Using Undersampling with Ensemble Learning to Identify Factors Contributing to Preterm Birth

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Abstract—In this paper, we propose Ensemble Learning models to identify factors contributing to preterm birth. Our work leverages a rich dataset collected by a NIEHS P42 Center that is trying to identify the dominant factors responsible for the high rate of premature births in northern Puerto Rico. We investigate analytical models addressing two major challenges present in the dataset: 1) the significant amount of incomplete data in the dataset, and 2) class imbalance in the dataset. First, we leverage and compare two types of missing data imputation methods: 1) mean-based and 2) similarity-based, increasing the completeness of this dataset. Second, we propose a feature selection and evaluation model based on using undersampling with Ensemble Learning to address class imbalance present in the dataset. We leverage and compare multiple Ensemble Feature selection methods, including Complete Linear Aggregation (CLA), Weighted Mean Aggregation (WMA), Feature Occurrence Frequency (OFA) and Classification Accuracy Based Aggregation (CAA). To further address missing data present in each feature, we propose two novel methods: 1) Missing Data Rate and Accuracy Based Aggregation (MAA), and 2) Entropy and Accuracy Based Aggregation (EAA). Both proposed models balance the degree of data variance introduced by the missing data handling during the feature selection process, while maintaining model performance. Our results show a 42% improvement in sensitivity versus fallout over previous state-of-the-art methods.

I. INTRODUCTION

Machine Learning (ML) techniques have shown significant potential in facilitating public health research. Unlike conventional data-driven statistics and analytics in the public health domain, ML can discover hidden patterns and knowledge in large amounts of data, significantly facilitating the research process. In recent years, public health research institutes have invested a vast amount of effort on: i) increasing the availability of research data, and ii) investigating and building accurate and reliable analytical models with ML and AI algorithms [1].

Although ML algorithms provide a rich collection of candidate methods, building a reliable ML-based analytical model can be challenging. For instance, ML algorithms provide a range of performance and accuracy based on the characteristics of the input dataset. Selecting the right ML algorithm requires a thorough understanding of the characteristics of the given dataset. After building the analytical model, we can apply the same model to all datasets with similar characteristics.

We have developed an analytical model that can identify the key factors tied to preterm birth. Our work is informed through access to a rich dataset from an National Institute for Environmental Health Sciences (NIEHS) P42 Center. Unfortunately, analyzing this dataset presents two major challenges. First, given the nature of the data collected, we encounter a significant number of missing data entries in a variety of fields, resulting in incomplete records. Second, given that the dataset includes a limited number of preterm birth outcomes (anomalies), the dataset is highly imbalanced in terms of class distribution, where the number of negative samples (term births) is several times larger than the number of positive samples (preterm births). If we attempt to use standard algorithms on this data, the outcome imbalance will clearly bias our results, misleading researchers, and resulting in poor model performance.

In this paper, we explore analytical models that address these two challenges present in the dataset. First, to impute the missing data, we leverage and compare two types of missing data imputation methods, i.e., mean-based and similarity-based algorithms. Next, we propose a model built using undersampling with ensemble learning [2], incorporating Decision Trees [3] and Recursive Feature Elimination (RFE) [4] to rank and select features. We also leverage multiple commonly used ensemble feature selection methods, including Complete Linear Aggregation (CLA), Weighted Mean Aggregation (WMA), Feature Occurrence Frequency (OFA), and Classification Accuracy Based Aggregation (CAA) [5]. As a further step, to address the high variance introduced by missing data handling, we propose two novel methods: 1) Missing Data Rate and Accuracy Based Aggregation (MAA) and 2) Entropy and Accuracy Based Aggregation (EAA), which are both based on CAA. The former leverages the missing data rate of each feature, selecting features that have a lower degree of missing data. The latter leverages the information entropy change of each feature after applying missing data handling, selecting features that preserve the original entropy.

We evaluate and compare these two methods, and provide insights when leveraging them during feature selection for the dataset. We use both accuracy and Area Under the ROC Curve (AUC) metrics to evaluate the performance, considering that
the data is highly imbalanced in terms of the outcome class distribution. The experiment results show a 42% improvement in AUC over previous state-of-the-art results on the same data, which pursued a hybrid approach [6].

II. RELATED WORK

The use of Machine Learning analytics in public and environmental health research has attracted global attention. Auffray et al. evaluate the potential benefits of leveraging Big Data techniques for public health in Europe [7]. Chen et al. present the work of leveraging cognitive computing to address Big Data challenges in medical research [1]. Zheng et al. present a study of air pollution using Big Data techniques [8]. Beam et al. elaborate the connections between Big Data technologies and health studies, clarifying their roles in each research domain [9].

Other than general work related to Big Data analytics, our work here focuses on addressing challenges of missing data and outcome class imbalance. Many previous studies have considered these issues. For missing data imputation, multiple techniques are available, including mean substitution, multiple imputation and full information maximum likelihood, and have been applied to survey and clinical data [10]–[14]. There are multiple techniques for tackling class imbalance as well. Prior studies adopted using a Decision Tree combined with AUC as training criterion to select features with the best AUC performance [6]. SMOTEBoost [15] and RUSBoost [16] are two similar approaches leveraging oversampling and ensemble of weak learners that are commonly used in boosting. Balanced random forest [17] and EasyEnsemble [2], on the other hand, leverage undersampling and ensemble learning, on which our work is based.

Ensemble feature selection methods have long been investigated and applied in bioinformatics, especially given their robustness when identifying biomarkers. Brahim et al. analyze and compare multiple commonly used ensemble feature selection methods [5]. Abeel et al. propose two aggregation methods, Complete Linear Aggregation (CLA) and Weighted Mean Aggregation (WMA) [18]. The former aggregates the complete feature rank lists across each instance of ensemble and selects a set of features with the highest ranks. The latter uses the mean of the AUC as a weight for ranking. Chan et al. propose classification Accuracy Based Aggregation (CAA) [19]. In this paper, we investigate these various ways for performing feature selection from ensemble methods. Furthermore, since our data has a high percentage of missing data entries, we introduced a novel strategy for taking data missing rate into account (via the effect of data imputation) in ensemble feature selection.

The previous work relied on correlation, SVMs and PCA [18], [19] to select features. In our work we used Decision Trees because they have shown to produce superior performance on this specific dataset [6], and intrinsically performs feature selection.

III. BACKGROUND

A. The NIEHS Dataset Targeting Preterm Birth

The NIEHS has established a Center to study environmental links to the high preterm birth rates observed in northern Puerto Rico. Over an 8 year period, the Center has collected both environmental data (e.g., water samples, air samples, product use surveys) and personal health data (e.g., chemicals in blood, urine, and placenta, health records, birth outcomes) for a cohort of over 2000 expectant mothers in this region, as well as the birth outcome (i.e., preterm birth or term birth) [20]. The dataset consists of a wide range of sources, including:

- Human subjects information - medical history, reproductive health records, product use data surveys, and birth outcomes
- Biological samples - blood, urine, hair and placenta samples
- Environmental samples and measurements - soil samples, well and tap water samples, historical Environmental Protection Agency (EPA) data, soil samples, superfund site data

The Center has made their data available to the broader research community. We have gained access to this large repository of environmental health data records, enabling us to explore analytical models to identify linkages between a large number of potential contributing factors and the high premature birth rate. We consider one subset of data in this paper, i.e., the human subject data. It contains forms capturing demographics and lifestyle information, as well as medical and health records. The dataset includes many key factors that can help guide domain-specific research. Additionally, the collection procedure for the human subject data is carried out in four phases chronically, associated with four visits according to different stages of the mother’s pregnancy. The data subset analyzed in this paper represents information collected at different pregnancy stages, defined as visit 1 (V1) to visit (V3), which correspond to the trimesters, and with visit (V4) included as a postpartum collection.

IV. MISSING DATA IMPUTATION

A. Data Preprocessing Model

Missing data entries pose significant challenges to machine learning algorithms. We developed a series of tools to preprocess the data stored in the database, generating valid inputs that can be readily consumed by the analytical model. Our tools provide an end-to-end processing pipeline.

Specifically, the tool chain consists of components including a form extractor, data filter, missing data handler and merger. Figure 1 shows the workflow of our tool chain. First, the form extractor obtains the selected forms for a set of selected patient IDs which correspond to pregnant mothers. Second, data is fed to the data filter and missing data handler. The data filter takes care of two things: 1) filtering out text-based data entries, and 2) filtering out features that have less than $P\%$ of the participants completing the field, and filtering out patients...
having less than the $P\%$ of their feature values completed, in which, $P\%$ is a configurable parameter, representing a data completion rate threshold. After data filtering, each form only contains numerical and categorical data values, and at least $P\%$ valid data values on both a feature basis and participant basis. The missing data handler implements the means-based or similarity-based imputation algorithm.

### B. Missing Data Imputation Algorithm

To address missing data handling, we leverage two commonly-used missing data imputation methods: 1) mean-based imputation and 2) similarity-based imputations [10], [11]. Applying mean-based algorithm is straightforward. The missing data entries in each feature are replaced with the mean value of the corresponding feature. For the categorical data, we replace the missing data entries using the category closest to the mean value. But applying a similarity-based imputation is challenging on the given dataset. The dataset contains mixed data types (i.e., textual, numerical and categorical data), restricting the use of standard data imputation methods which are commonly used when targeting a specific type of data. Alternatively, we propose a customized algorithm that can handle both categorical and numerical data, based on K-nearest Neighbor Imputation [21]. To calculate the similarity, we use the following equation:

$$S_{ij} = \frac{x_i \cdot y_j}{N}$$

where $S_{ij}$ stands for the similarity between sample vectors $x_i$ and $y_j$, $\cdot$ denotes a dot product operation and $N$ denotes the number of features involved in the dot product operation. Note that the numerical data is normalized and the categorical data is transformed to a one-hot encoding format.

### V. Undersampling Ensemble Feature Selection

#### A. Undersampling Ensemble Learning

Undersampling forcibly balances positive and negative samples to address the class imbalance issue in a straightforward manner. In a binary class problem, an instance of undersampled training data can be constructed with equal-sized samples from both the majority and minority classes. For example, one dataset contains two classes, positive and negative outcomes. The positive samples are assigned to subset $P$ and the negative samples are assigned to subset $N$. When the class distribution is highly imbalanced, the size of $P$ ($|P|$) is several times smaller than the size of $N$ ($|N|$). The undersampling randomly selects samples from $N$ to construct a subset $\bar{N}$ so that $|P| = |\bar{N}|$, and then combines them into a single training data instance, i.e., $P \cup \bar{N}$.

Although the new instance of the training data is balanced, we lose a lot of information from the majority class. Solely relying on this training data can lead to poor model performance. To deal with this problem, a form of ensemble learning can be applied using multiple instances of training data in a set, i.e. $\{P \cup \bar{N}_0, P \cup \bar{N}_1, \ldots, P \cup \bar{N}_{n-1}\}$, in which each $\bar{N}_i$ is a subset of the samples from the majority class obtained through undersampling, where $n$ is the number of ways.

In our proposed model, which is based on undersampling with ensemble learning, we elect to incorporate Decision Trees [3] as the base classifier and feature selector, considering that it can serve as both a feature ranker and a classifier.

Figure 2 shows the proposed analytical model leveraging undersampling ensemble learning with Decision Trees. The model consists of three steps: i) feature selection, ii) data selection and iii) performance evaluation. In the feature selection step, an ensemble of training data sets are produced by applying the undersampling procedure described above. Each training set will be used to build a Decision Tree to rank the entire feature set during the training process. Then an ensemble feature selection method (aggregation algorithm) is applied to produce the final selected feature set, according to a configurable number of features. We will explore a collection of commonly used ensemble feature selection methods in the next section. In the data selection step, a subset of data is selected according to the selected features. During performance evaluation, we follow a similar process. The only difference is that the Decision Tree built using each training data instance is used for prediction instead of ranking features. The classification results from each ensemble instance will be fed to an ensemble classifier to make a final prediction. In this work, we use a simple majority voting as the ensemble classifier. Based on the prediction, we calculate performance metrics such as accuracy and AUC.

#### B. Ensemble Feature Selection

In this section, we describe four commonly-used ensemble feature selection methods used as the aggregation algorithms shown in figure 2.

Complete Linear Aggregation (CLA) and Weighted Mean Aggregation (WMA) [18] are representative of the class of algorithms that perform aggregation of list rankings, using the numbers ranging from 1 to the number of total features. Here, 1 represents the highest rank. CLA simply sums up every list
ranking to obtain the final ranking results for every feature. Finally, a number of features with highest ranks (smallest scores) are selected. WMA is an advanced version of CLA, treating model performance metrics of each feature selector as weights applied to feature ranks. Before summing up ranks, each rank will be weighted by a performance metric from the associated feature selector.

In our proposed model, Recursive Feature Elimination (RFE) is used to generate ranks based on the feature ratings from the Sklearn Decision Tree [4]. For WMA, we use \(1 - \text{AUC}\) as the weight for each rank. The \(\text{AUC}\) is obtained through a 10-fold cross validation using the training data for each ensemble instance.

Feature Occurrence Frequency (OFA) and Classification Accuracy Based Aggregation (CAA) [5] represent a class of algorithms that count the occurrence of each selected feature. OFA is the most straightforward method, only counting the occurrence of selected features from each ensemble instance. CAA takes one step further, weighing the counts for each occurrence with an associated classification accuracy. When selecting features, the ones with the largest score (weighted count of occurrence in CAA) are selected. In our proposed model, for each feature selector, we select features with only positive ratings using a Decision Tree and validate the AUC and accuracy using 10-fold cross validation.

C. Missing Data Rate/Entropy And Accuracy Based Aggregation (MAA/EAA)

Next, we present two novel ensemble-based feature selection methods: 1) Missing Data Rate And Accuracy Based Aggregation (MAA) and 2) Entropy And Accuracy Based Aggregation (EAA). Both methods are based on CAA, using the frequency of occurrence as the score. In addition to leveraging classification accuracy, our proposed approaches also consider mitigating the potential impact of the potential increase in variance introduced by data imputation when ranking features. MAA uses the missing data rate of each feature to adjust the weight, whereas EAA uses the change in entropy for each feature, computed before and after applying missing data handling. Equations 2 and 3 show the adjusted weights for MAA and EAA, respectively.

\[
\text{Accuracy} = \frac{\text{Accuracy}}{(\text{Missing_rate} + \alpha)^\beta} \quad (2)
\]

\[
\text{Accuracy} = \frac{\text{Accuracy}}{(\Delta \text{Entropy} + \alpha)^\beta} \quad (3)
\]

In the above two equations, \(\text{Missing_rate}\) is the missing data rate for the associated features and \(\Delta \text{Entropy}\) is the information entropy change of the associated features, as computed before and after applying missing data handling. \(\text{Accuracy}\) is the classification accuracy, and the \(\alpha\) and \(\beta\) are hyper-parameters to adjust the importance of the \(\text{Missing_rate}\) and \(\Delta \text{Entropy}\).

The implications of using MAA and EAA are to adjust the classification accuracy based on the magnitude of data variance, represented by the missing data rate and information entropy change for MAA and EAA, respectively. Instead of estimating the actual value of the data variance, we estimate the magnitude of data variance based on two assumptions: 1) features with a lower \(\text{Missing_rate}\) will experience less data variance after applying missing data handling, and 2) a small \(\Delta \text{Entropy}\) preserves the information entropy, leading to lower data variance. MAA uses the \(\text{Missing_rate}\) to select features with a lower missing data rate. Based on equation 2, whenever the missing data rate is high, the associated weight value shrinks, reducing the overall contribution of the corresponding feature. Likewise, EAA uses the \(\Delta \text{Entropy}\), following the same procedure.

MAA is a straightforward method to select the features that exhibit low data variance. However, MAA also potentially eliminates important features that have a relatively large missing data rate, but small \(\Delta \text{Entropy}\) values. EAA, on the other hand, focuses on using the change in information entropy to adjust the classification accuracy, potentially being able to select important features MAA ignores.

VI. EXPERIMENTS

A. Dataset Specifications

We generate a series of multi-dimensional datasets, selected based on different data completion rate thresholds (50% to 80%, in steps of 10%), with participants (i.e., mothers) and features selected across all visits (V1 to V4).

Table I lists the details of the datasets. As indicated in the table, the dataset is highly imbalanced in terms of the outcome class distribution. Almost 90% of the samples are negative (i.e., term births) and only 10% are positive (i.e., preterm birth).

\begin{table}[h]
\centering
\caption{Details of data.}
\begin{tabular}{|c|c|c|c|c|}
\hline
Data completion rate threshold & 50\% & 60\% & 70\% & 80\% \\
\hline
Number of samples & 667 & 647 & 626 & 598 \\
Number of positive samples & 66 & 58 & 53 & 51 \\
Number of negative samples & 601 & 589 & 573 & 547 \\
Number of features & 1011 & 975 & 954 & 928 \\
\hline
\end{tabular}
\end{table}

\begin{table}[h]
\centering
\caption{Parameters and elements for the undersampling ensemble learning model.}
\begin{tabular}{|c|c|c|}
\hline
Decision Tree & CART [22] & CLA, WMA, OFA, CAA \\
Ensemble Feature Selection Methods & MAA, EAA (\(\alpha : 1, 0.5; \beta : 2\)) & Accuracy, AUC \\
Performance metrics & Number of ensemble ways & 91 \\
Number of selected features & Number of samples & 20 \\
\hline
\end{tabular}
\end{table}

B. Experimental Setup

Table II below lists the parameters and elements used to construct the undersampling ensemble learning model. As indicated in table II, we use CART from sklearn [4] as the Decision Tree model in our proposed undersampling...
ensemble learning model. We leverage and compare CLA, WMA, OFA and CAA, as well as our proposed MAA and EAA, and evaluate accuracy and compute the AUC. We use an odd number as the ensemble way size so that we can easily apply majority voting during the performance evaluation step shown in figure 2. In addition, we found that an ensemble way size larger than 91 produces negligible performance gains in our experiments. Therefore, we select 91 as the ensemble way size. We use the number of selected features reported in previous studies that analyzed this same dataset [6] to provide a fair comparison. We use 10-fold cross validation in the performance evaluation step.

VII. Results

In this section, we present results from our experiments. We use AUC and classification accuracy as the major measures of model performance.

Given the same experimental setup as used in a previous study using the same dataset [6] (where only data from visits V1 and V2 were evaluated), the undersampling ensemble model combined with CAA produces a 30% improvement in AUC over a hybrid approach, while achieving an accuracy of 72%. When using the dataset that combines all 4 visits, the baseline method [6] is equivalent to random guessing (the number of samples is too small to apply the method explored in previous work [6]). Given this behavior, we exclude this option in our results. Using undersampling with ensemble learning, on the other hand, overcomes this issue.

Tables III and IV list the AUC and accuracy of CLA, WMA, OFA and CAA, while using both mean-based and similarity-based missing data imputation methods. From the results, we find that mean-based missing data imputation produces better performance across all ensemble feature selection methods. We see that CAA achieves the best performance in terms of both accuracy and AUC. Considering that the hybrid approach presented in prior work [6] produced results close to random guessing (AUC was close to 50%), our proposed analytical model that includes CAA can improve AUC by 42%, while achieving a 79% accuracy, as seen in Table III.

Next, we present the results based on data selected across different data completion rates. As CLA and WMA do not have as good performance as compared to OFA and CAA, we only present the results from OFA, CAA, MAA and EAA, with mean-based missing data imputation.

Figures 3 and 4 show the AUC and accuracy with ensemble feature selection with OFA, CAA, MAA and EAA, while varying the data completion rate from >50% to >80%, in steps of 10%. From these results we can see that EAA has the best performance. Both MAA and EAA have slightly better performance than OFA and CAA. They are both effective in selecting features, while improving evaluation performance.

Figures 5 and 6 show the average missing data rate and
the change in information entropy of features selected by OFA, CAA, MAA and EAA. From Figure 5, the feature set selected by MAA presents a lower missing data rate as compared to OFA and CAA. This is expected, as MAA adjusts the classification accuracy using the missing data rate. EAA, surprisingly, performs more effectively in selecting features with a lower missing data rate, especially when the data completion rate is larger than 70%. The feature set selected by EAA, as expected, results in a lower $\triangle \text{Entropy}$, as shown in Figure 6. From the results, MAA and EAA are found to be more effective at providing low-variance feature selection, as measured using the missing data rate and the $\triangle \text{Entropy}$, respectively.

**VIII. Conclusion**

In this paper, we present an analytical model for identifying potential key factors contributing to the high rates of preterm birth in northern Puerto Rico, based on the dataset available from the NIEHS P42 Center. We highlight the challenges of analyzing a diverse dataset with class imbalance in the target variable, as well as a high missing rate in predictors.

To address these issues, we propose an undersampling ensemble learning approach leveraging many state-of-the-art ensemble feature selection methods, including CLA, WMA, OFA and CAA. We found that a undersampling ensemble model equipped with CAA can achieve a 42% improvement in AUC as compared to previous studies. We also propose two novel feature selection methods, MAA and EAA, limiting the variance introduced by missing data handling. With OFA and CAA as a baseline, we evaluate MAA and EAA. Our results show that MAA and EAA are effective at selecting features with lower data variance, while providing slightly better performance than CAA.

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