Nonparametric Empirical Bayes Estimation on Heterogeneous Data

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Abstract

The simultaneous estimation of many parameters based on data collected from corresponding studies is a key research problem that has received renewed attention in the high-dimensional setting. Many practical situations involve heterogeneous data where heterogeneity is captured by a nuisance parameter. Effectively pooling information across samples while correctly accounting for heterogeneity presents a significant challenge in large-scale estimation problems. We address this issue by introducing the “Nonparametric Empirical Bayes Structural Tweedie” (NEST) estimator, which efficiently estimates the unknown effect sizes and properly adjusts for heterogeneity via a generalized version of Tweedie’s formula. For the normal means problem, NEST simultaneously handles the two main selection biases introduced by heterogeneity: one, the selection bias in the mean, which cannot be effectively corrected without also correcting for, two, selection bias in the variance. We develop theory to show that NEST is asymptotically as good as the optimal Bayes rule that uniquely minimizes a weighted squared error loss. In our simulation studies NEST outperforms competing methods, with much efficiency gains in many settings. The proposed method is demonstrated on estimating the batting averages of baseball players and Sharpe ratios of mutual fund returns. Extensions to other members of the two-parameter exponential family are discussed.

Keywords: compound decision, double shrinkage estimation, kernelized Stein’s discrepancy, non-parametric empirical Bayes, Tweedie’s formula.

1 Introduction

Suppose that we are interested in estimating a vector of parameters $\mu = (\mu_1, \ldots, \mu_n)$ based on the summary statistics $Y_1, \ldots, Y_n$ from $n$ study units. The setting where $Y_i \mid \mu_i \sim N(\mu_i, \sigma^2)$ is the most well-known example, but the broader scope includes the compound estimation of Poisson parameters $\lambda_i$, Binomial parameters $p_i$, and other members of the exponential family.

In modern large-scale applications it is often of interest to perform simultaneous and selective inference (Benjamini and Yekutieli, 2011, Berk et al., 2013, Weinstein et al., 2013), which has

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called for solving the compound estimation problem in new ways and for new purposes. For example, there has been recent work on how to construct valid simultaneous confidence intervals of $\mu_i$’s after a selection procedure is applied (Lee et al., 2016). In multiple testing, as well as related ranking and selection problems, it is often desirable to incorporate estimates of the effect sizes $\mu_i$ in the decision process to prioritize the selection of more scientifically meaningful hypotheses (Basu et al., 2017, Benjamini and Hochberg, 1997, He et al., 2015, Henderson and Newton, 2016, Sun and McLain, 2012).

However, the simultaneous inference of thousands of means, or other parameters, is challenging because, as described in Efron (2011), the large scale of the problem introduces selection bias, wherein some data points are large merely by chance, causing traditional estimators to over-estimate the corresponding means. Shrinkage estimation, exemplified by the seminal work of James and Stein (1961), has been widely used in simultaneous inference. There are several popular classes of methods, including linear shrinkage estimators (Berger, 1976, Efron and Morris, 1975, James and Stein, 1961), non-linear thresholding–based estimators motivated by sparse priors (Abramovich et al., 2006, Donoho and Johnstone, 1994, Johnstone and Silverman, 2004), and both Bayes or empirical Bayes estimators with unspecified priors (Brown and Greenshtein, 2009, Castillo and van der Vaart, 2012, Jiang and Zhang, 2009). This article focuses on a class of estimators based on Tweedie’s formula (Dyson, 1926, Eddington, 1940, Robbins, 1956). The formula is an elegant shrinkage estimator, for distributions from the exponential family, that has recently received renewed interest (Brown and Greenshtein, 2009, Efron, 2011, Koenker and Mizera, 2014). Tweedie’s formula is simple and intuitive, and its implementation only requires estimating the marginal distribution of $Y_i$. This property is particularly appealing for large–scale estimation problems where such estimates can be easily constructed from the observed data. The resultant empirical Bayes estimator enjoys optimality properties (Brown and Greenshtein, 2009) and the work of Efron (2011) further convincingly demonstrates that Tweedie’s formula provides an effective bias correction tool when estimating thousands of parameters simultaneously.

### 1.1 Issues with heterogeneous data

Most of the research in this area has been restricted to models where the distribution of $Y_i$ is solely a function of the parameter of interest $\mu_i$. In situations involving a nuisance parameter $\tau_i$, it
is generally assumed to be known and identical for all $Y_i$. For example, homoskedastic Gaussian models of the form $Y_i \mid \mu_i, \sigma_i \sim N(\mu_i, \sigma_i^2)$ involve a common nuisance parameter $\tau_i = 1/\sigma_i^2$ for all $i$. However, in large-scale studies when the data are collected from heterogeneous sources, the nuisance parameters may vary over the $n$ study units. Perhaps the most common example, and the setting we concentrate most on, involves heteroskedastic errors, where $\sigma_i^2$ varies over $Y_i$. Microarray data (Chiaretti et al., 2004, Erickson and Sabatti, 2005), returns on mutual funds (Brown et al., 1992), and the state-wide school performance gaps (Sun and McLain, 2012) are all instances of large-scale data where genes, funds, or schools have heterogeneous variances. Heteroskedastic errors also arise in analysis of variance (Weinstein et al., 2018) and linear regression settings (Kou and Yang, 2017). Moreover, in compound binomial problems, heterogeneity arises through unequal sample sizes across different study units. Unfortunately, the conventional Tweedie’s formula assumes identical nuisance parameters across study units and so cannot eliminate selection bias for heterogeneous data. Moreover, various works show that failing to account for heterogeneity leads to inefficient shrinkage estimators (Weinstein et al., 2018), methods with invalid false discovery rates (Cai and Sun, 2009, Efron, 2008), unstable multiple testing procedures (Tusher et al., 2001) and suboptimal ranking and selection algorithms (Henderson and Newton, 2016), exacerbating the replicability crisis in large-scale studies. Few methodologies are available to address this issue.

For Gaussian data, a common goal is to find the estimator of the means $\mu_i$ that minimizes the expected squared error loss. A plausible–seeming solution might be to scale each sample mean $Y_i$ by its estimated standard deviation $S_i/\sqrt{m_i}$ so that a homoskedastic method could be applied to $X_i = \sqrt{m_i}Y_i/S_i$, before undoing the scaling on the final estimate of $\mu_i$. Indeed, this is essentially the approach taken whenever we compute standardized test statistics, such as $t$–values and $z$–values. However, this approach, which disregards important structural information, can be highly inefficient. More advanced methods have been developed, but all suffer from various limitations. For instance, the methods proposed by Xie et al. (2012), Tan (2015), Jing et al. (2016a), Kou and Yang (2017), and Zhang and Bhattacharya (2017) are designed for heteroscedastic data but assume a parametric Gaussian prior or semi-parametric Gaussian mixture prior, which leads to loss of efficiency when the prior is misspecified. Moreover, existing methods, such as Xie et al. (2012) and Weinstein et al. (2018), often assume that the nuisance parameters, the variances $1/\tau_i$, are known and use a consistent estimator for implementation. However, when a large number of
units are investigated simultaneously, traditional sample variance estimators may similarly suffer from selection bias, which often leads to severe deterioration in the MSE for estimating the means.

1.2 The proposed approach and main contributions

In the homogeneous setting, Tweedie’s formula for Gaussian data estimates $\mu_i$ using the score function at $Y_i$, which is the gradient of the log marginal density of the sufficient statistic $Y_i$. However, this approach does not immediately extend to heterogeneous data. A significant challenge in the heterogeneous setting is how to pool information from different study units effectively while accounting for the heterogeneity captured by the possibly unknown nuisance parameter $\tau_i$. In this article, we address this issue by proposing a double shrinkage method that simultaneously incorporates the structural information encoded in both the primary (e.g. $Y_i$) and auxiliary (e.g. $S_i$) data. We develop a two–step approach, “Nonparametric Empirical Bayes Structural Tweedie” (NEST), which first estimates the bivariate score function at the primary and auxiliary data, and second predicts $\mu_i$ under a weighted squared error loss function using a generalized version of Tweedie’s formula that effectively incorporates the structural information encoded in the nuisance parameter. Unlike the squared error loss, which has been widely used for large-scale shrinkage estimation, the weighted squared error loss provides a more suitable performance metric for analyzing heterogeneous data by utilizing a weighting scheme that captures the underlying heterogeneity across the study units. The corresponding shrinkage estimator that minimizes the expected weighted squared error loss naturally encompasses the ability to differentiate between the study units both with respect to $\mu_i$ and the heterogeneity captured by the unknown nuisance parameter $\tau_i$.

NEST has several clear advantages. First, it simultaneously handles the two main selection biases introduced by heterogeneity: one, selection bias in the primary data (e.g. sample means), which cannot be effectively corrected without also correcting for, two, selection bias in the auxiliary data (e.g. sample variances). By producing more accurate estimates for the nuisance parameters, NEST in general renders improved shrinkage factors for estimating the primary parameters. Second, NEST makes no parametric assumptions about the prior since it uses a nonparametric method to directly estimate the bivariate score function. Third, NEST exploits the structure of
the entire sample and avoids the information loss that occurs in the discretization step used in grouping methods (Weinstein et al., 2018). Fourth, we establish that NEST is asymptotically as good as the optimal Bayes rule that uniquely minimizes the weighted squared error loss. Finally, NEST provides a general estimation framework for members of the two-parameter exponential family and is robust against model mis-specification. We demonstrate numerically that NEST can provide high levels of estimation accuracy relative to a host of benchmark methods.

### 1.3 Connection to existing works

For empirical Bayes (EB) estimation of the means, recent works such as Jing et al. (2016a), Kou and Yang (2017), Tan (2015), Xie et al. (2012) assume a parametric prior distribution for the means. This is in contrast to the setting where such a prior is unknown and in those settings there are two main modeling strategies for EB estimation. They are known as the $g$-modeling and $f$-modeling strategies in the terminology of Efron (2014). The idea of $g$-modeling is to first obtain a deconvolution estimate of the unknown prior distribution of $\mu_i$, and then predict $\mu_i$ by plugging this estimate into Bayes rule for various loss functions. The deconvolution estimate can be constructed via the nonparametric maximum likelihood estimate (NPMLE; Kiefer and Wolfowitz, 1956, Laird, 1978), or by modeling the unknown prior as a low-dimensional exponential family distribution (Efron, 2016). Some notable works along this line include Jiang and Zhang (2009), Koenker and Mizera (2014), Gu and Koenker (2017a), Saha and Guntuboyina (2020), and Soloff et al. (2021). In contrast, the $f$-modeling strategy usually assumes that the nuisance parameter, such as the variance, is known and sidesteps the need of deconvolution estimation by directly predicting $\mu_i$ based on Tweedie’s formula (or its generalized version), which only depends on the score function at the sufficient statistic $Y_i$ and the known nuisance parameter. Notable works along this line include Brown and Greenshtein (2009) and Efron (2011), both of which rely on fixed and known variances. In particular, their approach involves kernel density estimation techniques for separately estimating the marginal density and its gradient, and then taking their ratio to construct an estimate of the score function. NEST adopts the $f$-modeling strategy. However, in contrast to existing $f$-modeling methods, we allow the variances to be unknown and develop a convex optimization approach that directly provides consistent estimate of the score function and is capable of incorporating various structural constraints in the data-driven NEST estimator.
The \(g\)-modeling approach via the NPMLE provides an excellent tool for EB estimation of heteroskedastic means. However, to the best of our knowledge, the asymptotic properties of the NPMLE are highly nontrivial to establish and often require strong assumptions. For instance, the analysis in Saha and Guntuboyina (2020) only works for a limited class of covariance structures, and the theory on the rate of convergence is applicable only when the degree of heteroskedasticity is “mild”; alternatively the analysis in Soloff et al. (2021) assumes that \(\mu_i\) are independent of \(\sigma_i\), which is often violated in practice (Weinstein et al., 2018). In contrast, we establish the asymptotic properties of NEST without assumptions on the degree of heteroskedasticity or independence between \(\mu_i\) and \(\sigma_i\). A key advantage of \(g\)-modeling is its capability to deal with a wider range of problems, particularly those in which direct use of the marginal density of the sufficient statistic itself cannot yield a solution. Meanwhile, the \(f\)-modeling approach, which often has a simple and intuitive form (e.g. Tweedie’s formula), is attractive when only the information on the marginal distribution is needed for solving the problem of interest.

### 1.4 Organization

The rest of the paper is structured as follows. In Section 2 we present our hierarchical Gaussian model where both the mean and variance parameters are unknown. In Section 2.1 we discuss the weighted squared error loss function and in Section 2.2 we rely on the natural-parameter Tweedie’s formula for our hierarchical model to introduce the oracle NEST estimator under the weighted squared error loss in Definition 1. We then develop a convex optimization approach in Section 3 for estimating the unknown shrinkage factors in the oracle NEST formula. In Section 4, we describe the theoretical setup, justify the optimization criterion, and finally establish asymptotic theories for the proposed NEST estimator. Simulation studies are carried out in Section 5 to compare NEST to competing methods. The article concludes with a discussion in Section 6. Two data applications are presented in Section D of the Appendix. The proofs, as well as additional theoretical and numerical results, are provided in the EC.

### 2 Double shrinkage estimation on heteroskedastic Normal data

In the main text of this article, we focus on the normal means problem. Extensions of the methodology to other members of the two-parameter exponential family are discussed in Section E of
Suppose we collect \(m_i\) observations for the \(i^{th}\) study unit, \(i = 1, \ldots, n\). The data are normally distributed obeying the following hierarchical model:

\[
Y_{ij} \mid \mu_i, \tau_i \overset{i.i.d.}{\sim} N(\mu_i, 1/\tau_i), \quad j = 1, \ldots, m_i, \quad i = 1, \ldots, n.
\]

Let \(Y_i = m_i^{-1} \sum_{j=1}^{m_i} Y_{ij}\) and \(S_i^2 = (m_i - 1)^{-1} \sum_{j=1}^{m_i} (Y_{ij} - Y_i)^2\) respectively denote the sample mean and sample variance under Model (1). Further, let \(Y = (Y_1, \ldots, Y_n)\) and \(S = (S_1^2, \ldots, S_n^2)\) be the vectors of summary statistics, and \(y = (y_1, \ldots, y_n)\) and \(s = (s_1^2, \ldots, s_n^2)\) the observed values. We view \(\mu_i\) and \(\tau_i\), both of which are unknown, as the primary and nuisance parameters, respectively. The prior distributions \(G_{\mu}(\cdot | \tau_i)\) and \(H_{\tau}(\cdot)\) are unspecified. When the precisions \(\tau_i\) are known in Model (1), compound estimation of the means \(\mu = (\mu_1, \ldots, \mu_n)^T\) under the squared error loss has received significant attention in recent years (see for example Cai et al. (2021), Weinstein et al. (2018), Xie et al. (2012) and the references therein) and the estimator minimizing the expected squared error loss is,

\[
\mathbb{E}(\mu_i \mid Y_i = y_i, \tau_i, m_i) = y_i + \frac{1}{m_i \tau_i} w_1(y_i; m_i, \tau_i),
\]

where \(w_1(y; m, \tau) := \frac{\partial}{\partial y} \log f_{m,\tau}(y)\) and \(f_{m,\tau}(\cdot)\) is the pdf of the marginal distribution of \(Y\). Equation (2) is the celebrated Tweedie’s formula with known variances (Efron, 2011) which forms the basis for \(f\)-modeling strategies for estimating \(\mu_i\), and only requires estimation of the score functions \(w_1(y_i; m_i, \tau_i)\) in order to compute the estimator. This is particularly appealing in large-scale studies where one observes thousands of \((Y_i, \tau_i)\), making it possible to obtain an accurate estimate of \(w_1(y_i; m_i, \tau_i)\) (see Section E of the Appendix for more details). Here, we develop a nonparametric empirical Bayes method for estimating \(\mu\) under the weighted squared error loss (Section 2.1) when \(\tau_i\) are unknown.
Figure 1: A toy illustration of the squared error and weighted squared error losses. Here $n = 100$ and $\mu_i = 0$ for all $i = 1, \ldots, n$. The variances $\sigma_i^2 = 1/\tau_i$ are $n$ equispaced numbers between 0.5 and 2. In the right panel $\delta_i$ are $n$ equispaced numbers between −0.5 and 0.5. The red dots represent the squared error losses while the green dots represent the corresponding weighted squared error losses for estimating $\mu_i$ using $\delta_i$ for $i = 1, \ldots, n$.

2.1 The weighted squared error loss

Let $\delta = (\delta_1, \ldots, \delta_n)^T$ be an estimator for $\mu$ based on $(Y, S)$. Consider a class of loss functions

$$\ell^{(p)}(\mu_i, \delta_i; \tau_i) = \tau_i^p (\mu_i - \delta_i)^2,$$

(3)

for $p \in \{0, 1\}$, where $\ell^{(0)}(\mu_i, \delta_i; \tau_i)$ represents the usual squared error loss and $\ell^{(1)}(\mu_i, \delta_i; \tau_i) = \tau_i (\mu_i - \delta_i)^2$ is the weighted squared error loss with weight given by the precision $\tau_i$. While the squared error loss treats any two study units equally with respect to their precisions $\tau_i$, the weighted squared error loss $\ell^{(1)}(\mu_i, \delta_i; \tau_i)$ is a particularly natural loss function which effectively captures the underlying heterogeneity across the study units and appropriately adjusts their contributions to the total estimation error. Under Model (1), and unlike the squared error loss, $\ell^{(1)}(\mu_i, \delta_i; \tau_i)$ is also invariant under a linear transformation of the data $Y_{ij}$ to $\tilde{Y}_{ij}$ i.e. $\ell^{(1)}(\mu_i, \delta_i; \tau_i) = \ell^{(1)}(\bar{\mu}_i, \bar{\delta}_i; \bar{\tau}_i)$ where $\bar{Y}_{ij} = a_i + b_i Y_{ij}$, $\bar{\mu}_i = a_i + b_i \mu_i$, $\bar{\tau}_i = \tau_i / b_i^2$ and $\bar{\delta}_i = a_i + b_i \delta_i$ for $(a_i, b_i) \in \mathbb{R}^2$ with $b_i \neq 0$. Moreover, the weighting scheme alleviates potential concerns on the fairness issue in terms of the estimation accuracy that each study unit achieves since estimating $\mu_i$ as accurately when $\tau_i$ is small is difficult compared to when $\tau_i$ is large. Intuitively, the weighted loss down-weights the impacts of observations with extremely large variances and allows a relatively larger estimation error when the precisions are small while penalizing the estimation error more when the precisions are large. In Figure 1 we present a toy illustration of this
phenomenon. We fix \( n = 100 \) and \( \mu_i = 0 \) for all \( i = 1, \ldots, n \). The variances \( \sigma_i^2 = 1/\tau_i \) are \( n \) equispaced numbers between 0.5 and 2. The red dots in Figure 1 represent the squared error losses while the green dots represent the corresponding weighted squared error losses for estimating \( \mu_i \) using \( \delta_i \) for \( i = 1, \ldots, n \). In the left panel of Figure 1 we see that when the precisions are large, the same \( \delta_i \), say \( \delta_i = 0.5 \), incurs a relatively larger loss under \( \ell(1)(\mu_i, \delta_i; \tau_i) \) than when the precisions are small. In the right panel, we pair each \( \sigma_i^2 \) with a \( \delta_i \), where \( \delta_i \) are \( n \) equispaced numbers between \(-0.5\) and \(0.5\), and calculate the two losses. Both the left and the right panels reveal that \( \ell(0)(\mu_i, \delta_i; \tau_i) \) does not adjust the estimation error in response to the underlying heteroskedasticity represented by the precisions.

In the following section, we present the oracle NEST estimator of \( \mu \) under this weighted squared error loss. Thereafter in Section A of the Appendix we present an analysis under the usual squared error loss.

### 2.2 The oracle NEST estimator under the weighted squared error loss

Denote the average loss \( l_n^{(p)}(\mu, \delta; \tau) = n^{-1} \sum_{i=1}^{n} \ell^{(p)}(\mu_i, \delta_i; \tau_i) \). The compound Bayes risk is

\[
r_p(\delta, G) = \mathbb{E}\{l_n^{(p)}(\mu, \delta; \tau)\} = \frac{1}{n} \sum_{i=1}^{n} \int \int \ell^{(p)}(\mu_i, \delta_i; \tau_i)f_{m_i}(y, s^2 | \psi_i)dyds^2dG(\psi_i),
\]

where \( \psi_i = (\mu_i, \tau_i) \), \( G(\psi_i) = G_\mu(\mu_i | \tau_i)H_\tau(\tau_i) \), and \( f_{m_i}(y, s^2 | \psi_i) \) is the likelihood function of \((Y_i, S_i^2)\). The Bayes estimator that uniquely minimizes Equation (4) is given by \( \delta_{(p)}^\pi = (\delta_{1,(p)}, \ldots, \delta_{n,(p)}) \), where

\[
\delta_{i,(p)}^\pi := \delta_{(p)}^\pi(y_i, s_i^2, m_i) = \frac{\mathbb{E}(\tau_i^p \mu_i | y_i, s_i^2, m_i)}{\mathbb{E}(\tau_i^p | y_i, s_i^2, m_i)}.
\]

Denote \( \zeta_i := \tau_i \mu_i \). In Definition 1 we present the oracle NEST estimator of \( \mu \) under the weighted squared error loss which is the Bayes estimator \( \delta_{(1)}^\pi \). The key idea is to exploit the exponential family representation of the posterior distribution of \((\mu_i, \tau_i)\) to construct a Tweedie-type formulae for \( \delta_{i,(1)}^\pi \).

**Definition 1.** Consider hierarchical Model (1) with \( m_i > 3 \). Then the oracle NEST estimator of
The natural parameter Tweedie’s formulae for the parameters \( (\zeta_i, \tau_i) \) under Model (1) and are significant for several reasons. First, they generalize Tweedie’s formula with known variances (Equation (2)), which can be recovered from Equation (7) by treating \( \tau_i \) as a known constant and dividing both sides by \( \tau_i \). Second, \( \hat{\tau}^\pi_i \) in Equation (8) has an interesting interpretation. Apart from being the Bayes estimate of \( \tau_i \) under the squared error loss, \( 1/\hat{\tau}^\pi_i \) is the Bayes estimate of \( 1/\tau_i \) under Stein’s loss (James and Stein, 1961). Third, Equations (7) and (8) can be used for compound estimation of the parameters \( \theta_i = T(\zeta_i, \tau_i, m_i) \) where \( T(\cdot) \) is a known function of \( (\zeta_i, \tau_i, m_i) \). For instance, in finance applications the Sharpe ratio for a portfolio \( i \) is given by \( T(\zeta_i, \tau_i, m_i) = \zeta_i / \sqrt{\tau_i} \) and one can construct an estimator \( \hat{\theta}_i := T(\hat{\zeta}^\pi_i, \hat{\tau}^\pi_i, m_i) \) using Equations (7) and (8) to estimate \( \theta_i \). In Section D.2 of the Appendix we present a real data
application that is dedicated to the compound estimation of Sharpe ratios for a large number of mutual fund portfolios. However we note that \( \hat{\theta}_i \) constructed in the aforementioned fashion may not, in general, be the optimal estimator of \( \theta_i \) under the squared error or the weighted squared error loss considered here. In Section B.7 of the Appendix we provide a proof of Equations (7) and (8).

3 The data-driven NEST estimator

In this section we discuss the estimation of the shrinkage factors and introduce the data-driven NEST estimator in Definition 2. We begin by introducing some notation. Let \( x = (y, s^2) \) be a generic pair of observations from the distribution with marginal density \( f_m(x) \), which we assume is continuously differentiable with support on \( \mathcal{X} \subseteq \mathbb{R} \times \mathbb{R}^+ \). Denote the score function

\[
\mathbf{w}(x; m) = \nabla_x \log f_m(x) := \{w_1(x; m), w_2(x; m)\}. 
\]

(9)

Next, for \( i = 1, \ldots, n \), let \( z^i = (x^i, m_i) \) where \( x^i \) is an observation from a distribution with density \( f_{m_i}(x) \).

3.1 Convex optimization

We first describe the methodology and then provide explanations. Let,

\[
K_\lambda(z^i, z^j) = \exp \left\{ -\frac{1}{2\lambda^2} (z^i - z^j)^T \Omega (z^i - z^j) \right\},
\]

(10)

be the Radial Basis Function (RBF) kernel with bandwidth parameter \( \lambda > 0 \) and \( \Omega_{3\times3} \) being the inverse of the sample covariance matrix of \( (z^1, \ldots, z^n) \). Denote

\[
\mathbf{W}_0^{n\times2} = \left\{ \mathbf{w}(x^1; m_1), \ldots, \mathbf{w}(x^n; m_n) \right\}^T,
\]

where

\[
\mathbf{w}(x^i; m_i) = \nabla_x \log f_{m_i}(x) \bigg|_{x=x^i} := \{w_1(x^i; m_i), w_2(x^i; m_i)\}.
\]

(11)

We denote \( w_k(x^i; m_i) \) by \( w_{k,i} \), \( k = 1, 2 \), for the remainder of this article.

Let \( \nabla_{z^i} K_\lambda(z^i, z^j) \) be the partial derivative of \( K_\lambda(z^i, z^j) \) with respect to the \( k^{th} \) component
of \( z^j \). The following matrices are needed in our proposed estimator:

\[
K_{\lambda}^{n \times n} = [K_{ij,\lambda}]_{1 \leq i \leq n, 1 \leq j \leq n}, \quad \nabla K_{\lambda}^{n \times 2} = [\nabla K_{ik,\lambda}]_{1 \leq i \leq n, 1 \leq k \leq 2},
\]

where \( K_{ij,\lambda} = K_{\lambda}(z_i, z_j) \) and \( \nabla K_{ik,\lambda} = \sum_{j=1}^{n} \nabla z^j K_{\lambda}(z^i, z^j) \). Next we formally define our proposed Nonparametric Empirical-Bayes Structural Tweedie (NEST) estimator.

**Definition 2.** Consider hierarchical Model (1) with \( m_i > 3 \). For a fixed bandwidth parameter \( \lambda > 0 \), let \( \hat{W}_n(\lambda) = \{ \hat{w}_{\lambda,n}(1), \ldots, \hat{w}_{\lambda,n}(n) \}^T \), where \( \hat{w}_{\lambda,n}(i) = \{ \hat{w}^{(1)}_{\lambda,n}(i), \hat{w}^{(2)}_{\lambda,n}(i) \}^T \), be the solution to the following quadratic optimization problem:

\[
\min_{W \in \mathbb{R}^{n \times 2}} \frac{1}{n^2} \text{trace} \left( W^T K_{\lambda} W + 2 W^T \nabla K_{\lambda} \right).
\]  

Then the NEST estimator for \( \mu_i \) is

\[
\delta_{i,n}^{ds}(\lambda) = y_i + \frac{s^2_i}{m_i} \hat{\gamma}_{i,n}(\lambda) \hat{w}^{(1)}_{\lambda,n}(i),
\]

where \( \hat{\gamma}_{i,n}(\lambda) = \frac{m_i - 1}{m_i - 3 - 2s^2_i \hat{w}^{(2)}_{\lambda,n}(i)} \).  

with the superscript \( \text{ds} \) denoting “double shrinkage”.

Although not immediately obvious, we show in Section 4 that, under the compound estimation setting, minimizing the objective function (12) is asymptotically equivalent to minimizing the kernelized Stein’s discrepancy (KSD; Chwialkowski et al., 2016, Liu et al., 2016). Roughly speaking, the KSD measures how far a given \( n \times 2 \) matrix \( W \) is from the true score matrix \( W_0 \). A key property of the KSD is that it is always non-negative and is equal to 0 if and only if \( W \) and \( W_0 \) are equal. Hence, solving the convex program (12) is equivalent to finding a \( \hat{W} \) that is as close as possible to \( W_0 \). Since the oracle NEST estimator in Definition 1 is constructed based on \( W_0 \), we can expect that the data-driven NEST estimator based on \( \hat{W}_n(\lambda) \) would be asymptotically close to its oracle counterpart. Theory underpinning this intuition are established in Section 4.

The bandwidth parameter \( \lambda \) in the RBF kernel \( K_{\lambda} \) (Equation (10)) controls the bias-variance trade-off in the score function estimate. For instance, a small value of \( \lambda \) allows unbiasedness but the resulting \( n \)-dimensional score function estimator has more variance compared to when \( \lambda \) is large, which forces the estimated shrinkage factors towards 0, and in the limit the NEST estimate
is simply the unbiased estimate $y_i$. In Section 3.2 we discuss a data-driven approach to choose $\lambda$ for practical implementation of the NEST estimator.

A key characteristic of $\hat{\delta}_{ds,i,n}(\lambda)$ in Equation (13) is that it exploits the joint structural information available in both $Y_i$ and $S_i^2$ through $\hat{W}_n(\lambda)$. Since the weighted loss function involves both the mean and the variance, we perform shrinkage on both these dimensions. Inspecting Equations (13) and (14), we expect that the improved accuracy achieved by $\hat{\hat{\gamma}}_{i,n}(\lambda)$ will lead to better shrinkage factors for $\hat{\delta}_{ds,i,n}(\lambda)$ and hence additional reduction in the estimation risk. Our numerical results in Sections 5 and D reveal that this is indeed true and the proposed NEST estimator dominates other shrinkage estimators across many settings. In Section E of the Appendix we adopt a similar strategy to extend the estimation framework presented in Definition 2 to other distributions in the two-parameter exponential family where the nuisance parameter is known.

We end this section with a discussion of the simpler case of equal sample sizes, i.e. $m_i = m$ for all $i$. For instance, the leukemia dataset analyzed in Jing et al. (2016a) consists of the expression levels of $n = 5,000$ genes for $m = 27$ acute lymphoblastic leukemia patients. The heterogeneity across the $n$ units in this case is due to the intrinsic variability instead of the varied number of replicates. When the $m_i$’s are equal, the RBF kernel $K_\lambda(\cdot, \cdot)$ needs to be modified to avoid a singular sample covariance matrix. Denote $K_\lambda(x^i, x^j) = \exp\{-0.5\lambda^{-2}(x^i - x^j)^T\Omega(x^i - x^j)\}$ the modified RBF kernel with $\Omega^{2\times2}$ being the inverse of the sample covariance matrix of $(x^1, \ldots, x^n)$. Correspondingly in Definition 2, $\hat{W}_n(\lambda)$ are the estimates of the shrinkage factors $W_{0}^{n\times2} = \{w(x^1; m), \ldots, w(x^n; m)\}^T$, where $w(x^i; m) = \{w_1(x^i; m), w_2(x^i; m)\}^T := (w_{1,i}, w_{2,i})^T$.

### 3.2 Details around implementation

In this section we discuss details around the implementation of NEST. First note that Equation (12) can be solved separately for the two columns of $W$, which respectively yield the estimates for $w_{1,i}$ and $w_{2,i}$. Next, the solution to Equation (12) is available in the closed form of $\hat{W}_n(\lambda) = -K_\lambda^{-1} \nabla K_\lambda$. However, in our implementation the closed form solution is replaced by a convex program that directly solves (12) with the following two constraints: (1) $W\mathbf{a} \leq \mathbf{b}$, where $\mathbf{a} = (0, 1)^T$ and $\mathbf{b} = (b_1, \ldots, b_n)$ with $b_i = \frac{1}{2}(m_i - 3)/s_i^2 - \kappa$ for some $\kappa > 0$, and (2) $1^T W\mathbf{e}_k = 0$, $k = 1, 2$, where $\mathbf{e}_k$ is the canonical basis vector with 1 at coordinate $k$. Inspecting Equation (14) shows
that adding the first constraint guarantees that $\hat{\gamma}_{i,n}(\lambda) < \infty$. This is desirable in both numerical and theoretical analyses. Similar ideas have been used in the seminal work of Koenker and Mizera (2014). The second constraint ensures that the sample mean of the $k^{th}$ estimated shrinkage factor is zero since the expectation of the gradient of the log marginal density is indeed zero.

The practical implementation requires a data-driven scheme for choosing $\lambda$. We propose to use a variation of the modified cross validation scheme of Brown et al. (2013), which involves splitting $Y_{ij}$ into two parts: $U_{ij} = Y_{ij} - (1/\alpha)e_{ij}$ and $V_{ij} = Y_{ij} + \alpha e_{ij}$, where $e_{ij} \sim N(0, \tau_i^{-1})$, and $U_{ij}$ and $V_{ij}$ are used to construct the estimator and to choose the tuning parameter, respectively. However in our setup $\tau_i$ is unknown; hence we define $U_{ij} = Y_{ij} - (1/\alpha)\sqrt{S_i} e_{ij}$, $V_{ij} = Y_{ij} + \alpha \sqrt{S_i} e_{ij}$ and sample $e_{ij}$ independently from $N(0, 1)$. Let $\bar{V}_i = m_i^{-1} \sum_{j=1}^{m_i} V_{ij}$, $\bar{U}_i = m_i^{-1} \sum_{j=1}^{m_i} U_{ij}$, $U = \{U_{ij} : 1 \leq i \leq n, 1 \leq j \leq m_i\}$ and $V = \{V_{ij} : 1 \leq i \leq n, 1 \leq j \leq m_i\}$. Then conditional on $(\mu_i, \tau_i)$, $\bar{U}_i$ and $\bar{V}_i$ are uncorrelated with mean $\mu_i$ and variances $(1 + \alpha^2)/(m_i \tau_i)$ and $(1 + \alpha^2)/(m_i \tau_i)$, respectively. Define

$$\bar{\vartheta}_n(\lambda; U, V) = \frac{1}{n} \sum_{i=1}^{n} \frac{\bar{V}_i - \hat{\delta}^{ds}_{i,n}(\bar{U}_i; U, \lambda)}{S_{V,i}^2},$$

where $\hat{\delta}^{ds}_{i,n}(\bar{U}_i; U, \lambda)$ is the NEST estimator of $\mu_i$ based on $U$ and $S_{V,i}^2$ is the sample variance of $\{V_{ij} : 1 \leq j \leq m_i\}$. The tuning parameter will be chosen as $\hat{\lambda} := \arg\min_{\lambda \in \Lambda} \bar{\vartheta}_n(\lambda)$. In our numerical studies of Section 5, we set $\alpha = 1/2$, $\Lambda = [0.5, 10^2]$. The tuning parameter $\hat{\lambda}$ is obtained from this scheme and then used to estimate $\mu_i$ via Equation (13).

**Remark 1.** Note that for large $m$, $\bar{V}_i$ and $S_{V,i}^2$ are independent and $E[\bar{\vartheta}_n(\lambda; U, V)]$ can be well approximated by the sum $n^{-1} \sum_{i=1}^{n} E[\hat{\mu}_i, \hat{\delta}^{ds}_{i,n}(\bar{U}_i; U, \lambda); \tau_i/(1 + \alpha^2)] + n^{-1} \sum_{i=1}^{n} E(\bar{V}_i - \mu_i)^2/S_{V,i}^2$ for any fixed $\lambda > 0$ and $n$. The first term depends on $\lambda$ while the second term is independent of it. We choose the value of $\lambda$ that minimizes $\bar{\vartheta}_n(\lambda; U, V)$ to construct the data-driven NEST estimator of $\mu$ based on the original sample $(Y, S)$. For $\alpha \to 0$, this scheme is equivalent to choosing a value of $\lambda$ that minimizes the true weighted squared error loss of $\hat{\delta}^{ds}_{i,n}(\bar{U}_i; U, \lambda)$, $i = 1, \ldots, n$. We note in Section 5 that when $m$ is small, the risk of the data-driven NEST estimator is generally higher than the risk of the NEST estimator that chooses $\lambda$ by minimizing the true weighted squared error loss of $\hat{\delta}^{ds}_{n}(\lambda)$ based on $(Y, S)$ (see for example the left panels in Figures 2 and 3). This is not unexpected since $\bar{\vartheta}_n(\lambda; U, V)$ is not an unbiased estimator of the true risk. However, $\bar{\vartheta}_n(\lambda; U, V)$ provides a practical criterion for choosing $\lambda$ and
works well empirically in our numerical studies for larger \( m \). The development of an unbiased risk estimate, such as a SURE-type criterion, for this setting is a challenging topic requiring further research. See also Ignatiadis and Wager (2019) for related discussions.

We are developing an \( R \) package, \texttt{nest}, to implement the NEST estimator in Definition 2. The \( R \) code that reproduces the numerical results in this paper can be downloaded from the link: https://www.dropbox.com/sh/5yptcj4epxdqbgq/AADAyHZNDCv4Hqagsi97vC2Da?dl=0.

4 Theory

In this section we introduce the Kernelized Stein’s Discrepancy (KSD) measure (Chwialkowski et al., 2016, Liu et al., 2016) and discuss its connection to the quadratic program (12). While the KSD has been used in various contexts including goodness of fit tests (Liu et al., 2016, Yang et al., 2018), variational inference (Liu and Wang, 2016) and Monte Carlo integration (Oates et al., 2017), its connections to compound estimation and empirical Bayes methodology was established only recently (Banerjee et al., 2021, Luo et al., 2023). The analysis in this and following sections is geared towards the case \( m_i = m \) for \( i = 1, \ldots, n \). Under this setting, \((x^1, \ldots, x^n)\) constitute an i.i.d random sample from \( f_m(x) \). The case of unequal \( m_i \)'s can be analyzed in a similar fashion by assuming that \( m_i \)'s are a random sample from a distribution with mass function \( q(\cdot) \). Then \( z = (x, m) \) has distribution with density \( p(z) := q(m)f_m(x) \), where \( z^i = (x^i, m_i), (z^1, \ldots, z^n) \) are realizations of an i.i.d random sample from \( p(z) \).

4.1 Kernelized Stein’s Discrepancy

Suppose \( X \) and \( X' \) are i.i.d. copies from the marginal distribution of \((Y, S^2)\) that has density \( f \) wherein the dependence on \( m \) is implicit. Denote \( w(X) \) and \( w(X') \), defined in Equation (9), to be the score functions at \( X \) and \( X' \) respectively. Suppose \( \tilde{f} \) is an arbitrary density function on the support of \((Y, S^2)\), for which we similarly define \( \tilde{w}(X) \). The KSD, formally defined as

\[
S_{\lambda}(f, \tilde{f}) = \mathbb{E}_{X, X' \sim f}
\left[
\left\{\tilde{w}(X) - w(X)\right\}^T K_{\lambda}(X, X') \left\{\tilde{w}(X') - w(X')\right\}
\right],
\]

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provides a discrepancy measure between $f$ and $\tilde{f}$ in the sense that $S_\lambda(f, \tilde{f})$ tends to increase when there is a bigger disparity between $w$ and $\tilde{w}$ (or equivalently, between $f$ and $\tilde{f}$), and

$$S_\lambda(f, \tilde{f}) \geq 0 \text{ and } S_\lambda(f, \tilde{f}) = 0 \text{ if and only if } f = \tilde{f}.$$ 

The direct evaluation of $S_\lambda(f, \tilde{f})$ is difficult because $w$ is unknown. Liu et al. (2016) introduced an alternative representation of the KSD that does not directly involve $w$:

$$S_\lambda(f, \tilde{f}) = E_f \kappa_\lambda[\tilde{w}(X), \tilde{w}(X')](X, X')$$

$$= E_f \left\{ \frac{1}{n(n-1)} \sum_{i=1}^{n} \sum_{j=1}^{n} \kappa_\lambda[\tilde{w}(X^i), \tilde{w}(X^j)](X^i, X^j) \mathbb{I}(i \neq j) \right\}$$

$$= E_f [\bar{M}_{\lambda,n}(\tilde{W})] := M_\lambda(\tilde{W}),$$

(15)

where $\{X^1, \ldots, X^n\}$ is a random sample from $f$, $E_f$ denotes expectation under $f$ and

$$\kappa_\lambda[\tilde{w}(x), \tilde{w}(x')](x, x') = \tilde{w}(x)^T \tilde{w}(x') \kappa(x, x') + \tilde{w}(x)^T \nabla_x \kappa(x, x') + \nabla_x \kappa(x, x')^T \tilde{w}(x) + \text{trace}(\nabla_x \nabla_x \kappa(x, x')),$$

is a smooth and symmetric positive definite kernel function associated with the U-statistic $\bar{M}_{\lambda,n}(\tilde{W})$. The implementation of the NEST estimator in Definition 2 boils down to the estimation of $W_0$ via the convex program (12), which corresponds to minimizing

$$\tilde{M}_{\lambda,n}(\tilde{W}) = \frac{1}{n^2} \sum_{i=1}^{n} \sum_{j=1}^{n} \kappa_\lambda[\tilde{w}(X^i), \tilde{w}(X^j)](X^i, X^j)$$

(16)

w.r.t. $\tilde{W}$. A key observation is that if the empirical criterion $\tilde{M}_{\lambda,n}(\tilde{W})$ is asymptotically equal to the population KSD criterion $M_\lambda(\tilde{W})$, then minimizing $\tilde{M}_{\lambda,n}(\tilde{W})$ with respect to $\tilde{W}$ is effectively the process of finding a $\tilde{W}$ that is as close as possible to $W_0$ in Equation (11). This intuitively justifies the NEST estimator in Definition 2.

In Theorem 1, we formally establish the asymptotic consistency of the sample criterion $\tilde{M}_{\lambda,n}(\tilde{W})$ around the population criterion $M_\lambda(\tilde{W})$ for any fixed $\lambda > 0$. We impose the following regularity condition where $f$ denotes the density function of the joint marginal distribution of $(Y, S^2)$ and $E_f$ denotes expectation under $f$. 16
Assumption 1. \( \mathbb{E} f | \kappa_\lambda [\tilde{w}(X_i), \tilde{w}(X_j)](X_i, X_j) |^2 < \infty \) for any \( (i, j) \in \{1, \ldots, n\} \).

Assumption 1 is a standard moment condition on \( \kappa_\lambda [\tilde{w}(X_i), \tilde{w}(X_j)](X_i, X_j) \) [see, for example, Section 5.5 in Serfling (2009)], which is needed for establishing the Central Limit Theorem for the U-statistic \( \tilde{M}_{\lambda,n}(\tilde{W}) \) in Equation (15).

**Theorem 1.** Under Assumption 1 and for any fixed \( \lambda > 0 \), we have

\[
\left| \tilde{M}_{\lambda,n}(\tilde{W}) - M_\lambda(\tilde{W}) \right| = O_p(n^{-1/2}).
\]

Moreover, along with the fact that \( M_\lambda(W_0) = 0 \), Theorem 1 justifies \( \tilde{M}_{\lambda,n}(\tilde{W}) \) as an appropriate optimization criterion. Next we show that the NEST estimator in Definition 2 is asymptotically close to its oracle counterpart.

### 4.2 Asymptotic Properties of NEST

This section studies the asymptotic properties of the NEST estimator under the setting where \( \lambda := \lambda(n) \) varies with \( n \). We begin by recalling the oracle NEST estimator \( \delta_{\pi,1} = (\delta_{\pi,1(1)}, \ldots, \delta_{\pi,1(n)}) \) for \( \mu \) in Equation (6), where

\[
\delta_{\pi,1(i)} := \delta_{\pi}(y_i, s_i^2, m) = \frac{\tilde{\zeta}(y_i, s_i^2, m)}{\tilde{\tau}(y_i, s_i^2, m)} = y_i + \frac{s_i^2}{m} \gamma(y_i, s_i^2, m_i)w_1(y_i, s_i^2, m),
\]

and \( \tilde{\zeta} \) and \( \tilde{\tau} \) are respectively the Bayes estimators of \( \tau \mu \) and \( \tau \) as defined in Equations (7) and (8). Viewing the proposed NEST estimator \( \delta_{\pi,n}(\lambda) \) as a data-driven approximation to \( \delta_{\pi,1(1)} \), we study the quality of this approximation for large \( n \) and fixed \( m \).

In Theorem 2 we first establish the consistency of the estimated score functions \( \tilde{W}_n(\lambda) \) obtained as the solution to the quadratic optimization problem in Equation (12). To establish such consistency, we rely on the tools developed in Luo et al. (2023) for analyzing the KSD. In what follows, we denote the Frobenius norm of a matrix \( \mathbf{A} \) by \( \|\mathbf{A}\|_F \) and for two sequences \( a_n \) and \( b_n \), we denote \( a_n \asymp b_n \) to mean \( d_1 a_n \leq b_n \leq d_2 a_n \) for large \( n \) and some constants \( d_2 \geq d_1 > 0 \). The following regularity condition is needed in our technical analysis.

**Assumption 2.** The joint density \( f \) and its score functions are Lipschitz continuous.
**Theorem 2.** If $\lim_{n \to \infty} c_n (\log n)^4 n^{-1/3} = 0$ then under Assumption 2 and with $\lambda \asymp n^{-1/3}$, we have

$$
\lim_{n \to \infty} \mathbb{P} \left\{ \frac{1}{n} \| \hat{W}_n(\lambda) - W_0 \|_{\mathcal{F}}^2 \geq c_n^{-1} \epsilon \right\} = 0, \text{ for any } \epsilon > 0.
$$

We prove Theorem 2 for sub-exponential random variables $X$ whose density $f$ and the corresponding score functions are Lipschitz continuous (Assumption 2). The assumption of Lipschitz continuity is mild and allows a cleaner statement of our theoretical result. Theorem 2 reveals a slower than parametric $\sqrt{n}$ rate of convergence, barring the logarithmic terms. This is related to the observation, originally made by Luo et al. (2023), that the isometry between the reproducing kernel Hilbert space (RKHS) norm and the $L_2(\tilde{W}, W_0) = n^{-1} \| \tilde{W} - W_0 \|_{\mathcal{F}}^2$ metric that appears in statement of Theorem 2, may not exist. Specifically, the KSD approach for estimating the score functions essentially works by first mapping the data into an RKHS associated with the integrally strictly positive definite kernel $K_{\lambda}$ in Equation (10) and then estimating the scores by minimizing $\hat{M}_{\lambda,n}(\hat{W})$ with respect to $\tilde{W}$, where $\hat{M}_{\lambda,n}(\hat{W})$ is a data-driven approximation to the squared RKHS norm $M_{\lambda}(\hat{W})$. The statement of Theorem 2, however, is based on the $L_2(\tilde{W}, W_0)$ metric and not the RKHS norm. As such, it is not necessarily true that if $\hat{W}_n(\lambda)$ is a good estimate of $W_0$ in the RKHS norm then it is a good estimate of $W_0$ in the $L_2$ metric too for every $\lambda > 0$. Theorem 2 precisely tabulates how $\lambda$ must behave as $n \to \infty$ to establish consistency of the estimated scores in the $L_2$ metric and the slower than $\sqrt{n}$ rate of convergence is a consequence of the cost of inversion from the RKHS norm to this $L_2$ metric.

Next we provide decision theoretic guarantees on $\delta_{ds}(\lambda)$ in relation to $\delta^{\pi}_{(1)}$. We impose the following moment condition on the prior distributions of $\mu$ and $\tau$.

**Assumption 3.** For some $\epsilon_i \in (0, 1), i = 1, 2, 3, \mathbb{E}_G\{\exp(\epsilon_1|\mu|)\} < \infty, \mathbb{E}_H\{\exp(\epsilon_2/\tau)\} < \infty$ and $\mathbb{E}_H\{\exp(\epsilon_3\tau)\} < \infty$.

The moment conditions in Assumption 3 ensure that with high probability $|\mu| \leq \log n$ and $1/\log n \leq \tau \leq \log n$ as $n \to \infty$. This is formalized in Lemma 4 in Section B.6 of the EC. It is likely that Assumption 3 can be further relaxed but we do not seek the full generality here.

**Theorem 3.** Suppose Assumptions 2 – 3 hold. If $\lim_{n \to \infty} c_n n^{-1/3} (\log n)^8 = 0$, then as $n \to \infty$ with $\lambda \asymp n^{-1/3}$, we have $\frac{c_n}{n} \| \delta_{ds}(\lambda) - \delta^{\pi}_{(1)} \|_2 = o_p(1)$. Furthermore, under the same conditions, $c_n^{1/2} \left| \eta_n(1) \{\mu, \delta_{ds}(\lambda); \tau\} - \eta_n(1) \{\mu, \delta^{\pi}_{(1)}; \tau\} \right| = o_p(1)$ as $n \to \infty$. 

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Theorem 3 establishes the optimality theory of $\delta_n^{ds}(\lambda)$ by showing that (a) the average squared error between $\delta_n^{ds}(\lambda)$ and $\delta_n^{\pi}(1)$ is asymptotically small, and (b) the estimation loss of NEST converges in probability to that of its oracle counterpart as $n \to \infty$.

5 Numerical experiments

In this section we assess the performance of the NEST estimation framework for compound estimation of Normal means under the weighted squared error loss. We focus on the hierarchical Model of Equation (1) and compare the following four approaches for estimating $\mu$ when the precisions $\tau$ are unknown: (1) the proposed NEST method where the tuning parameter $\lambda$ is chosen using a modified cross-validation approach described in Section 3.2, (2) the NEST method which estimates $\lambda$ by minimizing the true loss (NEST Orc. $\lambda$), (3) the g-modelling approach of Gu and Koenker (2017a,b) which first estimates the joint prior distribution of $(\mu_i, \tau_i)$ using nonparametric maximum likelihood estimation (NPMLE) techniques (Kiefer and Wolfowitz, 1956, Laird, 1978) and then plugs them into the Bayes rule for estimating $\mu_i$ under the weighted squared error loss, and (4) the oracle Bayes estimator $\delta_n^{\pi}(1)$ in Equation (6) that has full knowledge of the prior distributions of $(\mu_i, \tau_i)$. For NPMLE we use the R-function WGLVmix, available within the R-package REBayes (Koenker and Gu, 2017), to estimate the joint prior distribution of $(\mu_i, \sigma^2_i)$. To compute $\delta_n^{\pi}(1)$ we rely on the R-package RStan (Carpenter et al., 2017, Stan Development Team, 2022).

The aforementioned four approaches are evaluated on nine different simulation settings, with the goal of assessing the relative performance of the competing estimators as the heterogeneity in the variances $\sigma^2_i$ is varied while keeping the sample sizes $m_i$ fixed at $m$. The nine simulation settings can be categorized into three types: two settings where mean and variances are independent, five settings where mean and variance are correlated and two settings that represent departures from the Normal data-generating model. For each setting we set $n = 1,000$ and compute the average Bayes risk for each competing estimator of $\mu$ across 50 Monte Carlo repetitions. Figures 2 to 5 plot the relative risk of each competing estimator, which is the ratio of the average Bayes risk of the competing estimator to that of the oracle Bayes estimator $\delta_n^{\pi}(1)$.

The first two settings, Figures 2a and 2b, correspond to the independent case. We describe the two settings below.
Setting 1: $\mu_i \overset{i.i.d.}{\sim} N(0, 3)$ and $\sigma_i^2 \overset{i.i.d.}{\sim} U(0.25, u)$ where we let $u$ to vary across five levels, \{1, 2, 3, 4, 5\}.

Setting 2: $\mu_i \overset{i.i.d.}{\sim}$ Laplace(location = 0, scale = 1) and $\sigma_i^2 \overset{i.i.d.}{\sim} N(u, 1)$ truncated below 0.1 where $u$ varies across \{0.25, 0.5, 0.75, 1, 1.25\}.

The three plots in Figures 2a and 2b show the relative risks as $u$ varies for $m = 10, 15$ and $20$ (left to right). We see that under Setting 1, NEST and NEST Orc. $\lambda$ have relative risks substantially closer to 1 than NPMLE. As $m$ increases, the performance of NEST and NEST Orc. $\lambda$ improves as their relative risks almost coincide with 1. The NPMLE, on the other hand, exhibits a slightly higher relative risk at $m = 20$ than at $m = 10$ and 15.

The next five settings, Figures 3 and 4, correspond to the correlated case. We describe these five settings below.

Setting 3: $\sigma_i^2 \overset{i.i.d.}{\sim} U(0.25, 3)$ and $\mu_i \mid \sigma_i^2 \overset{ind.}{\sim} 0.5 \ N(-\sigma_i^2, u^2) + 0.5 \ N(\sigma_i^2, u^2)$ where $u$
varies across \{0.25, 0.35, 0.45, 0.55, 0.65, 0.75\}.

- Setting 4: \( \sigma_i^2 \sim U(0.25, u) \) and \( \mu_i | \sigma_i^2 \overset{\text{i.i.d}}{\sim} 0.5 N(\sigma_i^2, \sigma_i^2) + 0.5 N(-\sigma_i^2, \sigma_i^2) \) where we let \( u \in \{1, 2, 3, 4, 5\} \).

- Setting 5: we sample \( \sigma_i^2 \) from an Inverse Gamma (IG) distribution with shape parameter fixed at 5 and rate at \( u \) where \( u \in \{1, 1.5, 2, 2.5, 3\} \). Conditional on the variances, the means \( \mu_i \) are sampled from the design described in Setting 4.

- Setting 6: \( \sigma_i^2 \sim U(0.1, u) \) and \( \mu_i | \sigma_i^2 \overset{\text{i.i.d}}{\sim} \text{Laplace(location = 0, scale = } \sqrt{\sigma_i^2/2}) \) where \( u \in \{1, 1.5, 2, 2.5, 3\} \).

- Setting 7: \( \sigma_i^2 \) are independently sampled from a two component mixture of truncated Normal distributions that are truncated below 0.1 with means 0.5 and \( u \), standard deviation 0.1 and mixing weights 0.5. Conditional on the variances, the means \( \mu_i \) are sampled from a Logistic distribution with location \( \sigma_i^2 \) and scale \( \sqrt{\sigma_i^2} \). We let \( u \) to vary across five levels, \( \{1, 1.5, 2, 2.5, 3\} \).

Setting 3 is presented in Figure 3a where \( u \) controls the spread of the prior distribution of \( \mu_i \) given \( \sigma_i^2 \) and reveals that when \( u \) is small, NPMLE has a relatively better risk performance than NEST. This is expected because when \( u \) is small, the prior distribution of \( \mu_i \) can be well approximated by a discrete distribution with just two mass points and the NEST estimator finds it particularly difficult to estimate the means well in these scenarios because such a discrete prior on \( \mu_i \) introduces potential multimodality of the underlying marginal distribution of \((Y_i, S_i^2)\) in which case the true score functions may no longer be Lipschitz continuous, an assumption that is needed for the KSD approach to produce consistent estimates of the true scores in Section 3. The NPMLE, on the other hand, has a near parametric rate of convergence in Gaussian denoising problems when the underlying prior on the means is discrete (Saha and Guntuboyina, 2020), which explains its improved performance in Setting 3, especially when \( u \to 0 \).

Settings 4 to 7, Figures 3b, 3c and 4, represent scenarios where NEST dominates NPMLE uniformly for all values of \( u \) and \( m \). The relative risk of NPMLE in these scenarios is sometimes as high as 1.6 while NEST and NEST Oracle \( \lambda \) exhibit relative risk profiles substantially closer to 1.
Figure 3: Comparison of relative risks when \((\mu_i, \sigma_i^2)\) are dependent - Settings 3, 4 and 5. Plots show \(m = 10, 15, 20\) left to right.

The eight and ninth settings, Figures 5a and 5b, correspond to the scenarios where the data \(Y_{ij}|(\mu_i, \sigma_i^2)\) are not normally distributed. In Figure 5a, \((\mu_i, \sigma_i^2)\) are sampled according to Setting 5 but \(Y_{ij}|(\mu_i, \sigma_i^2)\) are generated independently from a Laplace distribution with location \(\mu_i\) and scale \(\sqrt{\sigma_i^2/2}\). Similarly, in Figure 5b, \((\mu_i, \sigma_i^2)\) are sampled according to Setting 4 but \(Y_{ij}|(\mu_i, \sigma_i^2) = \mu_i + \xi_i\) where \(\xi_i\) are independently sampled from a central \(t\)–distribution with 4 degrees of
(a) Setting 6: $\sigma_i^2 \sim i.i.d. U(0.1, u)$ and $\mu_i | \sigma_i^2 \sim \text{Laplace}(\text{location} = 0, \text{scale} = \sqrt{\sigma_i^2/2})$

(b) Setting 7: $\sigma_i^2 \sim i.i.d. 0.5 \text{Trunc}.N(0.5, 0.1^2; 0.1) + 0.5 \text{Trunc}.N(u, 0.1^2; 0.1)$ where $\text{Trunc}.N(a, b; c)$ represents a truncated Normal distribution with mean $a$, variance $b$ and truncated below $c$. Here $\mu_i | \sigma_i^2 \sim \text{Logistic}(\text{location} = \sigma_i^2, \text{scale} = \sqrt{\sigma_i^2})$.

Figure 4: Comparison of relative risks when $(\mu_i, \sigma_i^2)$ are dependent - Settings 6 and 7. Plots show $m = 10, 15, 20$ left to right.

freedom. Across both these settings the proposed NEST method demonstrates robustness to departures from the Normal model. Proposition 7 in Barp et al. (2019) guarantees that, in general, the influence function of minimum KSD estimators, such as the NEST estimator, is bounded under data corruption and the behavior of the NEST estimator in Settings 8 and 9 is potentially an example of such a robustness property. The scenario in Setting 9 is particularly interesting because we note from Figure 5b that the NEST estimator exhibits a smaller estimation risk than the Bayes oracle across all values of $u$. This is not surprising because in Settings 8 and 9, the Bayes oracle estimator $\delta_{(i)}^{\pi}$ in Equation (6), while having full knowledge of the underlying prior distributions, continues to rely on the incorrect Gaussian data generation process for $Y_{ij}$ to estimate the means. The NEST estimator, on the other hand, exhibits potential robustness to such mis-specification for estimating the shrinkage factors from the data.
(a) Setting 8: \((\mu_i, \sigma_i^2)\) generated according to Setting 5 and \(Y_{ij}|(\mu_i, \sigma_i^2) \sim \text{Laplace}(\text{location} = \mu_i, \text{scale} = \sqrt{\sigma_i^2/2})\).

(b) Setting 9: \((\mu_i, \sigma_i^2)\) generated according to Setting 4 and \(Y_{ij}|(\mu_i, \sigma_i^2) = \mu_i + \xi_i\) where \(\xi_i\) are independently sampled from a central \(t\)–distribution with 4 degrees of freedom.

**Figure 5:** Comparison of relative risks when the data \(Y_{ij}|(\mu_i, \sigma_i^2)\) are not normally distributed. Plots show \(m = 10, 15, 20\) left to right.

Overall, the results of the preceding nine simulation settings reveal that when the variances are unknown, the NEST estimation framework, in general, enjoys a relatively better risk performance for estimating \(\mu\) under the weighted squared error loss than the NPMLE based \(g\)-modeling approach. However, as Setting 3 reveals, the NPMLE may exhibit a substantially better risk performance when the underlying prior distribution on the means can be well approximated by a discrete mixture with a finite number of mass points. In this scenario, the NEST estimator is unable to effectively account for the potential multimodality of the marginal distribution while estimating the shrinkage factors and therefore may return a poorer estimate of the means.

In Section C of the Appendix, we present additional numerical experiments for the following cases: compound estimation of Normal means under the squared error loss (Section C.1), under unequal and small sample sizes \(m_i\) (Section C.2) and compound estimation of ratios under the squared error loss (Section C.3). Two real data examples are discussed in Section D.
6 Discussion

In this article we propose a nonparametric empirical Bayes framework, NEST, for large scale estimation of normal means when the corresponding precisions are unknown. We show that the data-driven NEST estimator is asymptotically as good as the optimal Bayes rule for estimating the means under the weighted squared error loss. Under the usual squared error loss our analysis in Section A reveals that the oracle NEST estimator dominates Tweedie’s formula that relies on sample variances for shrinkage estimation of the means. Furthermore, in our numerical experiments and real data examples NEST delivers a better risk performance, often substantially, than competing shrinkage estimators under both the weighted squared error loss and the squared error loss.

A key component of the NEST framework is the consistent estimation of the bivariate shrinkage factors that appear in the oracle NEST estimator. Our approach for estimating these shrinkage factors relies on the KSD measure and we solve a convex optimization problem to derive the shrinkage factors. However, the advantage of the KSD approach for empirical Bayes methods extends far beyond the hierarchical Model (1) considered in this paper. For instance, a natural extension of Model (1) is in large-scale regression problems where the goal is to estimate a large number of low-dimensional regression coefficients when some potentially useful side information on these unknown regression coefficients are available. In this setting the KSD approach is a natural candidate for estimating the multivariate shrinkage factors that arise in this context. Our future research will be directed towards studying this problem and developing empirical Bayes methods for the same.

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Appendix for “Nonparametric Empirical Bayes Estimation On Heterogeneous Data”

This Appendix is organized as follows: in Section A we present an analysis of the oracle NEST estimator (Definition 1) under the squared error loss. The proofs of all theoretical results in the paper are presented in Section B. In Section C, we provide additional numerical experiments for the following cases: compound estimation of Normal means under the squared error loss (Section C.1), under unequal and small sample sizes $m_i$ (Section C.2) and compound estimation of ratios under the squared error loss (Section C.3). Two real data examples are presented in Section D. In Section E we discuss extensions of our methodology to several well known members in the two-parameter exponential family when the nuisance parameter is known. We conclude with a discussion on the computational complexity of NEST in Section F.

A Analysis under the squared error loss

In this section we present an analysis of the oracle NEST estimator $\delta^*_{(1)}$ with respect to the squared error loss. We first introduce a counterpart to Tweedie’s formula for $\mu_i$ (Equation (2)) that relies on the sample variances $S^2_i$ instead of the unknown variances. Thereafter, we compare these two shrinkage estimators under the squared error loss.

When the precisions are unknown existing shrinkage methods, such as Weinstein et al. (2018), Xie et al. (2012), usually rely on the sample variance $S^2_i$, a consistent estimator of the unknown population variance $1/\tau_i$, for practical implementation. For instance, in Definition 3 we present the oracle Pseudo-Tweedie’s formula with sample variance which is a natural counterpart to Equation (2) when the variance is unknown.

**Definition 3.** *(Oracle Pseudo-Tweedie’s formula with sample variances)* Consider the hierarchical Model (1). Then an estimator for $\mu_i$ is $\delta^{TF}_i$ where,

$$\delta^{TF}_i := \delta^{TF}(y_i, s^2_i, m_i) = y_i + \frac{s^2_i}{m_i} w_1(y_i, s^2_i; m_i),$$

and $w_1(y, s^2; m) := \frac{\partial}{\partial y} \log f_m(y, s^2)$. 

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The oracle Pseudo-Tweedie’s formula in Definition 3 has a striking similarity to the oracle NEST estimator \( \delta_{\pi, i} \) (Definition 1) in the sense that both \( \delta_{i, \pi} \) and \( \delta_{i, (1)} \) involve an unbiased estimate \( y_i \) of \( \mu_i \) plus a shrinkage factor. The key difference, however, is that (1) while the shrinkage factor in \( \delta_{\pi, i} \) relies on \( \gamma_i \), \( \delta_{i, \pi} \) directly uses sample variances \( s_i^2 \), and (2) while it is not immediately clear what is the underlying loss function that \( \delta_{i, \pi} \) is minimizing for estimating \( \mu_i \), \( \delta_{i, (1)} \) uniquely minimizes the expected weighted squared error loss.

Notwithstanding, Proposition 1 establishes that the oracle Pseudo-Tweedie’s estimator \( \delta_{TF} = (\delta_{TF, 1}, \ldots, \delta_{TF, n}) \) dominates the sample mean estimator under a squared error loss function.

**Proposition 1.** Denote \( w_1'(y, s^2; m) \) as the partial derivative of \( w_1(y, s^2; m) \) with respect to \( y \). Suppose \( f_m(y, s^2) \) is a log-concave density and \( w_1'(y, s^2; m) \) is a non-decreasing function of \( s^2 \). Then under Model (1) and for \( m_i > 3 \), we have,

\[
 r_0(\delta_{\pi, (0)}, \mathcal{G}) \leq r_0(\delta_{TF}, \mathcal{G}) < r_0(Y, \mathcal{G}).
\]

However, when a large number of units are investigated simultaneously, traditional sample variance estimators may suffer from selection bias (Jing et al., 2016a, Kwon and Zhao, 2022), and their direct use may lead to severe deterioration in the MSE for estimating the means. We present Proposition 2 which shows that under the squared error loss function the risk of the oracle NEST estimator is uniformly smaller than that of the Pseudo-Tweedie’s estimator \( \delta_{TF} \) (Definition 3).

We impose the following regularity conditions for comparing the estimation risks of \( \delta_{\pi, (1)} \) and \( \delta_{TF} \) in Proposition 2.

**Assumption 4.** The shrinkage factor \( \gamma(y, s^2, m) \) is non-increasing in \( s^2 \).

**Assumption 5.** Let \( \omega(y, s^2; m) := w_1(y, s^2; m) \frac{\partial}{\partial y} w_2(y, s^2; m) \). Then, \( \omega(y, s^2; m) \leq 0 \).

Assumptions 4 and 5 are regularity conditions on the behavior of the shrinkage factor \( \gamma \) and the score functions \( w_1, w_2 \). For instance, Assumptions 4 enforces monotonicity on shrinkage factor \( \gamma(y, s^2, m) \), which is satisfied, for example, when \( (\mu, \tau) \) have a conjugate prior under Model (1). Similarly, Assumption 5 holds under conjugate priors and is also true when the prior on \( \tau \) is discrete with just one mass point. In Remarks 2 and 3 we discuss these assumptions in more details. Proposition 1 can be extended as follows.
Proposition 2. Suppose \( f_m(y, s^2) \) is a log-concave density and \( w_1'(y, s^2; m) \) is a non-decreasing function of \( s^2 \). If Assumptions 4 – 5 hold then, under Model (1) and for \( m_i > 5 \), we have,

\[
r_0(\delta^\pi_{(0)}, \mathcal{G}) \leq r_0(\delta^\pi_{(1)}, \mathcal{G}) \leq r_0(\delta^{TF}, \mathcal{G}) < r_0(Y, \mathcal{G}).
\]

Note that the oracle NEST estimator \( \delta^\pi_{i,(1)} \) is, in general, different from \( \delta^\pi_{i,(0)} \); \( \delta^\pi_{i,(1)} \) is the Bayes estimator under the weighted squared error loss, while \( \delta^\pi_{i,(0)} \) is the Bayes estimator under the usual squared error loss. This follows since from Equation (3), and dropping subscript \( i \),

\[
\delta^\pi_{(1)} = \frac{\mathbb{E}(\tau \mu | y, s^2)}{\mathbb{E}(\tau | y, s^2)} = \frac{\mathbb{E}[\tau \mathbb{E}(\mu | y, s^2, \tau) | y, s^2]}{\mathbb{E}(\tau | y, s^2)} \neq \mathbb{E}(\mu | y, s^2) = \delta^\pi_{(0)}.
\]

However, \( r_0(\delta^\pi_{(0)}, \mathcal{G}) = r_0(\delta^\pi_{(1)}, \mathcal{G}) \) when \( \mu \) and \( \tau \) are conditionally independent given \( y \) and \( s^2 \), in which case \( \delta^\pi_{(1)} = \delta^\pi_{(0)} \). In particular, Corollary 1 shows that when \( \mathbb{E}(\mu_i | y_i, s_i^2, \tau_i) \) is independent of \( \tau_i \) then the data-driven NEST estimator is asymptotically close to \( \delta^\pi_{(0)} \).

Corollary 1. Consider hierarchical Model (1) and suppose \( \mathbb{E}(\mu_i | y_i, s_i^2, \tau_i) \) is independent of \( \tau_i \). Then, under the conditions of Theorem 3, we have \( c_n \| \delta_n^\pi(\lambda) - \delta^\pi_{(0)} \|_2 = o_p(1) \) as \( n \to \infty \). Furthermore, under the same conditions, \( c_n^{1/2} | l_n^{(0)}(\delta_n^{\pi}(\lambda), \mu; \tau) - l_n^{(0)}(\delta^\pi_{(0)}, \mu; \tau) | = o_p(1) \) as \( n \to \infty \).

Corollary 1 is a straightforward consequence of Theorem 3 and the fact that when \( \mathbb{E}(\mu_i | y_i, s_i^2, \tau_i) \) is independent of \( \tau_i \), \( \delta^\pi_{i,(1)} = \delta^\pi_{i,(0)} \). A popular setting where \( \mathbb{E}(\mu_i | y_i, s_i^2, \tau_i) \) is indeed independent of \( \tau_i \) is the scenario where \( G_\mu(\cdot | \tau) \) and \( H_\tau(\cdot) \) belong to the family of conjugate priors under Model (1). Under this setting the posterior expectation of \( \mu \) is a linear combination of the prior expectation of \( \mu \) and the maximum likelihood estimate \( y \) (Diaconis and Ylvisaker, 1979), and the weights in this linear combination are proportional to \( m \) and the prior sample size. For instance, if \( G_\mu(\cdot | \tau) \) is \( N(\mu_0, 1/\tau) \) and \( H_\tau(\cdot) = \Gamma(\text{shape} = \alpha, \text{rate} = \beta) \) in Model (1) then standard calculations give \( \delta^\pi_{(0)} = (\mu_0 + my)/(m + 1) \) where \( \mu_0 \) is the prior expectation of \( \mu \) and \( m_0 = 1 \) is the prior sample size. Moreover, in this setting we also have

\[
\log f_m(y_i, s_i^2) = c_0 + \frac{m - 3}{2} \log s_i^2 - (\alpha + m/2) \log \left\{ \beta + 0.5(m - 1)s_i^2 + 0.5 \frac{m}{m + 1}(y_i - \mu_0)^2 \right\},
\]
where \( c_0 \) is a constant independent of \((y_i, s_i^2)\). From the above display,

\[
\begin{align*}
 w_1(y_i, s_i^2) &= -\frac{(\alpha + m/2)\{m/(m+1)\}(y_i - \mu_0)}{\beta + 0.5(m-1)s_i^2 + 0.5\{m/(m+1)\}(y_i - \mu_0)^2} \\
 w_2(y_i, s_i^2) &= \frac{(m-3)}{2s_i^2} - \frac{0.5(\alpha + m/2)(m-1)}{\beta + 0.5(m-1)s_i^2 + 0.5\{m/(m+1)\}(y_i - \mu_0)^2}.
\end{align*}
\]

Substituting these expressions for \( w_1(y_i, s_i^2) \), \( w_2(y_i, s_i^2) \) in Equation (6) gives \( \delta_i^\pi = \delta_i^{\pi(0)} \). Another example where \( \delta_i^\pi = \delta_i^{\pi(0)} \) is when \( G_\mu(\cdot|\tau) \) and \( H_\tau(\cdot) \) are independent discrete distributions with just one mass point, respectively. These two examples are also the settings, among others, under which the first inequality in the left hand side of Proposition 1 is strict, that is, 
\[ r_0(\delta_i^{\pi(0)}, \mathcal{G}) < r_0(\delta_T^{\pi}, \mathcal{G}). \]

In Section C.1 we present several numerical experiments to compare the data-driven NEST estimator against other shrinkage estimators for compound estimation of the Normal means from Model (1) under the squared error loss. Our empirical results corroborate Proposition 2 and suggest that the efficiency gain of the data-driven NEST estimator over competing linear shrinkage methods (Jing et al., 2016b, Weinstein et al., 2018) and Pseudo-Tweedie’s formula (Definition 3) is substantial across many settings. We end this section with the following remarks:

**Remark 2.** Here we provide three examples of \( \mathcal{G} \) that satisfy Assumptions 4 and 5.

1. \( G_\mu(\cdot|\tau) \) and \( H_\tau(\cdot) \) belong to the family of conjugate priors under Model (1).
2. \( H_\tau(\cdot) \) is a distribution with just one mass point, say \( \tau_0 \), and \( G_\mu(\cdot|\tau) \) has a log-concave density.
3. \( H_\tau(\cdot) \) has a log-concave density and \( G_\mu(\cdot|\tau) \) is a discrete distribution with just one mass point, say \( \mu_0 \).

In Figure 6 we consider specific choices of \( \mathcal{G} \) for each of these three examples. In particular, the left panel of Figure 6 provides a contour plot of \( \mathcal{G} \) for example 1 where \( G_\mu(\cdot|\tau) \) is \( N(0, 1/\tau) \) and \( H_\tau(\cdot) \) is \( \Gamma(\text{shape} = 10, \text{rate} = 5) \). The center panel presents the contour plot of the marginal density \( f_m \) for example 2 where \( G_\mu(\cdot|\tau) \) is \( N(0, 1/\tau) \) and \( H_\tau(\cdot) \) has all its mass concentrated at \( \tau_0 = 1 \). Finally, the right panel presents the contour plot of \( f_m \) for example 3 where \( \mu_0 = 0 \) and \( H_\tau(\cdot) \) is \( \Gamma(\text{shape} = 10, \text{rate} = 5) \).
Remark 2 contour plots with $m = 5$. Left to Right: contour plot of $G$ for example 1 where $G_{\mu}(\cdot | \tau)$ is $N(0, 1/\tau)$ and $H_{\tau}(\cdot)$ is $\Gamma$ (shape = 10, rate = 5). Contour plot of the marginal density $f_m$ for example 2 where $G_{\mu}(\cdot | \tau)$ is $N(0, 1/\tau)$ and $H_{\tau}(\cdot)$ has all its mass concentrated at $\tau_0 = 1$. Contour plot of $f_m$ for example 3 where $\mu_0 = 0$ and $H_{\tau}(\cdot)$ is $\Gamma$ (shape = 10, rate = 5).

**Remark 3.** Here we provide two examples of $G$ that induce marginal densities $f_m$ which are no longer log-concave and, hence, do not satisfy Assumption 4. In these examples, log-concavity of $f_m$ is lost because $w'_1(y, s^2; m)$ may be positive on its domain.

1. $H_{\tau}(\cdot)$ is a distribution with just one mass point, say $\tau_0$, and $G_{\mu}(\cdot | \tau_0)$ has mass points at $2\tau_0$ and $-2\tau_0$ with equal probability.

2. $H_{\tau}(\cdot)$ has mass points at 1 and 2 with equal probability, and $\mu = 4/\tau$.

Figure 7 presents the contour plots of $f_m$ for these two examples.

**B Proofs**

**B.1 Proof of Proposition 1**

Recall from Definition 3 that

$$\delta_i^{TF} := \delta^{TF}(y_i, s_i^2, m_i) = y_i + \frac{s_i^2}{m_i} w_1(y_i, s_i^2; m_i),$$

where $w_1(y, s^2; m) := \frac{\partial}{\partial y} \log f_m(y, s^2)$. Since the $n$ study units are independent, we will focus on unit $i$.  

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Figure 7: Remark 3 contour plots. Left: contour plot of the marginal density $f_m$ for example 1 where $\tau_0 = 1$ and $m = 5$. Right: contour plot of $f_m$ for example 2 with $m = 5$.

We note that the inequality in the left hand side of Proposition 1 follows from the fact that $\delta_{\pi_i,(0)}$ is the Bayes estimator of $\mu_i$ that uniquely minimizes the expected squared error loss under Model (1). To prove the inequality in the right hand side of Proposition 1, we proceed as follows.

Denote $w_{1,i} := w_1(y_i, s^2_i; m_i)$. We have,

$$
\mathbb{E}(\mu_i - \delta_i^{TF})^2 = \frac{1}{m_i} \mathbb{E}(1/\tau_i) + \frac{1}{m_i} \mathbb{E}(S^2_i w_{1,i})^2 + \frac{2}{m_i} \mathbb{E}(Y_i - \mu_i) S^2_i w_{1,i} = \mathbb{E}(\mu_i - Y_i)^2 + \frac{1}{m_i} \mathbb{E}(S^2_i w_{1,i})^2 + \frac{2}{m_i} \mathbb{E} \left( \frac{S^2_i}{\tau_i} w'_{1,i} \right),
$$

(17)

where $w'_{1,i} := w'_1(y_i, s^2_i; m_i)$ and $w_1(y, s^2; m) = \frac{\partial}{\partial y} w_1(y, s^2; m)$. The equality in Equation (17) follows from integration by parts and the fact that $Y_i|\mu_i, \tau_i \sim N(\mu_i, 1/(m_i\tau_i))$. Consider the term

$$
T_1 := \frac{1}{m_i} \mathbb{E}(S^2_i w_{1,i})^2 + \frac{2}{m_i} \mathbb{E} \left( \frac{S^2_i}{\tau_i} w'_{1,i} \right)
$$

and note that

$$
T_1 = \frac{1}{m_i} \mathbb{E} \left( S^2_i w_{1,i} \right)^2 + \frac{1}{m_i} \mathbb{E} \left( (S^2_i)^2 w'_{1,i} \right) + \frac{2}{m_i} \mathbb{E} \left( \frac{S^2_i}{\tau_i} w'_{1,i} \right) - \frac{1}{m_i} \mathbb{E} \left( (S^2_i)^2 w'_{1,i} \right) = \frac{1}{m_i} \mathbb{E} \left( \frac{2S^2_i}{\tau_i} - (S^2_i)^2 \right) w'_{1,i}.
$$

(18)

The equality in Equation (18) follows because, dropping subscript $i$,

$$
\mathbb{E} \left[ (S^2)^2 (w_1^2 + w'_1) \right] = \mathbb{E} \left[ (S^2)^2 \frac{f''_{m_i}(1)}{f_m(Y, S^2)} \right] = 0,
$$
where \( f''_{m,1}(y,s^2) \) is the second order partial derivative of \( f_m(y,s^2) \) with respect to \( y \). Now, we can re-write Equation (18) as

\[
T_1 = \frac{1}{m_i^2} E_{\mu,\tau} E_{Y_i,S_i^2|\mu_i,\tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] w'_{1,i} \right\} = \frac{1}{m_i^2} E_{\mu_i,\tau_i} \left( T_2 \right) + \frac{1}{m_i^2} E_{\mu_i,\tau_i} \left( T_3 \right),
\]

where \( E_{\mu,\tau} \) is the expectation with respect to the joint distribution of \( (\mu, \tau) \), \( E_{Y_i,S_i^2|\mu_i,\tau_i} \) is the expectation with respect to the joint distribution of \( (Y_i, S_i^2) \) conditional on \( (\mu, \tau) \) and

\[
T_2 = E_{Y_i,S_i^2|\mu_i,\tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] w'_{1,i} \left| S_i^2 < \frac{2}{\tau_i} \right. \right\} \mathbb{P} \left( S_i^2 < \frac{2}{\tau_i} \right),
\]

\[
T_3 = E_{Y_i,S_i^2|\mu_i,\tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] w'_{1,i} \left| S_i^2 > \frac{2}{\tau_i} \right. \right\} \mathbb{P} \left( S_i^2 > \frac{2}{\tau_i} \right).
\]

Denote \( p := \mathbb{P}(S_i^2 > 2/\tau_i | \tau_i) \) and let \( c_i \) be the partial derivative of \( w_1(y, s^2; m) \) with respect to \( y \) and evaluated at \( (y_i, 2/\tau_i, m_i) \). Now, using equations (18) and (19) in Equation (17), we get

\[
E(\mu_i - \delta_i)^2 = E(\mu_i - Y_i)^2 + \frac{1}{m_i^2} E_{\mu_i,\tau_i} (T_2 + T_3).
\]

(20)

We will show that \( T_2 + T_3 < 0 \) which will be enough to prove the statement of Proposition 1 using Equation (20).

We first state a few results that are straightforward consequences of Model (1) and the assumptions of Proposition 1. We have,

1. Under Model 1, \((m_i - 1)S_i^2/\tau_i \sim \chi_{m_i-1}^2\).

2. Additionally,

\[
E_{Y_i,S_i^2|\mu_i,\tau_i} \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] = \frac{m_i - 3}{(m_i - 1)\tau_i^2} > 0,
\]

since \( m_i > 3 \) in the statement of Proposition 1.

3. Since \( f_m(y, s^2) \) is a log-concave density, \( w'_{1,i} \leq 0 \) and so \( T_2 \leq 0 \) while \( T_3 \geq 0 \).

Assume, without loss of generality, \( w'_{1,i} < 0 \). Since \( E_{Y_i,S_i^2|\mu_i,\tau_i} [2S_i^2/\tau_i - (S_i^2)^2] > 0 \), we have

\[
(1 - p)E_{Y_i,S_i^2|\mu_i,\tau_i} \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] S_i^2 < \frac{2}{\tau_i} > -pE_{Y_i,S_i^2|\mu_i,\tau_i} \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] S_i^2 > \frac{2}{\tau_i} > 0.
\]

(21)
Furthermore, as \( w'_{1,i} < 0 \) and \( w'_{1,i} \) is a non-decreasing function of \( s_i^2 \),

\[
T_2 \leq (1 - p)E_{Y_i, s_i^2 | \mu_i, \tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] c_i \big| S_i^2 < \frac{2}{\tau_i} \right\},
\]

(22)

Note that \( c_i < 0 \), and as defined earlier, it is the partial derivative of \( w_1(y, s^2; m) \) with respect to \( y \) and evaluated at \( (y_i, 2/\tau_i, m_i) \). Therefore, using equations (21) and (22),

\[
T_2 \leq (1 - p)E_{Y_i, s_i^2 | \mu_i, \tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] c_i \big| S_i^2 < \frac{2}{\tau_i} \right\} < -pE_{Y_i, s_i^2 | \mu_i, \tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] c_i \big| S_i^2 > \frac{2}{\tau_i} \right\} < 0.\quad(23)
\]

Now, we consider the term \( T_3 \). Recall that

\[
T_3 = pE_{Y_i, s_i^2 | \mu_i, \tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] w'_{1,i} \big| S_i^2 > \frac{2}{\tau_i} \right\} > 0.
\]

Since \( w'_{1,i} < 0 \) and \( w'_{1,i} \) is a non-decreasing function of \( s_i^2 \),

\[
0 < T_3 \leq -pE_{Y_i, s_i^2 | \mu_i, \tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] c_i \big| S_i^2 > \frac{2}{\tau_i} \right\}.
\]

(24)

So, using equations (23) and (24)

\[
T_2 + T_3 < -pE_{Y_i, s_i^2 | \mu_i, \tau_i} \left\{ \left[ \frac{2S_i^2}{\tau_i} - (S_i^2)^2 \right] c_i \big| S_i^2 > \frac{2}{\tau_i} \right\} = 0,
\]

Hence the desired result follows from the display above and Equation (20).

\[\blacksquare\]

### B.2 Proof of Proposition 2

We will first collect a few notations that will be used throughout the proof. Denote \( w_{1,i} := w_1(y_i, s_i^2; m_i) \), \( w_{2,i} := w_2(y_i, s_i^2; m_i) \) and \( \gamma_i := \gamma(y_i, s_i^2, m_i) \). Let \( w'_1(y, s^2; m) = \frac{\partial}{\partial y} w_1(y, s^2; m) \) and denote \( w'_{1,i} := w'_1(y_i, s_i^2; m_i) \). Similarly, \( \nu(y, s^2; m) = \frac{1}{f_m(y, s^2)} \frac{\partial^2}{\partial y^2} f_m(y, s^2) \) and denote \( \nu_i := \nu(y_i, s_i^2; m_i) \). Finally, let \( \gamma'(y, s^2, m) = \frac{\partial}{\partial y} \gamma(y, s^2, m) \) and denote \( \gamma'_i := \gamma'(y_i, s_i^2, m_i) \).

The proof of Proposition 2 will use the following two lemmata.

**Lemma 1.** Suppose \( f_m(y, s^2) \) is a log concave density and Assumption 4 holds. Then under
Model (1) and \( m_i > 5 \), we have,
\[
E\left\{ \left[ \frac{2S_i^2}{\tau_i} \left( 1 - \gamma_i \right) - \left( S_i^2 \right)^2 \left( 1 - \gamma_i^2 \right) \right] w_{1,i}' \right\} \geq 0.
\]

**Lemma 2.** Under Assumption 5 and Model (1), we have,
\[
\frac{1}{m_i^2} E\left[ S_i^2 \left( S_i^2 \gamma_i^2 \nu_i + 2 \frac{w_{1,i}}{\tau_i} \gamma_i' \right) \right] \leq 0.
\]

Lemmata 1 and 2 are proved in Sections B.2.1 and B.2.2 respectively. We now prove Proposition 2. Recall from Definition 1 that the oracle NEST estimator for \( \mu_i \) is
\[
\delta_{i,(1)}^{\pi} = y_i + \frac{s_i^2}{m_i} \gamma(y_i, s_i^2, m_i) w_1(y_i, s_i^2, m_i),
\]
where
\[
\gamma(y_i, s_i^2, m_i) = \frac{m_i - 1}{m_i - 3 - 2s_i^2 w_2(y_i, s_i^2, m_i)}.
\]

Since the \( n \) study units are independent, we will focus on unit \( i \).

Under the squared error loss, \( \delta_{i,(0)}^{\pi} \) is the Bayes estimator of \( \mu_i \) in Model (1). This establishes the first inequality on the left hand side of Proposition 2. Together with Proposition 1, we only need to show \( r_0(\delta_{i,(1)}^{\pi}, G) \leq r_0(\delta_{TF}, G) \). First note that
\[
\begin{align*}
& r_0(\delta_{TF}, G) - r_0(\delta_{i,(1)}^{\pi}, G) = \frac{1}{m_i^2} E\left[ \left( S_i^2 \right)^2 \left( 1 - \gamma_i^2 \right) w_{1,i}' \right] + \frac{2}{m_i} E\left[ (Y_i - \mu_i) S_i^2 \left( 1 - \gamma_i \right) w_{1,i} \right] \\
& \quad = \frac{1}{m_i^2} E\left[ \left( S_i^2 \right)^2 \left( 1 - \gamma_i^2 \right) w_{1,i}' \right] + \frac{2}{m_i} E\left\{ \frac{S_i^2}{\tau_i} \left[ (1 - \gamma_i) w_{1,i}' y_i - w_{1,i}' \gamma_i \right] \right\}. \quad (25)
\end{align*}
\]

The equality in equation (25) follows from integration by parts and the fact that \( Y_i \sim N \left( \mu_i, \frac{1}{m_i \tau_i} \right) \).

We can re-write Equation (25) as,
\[
\begin{align*}
& \frac{1}{m_i^2} E\left[ \left( S_i^2 \right)^2 \left( 1 - \gamma_i^2 \right) w_{1,i}' \right] + \frac{1}{m_i^2} E\left[ \left( S_i^2 \right)^2 \left( 1 - \gamma_i^2 \right) w_{1,i}' \right] - \frac{2}{m_i} E\left\{ \frac{S_i^2}{\tau_i} w_{1,i}' \gamma_i \right\} \\
& \quad + \frac{1}{m_i^2} E\left\{ \left[ \frac{2S_i^2}{\tau_i} \left( 1 - \gamma_i \right) - \left( S_i^2 \right)^2 \left( 1 - \gamma_i^2 \right) \right] w_{1,i}' \right\}. \quad (26)
\end{align*}
\]

From Lemma 1, the last term in Equation (26) is non-negative. Let us consider the first three
terms in Equation (26) and denote them by,

\[ T := \frac{1}{m_i^2} \mathbb{E} \left[ (S_i^2)^2 \left( 1 - \gamma_i^2 \right) w^2_{i,i} \right] + \frac{1}{m_i^2} \mathbb{E} \left[ (S_i^2)^2 \left( 1 - \gamma_i^2 \right) w'_{i,i} \right] - \frac{2}{m_i^2} \mathbb{E} \left( \frac{S_i^2}{\tau_i} w_{1,i} \gamma_i' \right). \]

As shown in the proof of Proposition 1, \( \mathbb{E}[(S^2)^2(w^2_1 + w'_1)] = 0 \). The above display involving the term \( T \) can be written as

\[ T = -\frac{1}{m_i^2} \mathbb{E} \left[ (S_i^2)^2 \gamma_i^2 \nu_i \right] - \frac{2}{m_i^2} \mathbb{E} \left( \frac{S_i^2}{\tau_i} w_{1,i} \gamma_i' \right). \] (27)

From Lemma 2, the term \( T \) in Equation (27) is non-negative. This establishes the inequality on the right hand side of \( r_0(\delta_{(1)}, \mathcal{G}) \) in Proposition 2 and completes the proof. \( \blacksquare \)

### B.2.1 Proof of Lemma 1 -

Denote

\[ Z_i := \frac{2S_i^2}{\tau_i} (1 - \gamma_i) - (S_i^2)^2 (1 - \gamma_i^2) = \frac{Q_i}{\tau_i}, \]

where \( Q_i = 2S_i^2(1 - \gamma_i) - (S_i^2)^2(1 - \gamma_i^2)\tau_i \). Since \( f_m(y, s^2) \) is a log-concave density, \( w'_{1,i} \leq 0 \).

We will show that \( Q_i \leq 0 \) which will be sufficient to prove \( Z_i w'_{1,i} \geq 0 \) and hence the statement of Lemma 1.

Dropping subscript \( i \),

\[ \mathbb{E}(Q) = \mathbb{E}_{Y,S^2} \left[ 2S^2(1 - \gamma) - (S^2)^2(1 - \gamma^2)\hat{\tau}^\pi \right] = \mathbb{E}_{Y,S^2}[R(Y, S^2)], \]

where \( \hat{\tau}^\pi = \mathbb{E}(\tau|y, s^2, m) = (s^2 \gamma)^{-1} \) from Equation (8) and \( \mathbb{E}_{Y,S^2} \) is expectation with respect to the joint marginal distribution of \((Y, S^2)\). Suppose, if possible, \( Q_i > 0 \) for all \((\tau_i, y_i, s_i^2) \in \mathbb{R}^+ \times \mathbb{R} \times \mathbb{R}^+\). We will show that \( \mathbb{E}(Q) \leq 0 \) which will present a contradiction to \( Q > 0 \) for all \((\tau, y, s^2) \in \mathbb{R}^+ \times \mathbb{R} \times \mathbb{R}^+\). Fix a \( y \in \mathbb{R} \) and consider the following cases.

**Case 1** – Suppose \( 0 < \gamma(y, s^2, m) \leq 1/(2c) \) where \( c \geq 1 \) is a constant. Then, we have \( s^2 \hat{\tau}^\pi \geq 2c \) and consequently \( s^2(1 + \gamma)\hat{\tau}^\pi > 2 \). So \( R(y, s^2) < 0 \). Now, from Assumption 4, \( \gamma \) is a continuous and non-increasing function of \( s^2 \). Therefore, there exists \( c_1(y) \in \mathbb{R}^+ \), depending on \( y \), such that \( s^2 > c_1(y) \) whenever \( 0 < \gamma(y, s^2, m) \leq 1/(2c) \). Thus, \( R(y, s^2) < 0 \) for \( s^2 > c_1(y) \).

**Case 2** – Next, suppose \( 1/2 < \gamma(y, s^2, m) \leq (m - 1)/(2m - 6) \). Then \((2m - 6)/(m - 1) \leq
s^2 \hat{\tau}^\pi < 2 \text{ and } s^2(1 + \gamma)\hat{\tau}^\pi \geq (3m - 7)/(m - 1). Since m > 5, (3m - 7)/(m - 1) > 2 \text{ and so } s^2(1 + \gamma)\hat{\tau}^\pi > 2. Thus R(y, s^2) < 0 \text{ and using Assumption 4, } R(y, s^2) < 0 \text{ for } c_2(y) < s^2 \leq c_1(y), \text{ where } c_2(y) \text{ is such that } s^2 > c_2(y) \text{ whenever } \gamma(y, s^2, m) \leq (m - 1)/(2m - 6).

The remaining four cases proceed in a similar manner as follows:

**Case 3** – Suppose \((m - 1)/(2m - 6) < \gamma(y, s^2, m) \leq (2m - 6)/(3m - 11)\). Then \((3m - 11)/(2m - 6) \leq s^2 \hat{\tau}^\pi\). So \(s^2(1 + \gamma)\hat{\tau}^\pi \geq (5m - 17)/(2m - 6)\). Since \((5m - 17)/(2m - 6) > 2\) if \(m > 5\), we have \(s^2(1 + \gamma)\hat{\tau}^\pi > 2\) and so \(R(y, s^2) < 0\) for \(c_3(y) < s^2 \leq c_2(y)\). Similarly, we can show that \(R(y, s^2) \leq 0\) for \(c_4(y) < s^2 \leq c_3(y)\) where \(s^2 \in (c_4(y), c_3(y)]\) whenever \((2m - 6)/(3m - 11) < \gamma(y, s^2, m) \leq 1\).

**Case 4** – Now suppose, \(1 < \gamma(y, s^2, m) \leq (2m - 5)/(2m - 6)\). Then \((2m - 6)/(2m - 5) \leq s^2 \hat{\tau}^\pi < 1\) and \(s^2(1 + \gamma)\hat{\tau}^\pi < 2\). Note that here \(\gamma > 1\) as opposed to \(\gamma \leq 1\) in the earlier cases. Therefore, \(s^2(1 + \gamma)\hat{\tau}^\pi < 2\) implies \(R(y, s^2) < 0\) for \(c_5(y) < s^2 \leq c_4(y)\).

**Case 5** – Similarly, if \(\gamma\) is in the intervals \(((2m - 5)/(2m - 6), (m - 2)/(m - 3)]\), \(((m - 2)/(m - 3), (m - 1)/(m - 3)]\) and \(((m - 1)/(m - 3), (m + 1)/(m - 3)]\) then we have \(\gamma \geq 1\) and \(s^2(1 + \gamma)\hat{\tau}^\pi < 2\). Thus, on each of the corresponding intervals for \(s^2\), \(R(y, s^2) < 0\).

**Case 6** – Denote \(r_1 = 1\) and \(r_t = 2r_{t-1} + 3\) for \(t = 2, 3, \ldots\). Suppose \((m + r_{t-1})/(m - 3) < \gamma(y, s^2, m) \leq (m + r_t)/(m - 3)\). Then for each of these intervals indexed by \(t \geq 2\), we have \(\gamma \geq 1\), \(s^2(1 + \gamma)\hat{\tau}^\pi < 2\) and so \(R(y, s^2) < 0\) on the corresponding intervals for \(s^2\).

So from these six cases, \(R(y, s^2) \leq 0\) for all \(s^2 > 0\) and consequently \(\mathbb{E}_{Y, s^2}[R(Y, S^2)] \leq 0\). Therefore, \(\mathbb{E}(Q) \leq 0\) which contradicts that \(Q > 0\) for all \((\tau, y, s^2) \in \mathbb{R}^+ \times \mathbb{R} \times \mathbb{R}^+\). Now suppose that for some \(\Omega \subset \mathbb{R}^+ \times \mathbb{R} \times \mathbb{R}^+, Q > 0\) whenever \((\tau, y, s^2) \in \Omega\). However, Assumption 4 and the aforementioned six cases imply that \(\mathbb{E}(Q|\Omega) \leq 0\). Thus, \(Q \leq 0\) for all \((\tau, y, s^2) \in \mathbb{R}^+ \times \mathbb{R} \times \mathbb{R}^+\). So, we have \(Zw'_1 \geq 0\) and this completes the proof of Lemma 1.

**B.2.2  \textbf{Proof of Lemma 2 -}**

Dropping subscript \(i\), denote,

\[
T := -\frac{1}{m^2} \mathbb{E}\left[(S^2 \gamma(Y, S^2, m))^2 \nu(Y, S^2, m)\right] - \frac{2}{m^2} \mathbb{E}\left[\frac{S^2}{\tau} w_1(Y, S^2, m) \gamma'(Y, S^2, m)\right].
\]

We will show that \(T \geq 0\).

Let \(\gamma := \gamma(y, s^2, m), \nu := \nu(y, s^2, m), w_1 := w_1(y, s^2, m)\) and \(\gamma' := \gamma'(y, s^2, m)\). First note
that using standard integration by parts, we have

\[-\frac{1}{m^2} \mathbb{E}\left[(S^2 \gamma')^2 \nu\right] = \frac{2}{m^2} \mathbb{E}\left[(S^2) \gamma w_1 \gamma'\right].\]

So, we can write

\[T = \frac{2}{m^2} \mathbb{E}\left\{ S^2 w_1 \gamma' \left[ S^2 \gamma - \frac{1}{\tau}\right] \right\}.\]

Furthermore, from Definition 1,

\[\gamma' = \frac{2}{m-1} \gamma^2 \sigma^2 w_2',\]

where \(w_2' := \frac{\partial}{\partial y} w_2(y, s^2, m).\) So, we have

\[T = \frac{4}{m^2 (m-1)} \mathbb{E}\left\{ (S^2 \gamma)^2 w_1 w_2' \left[ S^2 \gamma - \frac{1}{\tau}\right] \right\}.\]

Now conditional on \((Y, S^2),\) note that from Jensen’s inequality,

\[s^2 \gamma = \frac{1}{\mathbb{E}(\tau \mid y, s^2)} \leq \mathbb{E}\left(\frac{1}{\tau} \mid y, s^2\right).\]

Furthermore, since \(w_1 w_2' \leq 0\) from Assumption 5, we have

\[T = \frac{4}{m^2 (m-1)} \mathbb{E}_{Y,S^2}\left\{ (S^2 \gamma)^2 w_1 w_2' \left[ S^2 \gamma - \mathbb{E}\left(\frac{1}{\tau} \mid Y, S^2\right)\right] \right\} \geq 0.\]

This completes the proof of Lemma 2. \[\blacksquare\]
B.3 Proof of Theorem 1

Consider the sample criterion \( \hat{M}_{\lambda,n}(\tilde{W}) \) (Equation (16)) and population criterion \( M_{\lambda}(\tilde{W}) \) (Equation (15)). We have

\[
\left| \hat{M}_{\lambda,n}(\tilde{W}) - M_{\lambda}(\tilde{W}) \right| \leq \left| \frac{1}{n^2} \sum_{i=1}^{n} \sum_{j=1}^{n} \kappa_{\lambda}[\tilde{w}(X^i), \tilde{w}(X^j)](X^i, X^j)I(i \neq j) - M_{\lambda}(\tilde{W}) \right| + \left| \frac{1}{n^2} \sum_{i=1}^{n} \kappa_{\lambda}[\tilde{w}(X^i), \tilde{w}(X^i)](X^i, X^i) \right| =: I_1 + I_2. \tag{28}
\]

Define \( \bar{M}_{\lambda,n}(\tilde{W}) = [n(n-1)]^{-1} \sum_{i=1}^{n} \sum_{j=1}^{n} \kappa_{\lambda}[\tilde{w}(X^i), \tilde{w}(X^j)](X^i, X^j)I(i \neq j) \). Then

\[
I_1 \leq |\bar{M}_{\lambda,n}(\tilde{W}) - M_{\lambda}(\tilde{W})| + n^{-1}|\bar{M}_{\lambda,n}(\tilde{W})|. \]

By Assumption 1, \( n^{-1}|\bar{M}_{\lambda,n}(\tilde{W})| \) is \( O_p(n^{-1}) \). Now, note that \( \bar{M}_{\lambda,n}(\tilde{W}) \) is an unbiased estimator of \( M_{\lambda}(\tilde{W}) \) and is a U-statistic with a symmetric kernel function \( \kappa_{\lambda}[\tilde{w}(X^i), \tilde{w}(X^j)](X^i, X^j) \).

From Assumption 1, \( \kappa_{\lambda}[\tilde{w}(X^i), \tilde{w}(X^j)](X^i, X^j) \) has finite second moments. Moreover, from theorem 4.1 of Liu et al. (2016), \( \bar{M}_{\lambda,n}(\tilde{W}) \) is a non-degenerate U-statistic whenever \( f \neq \tilde{f} \). Thus, from the CLT for U-statistics (Serfling (2009) section 5.5), \( |\bar{M}_{\lambda,n}(\tilde{W}) - M_{\lambda}(\tilde{W})| \) is \( O_p(n^{-1/2}) \).

Moreover, from Assumption 1, \( I_2 \) is \( O_p(n^{-1}) \). Theorem 1 is thus proved by combining these results.

\[ \blacksquare \]

B.4 Proof of Theorem 2

We begin by introducing some notations and a lemma that will be useful for proving Theorem 2.

Recall from Equation (11) that \( w_{k,i} \) denotes the \((i, k)\)th element of the true score matrix \( W_0 \) where \( k = 1, 2 \). Similarly, from Definition 2, \( \hat{w}_{\lambda,n}^{(k)}(i) \) denotes the \((i, k)\)th element of the estimated score matrix \( \hat{W}_n(\lambda) \). Let

\[
\Delta_{\lambda,n}^{(p)}(i, k) = E \left| \hat{w}_{\lambda,n}^{(k)}(i) - w_{k,i} \right|^p \quad \text{and} \quad \Delta_{\lambda,n}(i, k) = E \left\{ \left| \hat{w}_{\lambda,n}(i) - w_{k,i} \right|^2 f(X_i) \right\},
\]

where the expectation is taken with respect to the joint distribution of \( X_1, \ldots, X_n \) and \( X_i \)'s are
\[ \hat{M}_{\lambda,n}^{(k)}(w_{0}^{(k)}) = [w_{0}^{(k)}]^T K_{\lambda} w_{0}^{(k)} + \frac{1}{n^2} \sum_{i=1}^{n} \sum_{j=1}^{n} \{[\nabla x_i, \nabla x_j K_{\lambda}(X_i, X_j)] e_k \} , \]

where \( \nabla K_{\lambda}^{(k)} \) is the \( k \)th column of \( \nabla K_{\lambda} \) which was introduced in Section 3.1 and \( e_k \) is the canonical basis vector with 1 at coordinate \( k \). So from Equation (16), \( \hat{M}_{\lambda,n}^{(1)}(w_{0}^{(1)}) + \hat{M}_{\lambda,n}^{(2)}(w_{0}^{(2)}) \). Similarly, denote

\[ M_{\lambda}^{(k)}(\hat{w}_{\lambda,n}^{(k)}) = \mathbb{E}[D_{\lambda}(X_i, X_n)], \]

where \( D_{\lambda}(X_i, X_n) = \{w_{k,n} - \hat{w}_{\lambda,n}^{(k)}(n)\} K_{\lambda}(X_i, X_n)\{w_{k,i} - \hat{w}_{\lambda,n}^{(k)}(i)\} \). Notice that from the definition of the KSD in Equation (15), \( \hat{M}_{\lambda,n}^{(1)}(\hat{w}_{\lambda,n}^{(1)}) + \hat{M}_{\lambda,n}^{(2)}(\hat{w}_{\lambda,n}^{(2)}) \). Finally, for any two sequences \( a_n \) and \( b_n \), we will use \( a_n \lesssim b_n \) to denote \( a_n/b_n = O(1) \) as \( n \to \infty \).

**Lemma 3.** Under Assumption 2, the following statements hold:

(a) We have,
\[ \Delta_{\lambda,n}^{(1)}(i, k) \lesssim \sqrt{\{\log n\}^2 \Delta_{\lambda,n}(i, k)} + \frac{1}{n} \text{ as } n \to \infty. \]

(b) For any \( \lambda > 0 \) and some constant \( c > 0 \),
\[ \bar{\Delta}_{\lambda,n}(i, k) \lesssim \lambda^{-2} M_{\lambda}^{(k)}(\hat{w}_{\lambda,n+1}^{(k)}) + \lambda^2 (\log n)^4 + \lambda (\log n) \Delta_{\lambda,n}^{(2)}(i, k) + \frac{\lambda^{-2}}{n c \log n}. \]

(c) If \( K_{\lambda}(\cdot, \cdot) \) is the RBF kernel with bandwidth parameter \( \lambda \in \Lambda \) and \( \Lambda \) is a compact subset of \( \mathbb{R}^+ \) bounded away from zero, then
\[ M_{\lambda}^{(k)}(\hat{w}_{\lambda,n}^{(k)}) \leq \frac{\mathbb{E}\{w_{k,i}\}^2 - \mathbb{E}\{\hat{w}_{\lambda,n}^{(k)}(i)\}^2}{n - 1}. \]

The proof of Lemma 3 is available in Section B.5. We will now begin the proof of Theorem 2.

The main idea of the proof of Theorem 2 is to construct an appropriate bound for \( \Delta_{\lambda,n}^{(2)}(i, k) \). As we will see, this bound depends on \( \Delta_{\lambda,n}^{(1)}(i, k) \), \( \bar{\Delta}_{\lambda,n}(i, k) \) and \( M_{\lambda}^{(k)}(\hat{w}_{\lambda,n+1}^{(k)}) \). Once this bound is established, Markov’s inequality will provide the desired convergence in probability in the
statement of Theorem 2. Our approach for analyzing $\Delta^{(2)}_{\lambda,n}(i,k)$ relies on the technique developed in Luo et al. (2023) for proving their Theorem 1. We first note that since the density $f$ is a convolution with a Gaussian density, there exists some constant $c > 0$ such that for all large $\|X_i\|_2, |w_{k,i}|/\|X_i\|_2 \leq c$. Furthermore, we will consider a constrained version of our optimization problem such that the solution satisfies $|\hat{w}_{\lambda,n}(i)| = O(||X_i||_2)$. So, $|\hat{w}_{\lambda,n}(i) - w_{k,i}|$ is bounded above by $O(||X_i||_2)$. Thus,

$$E[\left\{\hat{w}_{\lambda,n}(i) - w_{k,i}\right\}^2 \|X_i\|_2 \leq 2c_0 \log n] \lesssim \Delta^{(1)}_{\lambda,n}(i,k) \log n,$$  \hspace{1cm} (29)

where $c_0 > 0$ is a constant. Similarly,

$$E[\left\{\hat{w}_{\lambda,n}(i) - w_{k,i}\right\}^2 \|X_i\|_2 > 2c_0 \log n] \lesssim E\left[||X_i||^2_2 \|X_i\|_2 > 2c_0 \log n\right].$$  \hspace{1cm} (30)

Since $X_i$ is sub-exponential, the right hand side of Equation (30) is bounded by $O(n^{-1})$. Therefore, from equations (29) and (30),

$$\Delta^{(2)}_{\lambda,n}(i,k) \lesssim \Delta^{(1)}_{\lambda,n}(i,k) \log n + \frac{1}{n}. \hspace{1cm} (31)$$

But, from statements (a) and (b) of Lemma 3

$$\Delta^{(1)}_{\lambda,n}(i,k) \lesssim \sqrt{(\log n)^2 \left[\lambda^{-2} M_{\lambda}^{(k)}(\hat{w}_{\lambda,n+1}) + \lambda^2 (\log n)^4 + \lambda \log n \Delta^{(2)}_{\lambda,n}(i,k) + \frac{\lambda^{-2}}{n^c \log n}\right] + \frac{1}{n}}. \hspace{1cm} (32)$$

So, using Equation (31) and $\lambda \asymp n^{-1/3}$ in Equation (32) we get

$$\Delta^{(1)}_{\lambda,n}(i,k) \lesssim \sqrt{(\log n)^2 \left[n^{2/3} M_{\lambda}^{(k)}(\hat{w}_{\lambda,n+1}) + n^{-2/3}(\log n)^4 + n^{-1/3}(\log n)^2 \Delta^{(1)}_{\lambda,n}(i,k)\right]}.$$

(33)

Now, from statement (c) of Lemma 3, we have

$$M_{\lambda}^{(k)}(\hat{w}_{\lambda,n+1}) \leq \frac{E\{w_{k,i}\}^2 - E\{\hat{w}_{\lambda,n}(i)\}^2}{n}. \hspace{1cm} (34)$$

Note that in Equation (34), the numerator $E\{w_{k,i}\}^2 - E\{\hat{w}_{\lambda,n}(i)\}^2 \lesssim \Delta^{(1)}_{\lambda,n+1}(i,k) \log n$ using the
arguments developed for equations (29) and (30). Therefore,

\[ M^{(k)}_\lambda(w^{(k)}_{\lambda,n+1}) \lesssim \frac{\log n}{n} \Delta^{(1)}_{\lambda,n+1}(i,k). \]

Substituting the above display in Equation (33) we get

\[ \Delta^{(1)}_{\lambda,n}(i,k) \lesssim \sqrt{(\log n)^2 \left[ n^{-1/3} \log n \Delta^{(1)}_{\lambda,n+1}(i,k) + n^{-2/3}(\log n)^4 + n^{-1/3}(\log n)^2 \Delta^{(1)}_{\lambda,n}(i,k) \right]}. \]

Since \( \Delta^{(1)}_{\lambda,n}(i,k) \) is bounded, the above display implies that that \( \Delta^{(1)}_{\lambda,n}(i,k) \to 0 \) as \( n \to \infty \).

We will now establish the rate of convergence of \( \Delta^{(1)}_{\lambda,n}(i,k) \). Let \( A_n = \max\{\Delta^{(1)}_{\lambda,n}(i,k), n^{-1/3}(\log n)^4\} \).

Using equations (33) and (34), we can derive the following inequality for all large \( n \):

\[ A_n \leq C (\log n)^{3/2} n^{-1/6} \sqrt{A_{n+1}}, \quad (35) \]

where the constant \( C \) is independent of \( n \). Recursively applying Equation (35) \( s > 0 \) times, we get

\[ A_n \leq \left[ C (\log n)^{3/2} n^{-1/6} \right]^{1 + \cdots + 2^{-s}} A_{n+s+1}. \]

But for large \( n \), \( A_n < 1 \) and this implies \( A_n \leq \left[ C (\log n)^{3/2} n^{-1/6} \right]^{1 + \cdots + 2^{-s}}. \) Therefore, letting \( s \to \infty \), we have \( A_n \leq C (\log n)^{3} n^{-1/3} \) and so \( \Delta^{(1)}_{\lambda,n}(i,k) \lesssim (\log n)^3 n^{-1/3}. \) Moreover, using Equation (31), \( \Delta^{(2)}_{\lambda,n}(i,k) \lesssim (\log n)^4 n^{-1/3}. \) The statement of Theorem 2 then follows using Markov’s inequality and by noting that \( \mathbb{E}\|W_n(\lambda) - W_0\|^2_F = \sum_{i=1}^{n} \{\Delta^{(2)}_{\lambda,n}(i,1) + \Delta^{(2)}_{\lambda,n}(i,2)\}. \)

\[ \Box \]

**B.5 Proof of Lemma 3**

**Proof of statement (a)** – From Cauchy-Schwartz inequality we get,

\[ \mathbb{E}\left[|\hat{w}^{(k)}_{\lambda,n}(i) - w_{k,i}|\mathbb{I}\{\|X_i\|_2 \leq 2c_0 \log n\}\right] \leq \left[c_1 (\log n)^2 \Delta_{\lambda,n}(i,k)\right]^{0.5}, \]

where \( c_1 \) is a constant depending only on \( c_0 \). Also, using arguments similar to that of Equation (30), we have

\[ \mathbb{E}\left[|\hat{w}^{(k)}_{\lambda,n}(i) - w_{k,i}|\mathbb{I}\{\|X_i\|_2 > 2c_0 \log n\}\right] = O(n^{-1}). \]

Thus, the proof of statement (a) then follows by combining the above two displays. \[ \Box \]
Proof of statement (b) – Suppose we have an i.i.d sample \((X_1, \ldots, X_{n+1})\) of size \(n+1\) from \(f\).

Now, the definition of \(D_\lambda(X_i, X_{n+1})\) implies

\[
D_\lambda(X_i, X_{n+1}) = K_\lambda(X_i, X_{n+1})\{w_{k,n+1} - \hat{w}_{\lambda,n+1}^{(k)}(n+1)\}\{w_{k,i} - \hat{w}_{\lambda,n+1}^{(k)}(i)\},
\]

where \(\hat{w}_{\lambda,n+1}^{(k)}(i)\) is the \((i, k)\)th element of the estimated score matrix \(\hat{W}_{n+1}(\lambda)\) that is constructed based on the sample of size \(n+1\). Denote \(\|X_i - X_n\| = \sqrt{(X_i - X_n)^T (X_i - X_n)}\) and fix \(\epsilon = \lambda \log n\). To prove statement (b), we will analyze the following four terms:

\[
I_1 := \mathbb{E}\left[D_\lambda(X_i, X_{n+1})\mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\}\right] \quad (36)
\]

\[
I_2 := \mathbb{E}\left[K_\lambda(X_i, X_{n+1})\{w_{k,i} - \hat{w}_{\lambda,n+1}^{(k)}(i)\}^2\mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\}\right] \quad (37)
\]

\[
I_3 := \mathbb{E}\left[K_\lambda(X_i, X_{n+1})\{w_{k,i} - \hat{w}_{\lambda,n+1}^{(k)}(i)\}^2\mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\}\right] \quad (38)
\]

\[
I_4 := \mathbb{E}\left[\int f(X_i)K_\lambda(X_i, X_{n+1})\{w_{k,i} - \hat{w}_{\lambda,n+1}^{(k)}(i)\}^2\mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\}dX_{n+1}\right]. \quad (39)
\]

For our analyses, we will assume that the eigen values of \(\Omega\) are finite and bounded away from zero. We begin by analyzing the term \(I_1\) in Equation (36). Now,

\[
\mathbb{E}[D_\lambda(X_i, X_{n+1})] - I_1 = \mathbb{E}[D_\lambda(X_i, X_{n+1})\mathbb{I}\{\|X_i - X_{n+1}\| \geq \epsilon\}].
\]

Since \(K_\lambda\) in Equation (10) is the RBF kernel and \(\epsilon = \lambda \log n\), we have \(K_\lambda(X_i, X_{n+1})\mathbb{I}\{\|X_i - X_{n+1}\| \geq \epsilon\} \lesssim n^{-c \log n}\) for some constant \(c > 0\). Thus \(\mathbb{E}[D_\lambda(X_i, X_{n+1})\mathbb{I}\{\|X_i - X_{n+1}\| \geq \epsilon\}] \lesssim \Delta_{\lambda,n+1}^{(2)}(i, k)n^{-c \log n}\). Since \(\Delta_{\lambda,n+1}^{(2)}(i, k)\) is bounded, we have established,

\[
I_1 \lesssim \mathbb{E}[D_\lambda(X_i, X_{n+1})] + O(n^{-c \log n}). \quad (40)
\]

Consider term \(I_2\) in Equation (37). We will show that

\[
I_2 \lesssim I_1 + O(\epsilon^2). \quad (41)
\]

From Assumption 2, the score function \(w_{k,i}\) is \(L_f\) – Lipschitz continuous. Moreover, for small \(\epsilon\)
and when $\|X_i - X_{n+1}\| < \epsilon$, $\hat{w}_{\lambda,n+1}^{(k)}(n + 1)$ is $L_{n,\epsilon}$-Lipschitz continuous since
\[
\left| \hat{w}_{\lambda,n+1}^{(k)}(n + 1) - \hat{w}_{\lambda,n+1}^{(k)}(i) \right| \leq L_{n,\epsilon} \epsilon,
\]
where $\mathbb{E}L_{n,\epsilon}^2 < \infty$. Then $I_2 - I_1$ is bounded above by
\[
\mathbb{E} \left[ \epsilon(L_f + L_{n,\epsilon}) K_{\lambda}(X_i, X_{n+1}) \left| w_{k,i} - \hat{w}_{\lambda,n+1}^{(k)}(i) \right| \mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\} \right].
\]
Using Cauchy-Schwartz inequality, the square of the above display has the following upper bound:
\[
\epsilon^2 \mathbb{E} \left[ (L_f + L_{n,\epsilon})^2 \mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\} \right] \mathbb{E} \left[ K_{\lambda}^2(X_i, X_{n+1}) \left| w_{k,i} - \hat{w}_{\lambda,n+1}^{(k)}(i) \right|^2 \mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\} \right].
\]
But, $\mathbb{E}[(L_f + L_{n,\epsilon})^2 \mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\}]$ is bounded above by
\[
C_f \frac{\pi}{\Gamma(2)} \epsilon^2 \mathbb{E} [(L_f + L_{n,\epsilon})^2],
\]
where $\Gamma(\cdot)$ is the Gamma function. Finally, noticing that $K_{\lambda}^2(X_i, X_{n+1}) \leq K_{\lambda}(X_i, X_{n+1})$, we have
\[
I_2 \lesssim I_1 + \epsilon \sqrt{\epsilon \Delta_{\lambda,n+1}^{(2)}(i, k)},
\]
which establishes Equation (41).

Similarly the difference between $I_3$ (Equation (38)) and $I_2$ can be bounded by $4\epsilon^2 \mathbb{E}[L_{n,\epsilon}^2 K_{\lambda}(X_i, X_{n+1}) \mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\}]$, which implies that
\[
I_3 \lesssim I_2 + O(\epsilon^4). \tag{42}
\]
Next, we will show that for the term $I_4$ in Equation (39), the following holds:
\[
\lambda^2 \Delta_{\lambda,n}(i, k) \lesssim I_4 \lesssim I_3 + \lambda^3 \Delta_{\lambda,n}^{(2)}(i, k) \log n. \tag{43}
\]
This inequality along with equations (42), (41) and (40) will establish statement (b).

First, note that
\[
I_4 = \mathbb{E} \left[ \left\{ w_{k,i} - \hat{w}_{\lambda,n}^{(k)}(i) \right\}^2 f(X_i) \right] \int K_{\lambda}(X_i, X_{n+1}) \mathbb{I}\{\|X_i - X_{n+1}\| < \epsilon\} dX_{n+1},
\]
since \( \hat{w}_{\lambda,n}(i) \) is independent of \( X_{n+1} \) for \( i < n + 1 \). Now, with \( \epsilon = \lambda \log n \), we have

\[
\int K_\lambda(X_i, X_{n+1})I\{\|X_i - X_{n+1}\| < \epsilon\}dX_{n+1} \geq c\lambda^2, \tag{44}
\]

for some constant \( c > 0 \) when \( n \) is large. So, Equation (44) implies \( \lambda^2 \Delta_{\lambda,n}(i,k) \leq I_4 \) and establishes the lower bound in Equation (43). To establish the upper bound we will assume that the density \( f \) is \( L_f - \)Lipschitz continuous (Assumption 2). We have,

\[
f(X_i) \int K_\lambda(X_i, X_{n+1})I\{\|X_i - X_{n+1}\| < \epsilon\}dX_{n+1} \lesssim \int K_\lambda(X_i, X_{n+1})I\{\|X_i - X_{n+1}\| < \epsilon\}f(X_{n+1})dX_{n+1} + L_f \epsilon \int K_\lambda(X_i, X_{n+1})I\{\|X_i - X_{n+1}\| < \epsilon\}dX_{n+1}.
\]

Then, pre-multiplying the above display with \( \{w_{k,i} - \hat{w}_{\lambda,n}(i)\}^2 \) and taking expectations, we get

\[
I_4 - I_3 \leq L_f \epsilon \Delta_{\lambda,n}^{(2)}(i,k) \int K_\lambda(X_i, X_{n+1})I\{\|X_i - X_{n+1}\| < \epsilon\}dX_{n+1} \leq L_f \epsilon \Delta_{\lambda,n}^{(2)}(i,k) \lambda^2 C,
\]

for some constant \( C \) and \( n \) large. This implies, \( I_4 \leq I_3 + \lambda^3 \Delta_{\lambda,n}^{(2)}(i,k) \) and establishes the upper bound in Equation (43). \( \blacksquare \)

Proof of statement (c) – We first note that \( \hat{M}^{(k)}(\hat{w}_{\lambda,n}(i)) \) and \( \hat{M}^{(k)}(w_0^{(k)}) \) are, respectively V-statistics of \( M^{(k)}_\lambda(\hat{w}_{\lambda,n}(i)) \) and \( M^{(k)}_\lambda(w_0^{(k)}) \). Moreover, we have

\[
\hat{M}^{(k)}(\hat{w}_{\lambda,n}(i)) \leq \hat{M}^{(k)}(w_0^{(k)}).
\]

So, taking expectations on both sides of the above display we get

\[
\frac{n^2 - n}{n^2} M^{(k)}_\lambda(\hat{w}_{\lambda,n}(i)) + \frac{1}{n} \mathbb{E}\{\hat{w}_{\lambda,n}(i)\}^2 \leq \frac{n^2 - n}{n^2} M^{(k)}_\lambda(w_0^{(k)}) + \frac{1}{n} \mathbb{E}\{w_{k,i}\}^2.
\]

But \( M^{(k)}_\lambda(w_0^{(k)}) = 0 \) and so from the above display, we have

\[
M^{(k)}_\lambda(\hat{w}_{\lambda,n}(i)) \leq \frac{\mathbb{E}\{w_{k,i}\}^2 - \mathbb{E}\{\hat{w}_{\lambda,n}(i)\}^2}{n - 1},
\]

which completes the proof of statement (c). \( \blacksquare \)
B.6 Proof of Theorem 3

We first state two lemmata that are needed for proving Theorem 3. Denote \( c_0, c_1, \ldots \) some generic positive constants which may vary in different statements.

**Lemma 4.** If Assumption 3 holds, then with probability tending to 1, \( C_1 / \log n \leq \tau \leq C_2 \log n \) and \( |\mu| \leq C_3 \log n \) for some positive constants \( C_1, C_2 \) and \( C_3 \).

**Lemma 5.** Consider Model (1). Suppose Assumption 3 holds. Then with probability tending to 1, \( \mathbb{E}(\tau_i|y_i, s_i^2) \geq c_0 \log n \).

Lemmata 4 and 5 are proved in Sections B.6.1 and B.6.2, respectively. We now prove Theorem 3.

To establish the first part of Theorem 3, note that

\[
\frac{1}{n} \left\| \delta_n^{ds}(\lambda) - \delta_{(1)} \right\|^2 = \frac{1}{nm^2} \sum_{i=1}^{n} \frac{w_{1,i}}{\tilde{\tau}_i} - \frac{\hat{w}_{\lambda,n}(i)}{\tilde{\tau}_{i,n}(\lambda)} \right\|^2 \leq \frac{2}{nm^2} \sum_{i=1}^{n} \left| w_{1,i} - \hat{w}_{\lambda,n}(i) \right|^2 + \frac{2}{nm^2} \sum_{i=1}^{n} \left| w_{1,i} \right|^2 \frac{1}{\tilde{\tau}_{i,n}(\lambda)} - \frac{1}{\tilde{\tau}_i} \right|^2.
\]

Consider the first term \( T_1 \). From the discussion in Section 3.2, there is a positive constant \( c_0 \) such that \( \tilde{\tau}_{i,n}(\lambda) > c_0 > 0 \) for all \( i = 1, \ldots, n \). It follows that for some constant \( c_1 > 0 \) depending on the fixed \( m \),

\[
T_1 \leq \frac{c_1}{n} \left\| w^{(1)}_0 - \hat{w}_{\lambda,n}^{(1)} \right\|^2 \leq \frac{c_1}{n} \left\| W_0 - \hat{W}_n(\lambda) \right\|^2_F, \tag{45}
\]

where \( \hat{w}_{\lambda,n}^{(1)} \) and \( w^{(1)}_0 \) are, respectively, the first column of \( \hat{W}_n(\lambda) \) and \( W_0 \). From Theorem 2 the last term on the right hand side of the inequality in equation (45) is \( O_p\{n^{-1/3}(\log n)^4\} \).

Next consider the second term \( T_2 \). We have

\[
T_2 \leq \frac{c_2}{n} \sum_{i=1}^{n} \left| \frac{w_{1,i}}{\tilde{\tau}_i} \right|^2 \left| w_{2,i} - \hat{w}_{\lambda,n}(i) \right|^2. \tag{46}
\]

We will use Lemma 5 to bound the terms \( |w_{1,i}/\tilde{\tau}_i| \) in equation (46). First note that since \( X_i \) is sub-exponential, \( |w_{1,i}| \leq c_1 \log n \) with high probability from the discussion in the proof of
Theorem 2. Next, Model (1), Assumption 3 and Lemma 4 imply that with probability tending to 1, \(|Y_i| \leq c_2 \log n\) and \((m - 1)n^{-1} \leq (m - 1)S_i^2 \tau_i \leq (m - 1) + 2\sqrt{(m - 1) \log n} + 2 \log n\) [cf. Lemma 1 of Laurent and Massart (2000)]. So conditional on these events, we have, from Lemma 5, \(|w_{1,i}/\hat{\pi}_i^\tau| \leq c_3 (\log n)^2\). Thus,

\[
T_2 \leq \frac{c_4 (\log n)^4}{n} \sum_{i=1}^{n} \left| w_{2,i} - w_{\lambda,n}(i) \right|^2,
\]

which is \(O_p \{n^{-1/3} (\log n)^8\}\) from Theorem 2. Thus \(n^{-1} \| \delta_n^{ds}(\lambda) - \delta_{(1)}^\tau \|_2^2\) is \(O_p \{n^{-1/3} (\log n)^8\}\).

Now we will prove the second part of Theorem 3. Observe that \(|l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau) - l_n^{(1)}(\mu, \delta_n^{ds}(\lambda); \tau)|\) equals

\[
\left| \sqrt{l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau)} - \sqrt{l_n^{(1)}(\mu, \delta_n^{ds}(\lambda); \tau)} \right| = \sqrt{l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau)} + \sqrt{l_n^{(1)}(\mu, \delta_n^{ds}(\lambda); \tau)}
\]

and Triangle inequality implies

\[
\left| \sqrt{l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau)} - \sqrt{l_n^{(1)}(\mu, \delta_n^{ds}(\lambda); \tau)} \right| \leq \sqrt{\frac{1}{n} \sum_{i=1}^{n} \tau_i \left| \delta_{i,(1)}^\tau - \delta_{i,n}^{ds}(\lambda) \right|^2} \leq c_0 \sqrt{\frac{\log n}{n}} \| \delta_n^{ds}(\lambda) - \delta_{(1)}^\tau \|_2,
\]

where the last inequality in Equation (47) follows from Lemma 4. Thus, from the first part of Theorem 3 and Lemma 4, the quantity on the right hand side of the inequality in equation (47) is \(O_p \{(\log n)^{9/2} n^{-1/6}\}\). Thus, it follows from equation (47) that

\[
\sqrt{l_n^{(1)}(\mu, \delta_n^{ds}(\lambda); \tau)} \leq \sqrt{l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau)} + O_p \{(\log n)^{9/2} n^{-1/6}\}, \quad \text{and}
\]

\[
\left| l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau) - l_n^{(1)}(\mu, \delta_n^{ds}(\lambda); \tau) \right| \leq 4 \sqrt{l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau)} \left| \sqrt{l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau)} - \sqrt{l_n^{(1)}(\mu, \delta_n^{ds}(\lambda); \tau)} \right| \left\{ 1 + o_p(1) \right\}.
\]

Now \(\delta_{(1)}^\tau\) is the Bayes estimator of \(\mu\) under the weighted squared error loss and so its risk \(\mathbb{E} l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau) < \infty\). This implies \(l_n^{(1)}(\mu, \delta_{(1)}^\tau; \tau)\) is \(O_p(1)\). Thus, from equation (47), the first part of Theorem 3 and the display above, we have the desired result.
B.6.1 Proof of Lemma 4

The proof of Lemma 4 follows directly from Assumption 3 and Markov’s inequality. For example, fix a $\nu > 0$ and note that, for $r = \epsilon_2^{-\nu} > 1$,

$$\mathbb{P}\left(\tau \leq \frac{\epsilon_2^{1+\nu}}{\log n}\right) \leq \frac{\mathbb{E}_H\left\{\exp(\epsilon_2/\tau)\right\}}{n^r}.$$  ■

B.6.2 Proof of Lemma 5

Recall that $f(y_i, s_i^2) = \int_{\mathbb{R}^+} f_1(y_i|\mu, \tau)f_2(s_i^2|\tau)g(\mu|\tau)h(\tau)d\mu d\tau$, where $g(\cdot|\tau)$ and $h(\cdot)$ are, respectively, the density functions associated with the distribution functions $G_\mu(\cdot|\tau)$ and $H_\tau(\cdot)$ in Model (1), $f_1$ is the density of a Gaussian random variable with mean $\mu$ and variance $1/(m\tau)$ and $f_2$ is the density of $S^2$ where $(m-1)S^2\tau \sim \chi^2_{m-1}$.

We will first analyze the behavior of $f(y_i, s_i^2)$. Gaussian concentration implies that with high probability $\{(Y_i - \mu_i)^2m\tau_i\} \leq 2\log n$ and so, conditional on this event,

$$f_1(y_i|\mu, \tau) \geq c_0\sqrt{\tau}/n.$$  (48)

Moreover, using the Chi-square concentration in Lemma 1 of Laurent and Massart (2000), $(m-1)S_i^2\tau_i \leq (m-1) + 2\sqrt{(m-1)\log n} + 2\log n$ and $S_i^2\tau_i \geq n^{-1}$ with high probability. It follows that

$$f_2(s_i^2|\tau) \geq c_1\tau a_n^{1/(m-1)/2},$$  (49)

conditional on this event, where $a_n = \exp\{-\sqrt{(m-1)\log n}\}$. Using equations (48) and (49), we have

$$f(y_i, s_i^2) \geq c_2\frac{a_n^{1/(m+1)/2}}{n^{(m+1)/2}} \int_{\mathbb{R}^+} \tau^{3/2}h(\tau)d\tau.$$  (50)

Now, use Assumption 3 and Lemma 4 on the quantity $\int_{\mathbb{R}^+} \tau^{3/2}h(\tau)d\tau$ in equation (50) to conclude that

$$\int_{\mathbb{R}^+} \tau^{3/2}h(\tau)d\tau \geq \frac{c_3}{(\log n)^{3/2}}\mathbb{P}\left(\tau \geq C_2/\log n\right).$$  (51)
So, with equations (51), (50) and Lemma 4, we have with high probability,

\[ f(y_i, s_i^2) \geq c_4 \frac{a_n}{n^{(m+1)/2} (\log n)^{3/2}}. \] (52)

Now we proceed to prove the statement of Lemma 5. Fix \( \nu > 0 \) such that \( \varepsilon^{-\nu} > 2m \) in Lemma 4 and let \( C_1 = \varepsilon^{1+\nu} \). First note that from Assumption 3, \( \mathbb{P}(\tau \leq C_1/\log n) \leq c_0/n^{2m} \).

Now, Markov’s inequality implies,

\[ \mathbb{E}(\tau | y_i, s_i^2) \geq \frac{C_1}{\log n} \left\{ 1 - \mathbb{P}\left[ \tau \leq C_1(\log n)^{-1} | y_i, s_i^2 \right] \right\}. \]

Moreover,

\[ \mathbb{P}(\tau \leq \frac{C_1}{\log n} | y_i, s_i^2) = \left\{ f(y_i, s_i^2) \right\}^{-1} \int_0^{C_1/\log n} h(\tau) \int_{\mathbb{R}} g(\mu | \tau) f_1(y_i | \mu, \tau) f_2(s_i^2 | \tau) d\mu d\tau. \]

Now \( f_1(y_i | \mu, \tau) \leq c_0 \sqrt{\tau} \) and \( f_2(s_i^2 | \tau) \leq c_1 \tau \) where \( c_0, c_1 > 0 \) are constants. So for some positive constant \( c_2 \),

\[ \mathbb{P}(\tau \leq C_2/\log n | y_i, s_i^2) \leq \frac{c_2}{f(y_i, s_i^2)} \int_0^{C_2/\log n} \tau^{3/2} h(\tau) \int_{\mathbb{R}} g(\mu | \tau) d\mu d\tau \leq \frac{c_3}{f(y_i, s_i^2)(\log n)^{3/2}} \int_0^{C_2/\log n} h(\tau) d\tau = \frac{c_3}{f(y_i, s_i^2)(\log n)^{3/2}} \mathbb{P}(\tau \leq C_2/\log n). \]

Thus, from the above display, Assumption 3 and Lemma 4,

\[ \mathbb{E}(\tau | y_i, s_i^2) \geq \frac{C_2}{\log n} \left\{ 1 - \frac{c_3}{f(y_i, s_i^2)(\log n)^{3/2}} \right\}. \]

Finally, equation (52) and the above display prove the statement of Lemma 5. \( \blacksquare \)

**B.7 Proof of Equations (7) and (8)**

The proof follows by first recalling that under the hierarchical model of Equation (1),

\[ f_{m_i}(y_i, s_i^2 | \mu_i, \tau_i) \propto \exp\left\{-\frac{\tau_i}{2} [m_i y_i^2 + (m_i - 1) s_i^2] + m_i \tau_i y_i - \frac{m_i}{2} \tau_i \mu_i^2 + \frac{m_i - 3}{2} \log s_i^2 \right\}. \]
Therefore from Bayes theorem,
\[
f_{m_i}(\mu_i, \tau_i|y_i, s_i^2) = \frac{f_{m_i}(y_i, s_i^2|\mu_i, \tau_i)}{f_{m_i}(y_i, s_i^2)} g(\mu_i|\tau_i) h(\tau_i) \propto \exp \left\{ \eta_i^T T(\mu_i, \tau_i) - A(\eta_i) \right\} g(\mu_i|\tau_i) h(\tau_i).
\]

Here \( \eta_i = (m_iy_i, -m_iy_i^2 - (m_i - 1)s_i^2) := (\eta_{1i}, \eta_{2i}) \), \( T(\mu_i, \tau_i) = (\tau_i\mu_i, \tau_i/2) \) and
\[
A(\eta_i) = -0.5(m_i - 3) \log \gamma(\eta_{1i}, \eta_{2i}) + \log f_{m_i} \{ m_i^{-1} \eta_{1i}, \gamma(\eta_{1i}, \eta_{2i}) \}, \\
\gamma(\eta_{1i}, \eta_{2i}) = \frac{-\eta_{2i} - m_i^{-1} \eta_{1i}^2}{m_i - 1},
\]
with \( f_m(y, s^2) = \int \int f_m(y, s^2|\mu, \tau) g_\mu(\mu|\tau) h_\tau(\tau)d\mu d\tau \) being the marginal density function of \( (Y, S^2) \). So the posterior distribution of \( (\mu_i, \tau_i) \) belongs to a 2-parameter exponential family and using the properties of exponential family distributions we have, dropping subscript \( i \),
\[
\hat{\tau}_i^\pi := \hat{\tau}_i^\pi(y, s^2, m) = \mathbb{E}(\tau|y, s^2, m) = 2 \frac{\partial A(\eta)}{\partial \eta_2} = \frac{m - 3}{(m - 1)s^2} - \frac{2}{m - 1} w_2(y, s^2; m).
\]
Furthermore, with \( \zeta = \tau \mu \),
\[
\hat{\zeta}_i^\pi := \hat{\zeta}_i^\pi(y, s^2, m) = \mathbb{E}(\zeta|y, s^2, m) = \frac{\partial A(\eta)}{\partial \eta_1} = \frac{(m - 3)y}{(m - 1)s^2} + m^{-1} w_1(y, s^2; m) - \frac{2}{m - 1} y w_2(y, s^2; m)
\]
\[
= y \mathbb{E}(\tau|y, s^2, m) + m^{-1} w_1(y, s^2; m).
\]

\[\blacktriangle\]

C Additional Numerical Experiments

C.1 Compound estimation of Normal means under squared error loss

We focus on the hierarchical Model of Equation (1) and compare six approaches for estimating \( \mu \) under the squared error loss when the variances \( \sigma_i = 1/\tau_i \) are assumed to be unknown. These approaches can be categorized into three types: the first consists of the NEST method (NEST Orc. \( \lambda \)), which estimates \( \lambda \) by minimizing the true loss, the proposed data-driven NEST method and Tweedie’s formula (TF) that uses sample variances. For both NEST and TF, \( \lambda \) is chosen using the modified cross-validation approach described in Section 3.2 with \( \vartheta_n(\lambda; U, V) \) defined
as
\[ \vartheta_n(\lambda; \mathcal{U}, \mathcal{V}) = \frac{1}{n} \sum_{i=1}^{n} \left\{ \bar{V}_i - \delta_i^{ds}(\bar{U}_i; \mathcal{U}, \lambda) \right\}^2, \]
to reflect the squared error loss. The second are linear shrinkage methods: the group linear estimator (Grp Linear) of Weinstein et al. (2018) and the semi-parametric linear shrinkage rule (Jing.SM) from Jing et al. (2016b). Finally, the third type is the g-modelling approach of Gu and Koenker (2017a,b). For Grp. Linear we use code provided by Weinstein et al. (2018) while for Jing.SM we write our own routine in R. We continue to rely on the function WGLVmix in the R package REBayes (Koenker and Gu, 2017) for NPMLE. We note that amongst the six methods considered here, NEST and NEST Orc. \( \lambda \) are designed to estimate the means under the weighted squared error loss while the remaining four approaches target the squared error loss. Furthermore, with the exception of Grp. Linear, all other methods considered here estimate \( \mu \) when the variances are unknown. Grp. Linear, on the other hand, assumes full knowledge of the unknown variances for shrinkage estimation of the means and here we use sample variances for its implementation.

The aforementioned six approaches are evaluated on five different simulation settings, with the goal of assessing the relative performance of the competing estimators as the heterogeneity in the variances \( \sigma_i^2 \) is varied while keeping the sample sizes \( m_i \) fixed at \( m \). The five simulation settings can be categorized into three types: a setting where mean and variances are independent; three settings where mean and variance are correlated; and a setting that represents departure from the Normal data-generating model. For each setting we set \( n = 1,000 \) and compute the average squared error risk for each competing estimator of \( \mu \) across 50 Monte Carlo repetitions. Figures 8 to 12 plot the relative risk which is the ratio of the average squared error risk for any competing estimator to that of oracle Bayes estimator \( \delta^{\pi}_{(0)} \) of \( \mu \) (Equation (5)) so that a ratio bigger than 1 represents a poorer risk performance of the competing estimator relative to the Bayes oracle.

The first setting, Figure 8, corresponds to the independent case. Here, for each \( i = 1, \ldots, n \), \( \mu_i \overset{i.i.d.}{\sim} 0.7 N(0, 0.1) + 0.15 N(1, 3) + 0.15 N(-1, 3) \) and \( \sigma_i^2 \overset{i.i.d.}{\sim} U(0.5, u) \) where we let \( u \) vary across five levels, \( \{1, 2, 3, 4, 5\} \). The three plots in Figure 8 show the relative risks as \( u \) varies for \( m = 10, 15 \) and 20 (left to right). We see that for \( m = 10 \), the competing methods split into two levels of performance. The group with the lowest relative risks consists of NPMLE,
Figure 8: Comparison of relative risks when \((\mu_i, \sigma^2_i)\) are independent. Here \(\mu_i \overset{i.i.d}{\sim} 0.7 N(0, .1) + 0.15 N(1, 3) + 0.15 N(-1, 3)\) and \(\sigma^2_i \overset{i.i.d}{\sim} U(0.5, u)\). Plots show \(m = 10, 15, 20\) left to right.

TF and NEST while the two linear shrinkage methods exhibit substantially higher relative risks. Moreover, we also see that as heterogeneity increases with increasing \(u\), the gap between the two groups’ relative risks increases, indicating that NPMLE, TF and the proposed NEST method are particularly useful for compound estimation of normal means when the variances are unknown and heterogeneous, and the sample size for estimating those variances are themselves small. As \(m\) increases, the performance of the two linear shrinkage methods and TF improve which is expected as there are now more replicates per unit of study to construct a relatively reliable estimate of the unknown variances. However, the performance of NEST improves too and particularly at \(m = 20\) (Figure 8 right), NPMLE exhibits a slightly higher relative risk than NEST and TF.

The second setting, Figure 9, corresponds to the correlated case. The precisions \(\tau_i = 1/\sigma^2_i\) are generated independently from a gamma mixture, with an even chance of drawing \(\Gamma(20, \text{rate} = 20)\) or \(\Gamma(20, \text{rate} = u)\) and given \(\tau_i\), the means \(\mu_i\) are independently \(0.5 \ N(0.5/\tau_i, 0.5^2) + 0.5 \ N(-0.5/\tau_i, 0.5^2)\). In this setting, the magnitude of the variances increase with \(u\) and the means grow with the variances. We note from Figure 9 that Grp, Linear and Jing.SM exhibit improved performance particularly for small values of \(u\). As \(u\) increases, TF, NPMLE and NEST perform well although their relative risk profiles are substantially away from 1 as \(u\) increases. For TF and NEST this behavior is expected given the statements of Propositions 1 and 2. The improved performance of Jing.SM in this setting is potentially related to the observation that when \(u\) is small, the rate mixture of Gamma distributions on \(\tau_i\) can be well approximated by a single Gamma distribution and that coincides with the parametric prior that Jing et al. (2016b) use on the precision to derive their empirical Bayes estimator for the means.
In the third setting, Figure 10, \((\mu_i, \tau_i)\) continue to be correlated and have a conjugate prior distribution under Model (1). The precisions \(\tau_i\) are drawn from \(\Gamma(20, \text{rate} = u)\) and conditional on \(\tau_i\), \(\mu_i\) are independently \(N(0, 0.5/\tau_i)\). Under this data generating scheme, the posterior mean of \(\mu_i\) is \(my_i/(m+2)\) which is independent of \(u\). This is the reason that the relative risks of the competing estimators in Figure 10 do not vary with the heterogeneity in the variances. Compared to the first two settings, we see that the linear shrinkage estimators have a relatively better performance and Jing.SM dominates all other shrinkage estimators. This is expected because in this setting the posterior mean of \(\mu_i\) is indeed a linear function of the sample mean \(y_i\). For \(m = 10\) and \(15\), we notice that the relative risk of NEST is marginally better than the competing estimators while at \(m = 20\) Grp Linear and NEST have similar risk performance.

Figure 11 presents the fourth setting where the precisions \(\tau_i\) are drawn from the gamma
Figure 11: Comparison of relative risks when \( \mu \) is sparse. Here \( \tau_i \overset{i.i.d.}\sim 0.5 \Gamma(20, \text{rate} = 20) + 0.5 \Gamma(20, \text{rate} = u) \) and \( \mu_i|\tau_i \overset{\text{ind.}}\sim 0.7 N(0, 0.01) + 0.15 N(0.5/\tau_i, 1) + 0.15 N(-0.5/\tau_i, 1) \). Plots show \( m = 10, 15, 20 \) left to right.

Figure 12: Comparison of relative risk for non-normal data. Here \( Y_{ij}|\mu_i, \sigma_i^2 \overset{i.i.d.}\sim U(\mu_i - \sqrt{3}\sigma_i^2, \mu_i + \sqrt{3}\sigma_i^2), \sigma_i^2 \) are sampled independently from \( N(u, 1) \) truncated below at 0.1 and \( \mu_i|\sigma_i^2 \overset{\text{ind.}}\sim 0.8 N(0.25\sigma_i^2, 0.25) + 0.2 N(\sigma_i^2, 1) \). Plots show \( m = 10, 15, 20 \) left to right.

A mixture of Setting 2 and \( \mu_i|\tau_i \overset{\text{ind.}}\sim 0.7 N(0, 0.01) + 0.15 N(0.5/\tau_i, 1) + 0.15 N(-0.5/\tau_i, 1) \). We see a similar pattern to that in Figure 8 at \( m = 10 \). For \( m = 15 \) and 20, we notice that the relative risks of the linear shrinkage methods are now higher than their levels at \( m = 10 \). This is not surprising for in this setting, while the risk performance of all methods have improved with larger sample sizes, NPMLE, TF and NEST exhibit a bigger improvement in risk than those of Grp Linear and Jing.SM.

The fifth setting, Figure 12, corresponds to the setting where the data \( Y_{ij}|(\mu_i, \sigma_i^2) \) are not normally distributed. Here the proposed NEST method demonstrates robustness to departures from the Normal model particularly in comparison to TF, Grp. Linear and Jing.SM.

Overall, the results of the preceding five simulation settings corroborate the statement of
(a) Simulation settings 4 (left) and 7 (right) from Section 5 where estimation is conducted under the weighted squared error loss.

(b) Simulation settings 2 (left) and 5 (right) from Section C.1 where estimation is conducted under the squared error loss.

Figure 13: Compound estimation of means under small and unequal sample sizes \( m \). Here \( m \) is a fixed vector of size \( n \) with elements sampled randomly from \((4, 5, 6)\) with replacement.

Proposition 2 and reveal that when the variances are unknown, the NEST estimation framework enjoys a relatively better risk performance for estimating the means under the squared error loss than the linear shrinkage methods and Tweedie’s formula that rely on sample variances.

C.2 Numerical Experiments with unequal sample sizes \( m_i \)

In this section, we present the risk performance of the competing approaches of sections 5 and C.1 when the sample sizes \( m_i \) are small and differ across the \( n = 1000 \) units of study. We use the following four simulation settings: Settings 4 and 7 from Section 5 and Settings 2 and 5 from Section C.1. However, we change how \( m = (m_1, \ldots, m_n) \) are generated in these settings.
Figures 13a and 13b present the relative risks of the competing estimators across these four scenarios. Here \( m \) is generated as a fixed vector of size \( n \) with elements sampled randomly from \( \{4, 5, 6\} \) with replacement. The case where \( m \in \{4, 5\} \) represents a particularly challenging scenario for NEST which is based on the following observation: in Equation (8) \( (m_i - 3) ((m_i - 1)S_i^2)^{-1} \) is an unbiased estimator of \( \tau_i \) which follows from the fact that \( ((m_i - 1)S_i^2 \tau_i)^{-1} \) has an inverse Chi-square distribution with \( m_i - 1 \) degrees of freedom. However the variance of this distribution does not exist unless \( m_i > 5 \).

Figure 13a represents Settings 4 (left) and 7 (right) from Section 5. We note that NEST continues to dominate NPMLE under the weighted squared error loss even when \( m_i \) are small. However, at such small sample sizes the risk of the NEST estimator is relatively larger than that of \( \delta_{(1)}^\pi \).

Figure 13b exhibits Settings 2 (left) and 5 (right) from Section C.1 where the estimation is conducted under the squared error loss. Here, we use the semi-parametric monotonically constrained SURE estimator that shrinks towards the grand mean, XKB.SG, from Xie et al. (2012) in place of Jing.SM from Section C.1 as the latter was originally designed for the case \( m_i = m \). While the extension of Jing.SM to unequal \( m_i \) is straightforward, we do not pursue that direction in this article and, instead, use XKB.SG in its place.

In Figure 13b we note that NPMLE dominates NEST and NEST exhibits a relatively better risk performance than TF, Grp. Linear and XKB.SG. However, in Setting 5 (right panel of Figure 13b), where the data \( Y_{ij} | (\mu_i, \sigma^2_i) \) are not normally distributed, the performance of NEST is substantially poorer than NPMLE when \( u \) is large. The two main reasons for this behavior are related to (1) Proposition 2 where the oracle NEST estimator \( \delta_{(1)}^\pi \) is not, in general, the optimal estimator of the means under the squared error loss, and (2) when the sample size \( m_i \) is less than 6 the variance of \( ((m_i - 1)S_i^2 \tau_i)^{-1} \) under Model (1) does not exist.

C.3 Compound Estimation of Ratios

In this section we demonstrate the use of the NEST estimation framework for compound estimation of \( n \) ratios \( \theta_i = \sqrt{m_i \mu_i} / \sigma_i \) which represent a popular financial metric for assessing mutual fund performance (see Section D.2 for a related real data application involving compound estimation of mutual fund Sharpe ratios.). We evaluate the performance of the same six
methods under the squared error loss with $n$ fixed at 1000 and $m_i = m$ for $i = 1, \ldots, n$. We consider two simulation settings and plot the risk performance of the competing estimators of $\theta = (\theta_1, \ldots, \theta_n)$ relative to the optimal Bayes estimator that estimates $\theta$ using the vector of posterior means $\mathbb{E}(\theta_i | y_i, s_i^2)$ for $i = 1, \ldots, n$. Note that amongst the six methods considered here, NPMLE is the only method that is designed to estimate these posterior means. The other methods estimate $\theta_i$ by separately estimating $\mu_i$ and $\sigma_i$, and then take their ratio to construct an estimate. Grp. Linear and TF, in particular, rely on the sample standard deviation for estimating $\sigma_i$ while NEST and Jing.SM employ their respective empirical Bayes estimators of $\sigma_i^2$.

Setting 1 is presented in Figure 14. The data $Y_{ij}$ are generated independently from $N(\mu_i, \sigma_i^2)$, and the variances $\sigma_i^2$ are simulated uniformly between 0.1 and $u$, and the means are independently drawn from a mixture model with half chance $N(-\sigma_i^2/2, 0.5^2)$ and the other half $N(\sigma_i^2/2, 0.5^2)$. We continue to see that NEST has a lower relative risk than Grp. Linear and TF, both of which use sample variances. NEST also dominates Jing.SM for almost all values of $u$ while NPMLE dominates NEST for small values of $u$. As $u$ increases the heterogeneity in the data grows and we see that the relative risks of Grp. Linear and Tweedie’s Formula across all $m$ first decrease and then increase. The shift in the behavior of these estimators is related to the observation that as $u$ increases, the centers of the mixture model that generates $\mu_i$, are on average, further away from one another. This makes estimating the numerator of the ratio easier for all methods up until a point. As heterogeneity increases further, the risks of these methods that use the sample standard deviation in the denominator of $\theta_i$ are relatively worse than the risk of NEST and NPMLE.

Setting 2 is presented in Figure 15 where the means are generated according to Setting 1.
Figure 15: Comparison of relative risk for estimating $\theta$. Here $\tau_i \stackrel{i.i.d.}{\sim} 0.5 \Gamma(20, \text{rate} = 20) + 0.5 \Gamma(20, \text{rate} = u)$ and $\mu_i | \tau_i \sim 0.5 N(0.5 / \tau_i, 0.5^2) + 0.5 N(-0.5 / \tau_i, 0.5^2)$. Plots show $m = 10, 15, 20$ left to right.

but the precisions are independently drawn from a mixture model with half chance $\Gamma(\text{shape} = 20, \text{rate} = 20)$ and the other half $\Gamma(\text{shape} = 20, \text{rate} = u)$. We note that while NEST dominates TF and Grp. Linear, NPMLE and Jing.SM dominate NEST when the heterogeneity is relatively smaller and when $m = 10$. The improved performance of Jing.SM in this setting is potentially related to the observation that when $u$ is small, the rate mixture of Gamma distributions on $\tau_i$ can be well approximated by a single Gamma distribution and that coincides with the parametric prior that Jing et al. (2016b) use on the precision to derive their empirical Bayes estimator for the variances. In Figures 14 and 15 we note that for $m = 10$, the relative risk of NEST is substantially higher than NEST Orc. $\lambda$. This is not unexpected because NEST Orc. $\lambda$ estimates $\lambda$ by minimizing the true squared error loss involving $\theta$ while the data-driven NEST estimator relies on the modified cross-validation approach described in Section 3.2 with $\nu_n(\lambda; \mathcal{U}, \mathcal{V})$ defined as

$$\nu_n(\lambda; \mathcal{U}, \mathcal{V}) = \frac{1}{n} \sum_{i=1}^{n} \left( \bar{V}_i - \delta^d_i(\bar{U}_i; \mathcal{U}, \lambda) \right)^2,$$

to choose $\lambda$. 

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D Real Data Analyses

D.1 Baseball Data

We analyze the monthly data on the number of “at bats” and “hits” for all U.S. Major League baseball players over the regular seasons from 2002 until 2011. In this analysis we focus on both pitchers and non-pitchers using an approach similar to that of Gu and Koenker (2017a). The data are available from the R package REBayes and have been aggregated into half seasons to produce an unbalanced panel. It includes observations on 932 players who have at least ten at bats in any half season and appear in no fewer than five half-seasons (note that there are a total of 20 half-seasons that a player can appear in).

Following Brown (2008), let the transformed batting average \( Y_{ij} \) for player \( i = 1, \ldots, n \) at time \( j = 1, \ldots, m_i \) be denoted by \( Y_{ij} = \arcsin \left( \sqrt{H_{ij} + 0.25} \right) \) where \( H_{ij} \) denotes the number of “hits” and \( N_{ij} \) denotes the number of “at bats” at time \( j \) for player \( i \). We assume that \( Y_{ij} \sim N(\mu_i, v_{ij}^2/\tau_i) \) with \( \mu_i = \arcsin(\sqrt{p_i}) \), \( p_i \) being player \( i \)’s batting success probability, and \( v_{ij}^2 = 1/(4N_{ij}) \). Here \( 1/\tau_i \) are player specific scale parameters as described in Gu and Koenker (2017a). Under this setup, the sufficient statistics are

\[
\hat{\mu}_i = \left( \sum_{j=1}^{m_i} 1/v_{ij}^2 \right)^{-1} \sum_{j=1}^{m_i} Y_{ij}/v_{ij}^2 \sim N(\mu_i, v_i^2/\tau_i) \text{ with } v_i^2 = \left( 4 \sum_{j=1}^{m_i} N_{ij} \right)^{-1},
\]

\[
S_i^2 = \frac{1}{m_i - 1} \sum_{j=1}^{m_i} (Y_{ij} - \hat{\mu}_i)^2/v_{ij}^2 \text{ with } (m_i - 1)S_i^2\tau_i \sim \chi^2_{m_i-1}.
\]

In this analysis, the goal is to use the 2002-2011 data to predict the batting averages of the players in 2012. Players are divided into three categories: all, non-pitchers, and pitchers. We consider the following seven estimators of \( \mu_i \): two non-parametric maximum likelihood based estimators, denoted NPMLE-Indep and NPMLE-Dep which assume, respectively, independent and dependent priors on \((\mu_i, 1/\tau_i)\), the sufficient statistics \( \hat{\mu}_i = (\hat{\mu}_1, \ldots, \hat{\mu}_n) \) of \( \mu \), the grand mean across all players in the 2011 season \( Y^{2011}_i = \sum_{i=1}^{n} Y_{i,2011} \), the proposed NEST estimator and its counterpart with an oracle choice for the tuning parameter \( \lambda \) (NEST orc \( \lambda \)), and the naive estimator that uses 2011 batting averages \( Y^{2011}_i = (Y_{i,2011}, \ldots, Y_{i,2011}) \). To assess how well these methods predict 2012 batting averages \( Y^{2012}_i = (Y_{i,2012}, \ldots, Y_{i,2012}) \), we consider three
Table 1: Performance of the competing estimators relative to the performance of the naive estimator \( Y_{2011} \). Here \( R-TSE(\delta) = TSE(\delta)/TSE(Y_{2011}) \) with similar definitions for R-NSE and R-TSEp. The smallest two relative errors are bolded.

|                | NPMLE-Indep | NPMLE-Dep | \( \hat{\mu} \) | \( Y_{2011} \) | NEST | NEST orc \( \lambda \) |
|----------------|-------------|-----------|-----------------|----------------|------|----------------------|
| All            |             |           |                 |                |      |                      |
| \( n \) for estimation: 932 | R-TSE | 0.463 | 0.491 | 0.352 | 1.958 | **0.350** | **0.350** |
| \( n \) for prediction: 370        | R-NSE | **0.668** | 0.678 | 0.676 | 1.495 | **0.670** | **0.670** |
|                            | R-TSEp | 0.582 | 0.610 | 0.501 | 1.783 | **0.498** | **0.498** |
| Nonpitchers       |             |           |                 |                |      |                      |
| \( n \) for estimation: 792        | R-TSE | 0.535 | 0.559 | 0.503 | 0.551 | **0.488** | **0.484** |
| \( n \) for prediction: 325        | R-NSE | **0.656** | 0.665 | 0.679 | 0.973 | 0.667 | 0.659 |
|                            | R-TSEp | 0.651 | 0.677 | 0.619 | 0.682 | **0.604** | **0.601** |
| Pitchers          |             |           |                 |                |      |                      |
| \( n \) for estimation: 140        | R-TSE | 0.659 | 0.662 | 0.629 | 0.804 | **0.628** | **0.620** |
| \( n \) for prediction: 45         | R-NSE | 0.659 | 0.663 | 0.649 | 0.769 | **0.649** | **0.638** |
|                            | R-TSEp | **0.124** | 0.124 | 0.133 | 0.337 | 0.133 | **0.122** |

Criteria for evaluating any estimate \( \delta_i \) of \( \mu_i \): total squared error from Brown (2008) and defined as:

\[
TSE(\delta) = \sum_{i=1}^{n} \left\{ (Y_{i,2012} - \delta_i)^2 - (4N_{i,2012})^{-1} \right\},
\]

normalized squared error from Gu and Koenker (2017a) and defined as:

\[
NSE(\delta) = \sum_{i=1}^{n} \left\{ 4N_{i,2012}(Y_{i,2012} - \delta_i)^2 \right\},
\]

and total squared error on a probability scale from Jiang et al. (2010) which is defined as:

\[
TSE_p(\hat{\lambda}) = \sum_{i=1}^{n} \left\{ (p_{i,2012} - \hat{p}_i)^2 - p_{i,2012}(1 - p_{i,2012})(4N_{i,2012})^{-1} \right\}.
\]

Here \( \hat{p}_i = \sin^2(\delta_i) \) and \( p_{i,2012} = \sin^2(Y_{i,2012}) \).

In Table 1, we report the performance of the competing estimators relative to the performance of the naive estimator \( Y_{2011} \) wherein \( R-TSE(\delta) = TSE(\delta)/TSE(Y_{2011}) \) with similar definitions for R-NSE and R-TSEp. Thus, a smaller value of R-TSE, R-NSE or R-TSEp indicates a relatively better prediction error. Across “All” and “Nonpitchers”, NEST exhibits the best relative risk for two of the three performance metrics. It is interesting to note that the sufficient statistics \( \hat{\mu} \) are quite competitive in this example while NPMLE with independent priors dominate the one with dependent priors across “All”, “Nonpitchers” and “Pitchers”. The compound estimation problem for “Pitchers” is an example of a setting where \( n \) is relatively small and NEST demonstrates a better risk performance than NPMLE for total squared error and normalized squared error losses.

D.2 Mutual Fund Sharpe Ratios

In this section we analyze a dataset on \( n_1 = 5,000 \) monthly mutual fund returns spanning 12 months from January 2014 to December 2014. This data are sourced from the Wharton research.
data services (Wharton School, 1993). The goal in this analysis is to use Sharpe ratios constructed using the data on the first \( m_1 = 6 \) months, January 2014 - June 2014, to predict the corresponding Sharpe ratios for the next 6 months. Formally, let \( Y_{ij} \) denote the excess return of fund \( i (= 1, \ldots, n_1) \) in month \( j (= 1, \ldots, m_1) \) over the return on the 3 month treasury yield. Denote \( \bar{Y}_i = m_1^{-1} \sum_{j=1}^{m_1} Y_{ij}, \quad S_i^2 = (m_1 - 1)^{-1} \sum_{j=1}^{m_1} (Y_{ij} - \bar{Y}_i)^2 \) and \( \delta_i^{\text{naive}} = \bar{Y}_i / \sqrt{S_i^2} \) to be, respectively, the sample mean, the sample variance and the observed Sharpe ratio of the monthly excess returns.

Of the 5,000 funds available during these first 6 months, there are \( n_2 = 4,958 \) funds that appear in the next 6 months, July 2014 - December 2014, and have at least 3 months of returns available during this period. For our prediction, we consider these \( n_2 \) funds to assess the performance of various estimators for predicting \( \theta_i = \mu_i / \sigma_i \) where \( \mu_i \) and \( \sigma_i \) are the sample mean and sample standard deviation of the excess returns of the \( n_2 \) funds during the next 6 months.

We consider the following estimators of \( \theta = (\theta_1, \ldots, \theta_n) \): NEST, NEST orc. \( \lambda \), Tweedie’s formula (TF), Grp Linear, Jing.SM and NPMLE from Section C.1. Additionally, we consider the SURE estimators XKB.SG and XKB.G from Xie et al. (2012). Note that for predicting \( \theta_i \), TF, Grp Linear, XKB.SG and XKB.G rely on the sample variances \( S_i^2 \). To evaluate the performance of these estimators for predicting \( \theta \), we consider the following three criteria with \( m_{i,2} \in [3, 6] \):

- Total Squared Error: \( TSE(\delta) = \sum_{i=1}^{n_2} (\theta_i - \delta_i)^2 \);
- Weighted Squared Error: \( WSE(\delta) = \sum_{i=1}^{n_2} m_{i,2} (\theta_i - \delta_i)^2 \);
- Weighted Absolute Error: \( WAE(\delta) = \sum_{i=1}^{n_2} m_{i,2} |1 - \delta_i / \theta_i| \).

In Table 2, we present the performance of the competing estimators relative to the performance of the naive estimator \( \delta^{\text{naive}} = (\delta_i^{\text{naive}} : 1 \leq i \leq n) \) so that a smaller value of R-TSE, R-WSE or R-WAE indicates a relatively better prediction error.

Along the performance measures of Total Squared Error and Weighted Squared Error, NEST has the smallest relative risk among all competing estimators considered in this example. With respect to the Weighted Absolute Error, Jing.SM has a substantially smaller relative risk than NEST while Grp Linear appears to be doing relatively worse and exhibits an R-WAE bigger

| \( n_1 \) | \( n_2 \) | Grp Lin. | XKB.G | XKB.SG | NPMLE | Jing.SM | TF | NEST | NEST orc. \( \lambda \) |
|---|---|---|---|---|---|---|---|---|---|
| R - TSE | 5000 | 4958 | 0.896 | 0.931 | 0.922 | 0.842 | 0.720 | 0.997 | \textbf{0.688} | \textbf{0.686} |
| R - WSE | 5000 | 4958 | 0.893 | 0.930 | 0.920 | 0.837 | 0.716 | 0.997 | \textbf{0.686} | \textbf{0.684} |
| R - W AE | 5000 | 4958 | 1.067 | 0.974 | 0.983 | \textbf{0.479} | \textbf{0.171} | 0.999 | 0.782 | 0.780 |
than 1. When compared against the three linear shrinkage methods considered here, NEST and NPMLE demonstrate an overall value in joint shrinkage estimation of the means $\mu_i$ and the variances $\sigma^2_i$ for predicting $\theta$.

E Extensions

This section considers the extension of our methodology to several well known members in the two-parameter exponential family. We will focus on examples where the nuisance parameter is known. Our proposed estimation framework is motivated by the double shrinkage idea, but the approach nonetheless handles the case with known nuisance parameters. We discuss four examples, in each of which we derive the Bayes estimator of the natural parameter under the squared error loss. The Bayes estimator in these examples relies on the unknown score function (of the marginal density of the sufficient statistic), which can be consistently estimated using the ideas in Section 3.

Example 1 (Location mixture of Gaussians). Consider the following hierarchical model

$$Y_i \mid \mu_i, \tau_i \overset{i.i.d.}{\sim} N(\mu_i, 1/\tau_i), \quad \mu_i \overset{i.i.d.}{\sim} G_{\mu}(\cdot), \quad \text{for } i = 1, \ldots, n, \quad (53)$$

where $\tau_i$ are known and $G_{\mu}(\cdot)$ is an unspecified prior. Equation (53) represents the heteroskedastic normal means problem with known variances $1/\tau_i$ [see for example Weinstein et al. (2018)]. In this setting, the sufficient statistic for $\mu_i$ is $Y_i$ and the Bayes estimator of $\mu_i$ under the squared error loss is given by

$$\mu_i^\tau := \mathbb{E}(\mu_i | y_i, \tau_i) = y_i + \frac{1}{\tau_i} \frac{\partial}{\partial y_i} \log f(y_i | \tau_i),$$

where $f(\cdot | \tau_i)$ is the pdf of the distribution of $Y_i$ given $\tau_i$ marginalizing out $\mu_i$. From Section 3 and with $m_i = 1$, $x_i = (y_i, \tau_i)$, the NEST estimate of $\mu_i$ is given by $\delta_{nest}^{i,n}(\lambda) = y_i + \frac{1}{\tau_i} \hat{w}_{\lambda,n}^{(1)}(i)$.

Example 2 (Scale mixture of Gamma distributions). Consider the following model

$$Y_{ij} \mid \alpha_i, 1/\beta_i \overset{i.i.d.}{\sim} \Gamma(\alpha_i, 1/\beta_i), \quad 1/\beta_i \overset{i.i.d.}{\sim} G(\cdot),$$

where the shape parameters $\alpha_i$ are known and $G(\cdot)$ is an unspecified prior distribution on scale
parameters $1/\beta_i$. Here $T_i = \sum_{j=1}^m Y_{ij}$ is a sufficient statistic and $T_i|\alpha_i, \beta_i \overset{\text{ind.}}{\sim} \Gamma(m\alpha_i, 1/\beta_i)$. The posterior distribution of $1/\beta_i$ belongs to a one-parameter exponential family with density

$$f(1/\beta_i|T_i, \alpha_i) \propto \exp \left\{ -T_i\beta_i + (m\alpha_i - 1) \log T_i - \log f(T_i|\alpha_i) \right\}, \quad (54)$$

where $f(\cdot|\alpha_i)$ is the pdf of the distribution of $T_i$ given $\alpha_i$ (marginalizing out $1/\beta_i$). From Equation $(54)$, the Bayes estimator of $\beta_i$ under the squared error loss is given by

$$\beta^\pi_i := \mathbb{E}(\beta_i|T_i, \alpha_i) = \frac{m\alpha_i - 1}{T_i} - \frac{\partial}{\partial T_i} \log f(T_i|\alpha_i).$$

With $x_i = (T_i, \alpha_i)$, the NEST estimate of $\beta_i$ is given by

$$\delta^{\text{nest}}_{i,n}(\lambda) = \frac{m\alpha_i - 1}{T_i} - \hat{w}^{(1)}_{\lambda,n}(i).$$

**Example 3** (Shape mixture of Gamma distributions). We consider the following model:

$$Y_{ij} | \alpha_i, 1/\beta_i \overset{i.i.d.}{\sim} \Gamma(\alpha_i, 1/\beta_i), \quad \alpha_i \overset{i.i.d.}{\sim} G(\cdot),$$

where the scale parameters $1/\beta_i$ are known and $G(\cdot)$ is an unspecified prior distribution on the shape parameters $\alpha_i$. Let $Y_i = \sum_{j=1}^m Y_{ij}$. Then $Y_i|\alpha_i, \beta_i \overset{\text{ind.}}{\sim} \Gamma(m\alpha_i, 1/\beta_i)$ and $T_i = \log Y_i$ is a sufficient statistic. Moreover, the posterior distribution of $\alpha_i$ belongs to a one-parameter exponential family with density

$$f(\alpha_i|T_i, 1/\beta_i) \propto \exp \left\{ (m\alpha_i)T_i - \beta_i \exp (T_i) - \log f(T_i|1/\beta_i) \right\}, \quad (55)$$

where $f(\cdot|1/\beta_i)$ is the density of the distribution of $T_i$ given $1/\beta_i$ marginalizing out $\alpha_i$. From Equation $(55)$, the Bayes estimator of $\alpha_i$ under the squared error loss is given by

$$\alpha^\pi_i := \mathbb{E}(\alpha_i|T_i, 1/\beta_i) = \frac{\beta_i \exp (T_i)}{m} + \frac{1}{m} \frac{\partial}{\partial T_i} \log f(T_i|1/\beta_i).$$

With $x_i = (T_i, 1/\beta_i)$, the NEST estimate of $\alpha_i$ is

$$\delta^{\text{nest}}_{i,n}(\lambda) = \frac{\beta_i \exp (T_i)}{m} + \frac{1}{m} \hat{w}^{(1)}_{\lambda,n}(i).$$

**Example 4** (Scale mixture of Weibulls). We consider the following model:

$$Y_{ij} | k_i, \beta_i \overset{i.i.d.}{\sim} \text{Weibull}(k_i, \beta_i), \quad \beta_i \overset{i.i.d.}{\sim} G(\cdot).$$

(56)
We have \( f(y | k, \beta) = \beta ky^{k-1} \exp(-\beta y^k) \). In Equation (56) the shape parameters \( k_i \) are known, \( G(\cdot) \) is an unspecified prior distribution on the scale parameters \( \beta_i \), \( T_i = \sum_{j=1}^{m} \{Y_{ij}\}^{k_i} \) is a sufficient statistic, and \( T_i \mid k_i, 1/\beta_i \overset{ind.}\sim \Gamma(m, 1/\beta_i) \). From Example 2, the Bayes estimator of \( \beta_i \) is

\[
\beta_i^\pi := \mathbb{E}(\beta_i \mid T_i, k_i) = \frac{m - 1}{T_i} + \frac{\partial}{\partial T_i} \log f(T_i \mid k_i).
\]

With \( x_i = (T_i, k_i) \), the NEST estimate of \( \beta_i \) is given by \( \hat{\delta}_i^{\text{nest}}(\lambda) = \frac{m - 1}{T_i} - \hat{\omega}^{(1)}_{\lambda,n}(i) \).

The preceding examples present a setting with known nuisance parameter. When both parameters are unknown, extensions of our estimation framework to an arbitrary member of the two-parameter exponential family is difficult. The main reason is that in the Gaussian case the sufficient statistics are independent and their marginal distributions are known. However, for other distributions such as the Gamma and Beta, the joint distribution of the two sufficient statistics is generally unknown. This impedes a full generalization of our approach. We anticipate that an iterative scheme that conducts shrinkage estimation on the primary and nuisance coordinates in turn may be developed by combining the ideas in Examples 3 and 4 above. We do not pursue those extensions in this article.

**F Computational complexity**

Here we discuss the computational complexities of NEST and NPMLE, and provide a comparison of their running time.

In contrast to the linear shrinkage estimators, such as Group Linear Weinstein et al. (2018), SURE estimators of Xie et al. (2012) and Jing et al. (2016b), NEST and NPMLE are similar in the sense that both these approaches rely on solving a convex optimization problem to estimate the means. An implementation of the convex optimization problem for estimating the joint prior distribution of the means and the variances via the NPMLE is available in the R package REBayes (Koenker and Gu, 2017). This package relies on interior point methods for solving the convex problem. See Koenker and Mizera (2014) for more details. The main effort in solving this problem depends on computing a Hessian and the computational complexity of that is \( O(n^2 r^2) \) where \( r \) is the number of grid points at which the prior masses for each of the two prior distributions will be estimated. Usually \( n \gg r \). Recently, Kim et al. (2020) propose a fast sequential quadratic pro-
gramming based algorithm for estimating a one dimensional prior distribution using the NPMLE however their algorithm is not available for the case where the prior is two dimensional, such as the scenario when both mean and variances are unknown. For NEST, the convex quadratic optimization problem described in Equation (12) depends on the $n \times n$ matrix $K_\lambda$ and uses the interior-point optimizer in MOSEK. The worst case computational complexity for evaluating the hessian of the underlying objective function is $O(n^3)$, which is substantially larger than that of NPMLE. Figure 16 provides a comparison of the running time of NEST and NPMLE for increasing $n$ and $r$. In the left panel of Figure 16 we fix $r = 30$, the default choice implemented in the function $\text{WGLVmix}$ in REBayes, and vary $n$. In the right panel, $n$ is fixed at 5000 with $r$ varying. We see that for problems with large $n$ and $r$ fixed, NPMLE is substantially faster than NEST and NEST, in its current implementation, may not be as scalable as NPMLE is to large $n$ problems. Our future research efforts will be geared towards developing faster first-order methods to solve Equation (12) rather than relying on the interior-point solver in MOSEK.

Figure 16: Running time comparison of NEST and NPMLE. Left: $r$ fixed at 30. Right: $n$ fixed at 5000.