

Fast Bootstrapping Nonparametric Maximum Likelihood for Latent Mixture Models

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Abstract—Estimating the mixing density of a latent mixture model is an important task in signal processing. Nonparametric maximum likelihood estimation is one popular approach to this problem. If the latent variable distribution is assumed to be continuous, then bootstrapping can be used to approximate it. However, traditional bootstrapping requires repeated evaluations on resampled data and is not scalable. In this letter, we construct a generative process to rapidly produce nonparametric maximum likelihood bootstrap estimates. Our method requires only a single evaluation of a novel two-stage optimization algorithm. Simulations and real data analyses demonstrate that our procedure accurately estimates the mixing density with little computational cost even when there are a hundred thousand observations.

Index Terms—bootstrap/resampling, deep neural network, generative process, mixing density estimation, nonparametric maximum likelihood estimation, two-stage algorithm

I. INTRODUCTION

NONPARAMETRIC maximum likelihood estimation (NPMLE) is a popular methodology in signal processing applications such as pattern recognition [1, 2], channel estimation [3], signal recovery [4], and positron emission tomography [5, 6]. Suppose that we observe $\mathbf{y} = (y_1, \dots, y_n)$ where

$$y_i | \theta_i \sim f(y_i | \theta_i), \quad \theta_i \sim \pi(\theta), \quad i = 1, \dots, n, \quad (1)$$

and the density f is known but π is unknown. NPMLE aims to estimate the mixing density (or the prior) π .

The NPMLE estimator $\hat{\pi}$ in the latent mixture model (1) was first introduced by [7] and solves the optimization,

$$\begin{aligned} \hat{\pi} &= \operatorname{argmax}_{\pi \in \Pi} \sum_{i=1}^n \log[\mathbb{E}_{\pi}\{f(y_i | \theta_i)\}] \\ &= \operatorname{argmax}_{\pi \in \Pi} \sum_{i=1}^n \log \left\{ \int_{\Theta} f(y_i | \theta_i) d\pi(\theta_i) \right\}, \end{aligned} \quad (2)$$

where Π denotes the family of all probability distributions on the parameter space Θ . Under mild regularity conditions, [8] showed that the solution $\hat{\pi}$ of (2) exists and is unique, and even if the true π is continuous, $\hat{\pi}$ is almost surely discrete with a support of at most n points. Exploiting this property, there exist many algorithms to solve (2), including the EM algorithm [9, 10] and convex optimizers [11, 12].

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Nevertheless, a discrete NPMLE estimator $\hat{\pi}$ is unsatisfactory when the latent distribution is reasonably assumed to be continuous [5, 13]. As a result, smoothed variants of NPMLE have been proposed [2, 5, 14, 15]. However, these methods typically require careful tuning of smoothing parameters such as bandwidth [2, 15], roughness penalty term [5, 14], and/or spline degrees of freedom [2, 14].

As an alternative to smoothing, bootstrapping has been shown to be an effective way to simulate from both prior and posterior continuous densities [16]–[19]. However, bootstrapping is seldomly employed for NPMLE. This may be because the bootstrap traditionally involves resampling the data with replacement and repeatedly optimizing weighted objective functions, which can be very computationally intensive. In the context of NPMLE, the standard bootstrap procedure requires one to repetitively optimize

$$\hat{\pi}^{(b)} = \operatorname{argmax}_{\pi} \sum_{i=1}^n w_i^{(b)} \log[\mathbb{E}_{\pi}\{f(y_i | \theta_i)\}], \quad \mathbf{w}^{(b)} \stackrel{\text{iid}}{\sim} \mathbb{P}_{\mathbf{w}}, \quad (3)$$

for some probability measure $\mathbb{P}_{\mathbf{w}}$. For example, if $\mathbf{w} \sim \text{Multinomial}(n, \mathbb{1}_n/n)$, where $\mathbb{1}_n$ denotes an n -dimensional one vector, we have the nonparametric bootstrap [20]. If $\mathbf{w} \sim n \times \text{Dirichlet}(n, \mathbb{1}_n)$, we have the weighted likelihood bootstrap [18].

To resolve this computational bottleneck and broaden the applicability of bootstrapping, we build upon the generative bootstrap sampling (GBS) approach of [21]. We introduce a new generative framework called Generative Bootstrapping for NPMLE (GB-NPMLE). Our contributions are as follows:

- 1) We propose GB-NPMLE for estimating a continuous prior density via bootstrapping. Different from GBS which only requires weights \mathbf{w} as inputs [21], GB-NPMLE needs *additional* noise inputs \mathbf{z} to capture the latent representation of π in (1).
- 2) To optimize GB-NPMLE, we introduce a novel two-stage algorithm. In contrast to traditional bootstrapping, GB-NPMLE requires only a *single* evaluation of this two-stage algorithm instead of repetitive evaluations of (3).

II. PROPOSED METHOD

A. GB-NPMLE

Recent developments in generative learning such as variational autoencoders (VAEs) [22] and generative adversarial networks (GANs) [23]–[25] provide another perspective for NPMLE. Instead of estimating π , we can construct a generator T that matches to the target density π , i.e. $T(\mathbf{z}) \sim \pi$ where

$z \in \mathbb{R}^q$ follows a known distribution such as a Gaussian or a uniform. We then reformulate NPMLE optimization (2) via a generative framework,

$$\hat{T} = \operatorname{argmax}_{T \in \mathcal{T}} \sum_{i=1}^n \log[\mathbb{E}_z \{f(y_i | T(z))\}], \quad (4)$$

where $z \in \mathbb{R}^q$ follows (for example) a standard uniform. [26] considered (4) where the generator T was constructed by a VAE [22] with a Langevin bias-correction algorithm [27]. Unfortunately, both the classical NPMLE $\hat{\pi}$ in (2) and the generative NPMLE \hat{T} in (4) result in discrete solutions [8]. This discreteness is less attractive when the true latent π is continuous, as in many signal processing applications [5, 6].

To efficiently obtain a bootstrapped NPMLE distribution as a smooth estimator for π in (1), we introduce GB-NPMLE. Our framework (approximately) generates NPMLE bootstrap estimates via a generator function. The generator $G := G(\mathbf{w}, \mathbf{z})$ takes *both* noise $z \in \mathbb{R}^q$ and bootstrap weights $\mathbf{w} \in \mathbb{R}^n$ as inputs. We construct G using a feedforward neural network (FNN) [28, 29], because of the FNN's universal approximation properties for a large class of functions [28, 29].

As a fast approximation to bootstrapped NPMLE (3), GB-NPMLE incorporates the GBS approach of [21]. However, GBS is only geared towards uncertainty quantification of a single *point estimate*. Consequently, GBS takes *only* weights \mathbf{w} as inputs and *cannot* be directly used to learn an entire *probability density* π as in (3). In order to learn an unknown density π , GB-NPMLE *also* requires noise inputs \mathbf{z} , as in (4). In sum, GB-NPMLE approximates the bootstrap distribution for NPMLE with the objective function,

$$\hat{G} = \operatorname{argmax}_G \mathbb{E}_{\mathbf{w}} \left[\sum_{i=1}^n w_i \log[\mathbb{E}_z \{f(y_i | G(\mathbf{w}, \mathbf{z}))\}] \right], \quad (5)$$

where $G(\mathbf{w}, \mathbf{z}) : \mathbb{R}^{n+q} \mapsto \mathbb{R}$. Once we have optimized G (i.e. the weights and biases of the FNN), it is then effortless to generate novel NPMLE bootstrap estimates as $\hat{G}(\mathbf{w}_{\text{new}}, \mathbf{z}_{\text{new}})$.

To optimize (5) in practice, the expectations \mathbb{E}_z and $\mathbb{E}_{\mathbf{w}}$ are approximated by Monte Carlo averaging, similarly as in VAEs [22] and GANs [23]. Specifically, we independently sample $z \sim \text{Unif}(0, 1)$ and $\mathbf{w} \sim n \times \text{Dirichlet}(n, \mathbf{1}_n)$ a sufficiently large number of times and then approximate the expectations in (5) with sample averages.

Directly optimizing G in (5) using stochastic gradient descent (SGD) [30] is a nontrivial task. In practice, the Monte Carlo approximation for \mathbb{E}_z introduces high variance to the approximate gradient for SGD because the log functions in (5) greatly shatter the linearity of \mathbb{E}_z [31]. Consequently, this can lead to slow convergence when optimizing the GB-NPMLE objective function (5) [31]. We address these estimation difficulties with a novel two-stage algorithm.

B. GB-NPMLE Two-Stage Algorithm

To efficiently optimize G in (5), we generalize the generator output $\theta = G(\mathbf{w}, \mathbf{z})$ to be a multi-dimensional $\boldsymbol{\theta} \in \mathbb{R}^l$ where $l > 1$ is the number of realizations of θ . The resulting sequence $\boldsymbol{\theta} = \{\theta^{(j)}\}_{j=1}^l$ serves as the bootstrap sample *candidates*. As we explain shortly, multiple realizations of θ help to

stabilize the gradient approximation in SGD for the Monte Carlo estimate of \mathbb{E}_z in (5). Unfortunately, this generalization also leads to correlation across the bootstrap samples.

One way to overcome the correlation issue is to take a random draw of one $\theta^{(j)}$ from $\boldsymbol{\theta}$ according to mixing probabilities $\boldsymbol{\tau} = (\tau_1, \dots, \tau_l)$ for the entries of $\boldsymbol{\theta}$. Repeating this procedure B times produces B independent bootstrap samples. However, in addition to generator G , we now *also* need to estimate $\boldsymbol{\tau}$. Accordingly, we propose a two-stage algorithm where in Stage I, we train G , and in Stage II, we estimate $\boldsymbol{\tau}$.

GB-NPMLE Algorithm Stage I: Fix $\boldsymbol{\tau} = (\tau_1, \dots, \tau_l)$ where $\tau_1 = \dots = \tau_l = 1/l$, and solve the modified GB-NPMLE objective function,

$$\hat{G} = \operatorname{argmax}_G \mathbb{E}_{\mathbf{w}} \left[\sum_{i=1}^n w_i \log[\mathbb{E}_{z, \gamma} \{f(y_i | e_{\gamma}^{\top} G(\mathbf{w}, \mathbf{z}))\}] \right], \quad (6)$$

where $\gamma \sim \text{Multinomial}(1, \boldsymbol{\tau})$, e_{γ} is the γ -th unit vector, and $e_{\gamma}^{\top} G(\mathbf{w}, \mathbf{z}) \in \mathbb{R}^p$ is the γ -th entry of generator output $\boldsymbol{\theta}$.

To see how the modified GB-NPMLE objective (6) helps to relieve high Monte Carlo variance, let $\mathbb{E}_{z, \gamma} \{f(y_i | e_{\gamma}^{\top} G(\mathbf{w}, \mathbf{z}))\}$ in (6) be denoted by $\mathbb{E}_z \{ \mathbb{E}_{\gamma} \{f(\boldsymbol{\theta}[\mathbf{w}, \mathbf{z}, \gamma])\} \}$, and similarly, let $\mathbb{E}_z \{f(y_i | G(\mathbf{w}, \mathbf{z}))\}$ be $\mathbb{E}_z \{f(\boldsymbol{\theta}[\mathbf{w}, \mathbf{z}])\}$ in (5). By the law of total variance, $\text{Var}\{f(\boldsymbol{\theta}[\mathbf{w}, \mathbf{z}])\} \geq \text{Var}\{\mathbb{E}_{\gamma} f(\boldsymbol{\theta}[\mathbf{w}, \mathbf{z}, \gamma])\}$. Hence, the generalization of the generator output dimension to $l > 1$ reduces Monte Carlo variance.

GB-NPMLE Algorithm Stage II: Fix the well-trained generator \hat{G} from (6) and estimate $\boldsymbol{\tau}$ via the Monte Carlo EM (MCEM) algorithm [32]. In particular, the MCEM updates for each k th entry of $\boldsymbol{\tau}$ have the closed form,

$$\tau_k^{(t+1)} = \frac{1}{n} \sum_{i=1}^n \frac{\tau_k^{(t)} \cdot \mathbb{E}_{\mathbf{w}, \mathbf{z}} \{f(y_i | e_k^{\top} \hat{G}(\mathbf{w}, \mathbf{z}))\}}{\sum_{k=1}^l \tau_k^{(t)} \cdot \mathbb{E}_{\mathbf{w}, \mathbf{z}} \{f(y_i | e_k^{\top} \hat{G}(\mathbf{w}, \mathbf{z}))\}}. \quad (7)$$

Note that we *already* obtained an efficient estimator of $\boldsymbol{\theta}$ from Stage I, and hence, the MCEM algorithm tends to have very fast convergence. In the literature, slow convergence of EM [26] mostly lies in obtaining a sequence of estimators $\boldsymbol{\theta}$. We avoid this issue by *fixing* the generator from Stage I.

Once we have estimated both G (in Stage I) and $\boldsymbol{\tau}$ (in Stage II), we can easily generate a new bootstrap estimate by sampling $\mathbf{w}_{\text{new}}, \mathbf{z}_{\text{new}}, \gamma_{\text{new}}$, and then taking $\theta_{\text{new}} = e_{\gamma_{\text{new}}}^{\top} \hat{G}(\mathbf{w}_{\text{new}}, \mathbf{z}_{\text{new}})$. Repeating this procedure B times results in B independent bootstrap estimates. Thus, in contrast to bootstrapped NPMLE which requires B total repetitive evaluations of (3), GB-NPMLE requires only a *single* evaluation of the two-stage algorithm. The entire two-stage GB-NPMLE procedure is given in Algorithm 1.

III. EXPERIMENTS AND RESULTS

A. Simulation Studies

We investigate the performance of GB-NPMLE optimized by the two-stage algorithm in Section II-B. The generator G is an FNN with three hidden layers and 500 neurons per layer. The expectations $\mathbb{E}_{\mathbf{w}}, \mathbb{E}_z$ and \mathbb{E}_{γ} in the modified GB-NPMLE objective (6) are approximated by Monte Carlo averages

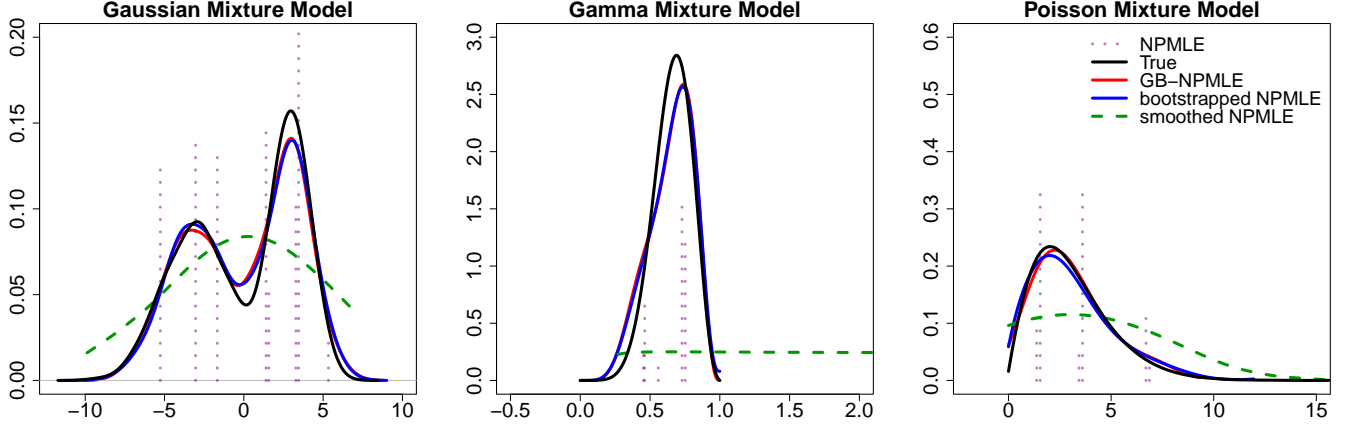


Fig. 1. Results from one replication of Simulations (i)-(iii). In addition to GB-NPMLE (solid red), bootstrapped NPMLE (solid blue), and smoothed NPMLE (dashed green), we also plot the true density (solid black) and the classical discrete NPMLE (dotted purple).

Algorithm 1 GB-NPMLE Two-stage Algorithm

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1: Set  $l = 100$ ,  $tol = 0.001$ ,  $T = 2000$ ,  $B = 1000$ .
2: Initialize neural network parameters  $\phi$  and  $\tau = (1/l, \dots, 1/l)^\top$ 
3: procedure: Stage I
4:   for  $t$  in  $1, \dots, \text{Epoch } T$  do
5:     Sample  $\mathbf{w} \sim n \times \text{Dirichlet}(n, \mathbb{1}_n)$ 
6:     Sample  $z \sim \text{Unif}(0,1)$ 
7:     Sample  $\gamma \sim \text{Multinomial}(1, \tau)$ 
8:     Update  $G_\phi$  by SGD using modified GB-NPMLE objective (6)
9:   end for
10:  return  $G_\phi$ 
11: end procedure
12: procedure: Stage II
13:  Initialize  $\tau^{new} = (1/l, \dots, 1/l)$  and  $\tau^{old} = (0, \dots, 0)$ 
14:  while  $\min(|\tau^{old} - \tau^{new}|) \geq tol$  do
15:    set  $\tau^{old} = \tau^{new}$ 
16:    for  $k$  in  $1, \dots, l$  do
17:      Sample  $\mathbf{w} \sim n \times \text{Dirichlet}(n, \mathbb{1}_n)$ 
18:      Sample  $z \sim \text{Unif}(0,1)$ 
19:      Update  $\tau^{new}$  as in (7)
20:    end for
21:  end while
22:  return  $\tau^{new}$ 
23: end procedure
24: procedure: Generate  $B$  bootstrap estimates
25:  for  $b$  in  $1, \dots, B$  do
26:    Sample  $\mathbf{w}^{(b)} \sim n \times \text{Dirichlet}(n, \mathbb{1}_n)$ 
27:    Sample  $z^{(b)} \sim \text{Unif}(0,1)$ 
28:    Sample  $\gamma^{(b)} \sim \text{Multinomial}(1, \tau^{new})$ 
29:    Generate  $\theta^{(b)} = e_{\gamma^{(b)}}^\top \widehat{G}(\mathbf{w}^{(b)}, z^{(b)})$ 
30:  end for
31:  return  $\theta^{(1)}, \dots, \theta^{(B)}$ 
32: end procedure

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of 100 independent samples of $\mathbf{w} \sim n \times \text{Dirichlet}(n, \mathbb{1}_n)$, $z \sim \text{Unif}(0,1)$ and $\gamma \sim \text{Multinomial}(1, \tau)$, and the length of τ is set to be $l = 100$. To train G , we use SGD with the Adam optimizer [33]. All experiments are conducted on a single NVIDIA GeForce RTX 2080 Ti graphics processing unit (GPU) with 11GB RAM.

We consider three simulation settings for the prior π in the mixture model (1): (i) bimodal, (ii) unimodal and bounded on $(0, 1)$, and (iii) skewed right. Specifically, we consider:

- (i) Gaussian mixture model (**GMM**): $y | \theta \sim \mathcal{N}(\theta, 1)$ and $\theta = 0.5\mathcal{N}(-3, 2) + 0.5\mathcal{N}(3, 1)$;

- (ii) Gamma mixture model (**GaMM**): $y | \theta \sim \text{Gamma}(10, \theta)$ and $\theta \sim \text{Beta}(10, 5)$;
- (iii) Poisson mixture model (**PMM**): $y | \theta \sim \text{Poisson}(\theta)$ and $\theta \sim \text{Gamma}(3, 1)$.

We compare GB-NPMLE to bootstrapped NPMLE (3) and a smoothed version of the discrete NPMLE (2). We use $B = 1000$ bootstrap estimates to estimate the prior for GB-NPMLE and bootstrapped NPMLE. For bootstrapped NPMLE, we use the nonparametric bootstrap, i.e. we sample n instances of the data with replacement and then optimize (2) a total of B times using the R package REBayes [34]. For the smoothed NPMLE, we use kernel smoothing of $\widehat{\pi}$ in (2) with the `KWsmooth` function in REBayes [34], where we select the optimal bandwidth from an equispaced grid from 0.1 to 20 using leave-one-out cross-validation (LOOCV).

For performance comparison of the three NPMLE methods, we generate artificial datasets of size $n = 1000$. We look at: a) Wasserstein-1 distance between π and $\widehat{\pi}$, defined as $W_1(\pi, \widehat{\pi}) = \int |\mathbb{F}_\pi(x) - \mathbb{F}_{\widehat{\pi}}(x)| dx$, where \mathbb{F} is the cumulative density function (CDF) of π , and b) Integrated Squared Error (ISE), defined as $\int (\widehat{\pi}(x) - \pi(x))^2 dx$. Lower $W_1(\pi, \widehat{\pi})$ and ISE indicate better performance. Our results for simulations (i)-(iii) averaged across 20 replications are shown in Table I.

TABLE I
COMPARISONS OF PERFORMANCE OF DIFFERENT NPMLE METHODS ON SIMULATED DATASETS

Model	GB-NPMLE		Bootstrapped NPMLE		Smoothed NPMLE	
	$W_1(\pi, \widehat{\pi})$	ISE	$W_1(\pi, \widehat{\pi})$	ISE	$W_1(\pi, \widehat{\pi})$	ISE
GMM	0.348	0.008	0.298	0.008	0.936	0.034
GaMM	0.003	0.263	0.037	0.497	0.223	1.731
PMM	0.400	0.045	0.389	0.036	1.763	0.059

From Table I, we first observe that GB-NPMLE has a very similar performance to bootstrapped NPMLE. On the other hand, smoothed NPMLE (with bandwidth parameter chosen by LOOCV) performs much worse than the bootstrapping approaches. This demonstrates the difficulty of optimally tuning smoothing parameters for smoothed NPMLE.

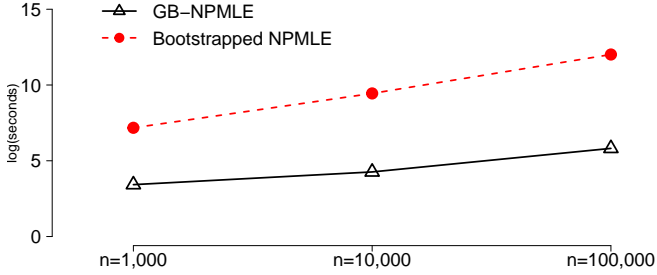


Fig. 2. Mean computation time in log(seconds) for the three simulations vs. sample size $n \in \{1000, 10,000, 100,000\}$. For $n = 100,000$, GB-NPMLE takes five minutes, whereas bootstrapped NPMLE takes almost two days.

The results from one replication of simulations (i)-(iii) are plotted in Fig. 1, which also plots the true latent density π (solid black) and the discrete NPMLE solution (2) (dotted purple). Fig. 1 shows that the classical (discrete) NPMLE (2) is inadequate for capturing π when π is truly continuous. For example, in the GaMM model, the classical NPMLE does not clearly indicate that π is unimodal. On the other hand, GB-NPMLE (solid red) and bootstrapped NPMLE (solid blue) are quite close to the true π , capturing all inherent aspects of the true π like bimodality, boundedness and unimodality, and skewness respectively in the GMM, GaMM, and PMM models. The smoothed NPMLE appears to be oversmoothed and is unable to capture bimodality of π in the GMM model or the unimodality and boundedness of π in the GaMM model. In the PMM model, the sharp peak of the skewed density π is also not adequately captured by smoothed NPMLE.

GB-NPMLE and bootstrapped NPMLE both give similarly good estimation performance, but GB-NPMLE is much faster because it only requires a single evaluation. We compare the average computational time of GB-NPMLE and bootstrapped NPMLE for sample sizes $n \in \{1000, 10,000, 100,000\}$ across all replicates of simulations (i)-(iii). Our results are plotted (on the log scale) in Fig. 2.

Fig. 2 shows that GB-NPMLE is much more scalable. In particular, GB-NPMLE takes only about five minutes to complete when $n = 100,000$; on the other hand, bootstrapped NPMLE requires almost two days to finish. Even if we were to parallelize bootstrapped NPMLE across 100 GPU cores with no competing processes and minimal communication cost, GB-NPMLE on a single GPU core would still be faster.

B. Evaluation on Real Datasets

We evaluate GB-NPMLE on three popular count datasets where the counts are assumed to follow a Poisson mixture model, $y_i | \theta_i \sim \text{Poisson}(\theta_i)$ and $\theta_i \sim \pi, i = 1, \dots, n$. The *Norberg* dataset consists of 1125 group life insurance statistics in Norway [35], the *Thailand* dataset consists of 602 pre-school children’s health condition, and the *Mortality* dataset contains the number of deaths for women aged greater than eighty [36]. Fig. 3 compares GB-NPMLE (top panel) with bootstrapped NPMLE (bottom panel). We see that GB-NPMLE approximates bootstrapped NPMLE very well. Fur-

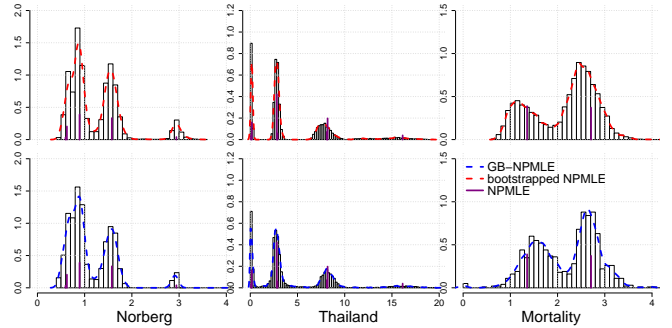


Fig. 3. Results for real datasets for GB-NPMLE (dashed red, top panel) and bootstrapped NPMLE (dotted blue, bottom panel). The pink vertical lines in both plots represent the discrete NPMLE solution.

thermore, both methods identify the same number of clusters (three in *Norberg*, four in *Thailand*, and two in *Mortality*).

We also examine the out-of-sample prediction $\mathbb{E}_{\theta}[-\log f(\mathbf{y}_{\text{new}}; \theta) | \mathbf{y}_{\text{obs}}]$, reported as “Likelihood” in Table II. Specifically, using K -fold cross validation (CV),

$$\text{Likelihood} = K^{-1} \sum_{k=1}^K \sum_{i \in I_k} -\log \left[B^{-1} \sum_{b=1}^B f(y_i; \hat{\theta}_{(-k)}^b) \right],$$

where $\hat{\theta}_{(-k)}^b \sim \hat{\pi}_{(-k)}$ and $\hat{\pi}_{(-k)}$ is the bootstrap distribution obtained by excluding the k th fold I_k . We set $K = 10$ and $B = 500$. From Table II, we observe that the GB-NPMLE predictions are quite close to their standard bootstrap counterparts. Table II also shows that GB-NPMLE is much more scalable, with only a modest increase in computation time for 10-fold CV when n increases from 72 to 1096. When $n = 1096$, GB-NPMLE is about 34 times faster than bootstrapped NPMLE.

TABLE II
COMPARISONS OF PERFORMANCE OF GB-NPMLE AND BOOTSTRAPPED NPMLE ON REAL COUNT DATASETS

Dataset	n	GB-NPMLE		Bootstrapped NPMLE	
		Likelihood	Time (sec)	Likelihood	Time (sec)
Norberg	72	20.35	126.1	20.83	732.8
Thailand	602	156.26	140.9	156.21	3439.3
Mortality	1096	199.30	153.3	199.31	5279.2

IV. CONCLUSION

In this letter, we introduced GB-NPMLE, a generative framework for estimating a mixing density through bootstrapping. A novel two-stage algorithm was proposed to rapidly and accurately obtain bootstrap estimates. Bootstrapping has rarely been employed for NPMLE in the past, and our work paves the way for bootstrapping to serve as an attractive alternative to smoothing NPMLE. Simulations and real data analyses demonstrated the effectiveness and scalability of our approach. Code to implement GB-NPMLE is available on GitHub at <https://github.com/shijiew97/GBnmpml>.

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