1 | y = 1)P(y = 1 | \mathbf{x})} = \\ &= \frac{p}{p + \beta(1 - p)}, \end{aligned}$$

where p and p' denote the posterior probability of encountering a sample from the minority class when employing the original and the undersampled dataset respectively. Whereas β denotes the probability of keeping a sample from the majority class.

The posterior is highly affected by the rate of the sampling. As more samples are removed, the classification is more biased towards the minority class. Figure ?? shows the decision region of a naive Bayes classifier. As the training set is undersampled the region of points that are labeled as the minority class grows. The rate of undersampling not only influences the posterior bias, but also the algorithm's ability to generalize. Thus, β should be chosen with care. Figure ?? presents the average F1-score over different training sets. We

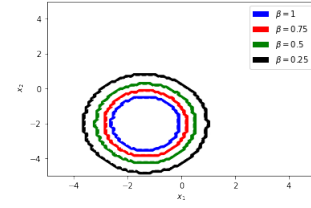


Fig. 2: Influence of undersampling on the classification region of a naive Bayes classifier trained with the Gaussian dataset. The area within each circle trained corresponds to the cluster of points that are classified as the minority class for a given undersampling rate.

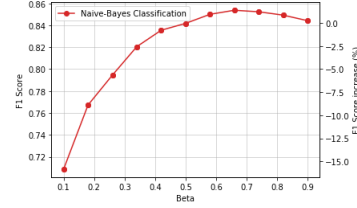


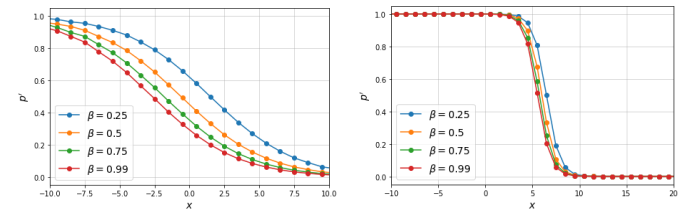
Fig. 3: F1-score vs. undersampling rate.

observe that the score is concave and in this case the optimum occurs with $\beta = 0.82$.

Another factor that strongly influences the posterior bias is class separability. The bias is higher when conditional distributions are similar across the classes [4]. To analyze this behaviour we reduced the problem to a one-dimensional setting, the results are depicted in Figure 4. We confirm that undersampling shifts the posterior distribution in favor of the minority class. Nevertheless, the shift caused by β is lower under the configuration with lower overlap.

B. Oversampling

Another technique to balance datasets is oversampling, which is artificially creating new instances of the minority class. This can be done just by cloning existing samples, but this is not really adding new information to the model. More elaborated methods create samples by interpolation. The most widely used is Synthetic Minority Oversampling Technique (SMOTE). It works by selecting a random sample, taking one near sample and creating points in between. This procedure



(a) $|\mu_0 - \mu_1| = 3$.

(b) $|\mu_0 - \mu_1| = 13$.

Fig. 4: Influence of undersampling on posterior probability of the minority class.

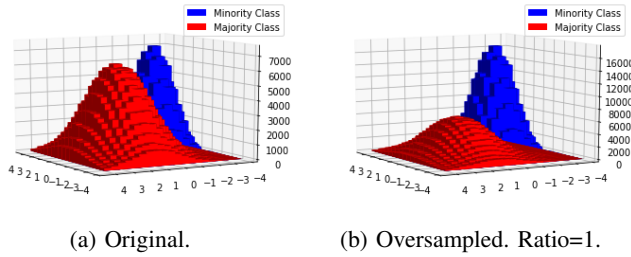


Fig. 5: Gaussian dataset.

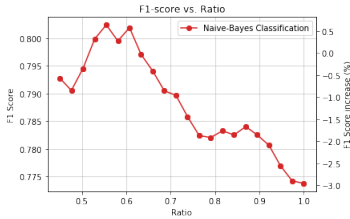


Fig. 6: F1-score vs. oversampling rate.

is repeated multiple times until the desired amount of new samples is created.

In Figure 5, it is possible to observe the effect of oversampling on the 2D dataset. For the chosen ratio of 1, the number of samples of both classes is the same.

A sensitivity analysis for several ratios has been conducted and the optimum ratio is determined to be around 0.55, which means that the minority class is oversampled until the number of minority samples is the 55% of the samples of the majority class. The increase in F1 score for this ratio is almost 0.6%. This can be appreciated in Figure 6.

C. Cost-Sensitive Techniques

Blabla

III. CLASSIFICATION IMPACT ON REAL DATA

Ensemble Classifiers are known to improve the performance of single classifiers in imbalanced datasets [A Review on Ensembles for the Class Imbalance] by combining separate weak learners into a composite whole. Both bagging and boosting algorithms have been implemented from scratch to explore their advantages and compare them with highly-effective classifiers such as XGBoost. [Check this](#) As we have stated, in recent years, Ensemble Classifiers have arisen as a possible solution to the class imbalance problem attracting great interest among researchers [45], [47], [50], [62].

Cita Fondo Flamenco: [5]

A. Bagging

Baggings algorithms are composed by two parts: bootstrapping and aggregation. The first part consist of selecting a subsample of the dataset. In bagging algorithms, weak learners are not trained over the whole dataset, but only over a part. For this task, decision tree classifiers are employed. Once all the classifiers are trained, the predictions of all models are

then aggregated to provide the final prediction. Important to mention that these algorithms are purely 'variance' reducing procedures intended to mitigate instability (especially associated to decision trees).

B. AdaBoost

Unlike bagging, boosting fits the weak learners in an adaptive way: a different significance is assigned to each of the base learners based on the samples that were misclassified in the previous iteration. It was first introduced by Schapire in 1990, who proved in [\[The strength of weak learnability\]](#) that a weak learner can be turned into a strong learner in the sense of probably approximately correct (PAC) learning framework. In particular, the AdaBoost algorithm stands out in the field of ensemble learning as one of the top ten data mining algorithms [\[Top 10 algorithms in data mining\]](#). One its main advantages is the versatility to incorporate cost-sensitive techniques. We implemented a custom AdaBoost classifier that enables us to gain control over the algorithm at a lower level, make adjustments when necessary and create other Adaboost-based classifiers such as AdaCost or Boosted SVM. The algorithm is fully detailed below.

Algorithm 1: AdaBoost Algorithm

Input: Training set $D = \{x_i, y_i\}, i = 1, \dots, N$; and $y_i \in \{-1, +1\}$; T : Number of iterations; I : Weak learner

Output: Boosted Classifier:
 $H(x) = \text{sign}\left(\sum_{t=1}^T \alpha_t h_t(x)\right)$ where h_t, α_t are the induced classifiers and their significance, respectively

```

1  $W_1(i) \leftarrow 1/N$  for  $i = 1, \dots, N$ 
  /* Create a weak learner in each iteration */
2 for  $t=1$  to  $T$  do
3    $h_t \leftarrow I(D, W)$ 
4    $\epsilon_t \leftarrow \sum_{i=1}^N W_t(i) [h_t(x_i) \neq y_i]$ 
5   if  $\epsilon_t > 0.5$  then
6      $T \leftarrow t - 1$ 
7   return
8    $\alpha_t = \frac{1}{2} \ln\left(\frac{1-\epsilon_t}{\epsilon_t}\right)$ 
  /* Update Weights */
9    $W_{t+1}(i) = W_t(i) e^{(-\alpha_t h_t(x_i) y_i)}$  for  $i = 1, \dots, N$ 
10  Normalize  $W_{t+1}$  such that  $\sum_{i=1}^N W_{t+1}(i) = 1$ 

```

This custom implementation uses single decision trees with 2 leaf nodes. It is given the number of maximum iterations (weak learners) that will be fitted, and a dataset D formed by N training samples, each constituted by a feature vector x_i and an associated binary label y_i . For each iteration, a weak classifier is created using the weighted training samples. For the first iteration, we allocate the same weight for all the samples ($1/N$). Once the decision tree has been created, we predict the class of the training samples and compute the weighted misclassification error ϵ_t . Then, the significance α_t

is calculated based on that error. Finally, the weights of each training samples are updated depending on the significance and whether the sample was correctly classified or not. Note that the weights are increased by e^α if the sample is misclassified or decreased by $e^{-\alpha}$ if correctly classified, with $\alpha_t \in [0, 1]$ because $\epsilon_t \in [0, 0.5]$.

AdaBoost uses the whole training samples to create each weak learner serially –in contrast to *Bagging*, where *bootstrap aggregating* is used to construct ensembles– giving more focus to challenging instances and turn the incorrect classifications into good predictions in the next iteration. Hence, AdaBoost can be considered as a “bias” reducing procedure, intended to increase the flexibility of stable (highly biased) weak learners when incorporated properly in a serial additive expansion. That is the reason behind why weak learners with low variance but high bias –such as decision trees– are well adapted for boosting. Lastly the overall literature suggest AdaBoost algorithms are generally resistant to overfitting regardless the number of weak estimators we use, and there are very few cases reported to overfit the training data [[Additive Logistic Regression: A Statistical View of Boosting]. This fascinating issue is explained due to 1) as the iterations proceed the weak learners tend to have less significance and 2) similarly to SVM, ensemble classifiers maximize the margin which allows promoting good generalization [The dynamics of AdaBoost: Cyclic behavior and convergence of margin]. The impact of its complexity on overfitting is covered in the result section.

C. AdaCost

The AdaBoost algorithm is an accuracy-oriented algorithm. In other words, it assumes each class has an even distribution. In case the class distribution is uneven, the algorithm may incur in systematic biases toward the majority class. Therefore, several methods have been proposed to incorporate an asymmetric weight update and eliminate these biases. The literature suggest two different strategies to implement this changes [Cost-sensitive boosting algorithms: Do we really need them?]: 1) modify the model learnt from data or 2) modifying how the model is used to make a decision. AdaCost falls into the former category, and it aims at modifying the update of the weights based on a slight modification:

$$W_{t+1}(i) = W_t(i) \exp\left(-\alpha_t y_i h_t(x_i) \boxed{\phi(i)}\right) \text{ with } W_1(i) = \frac{1}{\sum_{i=1}^n C_i}$$

Where the new boxed term is called cost-adjustment function. We use this function to allocate different penalizations across different classes. This function is left for design since the literature suggest more than one alternatives. In this project we used the function suggested in [A Review on Ensembles for the Class Imbalance]:

$$\phi(i) = 0.5 C_i (\mathbb{1}\{h_t(\mathbf{x}) = 1\} - \mathbb{1}\{h_t(\mathbf{x}) = -1\}) + 0.5$$

Where C_i is a hyper-parameter that establishes the misclassification cost of the sample i , which ultimately depends

upon the class of that sample. This cost-adjustment function yields to an upper bound cumulative misclassification cost equal to $d \prod_{t=1}^T Z_t$, where $d = \sum C_i$ and Z_t is the sum of the costs calculated for W_{t+1} , i.e. the coefficient that we use to normalize the weights in AdaBoost [AdaCost]. Due to the fact that boosting minimizes Z_t , the significance ultimately needs to be updated in a different way [Improved boosting algorithms using confidence-rated predictions.]:

$$\alpha = \frac{1}{2} \ln \frac{1+r}{1-r} \quad \text{where } r = \sum_i D(i) u_i, \quad u_i = y_i h(x_i) \phi(i)$$

All in all, the AdaCost algorithm follows the same procedure as AdaBoost with changes in the significance and weight update rule in order to incorporate an asymmetric penalization cost.

D. AdaMEC

The AdaMEC algorithm falls into the second category introduced in the previous part. The model is trained with the original AdaBoost, but modifies the decision rule by exploiting the bayesian decision theory:

$$H_{AdaMEC}(\mathbf{x}) = \text{sign} \left(\sum_{y \in \{-1, +1\}} c(y) \sum_{\tau: h_\tau(\mathbf{x}) = y} \alpha_\tau h_\tau(\mathbf{x}) \right)$$

This is called Minimum Expected Cost rule and it assumes that the wighted votes of each weak learners are proportional to the class probabilities. We have incorporated this classifier in this study as it has been reported as one of the best cost-sensitive AdaBoost classifiers [Cost-sensitive boosting algorithms: Do we really need them?].

E. Boosting SVM

One of the practical concerns about AdaBoost is the accuracy/diversity dilemma [A study of adaboost with SVM based on weak learners], which means that when the weak learners are actually strong classifiers and they do not disagree too much on their vote, the performance of boosting is weakened. Support vector machines (SVM) are known to be strong learners that minimize structural risk [12], so in a first glance they should not work well when ensemble. However, the literature [Boosting Support Vector Machines Successfully][A Study of AdaBoost with SVM Based Weak Learners][Boosting Support Vector Machines] suggests that if properly weakened, Boosted SVM can exceed the classification performance of traditional SVMs or AdaBoost on imbalanced datasets, and gain control over the accuracy/diversity issue. To do that, we have used radial basis function kernel (RBF), which uses the Gaussian width σ to map the features on a high dimensional feature space. The performance of RBF SVM can be conveniently controlled by adjusting σ . Large values of sigma will lead to very weak classifiers. The full algorithm description is detailed below:

The main disadvantage of SVM is that the running time to train the model does not scale well with the size of the

Algorithm 2: Boosted SVM Algorithm

Input: Training set $D = \{x_i, y_i\}, i = 1, \dots, N$; and $y_i \in \{-1, +1\}$; T : Maximum number of iterations; The initial $\sigma = \sigma_{ini}, \sigma_{min}, \sigma_{step}$

Output: Boosted Classifier:

$H(x) = \text{sign}\left(\sum_{t=1}^T \alpha_t h_t(x)\right)$ where h_t, α_t are the induced classifiers and their significance, respectively

```

1  $W_1(i) \leftarrow 1/N$  for  $i = 1, \dots, N$ 
  /* Create a weak learner in each iteration */
2 while  $\sigma > \sigma_{min}$  and  $t < T$  do
3    $t \leftarrow t + 1$ 
4    $h_t \leftarrow \text{RBFSVM}(D, W, \sigma)$ ; // Train RBFSVM
5    $\epsilon_t \leftarrow \sum_{i=1}^N W_t(i)[h_t(x_i) \neq y_i]$ 
6   if  $\epsilon_t > 0.5$  then
7      $\sigma \leftarrow \sigma - \sigma_{step}$ 
8   else
9      $\alpha_t = \frac{1}{2} \ln\left(\frac{1-\epsilon_t}{\epsilon_t}\right)$ 
10    /* Update Weights */
11     $W_{t+1}(i) = W_t(i)e^{(-\alpha_t h_t(x_i) y_i)}$  for  $i = 1, \dots, N$ 
    Normalize  $W_{t+1}$  such that  $\sum_{i=1}^N W_{t+1}(i) = 1$ 

```

training set [Elkin]. Our training set has more than 30,000 samples so the training would take a significant amount of time. Therefore, a preprocessing step was applied to the training data, whose dimensionality is reduced using PCA and ultimately under-sampled to speed-up the training process.

[TO-DO: Add comparison with SVM somehow]

F. XGBoost

XGBoost is one of the most effective classification algorithms nowadays, and it has been the winner of many competitions in the last years. The working principle is additive tree learning, which means that what it is learned on an iteration is used to train the next tree of the model. This new tree has to minimize the objective function, that includes a training loss and a regularization term. The second order Taylor expansion is employed, as the actual function can be very complex. The structure of the new tree is going to be optimized to produce the maximum reduction of the objective function. The tree split that gives the maximum similarity between its prediction and the output is the one that it is going to be added.

In practice, it is the classification algorithm that gives the highest F1-score and accuracy. The XGBoost library is used to train this algorithm for our dataset.

IV. RESULTS AND DISCUSSION

This section discusses the performance of each classifier in terms of imbalance impact, training error and diversity.

A. Performance Results

The best performance results of each methodology are shown in the table below. This table shows the metrics of the best F-1 score so it is possible to compare the algorithms from multiple perspectives.

TABLE I: Best Classification Performances

Type of Method	Classifier	Scores				
		F-1	Acc.	Prec.	Rec.	AUC
Ensemble	AdaBoost*	84.14	84.55	84.05	84.55	77.91
	Bagging*	76.32	78.49	76.91	78.49	65.58
	R. Forest	70.73	76.96	78.39	76.96	57.71
	BoostSVM*	62.18	73.42	53.91	73.42	0.5
	XGBoost	88.28	88.63	88.38	88.63	82.82
Cost Sensitive	AdaCost*	77.98	79.86	78.65	79.86	67.71
	AdaMEC					
Hybrid (using under-sampling)	AdaBoost*	83.12	82.89	83.45	82.89	79.52
	Bagging*	74.91	73.78	77.48	73.78	72.45
	R. Forest	80.14	80.87	79.97	80.87	72.32
	BoostSVM*					
	XGBoost	87.57	87.51	87.65	87.51	84.53
	AdaCost*	78.67	79.22	78.39	79.22	71.04
	AdaMEC					

*Custom implementation.

In the category of ensemble algorithms, XGBoost outperforms every other algorithm in all the metrics. F1-score is close to 90%. Next algorithm is AdaBoost, whose F1-score is above 84%. The rest of ensemble classifiers are clearly inferior. Regarding cost sensitive algorithm, for this particular dataset it doesn't give any advantage. For hybrid algorithms, which combine ensemble and cost-sensitive algorithms plus undersampling, a beta of 0.55 is chosen. A few of the classifiers improve, as it is the case of Random Forest and AdaCost, while the others become a little bit more. However, XGBoost still has the largest F1-score.

B. Impact of Imbalance Data on Classification

Chart of F1-score vs. Imbalance

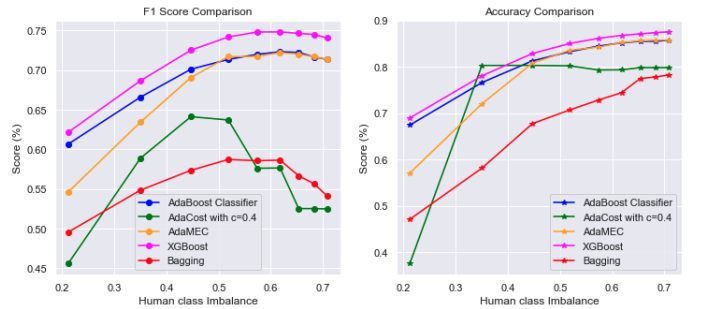


Fig. 7: Impact of imbalanced training set on classification.

– Comments –

C. Overfitting Analysis

Chart of risk (test and training errors) vs. Complexity

D. Diversity Analysis

Chart of Accuracy vs. Diversity

Blabla

A. Abbreviations and Acronyms

Define abbreviations and acronyms the first time they are used in the text, even after they have been defined in the abstract. Abbreviations such as IEEE, SI, MKS, CGS, ac, dc, and rms do not have to be defined. Do not use abbreviations in the title or heads unless they are unavoidable.

B. Units

- Use either SI (MKS) or CGS as primary units. (SI units are encouraged.) English units may be used as secondary units (in parentheses). An exception would be the use of English units as identifiers in trade, such as “3.5-inch disk drive”.
- Avoid combining SI and CGS units, such as current in amperes and magnetic field in oersteds. This often leads to confusion because equations do not balance dimensionally. If you must use mixed units, clearly state the units for each quantity that you use in an equation.
- Do not mix complete spellings and abbreviations of units: “Wb/m²” or “webers per square meter”, not “webers/m²”. Spell out units when they appear in text: “. . . a few henries”, not “. . . a few H”.
- Use a zero before decimal points: “0.25”, not “.25”. Use “cm³”, not “cc”.)

C. Equations

Number equations consecutively. To make your equations more compact, you may use the solidus (/), the exp function, or appropriate exponents. Italicize Roman symbols for quantities and variables, but not Greek symbols. Use a long dash rather than a hyphen for a minus sign. Punctuate equations with commas or periods when they are part of a sentence, as in:

$$a + b = \gamma \quad (1)$$

Be sure that the symbols in your equation have been defined before or immediately following the equation. Use “(1)”, not “Eq. (1)” or “equation (1)”, except at the beginning of a sentence: “Equation (1) is . . .”

D. \LaTeX -Specific Advice

Please use “soft” (e.g., `\eqref{Eq}`) cross references instead of “hard” references (e.g., (1)). That will make it possible to combine sections, add equations, or change the order of figures or citations without having to go through the file line by line.

Please don’t use the `{eqnarray}` equation environment. Use `{align}` or `{IEEEeqnarray}` instead. The `{eqnarray}` environment leaves unsightly spaces around relation symbols.

Please note that the `{subequations}` environment in \LaTeX will increment the main equation counter even when there are no equation numbers displayed. If you forget that, you might write an article in which the equation numbers skip from (17) to (20), causing the copy editors to wonder if you’ve discovered a new method of counting.

\BibTeX does not work by magic. It doesn’t get the bibliographic data from thin air but from .bib files. If you use \BibTeX to produce a bibliography you must send the .bib files.

\LaTeX can’t read your mind. If you assign the same label to a subsubsection and a table, you might find that Table I has been cross referenced as Table IV-B3.

\LaTeX does not have precognitive abilities. If you put a `\label` command before the command that updates the counter it’s supposed to be using, the label will pick up the last counter to be cross referenced instead. In particular, a `\label` command should not go before the caption of a figure or a table.

Do not use `\nonumber` inside the `{array}` environment. It will not stop equation numbers inside `{array}` (there won’t be any anyway) and it might stop a wanted equation number in the surrounding equation.

E. Some Common Mistakes

- The word “data” is plural, not singular.
- The subscript for the permeability of vacuum μ_0 , and other common scientific constants, is zero with subscript formatting, not a lowercase letter “o”.
- In American English, commas, semicolons, periods, question and exclamation marks are located within quotation marks only when a complete thought or name is cited, such as a title or full quotation. When quotation marks are used, instead of a bold or italic typeface, to highlight a word or phrase, punctuation should appear outside of the quotation marks. A parenthetical phrase or statement at the end of a sentence is punctuated outside of the closing parenthesis (like this). (A parenthetical sentence is punctuated within the parentheses.)
- A graph within a graph is an “inset”, not an “insert”. The word alternatively is preferred to the word “alternately” (unless you really mean something that alternates).
- Do not use the word “essentially” to mean “approximately” or “effectively”.
- In your paper title, if the words “that uses” can accurately replace the word “using”, capitalize the “u”; if not, keep using lower-cased.
- Be aware of the different meanings of the homophones “affect” and “effect”, “complement” and “compliment”, “discreet” and “discrete”, “principal” and “principle”.
- Do not confuse “imply” and “infer”.
- The prefix “non” is not a word; it should be joined to the word it modifies, usually without a hyphen.
- There is no period after the “et” in the Latin abbreviation “et al.”.
- The abbreviation “i.e.” means “that is”, and the abbreviation “e.g.” means “for example”.

An excellent style manual for science writers is [?].

F. Authors and Affiliations

The class file is designed for, but not limited to, six authors. A minimum of one author is required for all conference articles. Author names should be listed starting from left

to right and then moving down to the next line. This is the author sequence that will be used in future citations and by indexing services. Names should not be listed in columns nor group by affiliation. Please keep your affiliations as succinct as possible (for example, do not differentiate among departments of the same organization).

G. Identify the Headings

Headings, or heads, are organizational devices that guide the reader through your paper. There are two types: component heads and text heads.

Component heads identify the different components of your paper and are not topically subordinate to each other. Examples include Acknowledgments and References and, for these, the correct style to use is “Heading 5”. Use “figure caption” for your Figure captions, and “table head” for your table title. Run-in heads, such as “Abstract”, will require you to apply a style (in this case, italic) in addition to the style provided by the drop down menu to differentiate the head from the text.

Text heads organize the topics on a relational, hierarchical basis. For example, the paper title is the primary text head because all subsequent material relates and elaborates on this one topic. If there are two or more sub-topics, the next level head (uppercase Roman numerals) should be used and, conversely, if there are not at least two sub-topics, then no subheads should be introduced.

H. Figures and Tables

a) *Positioning Figures and Tables:* Place figures and tables at the top and bottom of columns. Avoid placing them in the middle of columns. Large figures and tables may span across both columns. Figure captions should be below the figures; table heads should appear above the tables. Insert figures and tables after they are cited in the text. Use the abbreviation “Fig. 8”, even at the beginning of a sentence.

TABLE II: Table Type Styles

Table Head	Table Column Head		
	Table column subhead	Subhead	Subhead
copy	More table copy ^a		

^aSample of a Table footnote.



Fig. 8: Example of a figure caption.

Figure Labels: Use 8 point Times New Roman for Figure labels. Use words rather than symbols or abbreviations when writing Figure axis labels to avoid confusing the reader. As an

example, write the quantity “Magnetization”, or “Magnetization, M”, not just “M”. If including units in the label, present them within parentheses. Do not label axes only with units. In the example, write “Magnetization (A/m)” or “Magnetization {A[m(1)]}”, not just “A/m”. Do not label axes with a ratio of quantities and units. For example, write “Temperature (K)”, not “Temperature/K”.

ACKNOWLEDGMENT

The preferred spelling of the word “acknowledgment” in America is without an “e” after the “g”. Avoid the stilted expression “one of us (R. B. G.) thanks . . .”. Instead, try “R. B. G. thanks. . .”. Put sponsor acknowledgments in the unnumbered footnote on the first page.

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