Master Research Report: Multilevel Multivariate Imputation by Chained Equations through Bayesian Additive Regression Trees

Methodology and Statistics for the Behavioural, Biomedical and Social Sciences

Heleen Brüggen

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MSc. T. Volker Dr. G. Vink MSc. H. Oberman 1384 Computational Statistics & Data Analysis 23-1778

> Utrecht University Utrecht University Utrecht University

1 Introduction

Incomplete data is a common challenge in many fields of research. A frequent approach for dealing with incomplete data is listwise deletion, also known as complete-case analysis, which is to remove all incomplete cases from the data. However, this could possibly lead to biased results if the data is not Missing Completely At Random (MCAR), meaning the cause of the missing data is unrelated to the data [Austin et al., 2021, Enders, 2017, Kang, 2013, Rubin, 1976, van Buuren, 2018]. Furthermore, other approaches to dealing with incomplete data include: pairwise deletion, mean imputation and regression imputation, which also yield biased results [van Buuren, 2018]. Pairwise deletion, also known as available-case analysis, is to remove all incomplete cases from the analysis when considering a specific pair of variables. Pairwise deletion leads to unbiased results when correlation between variables are low and the data is MCAR. Mean imputation is to replace missing values with the mean of the observed values. Mean imputation will bias almost all estimates except the mean when the data is not MCAR. Regression imputation is to replace missing values with the predicted values from a regression model and is unbiased when the data is Missing At Random (MAR), meaning that the missing data is related to the observed data [Rubin, 1976], and the factor influencing the missingness is present in the data [van Buuren, 2018]. So, in order to use these ad hoc strategies to deal with incomplete data correctly, the missing data mechanism should be carefully determined. However, determining the missing data mechanism is often difficult and MCAR is in practise an unrealistic assumption [van Buuren, 2018]. We can imagine that when a system becomes more complex, thus increases the amount of parameters, testing these assumptions also increases in complexity.

One of these complex systems are multilevel data structures. Multilevel data is hierarchically structured, where, for example, students are nested within schools, or patients are nested within hospitals [Hox and Roberts, 2011, Hox et al., 2017]. Thus, in these types of data sets there are level-1 and level-2 variables. Level-1 variables relate to the individual within a class and level-2 variables relate to the class as a whole. The recommended statistical technique to analyzing these models are multilevel models as it accounts for the specific dependencies in the multilevel data sets [Hox and Roberts, 2011, Hox et al., 2017, Lüdtke et al., 2017]. It can contain both level-1 and level-2 variables, random intercepts, random slopes, and cross-level interactions [Hox and Roberts, 2011, Hox et al., 2017].

Multiple imputation (MI) is considered a valid method for dealing with incomplete data and allows us the separate the missing data problem from the analysis problem [Audigier et al., 2018, Austin et al., 2021, Burgette and Reiter, 2010, Enders, 2017, Grund et al., 2021, Hughes et al., 2014, Mistler and Enders, 2017, Van Buuren, 2007, van Buuren, 2018]. MI imputes each missing value in the data set more than once given the observed data, thereby considering necessary variation associated with the missingness problem. The multiply imputed data sets are analyzed, and the corresponding inferences are pooled according to Rubin's rules [Austin et al., 2021, Carpenter and Kenward, 2013, Rubin, 1987, van Buuren, 2018]. Generally, multiple imputation operates under two frameworks: joint modeling (JM) and fully conditional specification (FCS) [Enders et al., 2018a,b, Hughes et al., 2014, Mistler and Enders, 2017, van Buuren, 2018]. JM employs a multivariate data distribution and regresses incomplete variables on all complete variables to impute missing values. FCS, or chained equations, iteratively imputes one variable with missing values at a time through conditional univariate distributions regressing an incomplete variable on complete variables and previously imputed variables [Enders et al., 2018a,b, 2016, Grund et al., 2018a, Hughes et al., 2014, Mistler and Enders, 2017, van Buuren, 2018].

JM and FCS use multivariate and univariate linear mixed models respectively in a multilevel context [Enders et al., 2018a, Mistler and Enders, 2017]. The implementation of JM and FCS in a multilevel context are equivalent in a two-level random intercept context with normally distributed data but differ beyond that [Enders et al., 2018b, 2016]. Overall, FCS is believed to be more flexible than JM [Audigier et al., 2018, Burgette and Reiter, 2010, Grund et al., 2018a, Van Buuren, 2007] and, thus, may be better suited for multilevel data. In FCS, one needs to define conditional models for all variables with missing values [Enders et al., 2018a,b, 2016, Grund et al., 2018a, Hughes et al., 2014, Mistler and Enders, 2017, van Buuren, 2018] and the imputation models should at least be as general as the analysis model and preferably all-encompassing [Grund et al., 2018b]. However, the complexity of the multilevel analysis model is built step-wise with non-linearities [Hox and Roberts, 2011, Hox et al., 2017]. Thus, defining imputation models for a multilevel data set is quite challenging [Burgette and Reiter, 2010, Hox and Roberts, 2011, van Buuren, 2018].

Using non-parametric tree-based models as imputation models might solve this problem. Tree-based models use recursive partitioning to split the data into smaller subgroups based on the predictor variables maximizing the homogeneity of the subgroups. Tree-based models are non-parametric, which means that

they do not assume a specific distribution of the data. Thus, they can handle non-linear relationships and interactions between the predictor variables well. Furthermore they handle continuous and catergorical variables simultaneously [Breiman et al., 1984, Burgette and Reiter, 2010, Chipman et al., 2010, Hill et al., 2020, James et al., 2021, Lin and Luo, 2019, Salditt et al., 2023].

In a single-level imputation context, the use of tree-based, non-parametric models like regression trees, random forests or Bayesian Additive Regression Trees (BART) simplified imputation models and performed better than parametric methods: the imputations showed better confidence interval coverage of the population parameters, lower variance and lower bias, especially in non-linear and interactive contexts [Burgette and Reiter, 2010, Silva and Gutman, 2022, Xu et al., 2016]. Others have also found lower normalized root mean squared error (NRMSE), which in essence encapsulates the bias of the imputations, when imputing with a random forest algorithm compared to MICE and KNN imputation [Stekhoven and Bühlmann, 2012, Waljee et al., 2013]. Furthermore, they also found that the algorithm reduced computational time and could handle multivariate data consisting of both continuous and categorical data simultaneously.

BART models have also been implemented in a multilevel prediction context. However, multilevel-BART models (M-BART) have predominantly been implemented with random intercepts and no random slopes and cross-level interactions [Chen, 2020, Tan et al., 2016, Wagner et al., 2020, Wundervald et al., 2022]. Wagner et al. [2020] have found that this random intercept M-BART model provided better predictions with a lower Mean Squared Error (MSE) compared to a parametric multilevel model, Tan et al. [2016] found higher Area Under the Curve values, and Chen [2020] found better predictions and better coverage compared to parametric models and a single-level BART model. Other researchers modeled the random intercept as an extra split on each terminal node within the BART algorithm and found a lower MSE compared to a standard BART model and parametric multilevel models [Wundervald et al., 2022]. Dorie et al. [2022] developed a multilevel BART model that included random intercepts and random slopes by combining BART with the Stan algorithm. However, the random intercept and slope are modeled by Stan, which is a parametric method. Their results showed that their algorithm 'stan4bart' showed better coverage of the population value and lower Root Mean Squared Error (RMSE) compared to BART models with varying intercept, BART models ignoring the multilevel structure, Bayesian Causal Forests (BCF), and parametric multilevel models.

In spite of these promising findings: tree-based model performing well in single-level imputation context [Burgette and Reiter, 2010, Silva and Gutman, 2022, Stekhoven and Bühlmann, 2012, Waljee et al., 2013, Xu et al., 2016] and M-BART models performing well in a multilevel prediction context [Chen, 2020, Dorie et al., 2022, Tan et al., 2016, Wagner et al., 2020, Wundervald et al., 2022], M-BART models have yet to be implemented in a multilevel multiple imputation context. Thus, my research question will be: Can multivariate imputation by chained equations through a multilevel bayesian additive regression trees model improve the bias, variance and coverage of the estimates in a multilevel context compared to current practices? Given the success of non-parametric models in single-level multiple imputation, I anticipate that employing multilevel BART models in a multilevel missing data context will reduce bias, accurately model variance, and improve estimate coverage compared to classical multilevel imputation through 2l.pmm, 2l.lmer, 2l.pan, 2l.jomo, rf and single-level pmm and complete case analysis in the R-package MICE [Buuren and Groothuis-Oudshoorn, 2011].

This research report is organised as follows: in section 2, will contain some theoretical background and describe the methods in which I will implement the M-BART model in a multilevel imputation context and Section 3 will provide some preliminary results.

2 Method

2.1 Theoretical background

Bayesian Additive Regression Trees (BART) is a sum-of-trees model proposed by Chipman et al. [Chipman et al., 2010]. Regression trees are its building blocks [Chipman et al., 2010, Hill et al., 2020, James et al., 2021]. Regression trees model non-linearities well and automatically through recursive binary partitioning of the predictor space [Burgette and Reiter, 2010, Hill et al., 2020]. Recursive binary partitioning doesn't assume a specific data form; it divides the predictor space to maximize variance explanation by automatically identifying best fitting splits [Hastie, 2017, James et al., 2021, Salditt et al., 2023]. BART models can be described as:

$$y_i = f(\mathbf{x}_i) + \epsilon_i, \tag{1.1}$$

$$y_i = g(\mathbf{x}_i, T_1, M_1) + g(\mathbf{x}_i, T_2, M_2) + \dots + g(\mathbf{x}_i, T_k, M_k) + \epsilon_i,$$
 (1.2)

where y_i is the outcome variable for person i, $f(\mathbf{x}_i)$ is the sum-of-trees many regression trees, and ϵ_i is the error term; $\epsilon \sim \mathcal{N}(0, \sigma^2)$. \mathbf{x} are the predictors included in the model, T_k is the kth tree and M_k is the collection of leaf parameters within the kth tree [Chipman et al., 2010, Hill et al., 2020, James et al., 2021]. Next to the sum-of-trees model, BART also includes a regularization prior that constrains the size and fit of each tree so that each contributes only a small part of the overall fit to prevent overfitting [Chipman et al., 2010, Hill et al., 2020, James et al., 2021]. The Bayesian back-fitting Markov Chain Monte Carlo (MCMC) algorithm is used to obtain estimates from BART. It updates each tree, conditional on the remaining trees, their associated parameters and σ , by fitting a new tree to the partial residuals, r_i , perturbing the tree from the previous iteration. The partial residuals, r_i , are defined as:

$$r_i = y_i - \sum_{k' < k} \hat{f}_{k'}^b(x_i) - \sum_{k' > k} \hat{f}_{k'}^{b-1}(x_i), \tag{2}$$

where $\hat{f}_{k'}^b(x_i)$ is the prediction of the k'^{th} tree in the b^{th} iteration for person i. The M-BART model including a random intercept can be identified as:

$$y_{ij} = \sum_{k=1}^{m} f(\mathbf{X}_{ij}; T_k, M_k) + \alpha_j + \epsilon_{ij}, \tag{3}$$

where, now, y_{ij} is the outcome variable for person i in cluster j and α_j is the random intercept for cluster j.

2.2 Simulation study

I will conduct a simulation study to examine the performance of the M-BART model in a multilevel multiple imputation context through FCS. I will compare the M-BART model to the following multilevel and single level imputation methods: 2l.pmm, 2l.lmer, 2l.pan, 2l.jomo, rf and pmm in MICE [Buuren and Groothuis-Oudshoorn, 2011]. The following five factors will be varied:

- 1. Intraclass Correlation (ICC = 0, .05, .3 and .5)
- 2. Number of clusters (J = 30 and 50)
- 3. Within-cluster sample size $(n_i = 5, 15, 25 \text{ and } 50)$
- 4. The Missing At Random (MAR) and Missing Completely At Random (MCAR) data rate (0%, 25% and 50%)
- 5. Within-group effect size: (size of the regression coefficients $\beta = .2, .5$ and .8)

All these values are realistic in practice and/or previously proposed [Enders et al., 2020, 2018b, Grund et al., 2018b, Gulliford et al., 1999, Hox et al., 2017, Murray and Blitstein, 2003]. The ICC can be interpreted as the expected correlation between two randomly sampled individuals from the same group or the proportion of the total variance at level-2 [Gulliford et al., 2005, Hox and Roberts, 2011, Shieh, 2012]. For each combination of varying parameters, a 1000 replicated data sets will be generated.

The simulation study will be performed in R [R Core Team, 2023] with the package MICE [Buuren and Groothuis-Oudshoorn, 2011] to perform the FCS imputations, which I will enchance by integrating a multilevel BART model. FCS multilevel imputation methods 2l.pmm, 2l.lmer, 2l.pan, 2l.jomo, rf and pmm in MICE and complete case analysis will serve as a benchmark. The population data-generating

mechanism will be based on the following multilevel model:

$$y_{ij} = \beta_{0j} + \beta_{1j}X_{1ij} + \beta_{2j}X_{2ij} + \beta_{3j}X_{3ij} + \beta_{4j}X_{4ij} + \beta_{5j}X_{5ij} + \beta_{6j}X_{6ij} + \beta_{7j}X_{7ij} + \epsilon_{ij},$$
(4)

$$\beta_{0j} = \gamma_{00} + \gamma_{01} Z_{1j} + v_{0j}, \tag{4.1}$$

$$\beta_{1j} = \gamma_{10} + \gamma_{11} Z_{1j} + \nu_{1j}, \tag{4.2}$$

$$\beta_{2j} = \gamma_{20} + \gamma_{21} Z_{1j} + v_{2j}, \tag{4.3}$$

$$\beta_{3j} = \gamma_{30} + \gamma_{32} Z_{2j} + v_{3j}, \tag{4.4}$$

$$\beta_{4j} = \gamma_{40} + v_{4j},\tag{4.5}$$

$$\beta_{5j} = \gamma_{50} + v_{5j},\tag{4.6}$$

$$\beta_{6j} = \gamma_{60} + \nu_{6j},\tag{4.7}$$

$$\beta_{7j} = \gamma_{70},\tag{4.8}$$

where y_{ij} is a continuous level 1 outcome variable for person i in group j and Z_{1j} and Z_{2j} are continuous level 2 variables. The random intercept β_{0j} is determined by the grand mean γ_{00} , the group effect $\gamma_{01}Z_{1j}$ and the group-level random residuals v_{0j} .

The regression coefficients β_{1j} , β_{2j} , and β_{3j} for the continuous variables X_{1ij} , X_{2ij} , and X_{3ij} depend on the the intercepts γ_{10} , γ_{20} , and γ_{30} , the cross-level interactions $\gamma_{11}Z_{1j}$, $\gamma_{21}Z_{1j}$, and $\gamma_{32}Z_{2j}$, and the random slopes v_{1j} , v_{2j} , and v_{3j} . The regression coefficients β_{4j} , β_{5j} and β_{6j} are determined by the intercepts γ_{40} , γ_{50} and γ_{60} and the random slopes v_{4j} , v_{5j} and v_{6j} . The regression coefficient β_{7j} is determined by the intercept γ_{70} . The residuals and random slopes v_{0j} , v_{1j} , v_{2j} , v_{3j} , v_{4j} , v_{5j} , v_{6j} and ϵ_{ij} follow a zero-mean normal distribution. v_{1j} , v_{2j} , v_{3j} , v_{4j} , v_{5j} , v_{6j} and v_{6j} are multivariate normally distributed. The level-2 variables v_{1j} and v_{2j} are also multivariate normally distributed.

The estimates will be evaluated on their relative bias (the difference between the average estimate and the true value), modeled variance and the 95% confidence interval coverage [Oberman and Vink, 2023], which will be calculated as follows [Morris et al., 2019]:

$$Bias = \frac{1}{n_{\text{sim}}} \sum_{i=1}^{n_{\text{sim}}} (\hat{\theta}_i - \theta), \tag{5}$$

$$Coverage = \frac{1}{n_{\text{sim}}} \sum_{i=1}^{n_{\text{sim}}} I(\theta \in [\hat{\theta}_{\min,i}, \hat{\theta}_{\max,i}]), \tag{6}$$

$$SE = \frac{1}{n_{\text{sim}}} \sum_{i=1}^{n_{\text{sim}}} \sqrt{\frac{\sigma_{\hat{\theta}_i}^2}{N}},\tag{7}$$

where $\hat{\theta}_i$ is the estimated parameter, θ is the true value, n_{sim} is the number of simulated data sets, $\hat{\theta}_{\min,i}$ and $\hat{\theta}_{\max,i}$ are the lower and upper bounds of the 95% confidence interval, $\sigma_{\hat{\theta}_i}^2$ is the variance of the estimated parameter and N is the size of each simulated data set.

3 Results

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