TOWARDS A UNIFIED STRATEGY FOR THE PREPROCESSING STEP IN DATA MINING

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Abstract: Data-related issues represent the main obstacle in obtaining a high quality data mining process. Existing strategies for preprocessing the available data usually focus on a single aspect, such as incompleteness, or dimensionality, or filtering out “harmful” attributes, etc. In this paper we propose a unified methodology for data preprocessing, which considers several aspects at the same time. The novelty of the approach consists in enhancing the data imputation step with information from the feature selection step, and performing both operations jointly, as two phases in the same activity. The methodology performs data imputation only on the attributes which are optimal for the class (from the feature selection point of view). Imputation is performed using machine learning methods. When imputing values for a given attribute, the optimal subset (of features) for that attribute is considered. The methodology is not restricted to the use of a particular technique, but can be applied using any existing data imputation and feature selection methods.

1 INTRODUCTION

Machine learning has become one of the main sources of learning techniques in data mining. During the last years a great number of state-of-the-art methods have emerged, while older methods have been improved. Thus, data mining researchers possess a rich collection of robust techniques which tackle the learning phase in the knowledge extraction process. Despite the strength of existing learning schemes, on some problems they fail to achieve a satisfying performance. The reason behind such a behaviour can be found in the quality of the training data. In this paper we propose a generic preprocessing methodology, which combines feature selection with data imputation, to improve the quality of the training data. The end goal is to obtain an increased learning accuracy.

The rest of the paper is organized as follows: some of the most common data-related issues are discussed in section 2. A data imputation and a wrapper feature selection approach are discussed in the first part of section 3, followed by the proposed preprocessing methodology. Section 4 presents the evaluations performed using two different implementations of the methodology, followed by the conclusion section.

2 WHY WE NEED PREPROCESSING

One of the main issues involved in the learning step is algorithm selection and engineering. This usually implies the choice of the “right” algorithm for the given problem, together with parameter tuning. The “right” algorithm, in theory, is the one which matches the problem bias, e.g. the separation hyper-planes it can learn match the data separation into classes. Algorithm selection and engineering are empirical-intensive activities, since many combinations need to be considered. Practical experience has shown us that the success of the mining process does not always depend on the choice and engineering of the learning technique alone. An important factor can be found in the quality of the training data which is presented to the algorithm. There are several dimensions to this problem that need to be addressed. Some of the most common are:

- Data volume (number of instances)
- Incompleteness
- Irrelevant/redundant information
- Dimensionality (number of features/attributes)

Even tough data mining has emerged from the need of analyzing large amounts of data, there are some
cases where the data volume is insufficient for modelling the true hidden hypothesis. Incomplete data has proved to be the rule, rather than the exception, of a data mining process. Incomplete datasets could result from many causes, such as: the values are out of the range of the measuring device, may signal a don’t care value, is the result of a random event, a.s.o. Incomplete datasets usually bias the learning step, even for techniques which are able to learn from incomplete data. Although there exists a taxonomy for data incompleteness (Little & Rubin, 1987), with systematic and efficient approaches for dealing with each type of incompleteness, this issue is neither trivial nor easy to handle. The main difficulty is that the type of incompleteness is, in most cases, unknown, rendering impossible the choice of the appropriate technique. This makes its treatment more challenging, since informative missing values should be handled differently than random incompleteness.

Supposing that for a given problem the “right” classifier has been identified, the monotonic assumption – that more attributes improve the learning phase – is not generally valid. This is due to the irrelevant/redundant information in the data, which harms the learning process. This suggests that, theoretically, the data should be cleaned and only strongly relevant attributes should be preserved. In practice, however, the strongly relevant attributes (as defined by theory) are hard to establish. Furthermore, in some cases, data which is theoretically redundant may enhance the learning process (e.g. adding the product of the two variables in an XOR problem improves the accuracy of a neural network classifier, (Georgieva, 2008)). High dimensionality (i.e. a very large number of attributes) has also other implications than the inclusion of irrelevant/redundant information. It usually results in slow learning and a complex and difficult to interpret output model.

This list of data-related problems we have enumerated is far from exhaustive, but we consider that finding viable, widely-applicable solutions to these issues is the first step to improving the data mining process. Starting from these facts, we have explored a data imputation and a wrapper feature selection strategy, and performed evaluations on benchmark data. The aim was to obtain a better accuracy after the learning process. Starting from the results obtained, we propose a combined preprocessing methodology for the data mining process. The final goal is to provide general approaches to handle different data-related problems in a unified manner.

3 PRE-PROCESSING APPROACHES

We have come across the need to employing more elaborate preprocessing strategies while mining a prostate cancer dataset using complex classification approaches. On medical benchmark data, the complex techniques obtained the best/at least similar results when compared to prominent learning techniques existing in literature (Vidrighin et al., 2007), in terms of classification accuracy or total cost. On a real prostate cancer dataset, the performance of the techniques remained at the highest level when compared to the performance of the same “competition” approaches. However, due to the sensitivity of this particular domain, the accuracy obtained is considered to be insufficiently high. Since the other techniques employed have failed to obtain a better accuracy, we have started to search for a preprocessing solution in order to boost the classification accuracy. A first approach started from the observation that the prostate cancer dataset contains a lot of missing values. Thus, we focused on methods for imputing the missing values. Also, even though in this case the data dimensionality is not an issue, we have employed a wrapper methodology to identify the optimal predictive subset for the problem. Given the domain particularities, this leads to a better model interpretability and also eliminates harmful irrelevant/redundant information, hence reducing the medical costs.

3.1 Data Imputation

Our method for data imputation is inspired from machine learning techniques. We found motivation in the increased accuracy rates obtained by an ensemble of artificial neural networks on various benchmark datasets. The technique involves training the ensemble on the available complete data for a given attribute. For each predictor attribute $A_i$, $i=1,n$, we split the training set $T$ into the training subset of complete data and the training subset of incomplete data: $T = CT_i + IT_i$. Given the attribute $A_i$ and an instance $t \in T$, $t \in CT_i$ if the value of $A_i$ is present and $t \in IT_i$ otherwise. The ensemble of neural networks is trained on $CT_i$, considering $A_i$ as the class, and the obtained model is employed to impute the values for $A_i$ in $IT_i$, resulting in $IT_i'$. The resulting training set $IT' = CT_i + IT_i'$ should have a higher quality, and thus improve the learning phase.
We have performed evaluations considering several different ratios of incompleteness. We have measured the classification accuracy of a decision tree learner. For each attribute we have compared results using three variants of the training set:

- **complete** training set: \( T_1 = \emptyset \), \( T = C T_i \)
- **incomplete**: \( p\% \) of the attribute values are considered missing, with \( p \) between 5 and 30, with a 5% increment
- **imputed**: \( p\% \) of the attribute values have been imputed by our method (\( p \) the same as before)

The evaluations were performed in a 10-fold cross validation loop (for the training/testing sets), 10 times for each incompleteness percentage, using different splits every time (for the training set).

The results suggest that the method could be used to improve the quality of the training set. The observations can be divided into three categories, considering each attribute’s correlation with the class (measured using the gain ratio): highly correlated attributes, mildly correlated attributes and weakly correlated attributes. The most stable improvements have been found for attributes which possessed a strong correlation with the class and for datasets with increased modelling power (measured in terms of the classification accuracy level) – (Vidrighin et al., 2008a). This result, together with the fact that the optimal subset contains the strongly relevant features, indicates the possibility of augmenting the mining process with a combined approach for preprocessing.

### 3.2 Wrapper Feature Selection

We have defined feature selection as the process of selecting the optimal subset of features for a dataset. Optimality may refer to: improving the classification accuracy, reducing the computation effort, improving model interpretability or avoiding costly features. Our target is to achieve the highest possible accuracy.

We have employed a 3-step methodology, based on an existing classification for feature selection methods (Kohavi, 1997). Therefore, we view the wrapper as a 3-tuple of the form \(<\text{generation procedure}, \text{evaluation function}, \text{validation function}>\). The **generation procedure** is a search procedure which selects a subset of features \((F_i)\) from the original feature set \((F)\), \( F_i \subseteq F \). There are many search methods available, from greedy hill climbing search, to genetic or random search methods. Each has its advantages and disadvantages. Previous work (Vidrighin et al., 2008b) has shown that greedy stepwise backward search and best first search constantly yield good results. We have employed these two search methods as generation procedures in our previous evaluations (Vidrighin et al., 2008c). The **evaluation function** measures the “quality” of a subset obtained from a given generation procedure. As the optimal features subset depends on the evaluation function, the process of selecting the appropriate evaluation function is dependent on the particular initial dataset. In the case of wrappers, the evaluation is performed by measuring the accuracy of a certain inducer on the projection of the initial dataset on the selected attributes. In our previous work we have employed three different learning schemes (inducers), representing three prominent classes of algorithms: decision trees (C4.5 – revision 8 – J4.8, as implemented by Weka (Witten, 2005)); Naïve Bayes (Cheeseman & Stutz, 1995) and ensemble methods (AdaBoost.M1 (Freund & Schapire, 1997)). For J4.8, we performed experiments both with and without pruning.

The **validation function** tests the validity of the selected subset through comparisons obtained from other feature selection and generation procedure pairs. The objective of the validation procedure is to identify the best performance that could be obtained in the first two steps of the method for a given dataset, i.e. to identify the selection method which is most suitable for the given dataset and classification method. In this phase we performed validations with all the three inducers employed in the evaluation phase. Again, J4.8 was considered both with pruning and without pruning. Validation is important in selecting the inducer for learning after feature selection has been performed.

Evaluations on 11 benchmark datasets (Vidrighin et al., 2008c) have shown that feature selection almost always improves the accuracy of any inducer. The most were found for combinations which included the Naïve Bayes classifier, but J4.8 combination also obtained good improvements. The best wrapper combinations obtained up to 13% relative improvements in accuracy (when compared to the initial accuracy of the inducer), while the second best obtained up to 7% relative improvements. Also, combinations using the initially best inducer for evaluation and validation have obtained relative improvements in accuracy up to 7%. Therefore, employing the reduced optimal feature set rather than the initial set almost always boosts the learning accuracy. For datasets having over 99% accuracy no improvement has been found.

### 3.3 A Combined Strategy

Starting from the results obtained by both preprocessing approaches explored, we propose a
unified strategy for preprocessing the data, which consists in two phases:
- a feature selection phase
- a data imputation phase

The novelty of the approach consists in enhancing the data imputation step with information from the feature selection step, and performing both operations jointly, as two phases in the same activity.

In the feature selection phase, we extract the class optimal feature subset (further referred as COS), i.e. the subset of features which best predicts the class. We have employed classification accuracy as optimality criterion. Also, for each attribute, \( A_i \), in the optimal subset, we perform feature selection on the entire training set, to obtain its optimal subset, \( AOS_i \). The optimality criterion should be accuracy here as well, but the methodology does not impose the use of a given method (e.g. filter ranking methods could be employed instead of wrappers).

In the data imputation phase we impute values for the incomplete attributes in the optimal subset of the class. For each attribute we will consider only the features in its optimal subset, \( AOS_i \), for imputation. By performing feature selection for each attribute in particular, we wish to eliminate any noise and harmful information.

After imputation, the resulting training set represents the improved training set, which is further used in the learning step.

This is a simple, generic strategy, which can be applied using any feature selection and data imputation techniques. It is neither restricted to wrapper feature selection in the feature selection phase, nor to a particular imputation method.

The next section presents the evaluations performed on two particular implementations of the methodology.

4 EVALUATIONS

We have performed evaluations using complete benchmark datasets from the UCI Machine Learning Data Repository (UCI), to check whether the imputation-feature selection combination can improve the accuracy of a classifier. We have experimented with two particularizations of the methodology: one that employs an ensemble of artificial neural networks for imputation, and one that utilizes kNN for imputation. Both strategies employ a wrapper method for feature selection.

4.1 Preprocessing with Ensembles of Artificial Neural Networks for Imputation

The first set of evaluations has been performed on a variant of the preprocessing methodology which utilizes the data imputation technique and the feature selection method already explored in (Vidrighin et. al, 2008a) and (Vidrighin et. al, 2008c).

The evaluation methodology is as follows:
1. select the best wrapper method for the given dataset
2. generate 10 random train/test sets pairs (from the original dataset), using a 80/20 percentage split.
3. for each training set, select the optimal feature subset (for the class attribute), using the wrapper method selected at 1.
4. for each training set, evaluate the imputation technique, as described in section 3.1

In step 3 of the evaluation methodology we have not considered the optimal attribute subset for each attribute separately, \( AOS_i \). Instead, we used COS for attribute imputation as well. This may lead to poorer results, since the attributes in COS are not necessarily predictive among each other. Therefore, when imputing values for attribute \( A_i \), only the features in its optimal subset should be employed.

Also, for each training/testing pairs, COS has been estimated on the complete training set, without taking into account any missing values. In a real world setting this should also have a different approach: either perform feature selection on \( CT_i \), for each attribute \( A_i \), or consider incompleteness in the evaluation function of the feature selection technique.

Table 1 presents the average classification accuracies obtained by training a multilayer perceptron classifier on different versions of the training set with respect to attribute GlucTest: preprocessed with the combined methodology, incomplete, and original complete set. GlucTest has been selected in COS as being strongly correlated with the class and imputation has been performed on \( A_i = \text{GlucTest} \), when \( AOS_i = \text{COS} \).

The improvement does not seem to be significant, but we believe this is due to the fact that we employed COS for imputing the values of \( A_i \), and not \( AOS_i \). Also, we consider that a more stable approach than artificial neural networks for data imputation would yield better results.
Table 1: Classification accuracies for different versions of the training set for attribute GlucTest, Pima dataset.

<table>
<thead>
<tr>
<th>% incomplete</th>
<th>5%</th>
<th>10%</th>
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<tbody>
<tr>
<td>Pre-processed (combined m.)</td>
<td>76.5</td>
<td>76.47</td>
<td>76.45</td>
<td>76.42</td>
<td>76.43</td>
<td>76.32</td>
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<tr>
<td>Incomplete</td>
<td>76.46</td>
<td>76.45</td>
<td>76.49</td>
<td>76.34</td>
<td>76.26</td>
<td>76.45</td>
</tr>
<tr>
<td>Original data</td>
<td>76.58</td>
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4.2 Preprocessing with kNN for Imputation

For evaluating this second approach we have employed the following datasets: Cleveland, Bupa, Pima and Cars. Two slightly different strategies have been considered:

- In the first one (FSAfterI), for each attribute Aᵢ in the training set (except the class) we have varied the percentage of incompleteness between 5% and 30%, with 5% increment. To impute the missing values for Aᵢ, we have employed kNN, k=3, using only the attributes in AOSᵢ (computed using a wrapper approach around kNN). The final classification was performed using J4.8, on COS (computed using a wrapper approach around J4.8).

- In the second approach (FSBeforeI), COS was computed initially, and only the attributes in COS were considered for imputation. In order to impute values for Aᵢ, AOSᵢ was extracted from the original training set.

A stratified 10-fold cross validation was performed for each experiment. The average classification accuracy and the standard deviation were computed. In the trials of the second strategy, for each attribute Aᵢ, averaging was performed only on the folds in which it was selected in COS. We have compared the accuracy of the model built using J4.8 on the entire training set against the accuracy of the model built on the pre-processed training set.

Tables 2-4 present the results obtained on the Pima dataset, for a strongly correlated attribute (with the class), a mildly and a weakly correlated attribute.

Table 2: Results of a strongly correlated attribute with the class, Pima dataset (GlucTest).

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<tbody>
<tr>
<td>FSAfterI</td>
<td>74.74</td>
<td>74.08</td>
<td>75.39</td>
<td>74.21</td>
<td>74.08</td>
<td>75.13</td>
</tr>
<tr>
<td>FSBeforeI</td>
<td>73.16</td>
<td>74.08</td>
<td>73.29</td>
<td>74.08</td>
<td>72.24</td>
<td>73.29</td>
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Table 3: Results of a mildly correlated attribute with the class, Pima dataset (Age).

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<tbody>
<tr>
<td>FSAfterI</td>
<td>73.95</td>
<td>73.95</td>
<td>74.61</td>
<td>74.21</td>
<td>74.16</td>
<td>73.55</td>
</tr>
<tr>
<td>FSBeforeI</td>
<td>72.63</td>
<td>72.63</td>
<td>71.84</td>
<td>72.11</td>
<td>72.11</td>
<td>72.11</td>
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Table 4: Results of a weakly correlated attribute with the class, Pima dataset (BloodPress).

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<tbody>
<tr>
<td>FSAfterI</td>
<td>74.21</td>
<td>74.21</td>
<td>73.82</td>
<td>74.21</td>
<td>73.16</td>
<td>73.55</td>
</tr>
<tr>
<td>FSBeforeI</td>
<td>72.63</td>
<td>72.63</td>
<td>71.84</td>
<td>72.11</td>
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We have compared the accuracy of the final classification with J4.8 on the complete set (74.74%) with its performance on the modified training sets obtained through preprocessing with the two versions of the proposed methodology (FSAfterI and FSBeforeI). Good results can be observed for strongly and mildly correlated attributes. There is no clear winner between the two approaches, and no success pattern can be identified.

The most remarkable improvements have been observed on the Cleveland data set, whose baseline accuracy lies somewhere around 50%. Figures 1-2 exemplify the results obtained for a strongly and a mildly correlated attribute with the class in the Cleveland data set. The combined preprocessing technique significantly boosts the performance of the model built on the pre-processed training set when compared to the performance of the model built on the original training set. For this dataset in particular, FSAfterI yields better results than FSBeforeI, for all attributes.

The results obtained on the other two datasets confirm the fact that preprocessing the training set using this combined methodology can help boost the classification accuracy. Also, there appears to be no absolute winner between the approaches.

Figure 1: Classification accuracy using different versions of the training set – attribute Thal.
5 CONCLUSIONS

Among the best known data preprocessing strategies are feature selection and procedures for handling incomplete data, with various existing techniques. Previous results on the data imputation step alone show that predicting strongly correlated attributes with the class can improve the learning accuracy. Wrapper feature selection has also been shown to boost the performance of an inducer.

In this paper we propose a new methodology for preprocessing the training set. Its novelty resides in the combination of the feature selection step with data imputation, in order to obtain an improved version of the training set. The main goal is to boost the classification accuracy (i.e., improve the learning step). The methodology is simple and generic, which makes it suitable for a wide range of application domains, where particular feature selection schemes/data imputation procedures may be preferred.

We have performed a number of evaluations of the combined methodology using benchmark datasets. The results indicate that performing preprocessing on the training set enhances the accuracy of the final model. The new methodology we have introduced is more successful than the individual steps it combines, producing similar or even superior results to the ones obtained with complete data.

However, just like in the case of classifiers (Moldovan et al., 2007), there is no absolute best preprocessing particularization for a given dataset. Therefore, there appears the need to assess a baseline performance using several approaches, and develop a semi-automated procedure for tuning the preprocessing method for a given problem. This is one of our current objectives. Another future development of the methodology is aimed at handling more complex patterns of incompleteness, closer to the ones encountered in real-life data sets.

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REFERENCES


