

Improving Imbalanced Learning in Land Cover Classification

A Heuristic Oversampling Method Based on K-Means and SMOTE

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Land cover maps are an important resource to make informed policy, development, planning and resource management decisions. The primary challenge for the development of accurate, timely and automated Land Use/Land Cover maps are technical skills. Specifically, remotely sensed data is often imbalanced, where the number of samples of a few classes is significantly greater the number of samples of the remaining classes. This asymmetric class distribution, impacts negatively the performance of classifiers and adds a new source of inaccuracy to the production of these maps. In this paper, we address this problem, known as the imbalanced learning problem, by using K-Means SMOTE, a recently proposed oversampling method. K-Means SMOTE is an oversampling algorithm that attempts to improve the quality of newly created artificial data by avoiding the generation of noisy data and effectively overcome data imbalance. The performance of K-Means SMOTE is compared to other popular oversampling methods using seven well known datasets and a variety of classifiers and evaluation metrics. The results show that the proposed method consistently outperforms the remaining oversamplers and produces higher quality land cover classifications.

1 Introduction

The increasing amount of remote sensing missions granted the access to dense time series (TS) data at a global level and provides up-to-date, accurate land cover information [Drusch et al., 2012]. This information is often materialized through Land Use/Land Cover (LULC) maps, which constitute an essential asset for various purposes, such as land cover change detection, urban planning, environmental monitoring and natural hazard assessment [Khatami et al., 2016]. However, the timely production of accurate and updated LULC maps is still a challenge within the remote sensing community [Wulder et al., 2018]. LULC maps are produced based on two main approaches: photo-interpreted by the human eye, or automatic mapping using remotely sensed data and classification algorithms.

While photo-interpreted LULC maps rely on human operators and can be more reliable, they also present some significant disadvantages. The most important disadvantage are the production costs, in fact photo-interpretation consumes significant resources, both money and time. Because of that, they

are not frequently updated and not suitable for operational mapping over large areas. Finally, there is also the issue of overlooking rare or small-area classes, due to factors such as the minimum mapping unit being used.

Automatic mapping with classification algorithms based on machine-learning (ML) have been extensively researched and used to speed up and reduce the costs of the production process. Improvements in classification algorithms are sure to have significant impact in the efficiency with which remote sensing imagery is used. Several challenges need to be tackled in order to improve automatic classification:

1. Improve the ability to handle high-dimensional datasets, in cases such as Multi-spectral TS composites high-dimensionality increases the complexity of the problem and creates a strain on computational power [Stromann et al., 2020].
2. Improve class separability, as the production of an accurate LULC map can be hindered by the existence of classes with similar spectral signatures, making these classes difficult to distinguish [Alonso-Sarria et al., 2019].
3. Resilience to mislabelled LULC patches, as the use of photo-interpreted training data poses a threat to the quality of any LULC map produced with this strategy, since factors such as the minimum mapping unit tend to cause the overlooking of small-area LULC patches and generates noisy training data that may reduce the prediction accuracy of a classifier [Pelletier et al., 2017].
4. Dealing with rare land cover classes, due to the varying levels of area coverage for each class. In this case using a purely random sampling strategy will amount to a dataset with a roughly proportional class distribution as the one on the landscape. On the other hand, the acquisition of training datasets containing balanced class frequencies is often unfeasible. This causes an asymmetry in class distribution, where some classes are frequent in the training dataset, while others have little expression [Wang et al., 2019, Feng et al., 2019].

The latter challenge is known as the imbalanced learning problem [Chawla et al., 2004]. It is defined as a skewed distribution of observations found in a dataset among classes in both binary and multi-class problems [Abdi and Hashemi, 2016]. This asymmetry in class distribution negatively impacts the performance of classifiers, especially in multi-class problems. During the learning phase, classifiers are optimized to maximize an objective function, with overall accuracy being the most common one [Maxwell et al., 2018]. This means that observations belonging to minority classes contribute less to the optimization process, translating into a bias towards majority classes. As an example, a trivial classifier can achieve 99% overall accuracy on a binary dataset where 1% of the observations belong to the minority class if it classifies all observations as belonging to the majority class. This is an especially significant issue in the automatic classification of LULC maps, as the distribution of the different land-use classes tends to be highly imbalanced. Therefore, improvements in the ability to deal with imbalanced datasets will translate into important progress in the automatic classification of LULC maps.

There are three different types of approaches to deal with the class imbalance problem [Fernández et al., 2013, Kaur et al., 2019]:

1. Cost-sensitive solutions. Introduces a cost matrix to the learning phase with misclassification costs attributed to each class. Minority classes will have a higher cost than majority classes, forcing the algorithm to be more flexible and adapt better to predict minority classes.
2. Algorithmic level solutions. Specific classifiers are modified to reinforce the learning on minority classes. Consists on the creation or adaptation of classifiers.

3. Resampling solutions. Rebalances the dataset’s class distribution by removing majority class instances and/or generating artificial minority instances. This is considered an external approach, where the intervention occurs before the learning phase, benefitting from versatility and independency from the classifier used.

It is important to note that resampling strategies have a significant advantage over the other approaches. By operating at the data level, resampling strategies allow the use of any off the shelf algorithm, without the need for any type of changes or adaptations to the algorithm. This is a significant advantage especially considering that most users in remote sensing are not expert machine learning engineers.

Within resampling approaches there are three subgroups of approaches [Fernández et al., 2013, Kaur et al., 2019, Luengo et al., 2020]:

1. Undersampling methods, which rebalance class distribution by removing instances from the majority classes.
2. Oversampling methods, which rebalance datasets by generating new artificial instances belonging to the minority classes.
3. Hybrid methods, which are a combination of both oversampling and undersampling, resulting in the removal of instances in the majority classes and the generation of artificial instances in the minority classes.

Resampling methods can be further distinguished between non-informed and heuristic (i.e., informed) resampling techniques [Fernández et al., 2013, Luengo et al., 2020, García et al., 2016]. The former consist of methods that duplicate/remove a random selection of data points to set class distributions to user-specified levels, and are therefore a simpler approach to the problem. The latter consists of more sophisticated approaches that aim to perform over/undersampling based on the points’ contextual information within their data space.

The imbalanced learning problem is not new in machine learning but its relevancy has been growing, as attested by [Haixiang et al., 2017]. The problem has also been addressed in the context of remote sensing [Douzas et al., 2019]. In this paper, we propose the K-means SMOTE (K-SMOTE) [Douzas et al., 2018] oversampler to address the imbalanced learning problem in a multiclass context for LULC classification using various remote sensing datasets. The K-SMOTE algorithm presents significant advantages over other oversamplers, by coupling two different procedures in the generation of artificial data. The algorithm starts by clustering the observations; next, the generation of the artificial observations is done taking into consideration the distribution of majority/minority cases in each individual cluster. The idea of starting with a clustering procedure before the data generation phase is important in remote sensing because the spectral signature of the different classes can change significantly based on the geographical area in which it is represented. In other words, the spectral signature of a specific class can vary greatly depending on the geography, meaning that often we will be facing within-class imbalance [Japkowicz, 2001].

In fact, we can decompose class imbalance into two different types: between-class imbalance and within-class imbalance [Douzas et al., 2018, Jo and Japkowicz, 2004]. While the first refers to the overall asymmetry between majority and minority classes, the second results from the fact that in different areas of the input space there might be different levels of imbalance. Depending on the complexity of the input space, different subclusters of minority and majority instances may be present. In order to achieve a balance between minority and majority instances, these subclusters should be treated separately. Assuming that the role of a classifier is to create rules in such a way that it is able to isolate the different relevant sub-concepts that represent both the majority and minority classes, the classifier will

create multiple disjunct rules that describe these concepts. If the input space is simple and the classes' instances are grouped together in a unique cluster, the classifier will only need to create (general) rules that comprise large portions of instances belonging to the same class. To the contrary, if the input space is complex and scatters through multiple small clusters, the classifier will need to learn a more complex set of (specific) rules, which can be seen in Figure 1. It is important to note that small clusters can happen both in the minority and majority class, although they will tend to be more frequent in the minority class due to its underrepresentation.

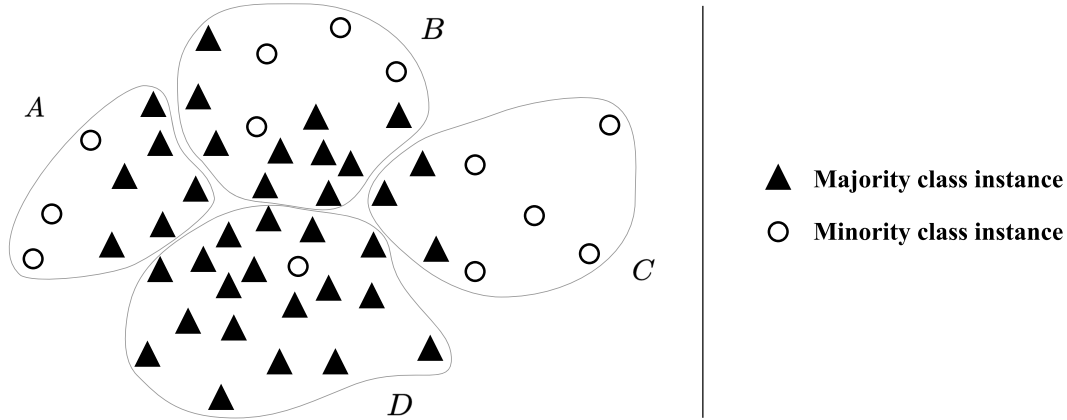


Figure 1: Example of a complex input space. In this example, a classifier would need to separate the minority class' samples across 4 distinguishable regions, or clusters.

The efficacy of K-SMOTE is tested using different types of classifiers. To do so, we employ both commonly used and state-of-the-art oversamplers as benchmarking methods: Random oversampling (ROS), Synthetic Minority Oversampling Technique (SMOTE) [Chawla et al., 2002] and Borderline-SMOTE (B-SMOTE) [Han et al., 2005]. Also as a baseline score we include classification results without the use of any resampling method.

This paper is organized in 5 sections: section 2 provides an overview of the state-of-art, section 3 describes the proposed methodology, section 4 covers the results and discussion and section 5 presents the conclusions taken from this study.

2 Imbalanced Learning Approaches

Imbalanced learning has been addressed in three different ways: over/undersampling, cost-sensitive training and changes/adaptations in the learning algorithms [Kaur et al., 2019]. These approaches impact different phases of the learning process, while over/undersampling can be seen as a pre-processing step, cost-sensitive and changes in the algorithm imply a more customized and complex intervention in the algorithms. In this section, we focus on previous work related with resampling methods, while providing a brief explanation of cost-sensitive and algorithmic level solutions.

All of the most common classifiers used for LULC classification tasks [Khatami et al., 2016, Gavade and Rajpurohit, 2019] are sensitive to class imbalance [Blagus and Lusa, 2010]. Algorithm-based approaches typically focus on adaptations based on ensemble classification methods [Mellor et al., 2015] or common non-ensemble based classifiers such as Support Vector Machines [Shao et al., 2014]. In [Lee et al., 2016], the reported results show that algorithm-based methods have comparable performance to resampling methods.

Cost-sensitive solutions refer to changes in the importance attributed to each instance through a cost matrix [Huang et al., 2016, Cui et al., 2019, Dong et al., 2017]. A common cost sensitive solution is found in [Huang et al., 2016]. The authors use the inverse class frequency (i.e., $1/|C_i|$) to give higher weight to minority classes. Cui et al. [Cui et al., 2019] extended this method by adding a hyperparameter β to class weights as $(1 - \beta)/(1 - \beta^{|C_i|})$. When $\beta = 0$, no re-weighting is done. When $\beta \rightarrow 1$, weights are the inverse of the frequency class matrix. Another method [Dong et al., 2017] explores adaptations of Cross-entropy classification loss by adding different formulations of class rectification loss.

Resampling (over/undersampling) is the most common approach to imbalanced Learning is most commonly addressed through data resampling in machine learning in general and remote sensing in particular [Feng et al., 2019]. The generation of artificial instances (i.e., augmenting the dataset), based on rare examples, is done independently of any other step in the learning process. Once the procedure is applied, any standard machine learning algorithm can be used. Its simplicity makes resampling strategies particularly appealing for any user (especially the non-sophisticated user) interested in applying several classifiers, while maintaining a simple approach. It is also important to notice that over/undersampling methods can also be easily applied to multiclass problems, common in LULC classification tasks.

2.1 Non-informed resampling methods

There are two main non-informed resampling methods. Random Oversampling (ROS) generates artificial observations through random duplication of minority instances. This method is used in remote sensing [Shariffar et al., 2019, Hounkpatin et al., 2018] for its simplicity, even though its mechanism makes the classifier prone to overfitting [Krawczyk, 2016]. Hounkpatin et al. [Hounkpatin et al., 2018] found that using ROS returned worse results than keeping the original imbalance in their dataset.

A few of the recent remote sensing studies employed Random Undersampling (RUS) [Ferreira et al., 2019], which randomly removes observations belonging to majority classes. Although it's not as prone to overfitting as ROS, it incurs into information loss by eliminating observations from the majority class [Feng et al., 2019].

Another disadvantage of non-informed resampling methods is their performance-wise inconsistency across classifiers. ROS' impact on the Indian Pines dataset was found inconsistent between Random Forest Classifiers (RFC) and Support Vector Machines (SVM) and lowered the predictive power of an artificial neural network (ANN) [Maxwell et al., 2018]. Similarly, RUS is found to generally lead to a lower overall accuracy due to the associated information loss [Maxwell et al., 2018].

2.2 Heuristic methods

The methods presented in this section appear as a means to overcome the insufficiencies found in non-informed resampling. They use either local or global information to generate new, relevant, non-duplicated instances to populate the minority classes and/or remove irrelevant instances from majority classes. In a comparative analysis between over- and undersamplers' performance for LULC classification [Feng et al., 2018] using the rotation forest ensemble classifier, authors found that oversampling methods consistently outperform undersampling methods. This result led us to exclude undersampling from our study.

SMOTE [Chawla et al., 2002] was the first heuristic oversampling algorithm to be proposed and has been the most popular one since then, likely due to its fair degree of simplicity and quality of generated data. It takes a random minority class sample and introduces synthetic examples along the line segment that

join a random k minority class nearest neighbor to the selected sample. Specifically, a single synthetic sample \vec{z} is generated within the line segment of a randomly selected minority class observation \vec{x} and one of its k nearest neighbors \vec{y} such that $\vec{z} = \alpha \vec{x} + (1 - \alpha) \vec{y}$, where α is a random floating point between 0 and 1, as shown in Figure 2.

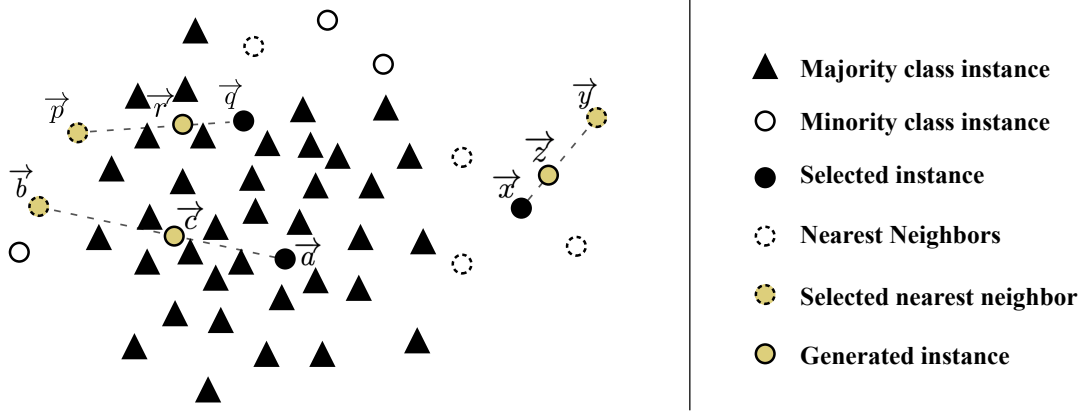


Figure 2: Example of SMOTE’s data generation process.

A number of studies implement SMOTE within the LULC classification context and reported improvements on the quality of the trained predictors [Jozdani et al., 2019, Bogner et al., 2018]. Another study proposes an adaptation of SMOTE on an algorithmic level for deep learning applications [Zhu et al., 2020]. This method combines both typical computer vision data augmentation techniques, such as image rotation, scaling and flipping on the generated instances to populate minority classes. Another algorithmic implementation is the variational semi-supervised learning model [Cenggoro et al., 2018]. It consists of a generative model that allows learning from both labeled and unlabeled instances while using SMOTE to balance the data.

Despite SMOTE’s popularity, its limitations have motivated the development of more sophisticated oversampling algorithms [Douzas and Bacao, 2019, Han et al., 2005, Ma and Fan, 2017, Douzas and Bacao, 2017, Douzas et al., 2018, Haibo He et al., 2008]. [Douzas and Bacao, 2019] identify four major weaknesses of the SMOTE algorithm, which can be summarized as:

1. Generation of noisy instances due to random selection of a minority observation to oversample. The random selection of a minority observation makes SMOTE oversampling prone to the amplification of existing noisy data. This has been addressed by variants such as B-SMOTE [Han et al., 2005] and ADASYN [Haibo He et al., 2008].
2. Generation of noisy instances due to the selection of the k nearest neighbors. In the event an observation (or a small number thereof) is not noisy but is isolated from the remaining clusters, known as the "small disjuncts problem" [Holte et al., 1989], much like sample \vec{b} from Figure 2, the selection of any nearest neighbor of the same class will have a high likelihood of producing a noisy sample.
3. Generation of nearly duplicated instances. Whenever the linear interpolation is done between two observations that are close to each other, the generated instance becomes very similar to its parents and increases the risk of overfitting. G-SMOTE [Douzas and Bacao, 2019] attempts to address both the k nearest neighbor selection mechanism problem as well as the generation of nearly duplicated instances problem.

4. Generation of noisy instances due to the use of observations from two different minority class clusters. Although an increased k could potentially avoid the previous problem, it can also lead to the generation of artificial data between different minority clusters, as depicted Figure 2 with the generation of point \vec{r} using minority class observations \vec{p} and \vec{q} . Cluster-based oversampling methods attempt to address this problem.

This last issue, the generation of noisy instances due to the existence of several minority class clusters, is particularly relevant in remote sensing. It is frequent that instances belonging to the same minority class can have different spectral signatures, meaning that they will be clustered in different parts of the input space. In this context, the use of SMOTE will lead to the generation of noisy instances of the minority class. This problem can be efficiently mitigated through the use of a cluster-based oversampling. According to our literature review cluster-based oversampling approaches have never been applied in the context of remote sensing. On the other hand, while there are several references of the application of cluster-based oversampling in the context of machine learning, there is no such application for the multiclass case, a fundamental requirement for the application of oversampling in the context of LULC.

Cluster-based oversampling approaches introduce an additional layer to SMOTE’s selection mechanism, which is done through the inclusion of a clustering process. This ensures that both between-class data balance and within each class balance is preserved. The self-organizing map oversampling (SOMO) [Douzas and Bacao, 2017] algorithm transforms the dataset into a 2-dimensional input, where the areas with the highest density of minority samples are identified. SMOTE is then used to oversample each of the identified areas separately. CURE-SMOTE [Ma and Fan, 2017] applies a hierarchical clustering algorithm (CURE) to discard isolated minority instances before applying SMOTE. Although it avoids noise generation problems, it ignores within-class data distribution. Another method [Santos et al., 2015] uses K-means to cluster the entire input space and applies SMOTE to clusters with the fewest observations, regardless of their class label. The label of the generated observation is copied from one of its parents. This method cannot ensure a balanced dataset since class imbalance is not specifically addressed, but rather dataset imbalance.

K-SMOTE [Douzas et al., 2018] avoids noisy data generation by modifying the data selection mechanism. It employs k -means clustering to identify safe areas using cluster-specific Imbalance Ratio (defined by $\frac{\text{count}(C_{\text{majority}})}{\text{count}(C_{\text{minority}})})$ and determine the quantity of generated samples per cluster based on a density measure. These samples are finally generated using the SMOTE algorithm. The K-SMOTE’s data generation process is depicted in Figure 3. Note that the number of samples generated for each cluster varies according to the sparsity of each cluster (the sparser the cluster is, the more samples will be generated) and a cluster is rejected if the cluster’s IR surpasses the threshold. Therefore, this method can be combined with any data generation mechanism, such as G-SMOTE. Also K-SMOTE includes the SMOTE algorithm as a special case when the number of clusters is set to one. Consequently, K-SMOTE is always guaranteed to return results as good as or better than SMOTE.

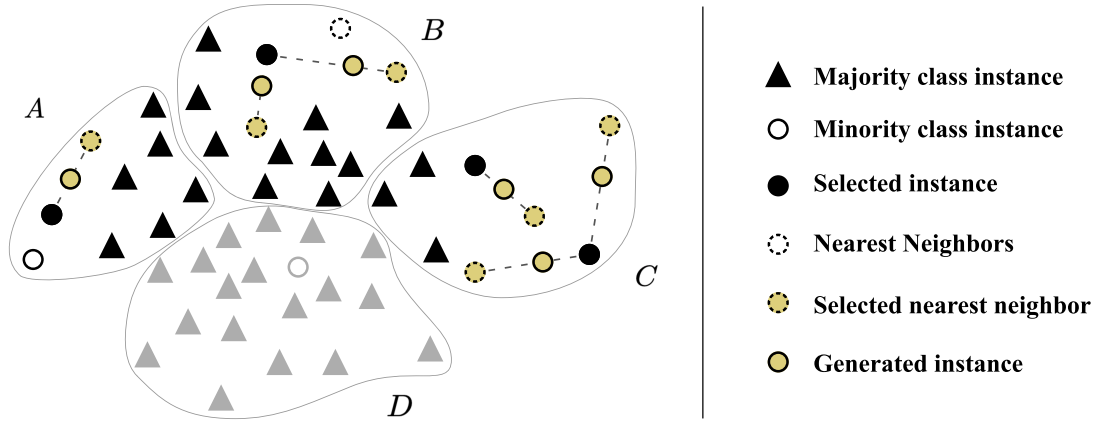


Figure 3: Example of K-SMOTE’s data generation process. Clusters *A*, *B* and *C* are selected for oversampling, whereas cluster *D* was rejected due to its high imbalance ratio. The oversampling is done using the SMOTE algorithm and the k nearest neighbors selection only considers observations within the same cluster.

Although no other study was found to implement cluster-based oversampling, another study [Douzas et al., 2019] compared the performance of SMOTE, ROS, ADASYN, B-SMOTE and G-SMOTE in a highly imbalanced LULC classification dataset. The authors found that G-SMOTE consistently outperformed the remaining oversampling algorithms regardless of the classifier used.

This paper’s main contributions are:

- Propose a cluster-based multiclass oversampling method appropriate for LULC classification;
- Compare its performance with the remaining oversamplers in a multiclass context with a set of widely used LULC classification datasets. Allows us to check for oversamplers’ performance statistical significance across datasets and report K-SMOTE’s performance in benchmark LULC datasets.
- Introducing a cluster-based oversampling algorithm within the remote sensing domain, as well as comparing its performance with the remaining oversamplers in a multiclass context.

3 Methodology

The purpose of this work is to understand the performance of K-SMOTE as opposed to other popular and/or state-of-the-art oversamplers for LULC classification. To do so, we employ 7 LULC datasets along with 3 evaluation metrics and 5 classifiers to evaluate the performance of oversamplers. In this section we describe the datasets, evaluation metrics, oversamplers, classifiers and software used as well as the procedure developed.

3.1 Datasets

The datasets used were extracted from publicly available hyperspectral scenes. Information regarding each of these scenes is provided in this subsection. A similar data preprocessing procedure was used for

each scene: 1) Conversion of each hyperspectral scene to a structured dataset and removal of instances with no associated LULC class, 2) random sampling to maintain similar class proportions on a sample of 10% of each dataset and 3) removal of instances belonging to a class with frequency lower than 20 or higher than 1000. This is done to maintain the datasets to a practicable size due to computational constraints, while conserving the relative LULC class frequencies and data distribution. Table 1 provides a description of the final datasets used for this work.

Dataset	Features	Instances	Min. Instances	Maj. Instances	IR	Classes
Botswana	145	288	20	41	2.05	11
Pavia Centre	102	3898	278	879	3.16	7
Kennedy Space Center	176	497	23	80	3.48	11
Salinas A	224	535	37	166	4.49	6
Pavia University	103	2392	89	679	7.63	8
Salinas	224	4236	91	719	7.9	15
Indian Pines	220	984	21	236	11.24	11

Table 1: Description of the datasets used for this experiment.

Indian Pines

The Indian Pines scene [Baumgardner et al., 2015] was collected on June 12, 1992 and consists of AVIRIS hyperspectral image data covering the Indian Pine Test Site 3, located in North-western Indiana, USA. As a subset of a larger scene, it is composed of 145×145 pixels (see Figure 4a) and 220 spectral reflectance bands in the wavelength range 400 to 2500 nanometers. Approximately two thirds of this scene is composed by agriculture and the other third is composed of forest and other natural perennial vegetation. Additionally, the scene also contains low density buildup areas.

Pavia Centre and University

Both Pavia Centre and University scenes were acquired by the ROSIS sensor. These scenes are located in Pavia, northern Italy. Pavia Centre is a 1096×1096 pixels image with 102 spectral bands, whereas Pavia University is a 610×610 pixels image with 103 spectral bands. Both images have a geometrical resolution of 1.3 meters and their ground truths are composed of 9 classes each (see Figures 4b and 4c).

Salinas and Salinas-A

These scenes were collected by the AVIRIS sensor over Salinas Valley, California and contain at-sensor radiance data. Salinas is a 512×217 pixels image with 224 bands and 16 classes regarding vegetables, bare soil and vineyard fields (see Figure 4d). Salinas-A, a subscene of Salinas, comprises 86×83 pixels and contains 6 classes regarding vegetables (see Figure 4e). These scenes have a geometrical resolution of 3.7 meters.

Botswana

The Botswana scene was acquired by the Hyperion sensor on the NASA EO-1 satellite over the Okavango Delta, Botswana in 2001-2004 at a 30m spatial resolution. Data preprocessing was performed by the

UT Center for Space Research. The scene comprises a 1476×256 pixels with 145 bands and 14 classes regarding land cover types in seasonal and occasional swamps, as well as drier woodlands (see figure 4f).

Kennedy Space Center

The Kennedy Space Center scene was acquired by the AVIRIS sensor over the Kennedy Space Center, Florida, on March 23, 1996. Out of the original 224 bands, water absorption and low SNR bands were removed and a total of 176 bands at a spatial resolution of 18m are used. The scene is a 512×614 pixel image and contains a total of 16 classes (see figure 4g).

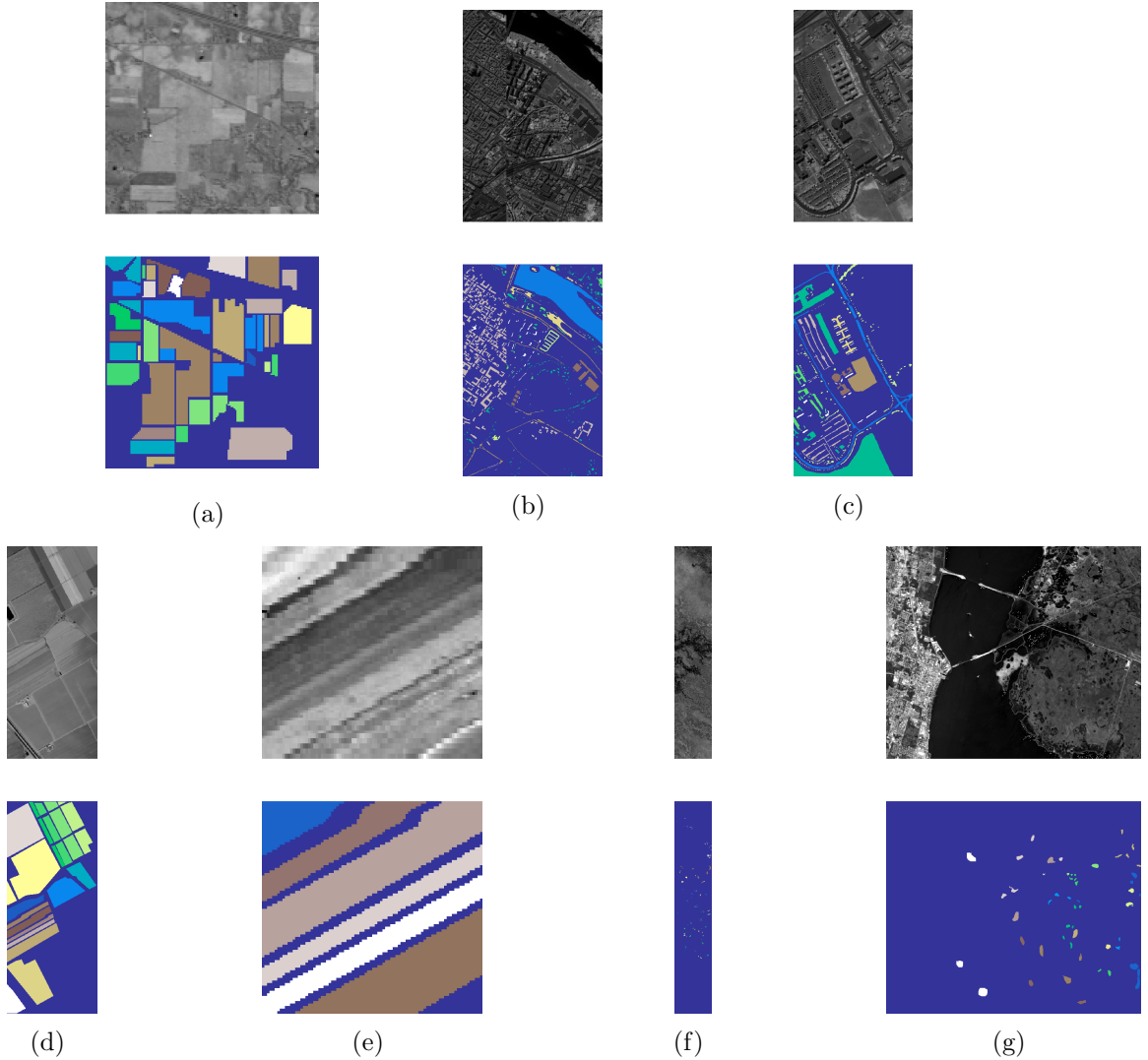


Figure 4: Gray scale visualization of a band (top row) and ground truth (bottom row) of each scene used in this study. (a) Indian Pines, (b) Pavia Centre, (c) Pavia University, (d) Salinas, (e) Salinas A, (f) Botswana, (g) Kennedy Space Center

3.2 Evaluation Metrics

Most of the satellite-based LULC classification studies (nearly 80%) employ *Overall Accuracy* (OA) and the *Kappa Coefficient* [Gavade and Rajpurohit, 2019]. Although, some authors argue that both evaluation metrics, even when used simultaneously, are insufficient to fully address the area estimation

and uncertainty information needs [Olofsson et al., 2013, Pontius and Millones, 2011]. Other metrics like User’s Accuracy (or *Precision*) and Producer’s Accuracy (or *Recall*) are also common metrics to evaluate per-class prediction power. These metrics consist of ratios employing the True and False Positives (TP and FP , number of correctly/incorrectly classified observations of a given class) and True and False Negatives (TN and FN , number of correctly/incorrectly classified observations as not belonging to a given class). These metrics are formulated as $Precision = \frac{TP}{TP+FP}$ and $Recall = \frac{TP}{TP+FN}$. While metrics like OA and *Kappa Coefficient* are significantly affected by imbalanced class distributions, *F-Score* is less sensitive to data imbalance and a more appropriate choice for performance evaluation [Jeni et al., 2013].

The datasets used present significantly high IRs (see Table 1). Therefore, it is especially important to attribute equal importance to the predictive power of all classes, which does not happen with OA and *Kappa Coefficient*. In this study, we employ 3 evaluation metrics: 1) *G-mean*, since it is not affected by skewed class distributions, 2) *F-Score*, as it proved to be a more appropriate metric for this problem when compared to other commonly used metrics [Jeni et al., 2013], and 3) *Overall Accuracy*, for discussion purposes.

- The *G-mean* consists of the geometric mean of $Specificity = \frac{TN}{TN+FP}$ and *Sensitivity* (also known as *Recall*). For multiclass problems, The *G-mean* is expressed as:

$$G-mean = \sqrt{Sensitivity \times Specificity}$$

- *F-score* is the harmonic mean of *Precision* and *Recall*. The *F-score* for the multi-class case can be calculated using their average per class values [He and Garcia, 2009]:

$$F-score = 2 \frac{\overline{Precision} \times \overline{Recall}}{\overline{Precision} + \overline{Recall}}$$

- *Overall Accuracy* is the number of correctly classified observations divided by the total amount of observations. Having c as the label of the various classes, *Accuracy* is given by the following formula:

$$Accuracy = \frac{\sum_c TP_c}{\sum_c (TP_c + FP_c)}$$

3.3 Machine Learning Algorithms

To assess the quality of the K-SMOTE algorithm, three other oversampling algorithms were used for benchmarking. ROS and SMOTE were chosen for their simplicity and popularity. B-SMOTE chosen as a popular variation of the SMOTE algorithm. We also include the classification results of no oversampling (NONE) as a baseline.

To assess the performance of each oversampler, we use the classifiers Logistic Regression (LR) [Nelder and Wedderburn, 1972], K-Nearest Neighbors (KNN) [Cover and Hart, 1967], Decision Tree (DT) [Salzberg, 1994], Gradient Boosting Classifier (GBC) [Friedman, 2001] and Random Forest (RF) [Liaw et al., 2002]. This choice was based on the classifiers’ popularity for LULC classification, learning type and training time [Maxwell et al., 2018, Gavade and Rajpurohit, 2019].

3.4 Experimental Procedure

The procedure for the experiment reported in this study is similar to the one proposed in [Douzas et al., 2019]. We start by defining a hyperparameter search grid, where a list of possible values for each relevant hyperparameter in both classifiers and oversamplers is stored. Based on this search grid, all possible combinations of oversamplers, classifiers and hyperparameters are formed. Finally, for each dataset, hyperparameter combination and initialization we use the evaluation strategy shown in Figure 5: k -fold cross-validation strategy where $k = 5$ to train each model defined and save the averaged scores of each split.

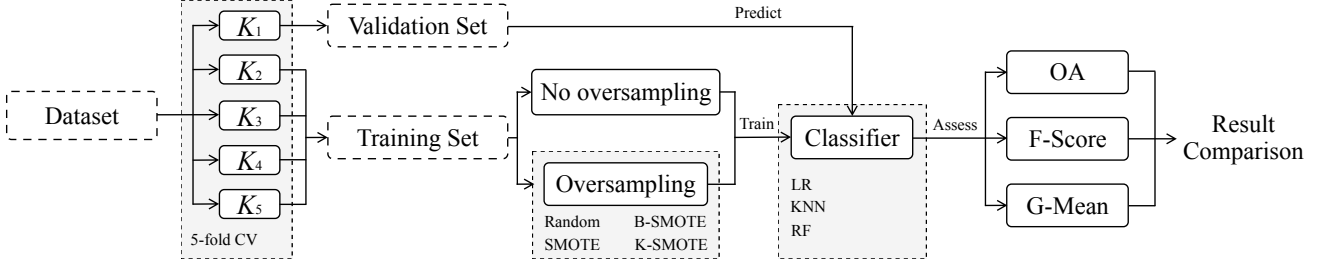


Figure 5: Experimental procedure. The performance metrics are averaged over the 5 folds across each of the 3 different initializations of this procedure for a given combination of oversampler, classifier and hyperparameter definition.

Each combination of oversampler, classifier and parameters definition is fit 5 times (once for each fold) per dataset. Each time, an oversampler will use the training set (80% of the dataset) to generate a set with artificial data, which is appended to the original training set in order to generate a training dataset with the exact same number of observations for each class. The newly formed training dataset is used to train the classifier and the test set (20% of the dataset, the remaining fold) is used to evaluate the performance of the classifier. The evaluation scores are then averaged over the 5 times the process is repeated. The range of hyperparameters used are shown in table 2.

Classifier	Hyperparameters	Values
LR	maximum iterations	10000
KNN	# neighbors	3, 5, 8
RF	maximum depth	None, 3, 6
	# estimators	50, 100, 200
Oversampler		
K-SMOTE	# neighbors	3, 5
	# clusters (as % of number of observations)	1*, 0.1, 0.3, 0.5, 0.7, 0.9
	Exponent of mean distance	auto, 2, 5, 7
	IR threshold	auto, 0.5, 0.75, 1.0
SMOTE	# neighbors	3, 5
BORDERLINE SMOTE	# neighbors	3, 5

Table 2: Hyper-parameters grid. * One cluster is generated in total, a corner case that mimics the behavior of SMOTE

3.5 Software Implementation

The experiment was implemented using the Python programming language, using the Scikit-Learn [Pedregosa et al., 2011], Imbalanced-Learn [Lemaître et al., 2017], Geometric-SMOTE, Cluster-Over-Sampling and Research-Learn libraries. All functions, algorithms, experiments and results are provided at the GitHub repository of the project.

4 Results

When evaluating the performance of an algorithm across multiple datasets, it is generally recommended to avoid direct score comparisons and use classification rankings instead [Demšar, 2006]. This is done by assigning a ranking to oversamplers based on the different combinations of classifier, metric and dataset used. These rankings are also used for the statistical analyses presented in Section 4.1.

The rank values are assigned based on the mean validation scores resulting from the experiment described in Section 3. The averaged ranking results are computed over 3 different initialization seeds and a 5 fold cross validation scheme, returning a float value within the interval $[1, 5]$. The mean rankings are presented in Table 3 and Figure 6.

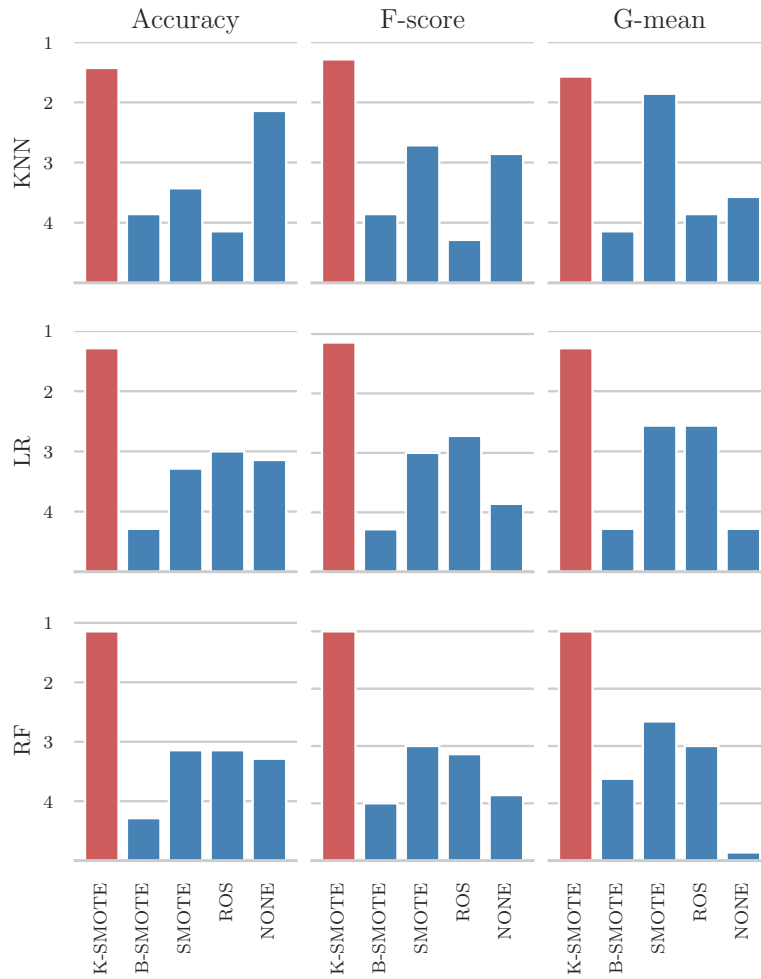


Figure 6: Mean ranking of oversamplers across datasets.

The mean ranking results show that K-SMOTE consistently presents the best results for every classifier and performance metric used. This is visually depicted in Figure 6. The quantitative results of this analysis is presented in Table 3. In addition to its better performance, in most cases K-SMOTE’s mean ranking has a lower standard deviation than any of the remaining methods, and particularly when opposed to SMOTE (the best performing benchmark method).

Classifier	Metric	NONE	ROS	SMOTE	B-SMOTE	K-SMOTE
LR	Accuracy	3.14 ± 0.47	3.00 ± 0.49	3.29 ± 0.42	4.29 ± 0.43	1.29 ± 0.18
LR	F-score	3.86 ± 0.26	2.71 ± 0.47	3.00 ± 0.44	4.29 ± 0.42	1.14 ± 0.14
LR	G-mean	4.29 ± 0.18	2.57 ± 0.48	2.57 ± 0.3	4.29 ± 0.42	1.29 ± 0.18
KNN	Accuracy	2.14 ± 0.55	4.14 ± 0.34	3.43 ± 0.43	3.86 ± 0.34	1.43 ± 0.2
KNN	F-score	2.86 ± 0.4	4.29 ± 0.29	2.71 ± 0.61	3.86 ± 0.34	1.29 ± 0.18
KNN	G-mean	3.57 ± 0.48	3.86 ± 0.4	1.86 ± 0.46	4.14 ± 0.26	1.57 ± 0.2
RF	Accuracy	3.29 ± 0.52	3.14 ± 0.4	3.14 ± 0.51	4.29 ± 0.29	1.14 ± 0.14
RF	F-score	3.86 ± 0.46	3.14 ± 0.51	3.00 ± 0.44	4.00 ± 0.22	1.00 ± 0.0
RF	G-mean	4.86 ± 0.14	3.00 ± 0.31	2.57 ± 0.3	3.57 ± 0.37	1.00 ± 0.0

Table 3: Results for mean ranking of oversamplers across datasets.

The mean percentage difference among K-SMOTE and SMOTE is presented in Figure 7. It is calculated as the score difference among the test (K-SMOTE) and control (SMOTE) oversampler, divided by the control oversampler’s score. K-SMOTE’s average performance improves classification performance of up to 1.9% and outperforms all other methods, with the exception of two situations when using the G-mean evaluation metric.

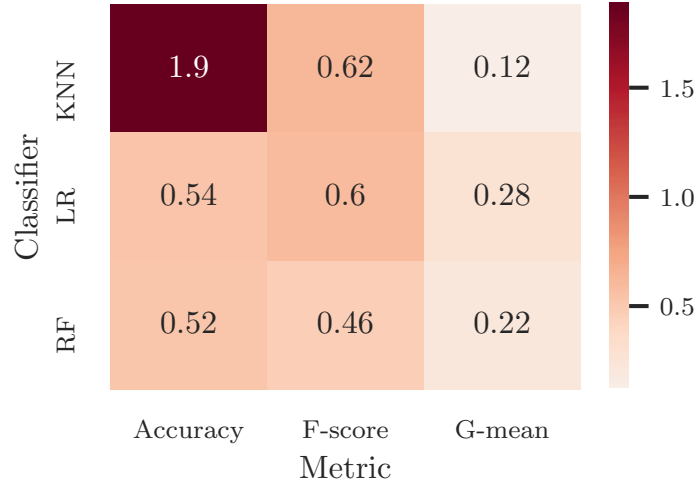


Figure 7: Mean score improvement (percentage difference) of the proposed method versus SMOTE across datasets.

The mean cross-validation scores are shown in Table 4. Considering the disparity of performance scores across datasets, the results presented in this table may not be as informative as the scores for each dataset, presented in Table 5. K-SMOTE’s performance is the highest in most classifier/metric combinations and datasets, showing more inconsistency on the Indian Pines and Kennedy Space Center datasets.

Classifier	Metric	NONE	ROS	SMOTE	B-SMOTE	K-SMOTE
LR	Accuracy	0.906 \pm 0.039	0.904 \pm 0.04	0.904 \pm 0.04	0.901 \pm 0.04	0.909 \pm 0.038
LR	F-score	0.891 \pm 0.041	0.893 \pm 0.042	0.893 \pm 0.042	0.890 \pm 0.042	0.898 \pm 0.04
LR	G-mean	0.936 \pm 0.025	0.940 \pm 0.025	0.940 \pm 0.025	0.937 \pm 0.025	0.943 \pm 0.024
KNN	Accuracy	0.879 \pm 0.043	0.865 \pm 0.048	0.867 \pm 0.05	0.862 \pm 0.054	0.881 \pm 0.045
KNN	F-score	0.859 \pm 0.05	0.853 \pm 0.049	0.861 \pm 0.047	0.851 \pm 0.053	0.866 \pm 0.048
KNN	G-mean	0.919 \pm 0.03	0.920 \pm 0.029	0.926 \pm 0.027	0.918 \pm 0.03	0.927 \pm 0.027
RF	Accuracy	0.898 \pm 0.032	0.901 \pm 0.031	0.900 \pm 0.031	0.898 \pm 0.032	0.905 \pm 0.031
RF	F-score	0.879 \pm 0.041	0.885 \pm 0.037	0.887 \pm 0.036	0.883 \pm 0.037	0.891 \pm 0.036
RF	G-mean	0.930 \pm 0.024	0.935 \pm 0.022	0.937 \pm 0.021	0.935 \pm 0.021	0.939 \pm 0.02

Table 4: Mean cross-validation scores of oversamplers.

The performance of both oversamplers and classifiers is generally dependent on the dataset being used. Although both absolute and relative scores between the different oversamplers are dependent on the choice of metric and classifier, K-SMOTE’s relative performance is consistent across datasets and generally outperforms the remaining oversampling methods.

Dataset	Classifier	Metric	NONE	ROS	SMOTE	B-SMOTE	K-SMOTE
Botswana	LR	Accuracy	0.920	0.917	0.920	0.921	0.927
Botswana	LR	F-score	0.913	0.909	0.913	0.914	0.921
Botswana	LR	G-mean	0.952	0.950	0.952	0.952	0.956
Botswana	KNN	Accuracy	0.875	0.862	0.881	0.869	0.889
Botswana	KNN	F-score	0.859	0.850	0.873	0.859	0.879
Botswana	KNN	G-mean	0.924	0.918	0.930	0.923	0.933
Botswana	RF	Accuracy	0.873	0.884	0.877	0.877	0.890
Botswana	RF	F-score	0.865	0.877	0.872	0.870	0.883
Botswana	RF	G-mean	0.925	0.933	0.929	0.928	0.936
IP	LR	Accuracy	0.687	0.681	0.680	0.678	0.692
IP	LR	F-score	0.662	0.663	0.659	0.659	0.674
IP	LR	G-mean	0.798	0.801	0.798	0.797	0.807
IP	KNN	Accuracy	0.644	0.602	0.589	0.557	0.632
IP	KNN	F-score	0.593	0.591	0.603	0.560	0.604
IP	KNN	G-mean	0.757	0.764	0.782	0.751	0.781
IP	RF	Accuracy	0.742	0.747	0.747	0.740	0.752
IP	RF	F-score	0.673	0.704	0.713	0.701	0.714
IP	RF	G-mean	0.806	0.826	0.835	0.831	0.838
KSC	LR	Accuracy	0.904	0.905	0.905	0.899	0.909
KSC	LR	F-score	0.868	0.873	0.874	0.862	0.877
KSC	LR	G-mean	0.928	0.932	0.932	0.924	0.934
KSC	KNN	Accuracy	0.855	0.859	0.862	0.857	0.865
KSC	KNN	F-score	0.808	0.819	0.827	0.810	0.826
KSC	KNN	G-mean	0.893	0.901	0.906	0.895	0.905
KSC	RF	Accuracy	0.860	0.859	0.863	0.859	0.868
KSC	RF	F-score	0.817	0.815	0.826	0.816	0.832
KSC	RF	G-mean	0.898	0.899	0.905	0.898	0.907
PC	LR	Accuracy	0.954	0.955	0.955	0.950	0.956
PC	LR	F-score	0.944	0.947	0.947	0.941	0.948
PC	LR	G-mean	0.968	0.972	0.972	0.966	0.973
PC	KNN	Accuracy	0.926	0.920	0.923	0.924	0.926
PC	KNN	F-score	0.915	0.909	0.913	0.913	0.915

Dataset	Classifier	Metric	NONE	ROS	SMOTE	B-SMOTE	K-SMOTE
PC	KNN	G-mean	0.953	0.955	0.957	0.954	0.957
PC	RF	Accuracy	0.938	0.941	0.940	0.938	0.942
PC	RF	F-score	0.928	0.932	0.931	0.928	0.933
PC	RF	G-mean	0.959	0.964	0.965	0.961	0.965
PU	LR	Accuracy	0.905	0.897	0.897	0.891	0.904
PU	LR	F-score	0.890	0.894	0.894	0.888	0.898
PU	LR	G-mean	0.932	0.947	0.947	0.942	0.949
PU	KNN	Accuracy	0.895	0.867	0.865	0.873	0.895
PU	KNN	F-score	0.891	0.868	0.868	0.874	0.891
PU	KNN	G-mean	0.940	0.935	0.936	0.936	0.941
PU	RF	Accuracy	0.912	0.908	0.907	0.908	0.911
PU	RF	F-score	0.909	0.906	0.906	0.908	0.909
PU	RF	G-mean	0.946	0.946	0.948	0.948	0.949
Salinas	LR	Accuracy	0.990	0.990	0.989	0.990	0.990
Salinas	LR	F-score	0.985	0.986	0.985	0.985	0.986
Salinas	LR	G-mean	0.992	0.993	0.992	0.992	0.993
Salinas	KNN	Accuracy	0.970	0.967	0.969	0.967	0.970
Salinas	KNN	F-score	0.959	0.957	0.960	0.957	0.960
Salinas	KNN	G-mean	0.977	0.978	0.981	0.976	0.981
Salinas	RF	Accuracy	0.984	0.983	0.983	0.983	0.985
Salinas	RF	F-score	0.979	0.979	0.977	0.978	0.980
Salinas	RF	G-mean	0.989	0.989	0.989	0.989	0.990
SA	LR	Accuracy	0.979	0.981	0.983	0.979	0.984
SA	LR	F-score	0.976	0.979	0.982	0.977	0.982
SA	LR	G-mean	0.985	0.988	0.990	0.987	0.989
SA	KNN	Accuracy	0.987	0.979	0.982	0.983	0.988
SA	KNN	F-score	0.986	0.979	0.981	0.982	0.987
SA	KNN	G-mean	0.992	0.989	0.990	0.991	0.993
SA	RF	Accuracy	0.980	0.983	0.984	0.979	0.985
SA	RF	F-score	0.979	0.982	0.983	0.978	0.984
SA	RF	G-mean	0.987	0.988	0.989	0.986	0.990

Table 5: Mean cross-validation scores of oversamplers for each dataset. Legend: IP Indian Pines, KSC Kennedy Space Center, PC Pavia Center, PU Pavia University, SA Salinas A.

4.1 Statistical Analysis

The experiment’s multi-dataset context was used to perform both a Friedman test [Friedman, 1937]. Table 6 shows the results obtained in the Friedman test performed, where the null hypothesis is rejected in all cases. Consequently, the Holm-Bonferroni comparison method (Holm’s method) [Holm, 1979] is used for post-hoc analysis.

Classifier	Metric	p-value	Significance
LR	Accuracy	9.8e-03	True
LR	F-score	2.3e-03	True
LR	G-mean	9.8e-04	True
KNN	Accuracy	4.3e-03	True

Classifier	Metric	p-value	Significance
KNN	F-score	4.3e-03	True
KNN	G-mean	3.0e-03	True
RF	Accuracy	5.5e-03	True
RF	F-score	2.9e-03	True
RF	G-mean	1.8e-04	True

Table 6: Results for Friedman test. Statistical significance is tested at a level of $\alpha = 0.05$. The null hypothesis is that there is no difference in the classification outcome across oversamplers.

The results of the Holm’s method are shown in Table 7. Even though K-SMOTE outperforms the remaining oversamplers, the datasets’ inherent high prediction scores make the rejection of this null hypothesis particularly difficult.

Classifier	Metric	NONE	ROS	SMOTE	B-SMOTE
LR	Accuracy	7.0e-02	7.0e-02	7.0e-02	2.6e-02
LR	F-score	1.5e-02	7.7e-02	7.7e-02	2.2e-02
LR	G-mean	5.1e-02	8.4e-02	8.4e-02	2.4e-02
KNN	Accuracy	5.7e-01	6.0e-02	2.1e-01	2.1e-01
KNN	F-score	1.5e-01	5.2e-02	1.5e-01	9.7e-02
KNN	G-mean	1.4e-01	8.8e-02	2.3e-01	1.4e-01
RF	Accuracy	4.4e-02	3.5e-02	4.4e-02	2.4e-02
RF	F-score	6.9e-02	6.9e-02	6.9e-02	3.6e-02
RF	G-mean	1.0e-01	1.0e-01	1.0e-01	3.8e-02

Table 7: Adjusted p-values using the Holm’s method. Bold values are statistically significant at a level of $\alpha = 0.05$. The null hypothesis is that the test method does not perform better than the control method.

5 Conclusion

This research paper was motivated by the difficulty posed in classifying rare classes in Land Use/Land Cover tasks. A number of existing methods to address this problem (known as imbalanced learning) was identified and their caveats were exposed. Typically, these methods are not only difficult to implement, they are also context dependent. We focused on oversampling methods due to their widespread usage, easy implementation and flexibility. Specifically, this paper demonstrated the efficacy of a recent oversampler, K-Means SMOTE, applied in a multi-class context for Land Cover Classification tasks. This was done with sampled data from seven well known and naturally imbalanced datasets: Indian Pines, Pavia Centre, Pavia University, Salinas, Salinas A, Botswana and Kennedy Space Center. The experiment comprised a hyper-parameter search in order to tune each algorithm to its specific use case. For each combination of dataset, oversampler and classifier, the results of every classification task was averaged across a 5 fold stratification strategy with 3 different initialization seeds, resulting in a mean validation score of 15 classification tasks. The optimal mean validation score of each combination was then used to perform the analyses presented in this report.

In most cases, classification tasks using K-SMOTE led to better results than using the original, unmodified, imbalanced data. More importantly, we found that K-Means SMOTE is always better or equal than the second best oversampling method. K-SMOTE’s performance was independent from both

the classifier and performance metric under analysis. In general, K-SMOTE shows a higher performance among the non tree-based classifiers employed, when compared to the remaining oversamplers. Although these findings are case dependent, they are consistent with the results presented in [Douzas et al., 2018]. The proposed method also had the most consistent results across datasets, since it had the lowest standard deviations across datasets in most cases for both analyses, either based on ranking or mean cross-validation scores.

The proposed algorithm is an extension of the original SMOTE algorithm. In fact, the SMOTE algorithm represents a corner case of K-SMOTE i.e. when the number of clusters equals to 1. Its data selection phase differs from the one used in SMOTE and Borderline SMOTE, providing artificially augmented datasets with less noisy data than the commonly used methods. This allows the training of classifiers with better defined decision boundaries, especially in the most important regions of the data space (the ones populated by a higher percentage of minority class instances).

As stated previously, the usage of this oversampler is technically simple. It can be applied to any classification problem relying on an imbalanced dataset, alongside any classifier. K-SMOTE is available as an open source implementation for the Python programming language (see Subsection 3.5). Consequently, it can be a useful tool for both remote sensing researchers and practitioners.

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