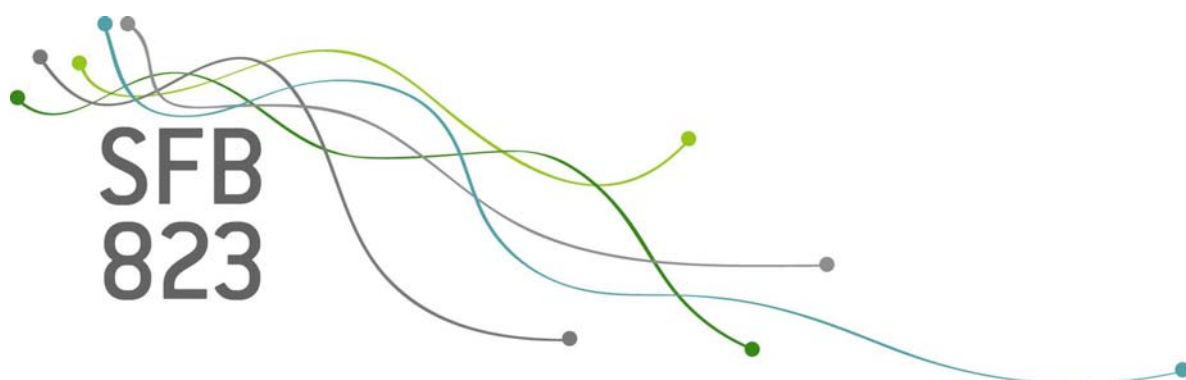


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Optimal designs for model averaging in non-nested models

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Discussion Paper

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Abstract

In this paper we construct optimal designs for frequentist model averaging estimation. We derive the asymptotic distribution of the model averaging estimate with fixed weights in the case where the competing models are non-nested and none of these models is correctly specified. A Bayesian optimal design minimizes an expectation of the asymptotic mean squared error of the model averaging estimate calculated with respect to a suitable prior distribution. We demonstrate that Bayesian optimal designs can improve the accuracy of model averaging substantially. Moreover, the derived designs also improve the accuracy of estimation in a model selected by model selection and model averaging estimates with random weights.

Keywords: Model selection, model averaging, model uncertainty, optimal design, Bayesian optimal designs

1 Introduction

There exists an enormous amount of literature on selecting an adequate model from a set of candidate models for statistical analysis. Numerous model selection criteria have been developed for this purpose. These procedures are widely used in practice and have the advantage of delivering a single model from a class of competing models, which makes them very attractive for practitioners. Exemplarily, we mention Akaike's information criterion (AIC), the Bayesian information criterion (BIC) and its extensions, Mallows's C_p , the generalized cross-validation and the minimum description length (see the monographs of Burnham and Anderson (2002), Konishi and Kitagawa (2008) and Claeskens and Hjort (2008) for more details). Different criteria have different properties, such as consistency, efficiency and parsimony (used in the sense of

Claeskens and Hjort (2008, Chapter 4)). Overall there seems to be no universally optimal model selection criterion and different criteria might be preferable in different situations depending on the particular application.

On the other hand, there exists a well known post-selection problem in this approach because model selection introduces an additional variance that is often ignored in statistical inference after model selection (see Pötscher (1991) for one of the first contributions discussing this issue). This post-selection problem is inter alia attributable to the fact, that estimates after model selection behave like mixtures of potential estimates. For example, ignoring the model selection step (and thus the additional variability) may lead to confidence intervals with coverage probability smaller than the nominal value, see for example Chapter 7 in Claeskens and Hjort (2008) for a mathematical treatment of this phenomenon.

An alternative to model selection is model averaging, where estimates of a target parameter are smoothed across several models, rather than restricting inference on a single selected model. This approach has been widely discussed in the Bayesian literature, where it is known as “Bayesian model averaging” (see the tutorial of Hoeting et al. (1999) among many others). For Bayesian model averaging prior probabilities have to be specified. This might not always be possible and therefore Hjort and Claeskens (2003) also proposed a “frequentist model averaging”, where smoothing across several models is commonly based on information criteria. Kapetanios et al. (2008) demonstrated that the frequentist approach is a worthwhile alternative to Bayesian model averaging. Stock and Watson (2003) observed that averaging predictions usually performs better than forecasting in a single model. Hong and Preston (2012) substantiate these observations with theoretical findings for Bayesian model averaging if the competing models are “sufficiently close”. Further results pointing in this direction can be found in Raftery and Zheng (2003), Schorning et al. (2016) and Buatois et al. (2018).

Independently of this discussion there exists a large amount of research how to optimally design experiments under model uncertainty (see Box and Hill (1967); Stigler (1971); Atkinson and Fedorov (1975) for early contributions). This work is motivated by the fact that an optimal design can improve the efficiency of the statistical analysis substantially if the postulated model assumptions are correct, but may be inefficient if the model is misspecified. Many authors suggested to choose the design for model discrimination such that the power of a test between competing regression models is maximized (see Ucinski and Bogacka (2005); López-Fidalgo et al. (2007); Tommasi and López-Fidalgo (2010) or Dette et al. (2015) for some more recent references). Other authors proposed to minimize an average of optimality criteria from different models to obtain an efficient design for all models under consideration (see Dette (1990), Zen and Tsai (2002); Tommasi (2009) among many others).

Although model selection or averaging are commonly used tools for statistical inference under model uncertainty most of the literature on designing experiments under model uncertainty does not address the specific aspects of these methods directly. Optimal designs are usually

constructed to maximize the power of a test for discriminating between competing models or to minimize a functional of the asymptotic variance of estimates in the different models. To the best of our knowledge Alhorn et al. (2019) is the first contribution, which addresses the specific challenges of designing experiments for model selection or model averaging. These authors constructed optimal designs minimizing the asymptotic mean squared error of the model averaging estimate and showed that optimal designs can yield a reduction of the mean squared error up to 45%. Moreover, they also showed that these designs improve the performance of estimates in models chosen by model selection criteria. However, their theory relies heavily on the assumption of nested models embedded in a framework of local alternatives as developed by Hjort and Claeskens (2003).

The goal of the present contribution is the construction of optimal designs for model averaging in cases where the competing models are not nested (note that in this case local alternatives cannot be formulated). Moreover, in contrast to most of the literature, we also consider the situation where all competing models misspecify the data underlying truth. In order to derive an optimality criterion, which can be used for the determination of optimal designs in this context, we further develop the approach of Hjort and Claeskens (2003) and derive an asymptotic theory for model averaging estimates for classes of competing models which are non-nested. Optimal designs are then constructed minimizing the asymptotic mean squared error of the model averaging estimate and it is demonstrated that these designs yield substantially more precise model averaging estimates. Moreover, these designs also improve the performance of estimates after model selection. Our work also contributes to the discussion of the superiority of model averaging over model selection. Most of the results presented in literature indicate that model averaging has some advantages over model selection in general. We demonstrate that conclusions of this type depend sensitively on the class of models under consideration. In particular we observe some advantages of estimation after model selection if the competing models are of rather different shape. Nevertheless, the optimal designs developed in this paper improve both estimation methods, where the improvement can be substantial in many cases. The remaining part of this paper is organized as follows. The pros and cons of model averaging and model selection are briefly discussed in Section 2 where we introduce the basic methodology and investigate the impact of similarity of the candidate models on the performance of the different estimates. In Section 3 we develop asymptotic theory for model averaging estimation in the case where the models are non-nested and all competing models might misspecify the underlying truth. Based on these results we derive a criterion for the determination of optimal designs. In Section 4 we study the performance of these designs by means of a simulation study. Finally, technical assumptions and proofs are given Section 6.

2 Model averaging versus model selection

In this section we introduce the basic terminology and also illustrate in a regression framework that the superiority of model averaging about estimation in a model chosen by model selection depends sensitively on the class of competing models.

2.1 Basic terminology

We consider data obtained at k different experimental conditions, say x_1, \dots, x_k chosen in a design space \mathcal{X} . At each experimental condition x_i one observes n_i responses, say y_{i1}, \dots, y_{in_i} ($i = 1, \dots, k$), and the total sample size is $n = \sum_{i=1}^k n_i$. We also assume that the responses y_{i1}, \dots, y_{in_i} are realizations of random variables of the form

$$Y_{ij} = \eta_s(x_i, \vartheta_s) + \varepsilon_{ij}, i = 1, \dots, k, j = 1, \dots, n_i, s = 1, \dots, r, \quad (2.1)$$

where the regression function η_s is a differentiable function with respect to the parameter ϑ_s and the random errors ε_{ij} are independent normally distributed with mean 0 and common variance σ^2 . Furthermore, the index s in η_s corresponds to different models (with parameters ϑ_s) and we assume that there are r competing regression functions η_1, \dots, η_r under consideration.

Having r different candidate models (differing by the regression functions η_s) a classical approach for estimating a parameter of interest, say μ , is to calculate an information criterion for each model under consideration and estimate this parameter in the model optimizing this information criterion. For this purpose, we denote the density of the normal distribution corresponding to a regression model (2.1) by $f_s(\cdot | x_i, \theta_s)$ with parameter $\theta_s = (\sigma^2, \vartheta_s)^\top$ and identify the different models by their densities f_1, \dots, f_r (note that in the situation considered in this sections these only differ in the mean). Using the observations $y_n = (y_{11}, \dots, y_{1n_1}, y_{21}, \dots, y_{kn_k})^\top$ estimate we calculate in each model the maximum likelihood estimate

$$\hat{\theta}_{n,s} = \arg \max_{\theta_s \in \Theta_s} \ell_{n,s}(\theta_s | y_n) \quad (2.2)$$

of the parameter θ_s , where

$$\ell_{n,s}(\theta_s | y_n) = \frac{1}{n} \sum_{i=1}^k \sum_{j=1}^{n_i} \log f_s(y_{ij} | x_i, \theta_s) \quad (2.3)$$

is the log-likelihood in candidate model f_s ($s = 1, \dots, r$). Note that we do not assume that the true data generating density is included in the set of candidate models f_1, \dots, f_r . Each

estimate $\hat{\theta}_{n,s}$ of the parameter θ_s yields an estimate

$$\hat{\mu}_s = \mu_s(\hat{\theta}_{n,s}), \quad (2.4)$$

for the quantity of interest, where μ_s is the target parameter in model s .

For example, regression models of the type (2.1) are frequently used in dose finding studies (see MacDougall (2006) or Bretz et al. (2008)). In this case a typical target function μ_s of interest is the “quantile” defined by

$$\mu_s(\theta_s) = \inf \left\{ x \in \mathcal{X} \mid \frac{\eta_s(x, \vartheta_s) - \eta_s(a, \vartheta_s)}{\eta_s(b, \vartheta_s) - \eta_s(a, \vartheta_s)} \geq \alpha \right\}. \quad (2.5)$$

The value defined in (2.5) is well-known as ED_α , that is, the effective dose at which $100 \times \alpha\%$ of the maximum effect in the design space $\mathcal{X} = [a, b]$ is achieved.

We now briefly discuss the principle of model selection and averaging to estimate the target parameter μ . For model selection we choose the model f_{s^*} from f_1, \dots, f_s , which maximizes Akaike’s information criterion (AIC)

$$\text{AIC}(f_s \mid Y_n) = 2\ell_{n,s}(\hat{\theta}_{n,s} \mid y_n) - 2p_s, \quad (2.6)$$

where p_s is the number of parameters in model f_s (see Claeskens and Hjort (2008), Chapter 2). The target parameter is finally estimated by $\hat{\mu} = \mu_{s^*}(\hat{\theta}_{n,s^*})$. Obviously, other model selection schemes, such as the Bayesian or focussed information criterion can be used here as well, but we restrict ourselves to the AIC for the sake of a transparent presentation.

Roughly speaking, model averaging is a weighted average of the individual estimates in the competing models. It might be viewed from a Bayesian (see for example Wassermann (2000)) or a frequentist point of view (see for example Claeskens and Hjort (2008)) resulting in different choices of model averaging weights. We will focus here on non-Bayesian methods. More explicitly, assigning nonnegative weights w_1, \dots, w_r to the candidate models f_1, \dots, f_r , with $\sum_{i=1}^r w_i = 1$, the model averaging estimate for μ is given by

$$\hat{\mu}_{\text{mav}} = \sum_{s=1}^r w_s \mu_s(\hat{\theta}_{n,s}). \quad (2.7)$$

Frequently used weights are uniform weights (see, for example Stock and Watson (2004), Kapetanios et al. (2008)). More elaborate model averaging weights can be chosen depending on the data. For example, Claeskens and Hjort (2008) define smooth AIC-weights as

$$w_s^{\text{smAIC}}(Y_n) = \frac{\exp\{\frac{1}{2}\text{AIC}(f_s \mid Y_n)\}}{\sum_{s=1}^r \exp\{\frac{1}{2}\text{AIC}(f_s \mid Y_n)\}}. \quad (2.8)$$

Model	Mean function η_s	Parameter specifications
Log-Linear (f_1)	$\eta_1(x_i, \vartheta_1) = \vartheta_{11} + \vartheta_{12} \log(x_i + \vartheta_{13})$	$\vartheta_1 = (0, 0.0797, 1)^\top$
Emax (f_2)	$\eta_2(x_i, \vartheta_2) = \vartheta_{21} + \frac{\vartheta_{22}x}{\vartheta_{23}+x}$	$\vartheta_2 = (0, 0.467, 25)^\top$
Exponential (f_3)	$\eta_3(x_i, \vartheta_3) = \vartheta_{31} + \vartheta_{32} \exp(x_i/\vartheta_{33})$	$\vartheta_3 = (-0.08265, 0.08265, 85)^\top$
Quadratic (f_4)	$\eta_4(x_i, \vartheta_4) = \vartheta_{41} + \vartheta_{42}x + \vartheta_{43}x^2$	$\vartheta_4 = (0, 0.00533, -0.00002)^\top$

Table 1: *Models and parameters used for the simulation study.*

Alternative data dependent weights can be constructed using other information criteria or model selection criteria. There also exists a vast amount of literature on determining optimal data dependent weights such that the resulting mean squared error of the model averaging estimate is minimal (see Hjort and Claeskens (2003), Hansen (2007) or Liang et al. (2011) among many others). For the sake of brevity we concentrate on smooth AIC-weights here, but similar observations as presented in this paper can also be made for other data dependent weights.

2.2 The class of competing models matters

In this section we illustrate the influence of the candidate set on the properties of model averaging estimation and estimation after model selection by means of a brief simulation study. For this purpose we consider four regression models of the form (2.1), which are commonly used in dose-response modeling and specified in Table 1 with corresponding parameters. Here we adapt the setting of Pinheiro et al. (2006) who model the dose-response relationship of an anti-anxiety drug, where the dose of the drug may vary in the interval $\mathcal{X} = [0, 150]$. In particular, we have $k = 6$ different dose levels $x_i \in \{0, 10, 25, 50, 100, 150\}$ and patients are allocated to each dose level most equally, where the total sample size is $n \in \{50, 100, 250\}$. We consider the problem of estimating the $ED_{0.4}$, as defined in (2.5).

To investigate the particular differences between both estimation methods we choose two different sets of competing models from Table 1. The first set

$$\mathcal{S}_1 = \{f_1, f_2, f_4\} \tag{2.9}$$

contains the log-linear, the Emax and the quadratic model, while the second set

$$\mathcal{S}_2 = \{f_1, f_2, f_3\} \tag{2.10}$$

contains the log-linear, the Emax and the exponential model. The set \mathcal{S}_1 serves as a prototype set of “similar” models while the set \mathcal{S}_2 contains models of more “different” shape. This is illustrated in Figure 1. In the left panel we show the quadratic model f_4 (for the parameters specified in Table 1) and the best approximations of this function by a log-linear model (f_1)

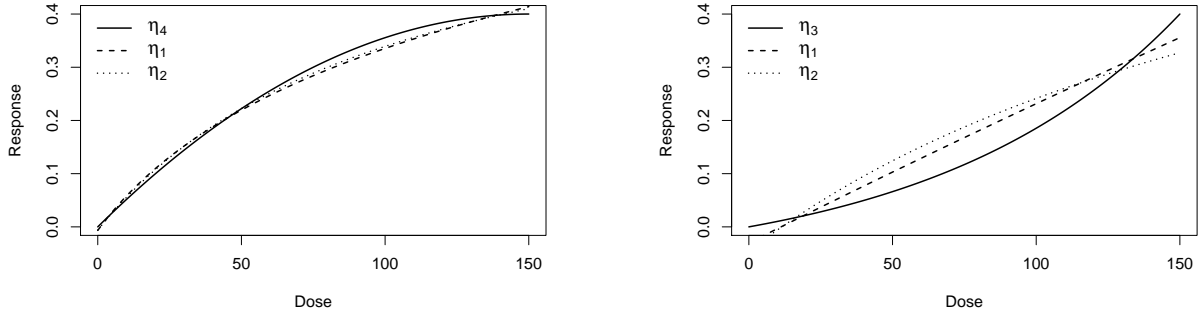


Figure 1: *Left panel: quadratic model (solid line) and its best approximations by the log-linear (dashed line) and the Emax model (dotted line) with respect to the Kullback-Leibler divergence (2.11). Right panel: exponential model (solid line) and its best approximations by the log-linear (dashed line) and the Emax model (dotted line).*

and an Emax model (f_2) with respect to the Kullback-Leibler divergence

$$\frac{1}{6} \sum_{i=1}^6 \int f_4(y | x_i, \theta_4) \log \left(\frac{f_4(y | x_i, \theta_4)}{f_s(y | x_i, \theta_s)} \right) dy, \quad s = 1, 2. \quad (2.11)$$

In this case, all models have a very similar shape and we obtain for the $ED_{0.4}$ the values 32.581, 32.261 and 33.810 for the log-linear (f_1), Emax (f_2) and quadratic model (f_4). Similarly the right panel shows the exponential model (f_3 , solid line) and its corresponding best approximations by the log-linear model (f_1) and the Emax model (f_2). Here we observe larger differences between the models in the candidate set and we obtain for the $ED_{0.4}$ the values 58.116, 42.857 and 91.547 for the models f_1 , f_2 and f_3 , respectively.

All results presented in this paper are based on 1000 simulations runs generating in each run n observations of the form

$$y_{ij}^{(l)} = \eta_s(x_i, \vartheta_s) + \varepsilon_{ij}^{(l)}, \quad i = 1, \dots, k, \quad j = 1, \dots, n_i, \quad (2.12)$$

where the errors $\varepsilon_{ij}^{(l)}$ are independent centered normal distributed random variables with $\sigma^2 = 0.1$ and η_s is one of the models η_1, \dots, η_4 (with parameters specified in Table 1). The parameter $\mu = ED_{0.4}$ is estimated by model averaging with uniform, smooth AIC weights in (2.8) and estimation after model selection by the AIC criterion.

In Table 2 and 3 we show the simulated mean squared errors of the model averaging estimates with uniform weights (left column), smooth AIC-weights (2.8) (middle column) and estimation after model selection (right column). Here, different rows correspond to different models. The numbers printed in bold face indicate the estimation method with the smallest mean squared error.

model	sample size	uniform weights	smooth AIC-weights	model selection
f_1	$n = 50$	437.045	498.323	758.978
	$n = 100$	223.291	218.99	285.062
	$n = 250$	111.973	82.713	78.371
f_2	$n = 50$	286.638	329.904	515.32
	$n = 100$	189.785	203.796	251.836
	$n = 250$	62.792	64.854	66.54
f_4	$n = 50$	276.037	361.101	669.873
	$n = 100$	190.662	244.558	391.443
	$n = 250$	92.653	109.852	139.859
f_3	$n = 50$	1503.903	1372.31	1381.033
	$n = 100$	1109.622	856.484	729.912
	$n = 250$	864.163	398.144	255.604

Table 2: *Simulated mean squared error of different estimates of the $ED_{0.4}$. The set of candidate models is $\mathcal{S}_1 = \{f_1, f_2, f_4\}$. Left column: model averaging with uniform weights; middle column: model averaging with smooth AIC-weights; right column: estimation after model selection.*

2.2.1 Models of similar shape

We will first discuss the results for the set of similar models in (2.9) (see Table 2). If the data generating model is an element of the set of candidate models, model averaging with uniform weights performs very well. Model averaging with smooth AIC-weights yields an about 10%-25% larger mean squared error (except for two cases, where it performs better than model averaging with uniform weights). On the other hand the mean squared error of estimation after model selection is substantially larger than that of model averaging, if the sample size is small. This is a consequence of the additional variability associated with data-dependent weights. For example, if the sample size is $n = 50$ and the data generating model is given by f_1 , the mean squared errors of the model averaging estimates with uniform and smooth AIC-weights and the estimate after model selection are given by 437.0, 498.3 and 759.0, respectively. The corresponding variances are given by 235.2, 337.6 and 599.7, respectively. For the squared bias the order is exactly the opposite, that is 201.9, 160.7, 159.3, but the differences are not so large. This means that the bias can be reduced by using random weights, because these put more weight on the “correct” model. As a consequence, compared to model averaging with uniform weights the performance of model averaging with smooth AIC-weights and the estimate after model selection improves with increasing sample size. Nevertheless, if the “true” model is an element of the candidate set and the functions in this set have a similar shape, model averaging performs better than estimation after model selection. In particular, model averaging with (fixed) uniform weights yield very reasonable results. These observation coincide with the findings of Schorning et al. (2016) and Buatois et al. (2018) who compared model averaging and model selection in the context of dose finding studies (see also Chen et al. (2018) for similar

model	sample size	estimation method		
		uniform weights	smooth AIC-weights	model selection
f_1	$n = 50$	834.295	553.427	776.311
	$n = 100$	712.404	340.254	353.707
	$n = 250$	524.518	48.587	38.591
f_2	$n = 50$	640.706	505.054	669.285
	$n = 100$	517.963	267.967	286.272
	$n = 250$	394.536	65.805	53.424
f_3	$n = 50$	1076.154	1141.476	1427.441
	$n = 100$	871.362	766.140	802.763
	$n = 250$	802.196	480.641	399.839
f_4	$n = 50$	288.091	486.501	852.377
	$n = 100$	208.628	298.315	419.651
	$n = 250$	162.689	138.331	142.673

Table 3: *Simulated mean squared error of different estimates of the $ED_{0.4}$. The set of candidate models is $\mathcal{S}_2 = \{f_1, f_2, f_3\}$. Left column: model averaging with uniform weights; middle column: model averaging with smooth AIC-weights; right column: estimation after model selection.*

results for the AIC in the context of ordered probit and nested logit models).

The situation changes if none of the candidate models from the set \mathcal{S}_1 is the “true” model. This is illustrated in the lower part of Table 2, where we show results if the exponential model f_3 is used for generating the data. We observe that model averaging with uniform weights is outperformed by model averaging with smooth AIC-weights. Moreover, the estimate after model selection is even better, if the sample size increases. These observations can be explained by the different shapes of the regression functions, as illustrated in Figure 1. By a suitable choice of parameters the quadratic model can adapt to the shape of the exponential model, whereas the log-linear and the Emax model still have different forms (see right panel of Figure 1 for the best approximations that are possible using the log-linear and the Emax model). Thus, incorporating these models in a model averaging estimate yields a large bias, that can be reduced substantially by data dependent weights or by model selection. For example, if $n = 100$ the squared bias of the model averaging estimate with uniform weights is 981.631, whereas the model averaging estimate with smooth AIC-weights and the estimate after model selection show a squared bias of 328.634 and 69.465, respectively.

2.2.2 Models of more different shape

We will now consider the candidate set \mathcal{S}_2 in (2.10), which serves as an example of more different models and includes the log-linear, the Emax and the exponential model. The simulated mean squared errors of the three estimates of the $ED_{0.4}$ are given in Table 3. The upper part of the table corresponds to cases, where data is generated from a model in the candidate set \mathcal{S}_2 used

for model selection and averaging. In contrast to Section 2.2.1 we observe only one scenario, where model averaging with uniform weights gives the smallest mean squared error (but in this case model averaging with smooth AIC-weights yields very similar results). If the sample size increases model averaging with smooth AIC-weights and estimation after model selection yield a substantially smaller mean squared error. An explanation of this observation consists in the fact that for a candidate set containing models with a rather different shape model averaging with uniform weights produces a large bias. On the other hand model averaging with smooth AIC-weights and estimation after model selection adapt to the data and put more weight on the “true” model, in particular if the sample size is large. As estimation after model selection has a larger variance and the variance is decreasing with increasing sample size, the bias is dominating the mean squared error for large sample sizes and thus estimation in the model selected by the AIC is more efficient for large sample sizes.

Finally, if the data is generated according to the quadratic model $f_4 \notin \mathcal{S}_2$ model averaging with uniform weights has the smallest mean squared error if the sample is $n = 50$ and $n = 100$. In this case estimation in the model selected by the AIC performs much worse (due to its large variance). However, the differences become smaller with increasing sample size. In particular for $n = 250$ model averaging with smooth AIC-weights and estimation after model selection show a substantially better performance than model averaging with uniform weights.

The numerical study in Section 2.2.1 and 2.2.2 can be summarized as follows. The results observed in the literature have to be partially relativized. The superiority of model averaging with uniform weights can only be observed for classes of “similar” competing models and a not too large signal to noise ratio. On the other hand if the models in the candidate set are of rather different structure, model averaging with data dependent weights (such as smooth AIC-weights) or estimation after model selection may show a better performance. For these reasons we will investigate optimal/efficient designs for all three estimation methods in the following sections. We will demonstrate that a careful design of experiments can improve the accuracy of these estimates substantially.

3 Asymptotic properties and optimal design

In this section we will derive the asymptotic properties of model averaging estimates with fixed weights in the case where the competing models are not nested. The results can be used for (at least) two purposes. On the one hand they provide some understanding of the empirical findings in Section 2, where we observed, that for increasing sample size the mean squared error of model averaging estimates is dominated by its bias. On the other hand, we will use these results to develop an asymptotic representation of the mean squared error of the model averaging estimate, which can be used in the construction of optimal designs.

3.1 Model averaging for non-nested models

Hjort and Claeskens (2003) provide an asymptotic distribution of frequentist model averaging estimates making use of local alternatives which require the true data generating process to lie inside a wide parametric model. All candidate models are sub-models of this wide model and the deviations in the parameters are restricted to be of order $n^{-1/2}$. Using this assumption results in convenient approximations for the mean squared error as variance and bias are both of order $O(1/n)$. However, in the discussion of this paper Raftery and Zheng (2003) pose the question if the framework of local alternatives is realistic. More importantly, frequentist model averaging is also often used for non-nested models (see for example Verrier et al. (2014)). In this section we will develop asymptotic theory for model averaging estimation in non-nested models. In particular, we do not assume that the “true” model is among the candidate models used in the model averaging estimate.

As we will apply our results for the construction of efficient designs for model averaging estimation we use the common notation of this field. To be precise, let Y denote a response variable and let x denote a vector of explanatory variables defined on a given compact design space \mathcal{X} . Suppose that Y has a density $g(y | x)$ with respect to a dominating measure. For estimating a quantity of interest, say μ , from the distribution g we use r different parametric candidate models with densities

$$f_1(y | x, \theta_1), \dots, f_r(y | x, \theta_r) \quad (3.1)$$

where θ_s denotes the parameter in the s th model, which varies in a compact parameter space, say $\Theta_s \subset \mathbb{R}^{p_s}$ ($s = 1, \dots, r$). Note, that in general we do not assume that the density g is contained in the set of candidate models in (3.1) and that the regression model (2.1) investigated in Section 2 is a special case of this general notation.

We assume that k different experimental conditions, say x_1, \dots, x_k , can be chosen in a design space \mathcal{X} and that at each experimental condition x_i one can observe n_i responses, say y_{i1}, \dots, y_{in_i} (thus the total sample size is $n = \sum_{i=1}^k n_i$), which are realizations of independent identically distributed random variables Y_{i1}, \dots, Y_{in_i} with density $g(\cdot | x_i)$. For example, if g coincides with f_s then the density of the random variables Y_{i1}, \dots, Y_{in_i} is given by $f_s(\cdot | x_i, \theta_s)$ ($i = 1, \dots, k$). To measure efficiency and to compare different experimental designs we will use asymptotic arguments and consider the case $\lim_{n \rightarrow \infty} \frac{n_i}{n} = \xi_i \in (0, 1)$ for $i = 1, \dots, k$. As common in optimal design theory we collect this information in the form

$$\xi = \{x_1, \dots, x_k; \xi_1, \dots, \xi_k\}, \quad (3.2)$$

which is called approximate design in the following discussion (see, for example, Kiefer (1974)). For an approximate design ξ of the form (3.2) and total sample size n a rounding procedure is applied to obtain integers n_i taken at each x_i ($i = 1, \dots, k$) from the not necessarily integer valued quantities $\xi_i n$ (see, for example Pukelsheim (2006), Chapter 12).

The asymptotic properties of the maximum likelihood estimate (calculated under the assumption that f_s is the correct density) is derived under certain assumptions of regularity (see the Assumptions (A1)-(A6) in Section 6). In particular, we assume that the functions f_s are twice continuously differentiable with respect to θ_s and that several expectations of derivatives of the log-densities exist. For a given approximate design ξ and a candidate density f_s we denote by

$$\text{KL}(g : f_s \mid \theta_s, \xi) = \int g(y \mid x) \log \left(\frac{g(y \mid x)}{f_s(y \mid x, \theta_s)} \right) dy d\xi(x), \quad (3.3)$$

the Kullback-Leibler divergence between the models g and f_s and assume that

$$\theta_{s,g}^*(\xi) = \arg \min_{\theta_s \in \Theta_s} \text{KL}(g : f_s \mid \theta_s, \xi) \quad (3.4)$$

is unique for each $s \in \{1, \dots, r\}$. For notational simplicity we will omit the dependency of the minimum on the density g , whenever it is clear from the context and denote the minimizer by $\theta_s^*(\xi)$. We also assume that the matrices

$$A_s(\theta_s, \xi) = \sum_{i=1}^k \xi_i \mathbb{E}_{g(\cdot \mid x_i)} \left(\frac{\partial^2 \log f_s(Y_{ij} \mid x_i, \theta_s)}{\partial \theta_s \partial \theta_s^\top} \right)_{s,t=1}^{p_s}, \quad (3.5)$$

$$B_{st}(\theta_s, \theta_t, \xi) = \sum_{i=1}^k \xi_i \mathbb{E}_{g(\cdot \mid x_i)} \left(\frac{\partial \log f_s(Y_{ij} \mid x_i, \theta_s)}{\partial \theta_s} \left(\frac{\partial \log f_t(Y_{ij} \mid x_i, \theta_t)}{\partial \theta_t} \right)^\top \right)_{s,t=1}^{p_s, p_t}, \quad (3.6)$$

exist, where expectations are taken with respect to the true distribution $g(\cdot \mid x_i)$.

Under standard assumptions White (1982) shows the existence of a measurable maximum likelihood estimate $\hat{\theta}_{n,s}$ for all candidate models which is strongly consistent for the (unique) minimizer $\theta_s^*(\xi)$ in (3.4). Moreover, the estimate is also asymptotically normal distributed, that is

$$\sqrt{n}(\hat{\theta}_{n,s} - \theta_s^*(\xi)) \xrightarrow{\mathcal{D}} \mathcal{N} \left(0, A_s^{-1}(\theta_s^*(\xi)) B_{ss}(\theta_s^*(\xi), \theta_s^*(\xi)) A_s^{-1}(\theta_s^*(\xi)) \right), \quad (3.7)$$

where we assume the existence of the inverse matrices, $\xrightarrow{\mathcal{D}}$ denotes convergence in distribution and we use the notations

$$A_s(\theta_s^*(\xi)) = A_s(\theta_s^*(\xi), \xi), \quad B_{st}(\theta_s^*(\xi), \theta_t^*(\xi)) = B_{st}(\theta_s^*(\xi), \theta_t^*(\xi), \xi) \quad (3.8)$$

($s, t = 1, \dots, r$). The following result gives the asymptotic distribution of model averaging estimates of the form (2.7).

Theorem 3.1. *If Assumptions (A1) - (A7) in Section 6.1 are satisfied, then the model averaging*

estimate (2.7) satisfies

$$\sqrt{n} \left(\hat{\mu}_{\text{mav}} - \sum_{s=1}^r w_s \mu_s(\theta_s^*(\xi)) \right) \xrightarrow{\mathcal{D}} \mathcal{N} \left(0, \sigma_w^2(\theta^*(\xi)) \right), \quad (3.9)$$

where the asymptotic variance is given by

$$\sigma_w^2(\theta^*(\xi)) = \sum_{s,t=1}^r w_s w_t \left(\frac{\partial \mu_s(\theta_s^*(\xi))}{\partial \theta_s} \right)^\top A_s^{-1}(\theta_s^*(\xi)) B_{st}(\theta_s^*(\xi), \theta_t^*(\xi)) A_t^{-1}(\theta_t^*(\xi)) \frac{\partial \mu_t(\theta_t^*(\xi))}{\partial \theta_t}. \quad (3.10)$$

Theorem 3.1 shows, that the model averaging estimate is biased for the true target parameter μ_{true} , unless we have $\sum_{s=1}^r w_s \mu_s(\theta_s^*(\xi)) = \mu_{\text{true}}$. Hence we aim to minimize the asymptotic mean squared error of the model averaging estimate. Note, that the bias does not depend on the sample size, while the variance is of order $O(1/n)$.

3.2 Optimal designs for model averaging of non-nested models

Alhorn et al. (2019) determined optimal designs for model averaging minimizing the asymptotic mean squared error of the estimate calculated in a class of nested models under local alternatives and demonstrated that optimal designs lead to substantially more precise model averaging estimates than commonly used designs in dose finding studies. With the results of Section 3.1 we can develop a more general concept of design of experiments for model averaging estimation, which is applicable for non-nested models and in situations, where the “true” model is not contained in the set of candidate models used for model averaging.

To be precise, we consider the criterion

$$\Phi_{\text{mav}}(\xi, g) = \frac{1}{n} \sigma_w^2(\theta^*(\xi)) + \left(\sum_{s=1}^r w_s \mu_s(\theta_s^*(\xi)) - \mu_{\text{true}} \right)^2 \approx \text{MSE}(\hat{\mu}_{\text{mav}}), \quad (3.11)$$

where μ_{true} is the target parameter in the “true” model with density g and $\sigma_w^2(\theta^*(\xi))$ and $\theta_s^*(\xi)$ are defined in (3.10) and (3.4), respectively. Note that this criterion depends on the “true” distribution via μ_{true} and the best approximating parameters $\theta_s^*(\xi) = \theta_{s,g}^*(\xi)$.

For estimating the target parameter μ via a model averaging estimate of the form (2.7) most precisely a “good” design ξ yields small values of the criterion function $\Phi_{\text{mav}}(\xi, g)$. Therefore, for a given finite set of candidate models f_1, \dots, f_r and weights $w_s, s = 1, \dots, r$, a design ξ^* is called *locally optimal design for model averaging estimation of the parameter μ* , if it minimizes the function $\Phi_{\text{mav}}(\xi, g)$ in (3.11) in the class of all approximate designs on \mathcal{X} . Here the term “locally” refers to the seminal paper of Chernoff (1953) on optimal designs for nonlinear regression models, because the optimality criterion still depends the unknown density $g(y | x)$.

A general approach to address this uncertainty problem is a Bayesian approach based on a class of models for the density g . To be precise, let \mathcal{G} denote a finite set of potential densities and let π denote a probability distribution on \mathcal{G} , then we call a design *Bayesian optimal design for model averaging estimation of the parameter μ* if it minimizes the function

$$\Phi_{\text{mav}}^{\pi}(\xi) = \int_{\mathcal{G}} \Phi_{\text{mav}}(\xi, g) d\pi(g) . \quad (3.12)$$

In general, the set \mathcal{G} can be constructed independently of the set of candidate models. However, if there is not much prior information available one can construct a class of potential models \mathcal{G} from the candidate set as follows. We denote the candidate set of models in (3.1) by \mathcal{S} . Each of these models depends on a unknown parameter θ_s and we denote by $\mathcal{F}_{f_s} \subset \Theta_s$ a set of possible parameter values for the model f_s . Now let π_2 denote a prior distribution on \mathcal{S} and for each $f_s \in \mathcal{S}$ let $\pi_1(\cdot | f_s)$ denote a prior distribution on \mathcal{F}_{f_s} . Finally, we define $\mathcal{G} = \{(g, \theta) : g \in \mathcal{S}, \theta \in \mathcal{F}_g\}$ and a prior

$$d\pi(g, \theta) = d\pi_1(\theta | g) d\pi_2(g) , \quad (3.13)$$

then the criterion (3.12) can be rewritten as

$$\Phi_{\text{mav}}^{\pi}(\xi) = \int_{\mathcal{S}} \int_{\mathcal{F}_g} \Phi_{\text{mav}}(\xi, g) d\pi_1(\theta | g) d\pi_2(g), \quad (3.14)$$

In the finite sample study of the following section the set \mathcal{S} and the set \mathcal{F}_g (for any $g \in \mathcal{S}$) are finite, which results in a finite set \mathcal{G} .

We conclude noting that the optimality criteria proposed in this section have been derived for model averaging estimates with fixed weights. The asymptotic theory presented here cannot be easily adapted to estimates using data-dependent (random) weights (as considered in Section 2), because it is difficult to get an explicit expression for the asymptotic distribution, which is not normal in general. Nevertheless, we will demonstrate in the following section that designs minimizing the mean squared error of model averaging estimates with fixed weights will also yield a substantial improvement in model averaging estimation with smooth AIC-weights and in estimation after model selection.

4 Bayesian optimal designs for model averaging

We will demonstrate by means of a simulation study that the performance of all considered estimates can be improved substantially by the choice of an appropriate design. For this purpose we consider the same situation as in Section 2, that is regression models of the form (2.1) with centred normal distributed errors. We also consider the two different candidate sets \mathcal{S}_1 and

\mathcal{S}_2 defined in (2.9) (log-linear, Emax and quadratic model) and (2.10) (log-linear, Emax and exponential model), respectively.

Using the criterion introduced in Section 3 we now determine a Bayesian optimal design for model averaging estimation of the $\text{ED}_{0.4}$ with uniform weights from $n = 100$ observations. We require a prior distribution for the unknown density g , and we use a distribution of the form (3.13) for this purpose. To be precise, let $f_s(y | x, \theta_s)$ denote the density of a normal distribution with mean $\eta_s(x, \vartheta_s)$ and variance $\sigma_s^2 = 0.1$ ($s = 1, \dots, r$), where the mean functions are given in Table 1. As the criterion (3.14) does not depend on the intercept ϑ_{s1} , these are not varied and taken from Table 1. For each of the other parameters we use three different values: the values specified in Table 1 and a 10% larger and smaller value of this parameter.

$$\begin{aligned}\mathcal{F}_{f_1} &= \{(0, \vartheta_{12}, \vartheta_{13}) : \vartheta_{12} = 0.0797 \pm 10\%, \vartheta_{13} = 1 \pm 10\%\}, \\ \mathcal{F}_{f_2} &= \{(0, \vartheta_{22}, \vartheta_{23}) : \vartheta_{22} = 0.467 \pm 10\%, \vartheta_{23} = 25 \pm 10\%\}, \\ \mathcal{F}_{f_3} &= \{(-0.08265, \vartheta_{32}, \vartheta_{33}) : \vartheta_{32} = 0.08265 \pm 10\%, \vartheta_{33} = 85 \pm 10\%\}, \\ \mathcal{F}_{f_4} &= \{(0, \vartheta_{42}, \vartheta_{43}) : \vartheta_{42} = 0.00533 \pm 10\%, \vartheta_{43} = -0.00002 \pm 10\%\}.\end{aligned}\tag{4.1}$$

4.1 Models of similar shape

We will first consider the candidate set $\mathcal{S}_1 = \{f_1, f_2, f_4\}$ consisting of the log-linear, the Emax and the quadratic model. For the definition of the prior distribution (3.13) in the criterion (3.14) we consider a uniform distribution π_2 on the set \mathcal{S}_1 and a uniform prior $\pi_1(\cdot | f_s)$ on each set \mathcal{F}_{f_s} in (4.1) ($s = 1, 2, 4$). The Bayesian optimal design for model averaging estimation of the $\text{ED}_{0.4}$ minimizing the criterion (3.14) has been calculated numerically using the COBYLA algorithm (see Powell (1994)) and is given by

$$\xi_{\mathcal{S}_1}^* = \{0, 14.447, 49.283, 150; 0.227, 0.167, 0.365, 0.240\} .\tag{4.2}$$

We will compare this design with the design

$$\xi_1 = \{0, 10, 25, 50, 100, 150; 1/6, 1/6, 1/6, 1/6, 1/6, 1/6\},\tag{4.3}$$

proposed in Pinheiro et al. (2006) for a similar setting (this design has also been used in Section 2) and the locally optimal design for estimation of the $\text{ED}_{0.4}$ in the log-linear model f_1 with parameter $\theta_1 = (0, 0.0797, 1)^\top$ given by

$$\xi_2 = \{0, 4.051, 150; 0.339, 0.5, 0.161\}\tag{4.4}$$

(see Dette et al. (2010)). Results for the locally optimal designs for the estimation of the $\text{ED}_{0.4}$ in the Emax and exponential model are similar and omitted for the sake of brevity. We use the

model	design	uniform weights	smooth AIC-weights	model selection
f_1	(4.2)	164.338	161.239	182.069
	(4.3)	223.291	218.99	285.062
	(4.4)	185.251	184.77	340.698
f_2	(4.2)	122.665	135.969	168.577
	(4.3)	189.785	203.796	251.836
	(4.4)	501.814	501.394	1162.654
f_4	(4.2)	174.535	212.746	331.552
	(4.3)	190.662	244.558	391.443
	(4.4)	404.716	427.548	1396.051
f_3	(4.2)	1073.859	804.385	630.458
	(4.3)	1109.622	856.484	729.912
	(4.4)	3184.11	3413.566	4102.964

Table 4: *Simulated mean squared errors of different estimates of the $ED_{0.4}$ for different experimental designs. The set of candidate models is $\mathcal{S}_1 = \{f_1, f_2, f_4\}$. Left column: model averaging estimate with uniform weights; middle column: model averaging estimate with smooth AIC-weights; right column: estimate after model selection.*

same setup as in Section 2. For the sake of brevity we only report results for the sample size $n = 100$. Other results are available from the authors.

The corresponding results are given in Table 4, where we use the models f_1, f_2, f_3 and f_4 from Table 1 to generate the data (note that the model f_3 is not in the candidate set used for model averaging and model selection). The different columns represent the different estimation methods (left column: model averaging with uniform weights; middle column: smooth AIC-weights, right column: model selection). The numbers printed in boldface indicate the minimal mean squared error for each estimation method obtained from the different experimental designs. First, we consider the situation, where the data generating model is contained in the set of candidate models $\mathcal{S}_1 = \{f_1, f_2, f_4\}$ corresponding to the upper part of the table. We observe that in this case model averaging yields better results than estimation after model selection and this superiority is independent of the design under consideration. Compared to the designs ξ_1 and ξ_2 the Bayesian optimal design $\xi_{\mathcal{S}_1}^*$ for model averaging with uniform weights improves the efficiency of all estimation techniques. For example, when data is generated using the log-linear model f_1 the mean squared error of the model averaging estimate with uniform weights is reduced by 26.4% and 11.3%, when the optimal design is used instead of the designs ξ_1 or ξ_2 , respectively. This improvement is remarkable as the design ξ_2 is locally optimal for estimating the $ED_{0.4}$ in the model f_1 and data is generated from this model. In other cases the improvement is even more visible. For example, if data is generated by the model f_2 the improvement in model averaging estimation with uniform weights is 35.4% and 75.6% compared to the designs ξ_1 and ξ_2 . Moreover, although the designs are constructed for model averaging

with uniform weights they also yield substantially more accurate model averaging estimates with smooth AIC-weights and a more precise estimate after model selection. For example, if the data is generated from model f_1 the mean squared error is reduced by 26.4% and by 12.7% for estimation with smooth AIC-weights and by 36.1% and 46.6% for estimation after model selection, respectively. Similar results can be observed for the models f_2 and f_4 .

Next, we consider the case where the data is generated from the exponential model f_3 , which is not contained in the candidate set \mathcal{S}_1 . The efficiency of all three estimates improves substantially by the use of the Bayesian optimal design $\xi_{\mathcal{S}_1}^*$. Interestingly, the improvement is less pronounced for model averaging with uniform weights (3.2% and 66.3% compared to the designs ξ_1 and ξ_2 , respectively) than for smooth AIC-weights (6.1% and 76.4%) and estimation after model selection (13.6% and 84.6%).

Summarizing, our numerical results show that the Bayesian optimal design for model averaging estimation of the $ED_{0.4}$ yields a substantial improvement of the mean squared error of the model averaging estimate with uniform weights (3.2%-75.6%), smooth AIC-weights (6.1%-76.4%) and the estimate after model selection (13.6%-85.5%) for all four models under consideration.

4.2 Models of different shape

We will now consider the second candidate set \mathcal{S}_2 consisting of the log-linear (f_1) the Emax (f_2) and the exponential model (f_3). For the definition of the prior distribution (3.13) in the criterion (3.14) we use a uniform distribution π_2 on the set \mathcal{S}_2 and a uniform prior $\pi_1(\cdot | f_s)$ on each set \mathcal{F}_{f_s} ($s = 1, 2, 3$) in (4.1). For this choice the Bayesian optimal design for model averaging estimation of the $ED_{0.4}$ is given by

$$\xi_{\mathcal{S}_2}^* = \{0, 5.529, 74.303, 77.186, 150; 0.179, 0.142, 0.274, 0.162, 0.243\}, \quad (4.5)$$

and has (in comparison to the design $\xi_{\mathcal{S}_1}^*$ in Section 4.2) five instead of four support points. The simulated mean squared errors of the three estimates under different designs are given in Table 5. We observe again that compared to the designs ξ_1 and ξ_2 the Bayesian optimal design $\xi_{\mathcal{S}_2}^*$ improves most estimation techniques substantially. However, if model averaging with uniform weights is used and data is generated by model f_1 , the mean squared error of the model averaging estimate from the optimal design is 1.7% larger than the mean squared error obtained by the design ξ_1 . For model averaging with smooth AIC-weights this difference is 2.8%. Overall, the reported results demonstrate a substantial improvement in efficiency by the use of the Bayesian optimal design independently of the estimation method. If the Bayesian optimal design is used, estimation after model selection yields the smallest mean squared error if the data is generated from a model of the candidate set \mathcal{S}_2 . On the other hand, if data is generated from model $f_4 \notin \mathcal{S}_2$ model averaging with equal weights shows the best performance. Summarizing, our numerical results show that compared to the designs ξ_1 and ξ_2 the design $\xi_{\mathcal{S}_2}^*$

model	design	estimation method		
		uniform weights	smooth AIC-weights	model selection
f_1	(4.5)	724.682	349.615	344.991
	(4.3)	712.404	340.254	353.707
	(4.4)	798.473	444.457	459.335
f_2	(4.5)	510.354	239.594	206.281
	(4.3)	517.963	267.967	286.272
	(4.4)	1155.793	1025.269	1835.355
f_3	(4.5)	834.284	665.07	646.381
	(4.3)	871.362	766.140	802.763
	(4.4)	1526.230	1842.633	2721.415
f_4	(4.5)	148.559	307.103	388.958
	(4.3)	208.628	298.315	419.651
	(4.4)	522.652	610.198	1907.066

Table 5: *Simulated mean squared errors of different estimates of the $ED_{0.4}$ for different experimental designs. The set of candidate models is $\mathcal{S}_2 = \{f_1, f_2, f_3\}$. Left column: model averaging estimate with uniform weights; middle column: model averaging estimate with smooth AIC-weights; right column: estimate after model selection.*

reduces the mean squared error of model averaging estimates with uniform weights up to 71.6%. Furthermore, for smooth AIC-weights and estimation after model selection the reduction can be even larger and is up to 76.6% and 88.8%, respectively. These improvements hold also for the quadratic model f_4 , which is not contained in the candidate set \mathcal{S}_2 used in the definition of the optimality criterion.

5 Conclusions

In this paper we derived the asymptotic distribution of the frequentist model averaging estimate with fixed weights from a class of not necessarily nested models. We neither assume that this class contains the “true” model. We use these results to determine Bayesian optimal designs for model averaging, which improve the estimation accuracy of the estimate substantially. Although these designs are constructed for model averaging with fixed weights, they also yield a substantial improvement of accuracy for model averaging with data dependent weights and for estimation after model selection.

We also demonstrate that the superiority of model averaging against estimation after model selection depends sensitively on the class of competing models, which is used in the model averaging procedure. If the competing models are similar (which means that a given model from the class can be well approximated by all other models), then model averaging should be preferred. Otherwise, we observe advantages for estimation after model selection, in particular,

if the signal to noise ratio is small.

Although, the new designs show a very good performance for estimation after model selection and for model averaging with data dependent weights, it is of interest to develop optimal designs, which address the specific issues of data dependent weights in the estimates. This is a very challenging problem for future research as there is no simple expression of the asymptotic mean squared error of these estimates. A first approach to solve this problem is an adaptive one and a further interesting and very challenging question of future research is to improve the accuracy of adaptive designs.

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6 Technical assumptions and proof of Theorem 3.1

6.1 Assumptions

Following White (1982) we assume:

- (A1) The random variables $Y_{ij}, i = 1, \dots, k, j = 1, \dots, n_i$ are independent. Furthermore, Y_{i1}, \dots, Y_{in_i} have a common distribution function with a measurable density $g(\cdot | x_i)$ with respect to a dominating measure ν .
- (A2) The distribution function of each candidate model $s \in \{1, \dots, r\}$ has a measurable density $f_s(\cdot | x, \theta_s)$ with respect to ν (for all $\theta_s \in \Theta_s$) that is continuous in θ_s .
- (A3) For all $x \in \mathcal{X}$ the expectation $\mathbb{E}(\log(g(Y | x)))$ exists (where expectation is taken with respect to $g(\cdot | x)$) and for each candidate model the function $y \mapsto |\log f_s(y | x, \theta_s)|$ is dominated by a function that is integrable with respect to $g(\cdot | x)$ and does not depend on θ_s . Furthermore the Kullback-Leibler divergence (3.3) has a unique minimum

$$\theta_{s,g}^*(\xi) = \arg \min_{\theta_s \in \Theta_s} \text{KL}(g : f_s | \theta_s, \xi), \quad (\text{A.1})$$

and $\theta_{s,g}^*(\xi)$ is an interior point of Θ_s .

(A4) For all $x \in \mathcal{X}$ the function $y \mapsto \frac{\partial \log f_s(y|x, \theta_s)}{\partial \theta_s}$ is a measurable function for all $\theta_s \in \Theta_s$ and continuously differentiable with respect to θ_s for all $y \in \mathbb{R}$.

(A5) The entries of the (matrix valued) functions

$$y \mapsto C^{(s)}(y) = \frac{\partial^2 \log f_s(y|x, \theta_s)}{\partial \theta_s \partial \theta_s^\top}, \quad y \mapsto D^{(st)}(y) = \frac{\partial \log f_s(y|x, \theta_s)}{\partial \theta_s} \left(\frac{\partial \log f_t(y|x, \theta_t)}{\partial \theta_t} \right)^\top$$

are dominated by integrable functions with respect to $g(\cdot|x)$ for all $x \in \mathcal{X}$ and $\theta_s \in \Theta_s$.

(A6) The matrices $B_{ss}(\theta_s^*(\xi), \theta_s^*(\xi), \xi)$ and $A_s(\theta_s^*(\xi), \xi)$ in (3.5) and (3.6) are nonsingular.

(A7) The functions $\theta_s \mapsto \mu_s(\theta_s)$ are once continuously differentiable.

6.2 Proof of Theorem 3.1

By equation (A.2) in White (1982) we have

$$\sqrt{n}(\hat{\theta}_{n,s} - \theta_s^*(\xi)) + A_s^{-1}(\theta_s^*(\xi)) \frac{1}{\sqrt{n}} \sum_{i=1}^k \sum_{j=1}^{n_i} \frac{\partial \log f_s(Y_{ij} | x_i, \theta_s^*(\xi))}{\partial \theta_s} \xrightarrow{p} 0, \quad (\text{A.2})$$

where \xrightarrow{p} denotes convergence in probability (note that the matrix $A_s(\theta_s^*) = A_s(\theta_s^*, \xi)$ is nonsingular by assumption). An application of the multivariate central limit theorem now leads to

$$\frac{1}{\sqrt{n}} \begin{pmatrix} \sum_{i=1}^k \sum_{j=1}^{n_i} \frac{\partial \log f_1(Y_{ij} | x_i, \theta_1^*(\xi))}{\partial \theta_1} \\ \vdots \\ \sum_{i=1}^k \sum_{j=1}^{n_i} \frac{\partial \log f_r(Y_{ij} | x_i, \theta_r^*(\xi))}{\partial \theta_r} \end{pmatrix} \xrightarrow{\mathcal{D}} \mathcal{N} \left(0, \begin{pmatrix} B_{11} & \dots & B_{1r} \\ \vdots & \ddots & \vdots \\ B_{r1} & \dots & B_{rr} \end{pmatrix} \right), \quad (\text{A.3})$$

where $B_{st} = B_{st}(\theta_s^*(\xi), \theta_t^*(\xi), \xi)$ is defined in (3.6). Combining (A.2) and (A.3) we obtain the weak convergence of the vector $\hat{\theta}_n = (\hat{\theta}_{n,1}^\top, \dots, \hat{\theta}_{n,r}^\top)^\top$, that is

$$\sqrt{n}(\hat{\theta}_n - \theta^*(\xi)) \xrightarrow{\mathcal{D}} \mathcal{N}(0, \Sigma), \quad (\text{A.4})$$

where $\Sigma = (\Sigma_{st})_{s,t=1,\dots,r}$ is a block matrix with $p_s \times p_t$ entries

$$\Sigma_{st} = A_s^{-1}(\theta_s^*(\xi)) B_{st}(\theta_s^*(\xi), \theta_t^*(\xi)) A_t^{-1}(\theta_t^*(\xi)), \quad s, t = 1, \dots, r \quad (\text{A.5})$$

and the vector $\theta_s^*(\xi)$ is given by

$$\theta^*(\xi) = ((\theta_1^*(\xi))^T, \dots, (\theta_r^*(\xi))^T)^T. \quad (\text{A.6})$$

Next, we define for the parameter vector $\theta^\top = (\theta_1^\top, \dots, \theta_r^\top) \in \mathbb{R}^{\sum_{s=1}^r p_s}$ the projection π_s by $\pi_s \theta := \theta_s$ and the vector

$$\tilde{\mu}(\theta) = (\mu_1(\pi_1 \theta), \dots, \mu_r(\pi_r \theta))^T \quad (\text{A.7})$$

with derivative

$$\mu'_\theta = \begin{pmatrix} \left(\frac{\partial \mu_1(\theta_1)}{\partial \theta_1}\right)^\top & 0 & \dots & 0 \\ 0 & \left(\frac{\partial \mu_2(\theta_2)}{\partial \theta_2}\right)^\top & 0 & \dots & 0 \\ 0 & \dots & 0 & \left(\frac{\partial \mu_r(\theta_r)}{\partial \theta_r}\right)^\top \end{pmatrix}. \quad (\text{A.8})$$

An application of the Delta method (see for example van der Vaart (1998, Chapter 3)) now shows that

$$\sqrt{n}(\tilde{\mu}(\hat{\theta}_n) - \tilde{\mu}(\theta^*(\xi))) \xrightarrow{\mathcal{D}} \mathcal{N}\left(0, \mu'_{\theta^*(\xi)} \Sigma (\mu'_{\theta^*(\xi)})^\top\right). \quad (\text{A.9})$$

The assertion finally follows from the continuous mapping theorem observing the representation

$$\hat{\mu}_{\text{mav}} = (w_1, \dots, w_r) \tilde{\mu}(\hat{\theta}_n). \quad (\text{A.10})$$

