

Adaptive Bayesian Procedures Using Random Series Priors

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ABSTRACT. We consider a general class of prior distributions for nonparametric Bayesian estimation which uses finite random series with a random number of terms. A prior is constructed through distributions on the number of basis functions and the associated coefficients. We derive a general result on adaptive posterior contraction rates for all smoothness levels of the target function in the true model by constructing an appropriate ‘sieve’ and applying the general theory of posterior contraction rates. We apply this general result on several statistical problems such as density estimation, various nonparametric regressions, classification, spectral density estimation and functional regression. The prior can be viewed as an alternative to the commonly used Gaussian process prior, but properties of the posterior distribution can be analysed by relatively simpler techniques. An interesting approximation property of B-spline basis expansion established in this paper allows a canonical choice of prior on coefficients in a random series and allows a simple computational approach without using Markov chain Monte Carlo methods. A simulation study is conducted to show that the accuracy of the Bayesian estimators based on the random series prior and the Gaussian process prior are comparable. We apply the method on Tecator data using functional regression models.

Key words: adaptation, B-spline, Gaussian process, nonparametric Bayes, posterior contraction rate, random series prior

1. Introduction

Bayesian methods have been widely used in the nonparametric statistical literature. Contraction rates of posterior distributions were studied in Ghosal *et al.* (2000), Shen & Wasserman (2001), Ghosal and van der Vaart (2007a,b) and van der Vaart & van Zanten (2008). The optimal contraction rate of estimating a univariate α -smooth function is typically $n^{-\alpha/(2\alpha+1)}$, where n is the sample size. Because the smoothness parameter α is usually unknown in practice, it is then of interest to investigate if a prior leads to optimal posterior contraction rates simultaneously for all values of α , possibly up to a logarithmic factor. If that holds, a procedure is called rate-adaptive.

Bayesian rate adaptation results are important for at least two reasons. Firstly, they guarantee maximum possible accuracy of the Bayesian estimation procedure within the given framework. Secondly, they assure that the same prior distribution can be used regardless of the smoothness of the underlying function being estimated. Bayesian adaptation results have been established for signal estimation by Belitser & Ghosal (2003) and Szabó *et al.* (2013), for density estimation by Ghosal *et al.* (2003, 2008), Scricciolo (2006) and Huang (2004), and for nonparametric regression by Huang (2004) using discrete mixtures. Alternatively, van der Vaart & van Zanten (2009) constructed a prior based on a randomly rescaled Gaussian process, which automatically adapts for a continuous range of smoothness parameters.

Gaussian processes have been widely used for constructing prior distributions (Lenk, 1988) and applications in spatial statistics (Banerjee *et al.*, 2008). Posterior computational methods

were developed in Choudhuri *et al.* (2007), Rasmussen & Williams (2006), Tokdar (2007) and Rue *et al.* (2009) among others. Posterior asymptotic properties, which are primarily driven by the structure of their reproducing kernel Hilbert space, were studied by Tokdar & Ghosh (2007), Ghosal & Roy (2006), Choi & Schervish (2007), van der Vaart and van Zanten (2007, 2008, 2009), Castillo (2008) & Castillo (2012), Castillo *et al.* (2014) and Bhattacharya *et al.* (2014).

Besides a Gaussian process, another common prior on functions, obtained by putting a prior on the the number of terms and the corresponding coefficients of a series expansion, has been used extensively in applications (Crainiceanu *et al.*, 2005). Study of posterior contraction rates for such finite random series priors have begun only recently. Rivoirard & Rousseau (2012b) considered univariate density estimation using an exponential link and wavelet or Fourier series basis; de Jonge & van Zanten (2012) considered a general approach for multivariate function estimations using tensor-product spline basis and Gaussian distributions on the coefficients; Arbel *et al.* (2013) proposed a class of sieve priors with general choice of basis functions and independent priors on the coefficients. A related work is Babenko & Belitser (2010), who obtained oracle inequalities for posterior contraction for the infinite dimensional normal mean problem by putting a prior on the number of non-zero entries and then independent normal priors on the resulting components.

In the present paper, our contributions are two-fold. Firstly, we obtain posterior contraction rates for finite random series priors for any curve estimation problems under both univariate and multivariate settings with arbitrary bases and arbitrary distributions on coefficients. Secondly, we show that for the B-splines basis and certain choices of priors on the coefficients, the posterior computation can be carried out by exploiting a conjugacy-like structure without using Markov chain Monte Carlo (MCMC) techniques. Inevitably, there are some overlap with Rivoirard & Rousseau (2012b), de Jonge & van Zanten (2012) and Arbel *et al.* (2013), but our goal is to emphasize the general properties of finite random series in all curve estimation problems and the possible MCMC-free calculation building on the conjugacy-like structure, which emerges only when one considers general prior distributions on the coefficients. We formulate one general theorem in an abstract setting suitable as a prelude for many different inference problems where we allow arbitrary basis functions and arbitrary multivariate distributions on the coefficients of the expansion. Thus, the resulting process induced on the function need not be Gaussian, and can accommodate a variety of functions starting from one with a bounded support to one with a heavy tail. The resulting rate obtained in the abstract theorem depends on the smoothness of the underlying function, approximation ability of the basis expansion used, tail of the prior distribution on the coefficients, prior on the number of terms in the series expansion, prior concentration and the metrics being used. We compute the rates for various combinations of these choices.

It may be noted that Gaussian process and random series priors are intimately related in two ways — a normal prior on the coefficients of a random series gives a Gaussian process, while the Karhunen–Loève expansion of a Gaussian process expresses itself as a random series with basis consisting of eigenfunctions of the covariance kernel of the Gaussian process. Thus, a random series prior may be regarded as a flexible alternative to a Gaussian process prior. It is interesting to note that the theory of posterior contraction for Gaussian process priors established in van der Vaart & van Zanten (2007), van der Vaart & van Zanten (2008), & van der Vaart & van Zanten (2009) use deep properties of Gaussian processes, while relatively elementary techniques lead to comparable posterior contraction rates for finite random series priors. Posterior computation for Gaussian process priors often need reversible jump MCMC procedures (Tokdar, 2007) typically with a large number of knots to approximate a given Gaussian process. For a random series prior based on B-spline expansion, for an appropriate prior on

the coefficients, the conjugacy-like structure model can avoid the use of MCMC altogether by representing the posterior mean analytically, although the number of terms in the representation may be large. When the sample size n is relatively small (e.g. $n = 10$), the number of terms is manageable and the exact values of posterior moments can be computed. When the sample size is large, we sample a few terms and estimate the sum. The Monte Carlo standard error of the expression can be estimated, and is often fairly controlled provided the terms are similar to each other.

The paper is organized as follows. In section 2, we present the main theorems of random series priors. In sections 3 and 4, we apply the theorems to a variety of statistical problems and derive the corresponding posterior contraction rates. Numerical results are presented in section 5.

2. General results

2.1. Notations

Let $\mathbb{N} = \{1, 2, \dots\}$, $\Delta_j = \{(x_1, \dots, x_j) : \sum_{i=1}^j x_i = 1, x_1, \dots, x_j \geq 0\}$, and δ_x stand for the degenerate probability distribution at a point x . Let the indicator function of a set A be denoted by $\mathbb{1}\{A\}$. For an open region Ω_0 in a Euclidean space, define the α -Hölder class $\mathcal{C}^\alpha(\Omega_0)$ as the collection of functions f on Ω_0 that has bounded derivatives up to the order α_0 , which is the largest integer strictly smaller than α , and the α_0 -th derivative of f satisfies the Hölder condition $|f^{(\alpha_0)}(x) - f^{(\alpha_0)}(y)| \leq C|x - y|^{\alpha - \alpha_0}$ for some constant $C > 0$ and any x, y in the support of f .

We use ' \lesssim ' to denote an inequality up to a constant multiple, where the underlying constant of proportionality is universal. By $f \asymp g$, we mean $f \lesssim g \lesssim f$. The packing number $D(\epsilon, T, d)$ is defined as the maximum cardinality of a subset of T whose elements are at least ϵ -separated out with respect to a distance d . Let $h^2(p, q) = \int (\sqrt{p} - \sqrt{q})^2 d\mu$, be the squared Hellinger distance, $K(p, q) = \int p \log(p/q) d\mu$, $V(p, q) = \int p \log^2(p/q) d\mu$, be the Kullback–Leibler (KL) divergences and $\mathcal{K}(p, \epsilon) = \{f : K(p, f) \leq \epsilon^2, V(p, f) \leq \epsilon^2\}$, be the KL neighbourhood. For a vector $\theta \in \mathbb{R}^d$, define $\|\theta\|_p = \left\{ \sum_{i=1}^d |\theta_i|^p \right\}^{1/p}$, $1 \leq p < \infty$, and $\|\theta\|_\infty = \max_{1 \leq i \leq d} |\theta_i|$. Similarly, we define $\|f\|_{p,G} = \left\{ \int |f(x)|^p dG \right\}^{1/p}$ and $\|f\|_\infty = \sup_x |f(x)|$ as the L_p -, $1 \leq p < \infty$, and L_∞ -norms of a function f with respect to a measure G .

2.2. Main results

We consider a random variable J taking values in \mathbb{N} . For each $J \in \mathbb{N}$, we consider a triangular array of linearly independent real-valued functions $\xi = (\xi_1, \xi_2, \dots, \xi_J)^T$ defined on a region Ω_0 . In applications, Ω_0 will be typically a bounded region. Note that the resulting basis functions may change from one stage to the next, although that is not made explicit in our notation. We use Π as a generic notation for priors assigned on J and the coefficients of basis functions $\theta = (\theta_1, \dots, \theta_J)^T$.

- (A1) For some $c_1, c_2 > 0$, $0 \leq t_2 \leq t_1 \leq 1$, $\exp\{-c_1 j \log^{t_1} j\} \leq \Pi(J = j) \leq \exp\{-c_2 j \log^{t_2} j\}$.
- (A2) Given J , we consider a J -dimensional joint distribution as the prior for $\theta = (\theta_1, \dots, \theta_J)^T$ satisfying $\Pi(\|\theta - \theta_0\|_2 \leq \epsilon) \geq \exp\{-c_3 J \log(1/\epsilon)\}$ for every $\|\theta_0\|_\infty \leq H$, where c_3 is some positive constant, H is chosen sufficiently large, and $\epsilon > 0$ is sufficiently small. Also, assume that $\Pi(\theta \notin [-M, M]^J) \leq J \exp\{-CM^{t_3}\}$ for some constants $C, t_3 > 0$.

Remark 1. Geometric, Poisson and negative binomial distributions on J satisfy Condition (A1) respectively with $t_1 = t_2 = 0$, $t_1 = t_2 = 1$ and $t_1 = t_2 = 0$. Examples of priors satisfying (A2) include independent gamma, exponential distributions assigned on each element of θ , and multivariate normal and Dirichlet distributions provided the parameters lie in a fixed compact set; see Lemma 6.1 of Ghosal *et al.* (2000) for the last conclusion.

We consider a distance metric d on functions belonging to Ω_0 satisfying the following condition for every $\theta_1, \theta_2 \in \mathbb{R}^J$, $J \in \mathbb{N}$, and some positive increasing function $a(\cdot)$:

$$d\left(\theta_1^T \xi, \theta_2^T \xi\right) \leq a(J)\|\theta_1 - \theta_2\|_2. \tag{2.1}$$

Now we state the main theorem, which gives unified conditions for posterior contraction rates for various inference problems, in a manner similar to Theorem 2.1 of van der Vaart & van Zanten (2008) and Theorem 3.1 of van der Vaart & van Zanten (2009).

Theorem 1. Let $\epsilon_n \geq \bar{\epsilon}_n$ be two sequence of positive numbers satisfying $\epsilon_n \rightarrow 0$ and $n\bar{\epsilon}_n^2 \rightarrow \infty$ as $n \rightarrow \infty$. For a function w_0 , suppose that a prior satisfies Conditions (A1) and (A2). Assume that there exist sequences of positive numbers J_n, \bar{J}_n and M_n , a strictly decreasing, nonnegative function $e(\cdot)$ and $\theta_{0,j} \in \mathbb{R}^j$ for any $j \in \mathbb{N}$, such that the following conditions hold for some positive constants $a_1 > 1$, a_2 , C_0 and H :

$$\|\theta_{0,j}\|_\infty \leq H, \quad d(w_0, \theta_{0,j}^T \xi) \leq e(j), \tag{2.2}$$

$$J_n \{\log J_n + \log a(J_n) + \log M_n + C_0 \log n\} \leq n\epsilon_n^2, \tag{2.3}$$

$$e(\bar{J}_n) \leq \bar{\epsilon}_n, \quad c_1 \bar{J}_n \log^{t_1} \bar{J}_n + c_3 \bar{J}_n \log(2a(\bar{J}_n)/\bar{\epsilon}_n) \leq a_2 n\bar{\epsilon}_n^2, \tag{2.4}$$

$$n\bar{\epsilon}_n^2 \leq C J_n \log^{t_2} J_n \text{ for any constant } C, \quad J_n \exp\{-CM_n^{t_3}\} \leq (a_1 - 1) \exp\{-n\bar{\epsilon}_n^2\}. \tag{2.5}$$

Let $\mathcal{W}_{J_n, M_n} = \{w = \theta^T \xi : \theta \in \mathbb{R}^j, j \leq J_n, \|\theta\|_\infty \leq M_n\}$. Then the following assertions hold:

$$\log D\left(n^{-C_0}, \mathcal{W}_{J_n, M_n}, d\right) \leq n\epsilon_n^2, \tag{2.6}$$

$$\Pi(W \notin \mathcal{W}_{J_n, M_n}) \leq a_1 \exp\{-bn\bar{\epsilon}_n^2\}, \tag{2.7}$$

$$-\log \Pi\{w = \theta^T \xi : d(w_0, w) \leq \bar{\epsilon}_n\} \leq a_2 n\bar{\epsilon}_n^2. \tag{2.8}$$

Proof. We first verify (2.6), using the definition of packing number, the assumptions on M_n , $J_n \geq 2$, the fact that $a(\cdot)$ is increasing and (2.1), we obtain

$$\begin{aligned} \log D\left(n^{-C_0}, \mathcal{W}_{J_n, M_n}, d\right) &\leq \log \left\{ \sum_{j=1}^{J_n} D\left(n^{-C_0}/a(j), \{\theta \in \mathbb{R}^j, \|\theta\|_\infty \leq M_n\}, \|\cdot\|_2\right) \right\} \\ &\leq \log \left[J_n \left\{ \sqrt{J_n} M_n a(J_n) n^{C_0} \right\}^{J_n} \right] \\ &\leq J_n (\log J_n + \log M_n + \log a(J_n) + C_0 \log n) \leq n\epsilon_n^2. \end{aligned} \tag{2.9}$$

Next, to verify (2.7), observe that for some $c'_2 > 0$,

$$\begin{aligned} \Pi(w \notin \mathcal{W}_{J_n, M_n}) &\leq \Pi(J > J_n) + \sum_{j=1}^{J_n} \Pi(\theta \notin [-M_n, M_n]^j) \Pi(J = j) \\ &\leq \exp(-c'_2 J_n \log^{t_2} J_n) + J_n \exp\{-CM_n^{t_3}\} \\ &\leq a_1 \exp\{-n\bar{\epsilon}_n^2\}. \end{aligned} \tag{2.10}$$

For (2.8), using (2.2), because $d(w_0, \theta_{0,j}^T \xi) \leq e(j) \leq \bar{\epsilon}_n$ for all $j \geq \bar{J}_n$, we have

$$\begin{aligned} \Pi\left\{w : d(w_0, \theta^T \xi) \leq 2\bar{\epsilon}_n\right\} &\geq \Pi(J = \bar{J}_n) \Pi(\|\theta - \theta_0\|_2 \leq \bar{\epsilon}_n/a(\bar{J}_n)) \\ &\geq \exp\{-c_1 \bar{J}_n \log^{t_1} \bar{J}_n\} \exp\left\{-c_3 \bar{J}_n \log\left(\frac{a(\bar{J}_n)}{\bar{\epsilon}_n}\right)\right\}. \end{aligned} \tag{2.11}$$

By taking the negative of the logarithm on both sides, and using (2.4), we obtain (2.8). □

Conditions (2.3)–(2.5) require sufficiently large J_n, \bar{J}_n in order to have sufficiently good approximation to w_0 while J_n, \bar{J}_n should not be too large if the complexity of the model is to be controlled. When studying Bayesian asymptotic properties, a balance between bias and complexity needs to be established to obtain the optimal posterior contraction rate.

Theorem 1 can be further simplified to obtain a posterior contraction rate at w_0 . We assume that the approximation error is of the form $e(J) \lesssim J^{-\alpha}$ for α -smooth functions. Such collections include B-splines, wavelets, Fourier series and many other commonly used bases. Let d be the L_2 -distance or the L_∞ -distance. For two groups of densities $p_{i,w_1}, p_{i,w_2}, i = 1, \dots, n$, we consider the root average squared Hellinger distance, defined by $\rho_n^2(w_1, w_2) = n^{-1} \sum_{i=1}^n h^2(p_{i,w_1}, p_{i,w_2})$. Note that when the observations are independent and identically distributed (i.i.d.), ρ_n reduces to the usual Hellinger distance. Then the following result gives the posterior contraction rate for various inference problems.

Theorem 2. *Suppose that we have independent observations X_i following some distributions with densities $p_{i,w}, i = 1, \dots, n$, respectively. Let $w_0 \in C^\alpha(\Omega_0)$ be the true value of w . Let r be either 2 or ∞ . Let $\epsilon_n \geq \bar{\epsilon}_n$ be two sequences of positive numbers satisfying $\epsilon_n \rightarrow 0$ and $n\bar{\epsilon}_n^2 \rightarrow \infty$ as $n \rightarrow \infty$. Assume that there exists a $\theta_0 \in \mathbb{R}^J, \|\theta_0\|_\infty \leq H$ and some positive constants C_1, C_2 and $K_0 \geq 0$ satisfying*

$$\|w_0 - \theta_0^T \xi\|_r \leq C_1 J^{-\alpha}, \tag{2.12}$$

$$\|\theta_1^T \xi - \theta_2^T \xi\|_r \leq C_2 J^{K_0} \|\theta_1 - \theta_2\|_2, \theta_1, \theta_2 \in \mathbb{R}^J. \tag{2.13}$$

Assume that the prior on J and θ satisfy Conditions (A1) and (A2). Let $J_n, \bar{J}_n \geq 2$ and M_n be sequences of positive numbers such that the following hold for positive constants a_3, a_4, c_3, c_4, C and any given constant $b > 0$:

$$bn\bar{\epsilon}_n^2 \leq J_n \log^{t_2} J_n, \log J_n + n\bar{\epsilon}_n^2 \leq M_n^{t_3}, \tag{2.14}$$

$$J_n\{(K_0 + 1) \log J_n + \log M_n + C_0 \log n\} \leq n\epsilon_n^2, \tag{2.15}$$

$$\bar{J}_n^{-\alpha} \leq \bar{\epsilon}_n, \bar{J}_n \{c_1 \log^{t_1} \bar{J}_n + c_3 K_0 \log(\bar{J}_n) + c_3 \log(1/\bar{\epsilon}_n)\} \leq 2n\bar{\epsilon}_n^2, \tag{2.16}$$

$$\rho_n(w_1, w_2) \lesssim n^{a_3} \|w_1 - w_2\|_r^{a_4} \text{ for any } w_1, w_2 \in \mathcal{W}_{J_n, M_n}, \tag{2.17}$$

$$\max \left\{ n^{-1} \sum_{i=1}^n K(p_{i, w_0}, p_{i, w}), n^{-1} \sum_{i=1}^n V(p_{i, w_0}, p_{i, w}) \right\} \leq C \|w_1 - w_2\|_r^2, \tag{2.18}$$

provided $\|w_1 - w_2\|_r$ is sufficiently small. Then the posterior distribution of w contracts at w_0 at the rate ϵ_n with respect to ρ_n .

Proof. In order to obtain the posterior contraction rate, we verify the following conditions as described in Theorem 4 of Ghosal & van der Vaart (2007a):

$$\log D(\epsilon_n, \mathcal{W}_{J_n, M_n}, \rho_n) \leq b_1 n \epsilon_n^2, \tag{2.19}$$

$$\Pi(w \notin \mathcal{W}_{J_n, M_n}) \leq b_3 \exp \left\{ -n \epsilon_n^2 \right\}, \tag{2.20}$$

$$\Pi(\mathcal{K}(w_0, \bar{\epsilon}_n)) \geq b_4 \exp \left\{ -b_2 n \bar{\epsilon}_n^2 \right\}, \tag{2.21}$$

where \mathcal{W}_{J_n, M_n} is defined in Theorem 1 and b_1, b_2, b_3, b_4 are some positive constants. Note that the conditions in Theorem 1 are satisfied for $a_1 = 1$ and $a_2 = 2$ in the following way: (2.2) is satisfied by the approximation assumption of ξ ; (2.15) implies (2.3); (2.16) implies (2.4); (2.5) holds because of (2.14). Using condition (2.17), we obtain

$$\log D(\epsilon_n, \mathcal{W}_{J_n, M_n}, \rho_n) \lesssim \log D(n^{-a_3} \epsilon_n^{a_4}, \mathcal{W}_{J_n, M_n}, \|\cdot\|_r) \lesssim n \epsilon_n^2$$

because $n^{-a_3} \epsilon_n^{a_4}$ is lower bounded by a polynomial in n^{-1} . Also, $\Pi(w \notin \mathcal{W}_{J_n, M_n}) \leq 2 \exp \left\{ -n \epsilon_n^2 \right\}$, therefore relation (2.20) holds for $b_3 = 2$. For (2.21), observe that $\Pi(\mathcal{K}(w_0, \bar{\epsilon}_n)) \geq \Pi(\|w - w_0\|_r \leq \bar{\epsilon}_n)$ so the conclusion holds for an appropriate adjustment of constants in the definitions of the rates $\bar{\epsilon}_n$ and ϵ_n . \square

Remark 2. For $r = 2$ or ∞ , relation (2.12) holds for polynomials, Fourier series, B-splines and wavelets. Relation (2.13) holds for B-splines, polynomials and Fourier series base with $K_0 = 1/2$ when $r = 2$ and $K_0 = 1$ when $r = \infty$. For wavelets, (2.13) holds with $K_0 = 1$ for $r = 2, \infty$. This is because $\|(\theta_1 - \theta_2)^T \xi\|_p \leq \sum_{j=1}^J |\theta_{1j} - \theta_{2j}| \max_{1 \leq j \leq J} \|\xi_j\|_p \leq \sqrt{J} \|\theta_1 - \theta_2\|_2 C_{p, J}$ for $C_{p, J} = \max_{1 \leq j \leq J} \|\xi_j\|_p$ and $1 \leq p \leq \infty$. For B-splines, polynomials and Fourier series bases, $C_{p, J} \asymp 1$ when $p = 2$ and $C_{p, J} \asymp \sqrt{J}$ when $p = \infty$. For wavelets, $C_{p, J} \asymp \sqrt{J}$ for $p = 2, \infty$.

Remark 3. It is possible to incorporate a finite-dimensional nuisance parameter η in our setup, such as a scale parameter in a normal regression model. In this case, the sieve will be defined as the product of \mathcal{W}_{J_n, M_n} with a suitable sieve for η whose metric entropy can be appropriately controlled and whose complement has exponentially small prior probability; see Remark 5 for a concrete analysis.

Theorem 2 suggests that in order to obtain adaptive posterior contraction rates, it is crucial to choose sequences $J_n, \bar{J}_n, \epsilon_n, M_n$ in the rate equations (2.14)–(2.16) and bound the KL divergences by the squared Euclidean distance $\|\cdot\|_r^2$. Bounding the KL divergence can be very different for various statistical problems, while the choices of J_n and \bar{J}_n are common for a set of basis functions. The following examples illustrate the use of the theorem.

Example 1. (Fourier trigonometric series). For a function $w_0 \in C^\alpha(0, 1)$, the best approximation has the error $e(J) \asymp J^{-\alpha}$ (Dai & Xu, 2013). Then the rate calculation proceeds in the following way: (2.16) implies $\bar{J}_n^{-\alpha} \lesssim \bar{\epsilon}_n$ and $\bar{J}_n \log n \lesssim n\bar{\epsilon}_n^2$, and hence $\bar{\epsilon}_n \asymp n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)}$ and $\bar{J}_n \asymp (n/\log n)^{1/(2\alpha+1)}$. Now use (2.14), we have $J_n \log^2 n \gtrsim n\bar{\epsilon}_n^2$, hence we choose $J_n \asymp n^{1/(2\alpha+1)}(\log n)^{2\alpha/(2\alpha+1)-t_2}$. Note that (2.15) implies $J_n \log n \lesssim n\epsilon_n^2$. As a result, we choose $\epsilon_n \asymp n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$.

Example 2. (Bernstein polynomials). We consider the Bernstein polynomial prior proposed by Petrone (1999). Consider a continuously differentiable density function w_0 with bounded second derivative, the approximation property of Bernstein polynomials to w_0 is $e(J) = C/J$ for some universal constant C and $r = 2$ (Lorenz, 2012). We can choose $\bar{J}_n = (n/\log n)^{1/3}$, $J_n = n^{1/3}(\log n)^{2/3-t_2}$, $\bar{\epsilon}_n = (n/\log n)^{-1/3}$ and $M_n = n^{1/t_3}$. The rate ϵ_n is $n^{-1/3}(\log n)^{1/3+(1-t_2)/2}$, which has the same polynomial power as given in Ghosal (2001). In fact, for any $0 \leq \alpha \leq 2$, the approximation rate of Bernstein polynomials is $J^{-\alpha/2}$, and the resulting posterior contraction rate is $n^{-\alpha/2(\alpha+1)}(\log n)^{\alpha/2(\alpha+1)}$ (Kruijer & van der Vaart, 2008). The poor contraction rate stems from the poor approximation rate of Bernstein polynomials. Kruijer & van der Vaart (2008) used coarsened Bernstein polynomials and showed that for any $f \in C^\alpha(0, 1)$ with $0 \leq \alpha \leq 1$, the approximation rate with J undetermined parameters is $J^{-\alpha}$. If we choose $\bar{J}_n \asymp (n/\log n)^{-\alpha/(2\alpha+1)}$, then the rate is $\epsilon_n = n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$, which adapts in the range $0 \leq \alpha \leq 1$.

Example 3. (Polynomial basis). Consider the orthogonal Legendre polynomials as the approximation tool for $w_0 \in C^\alpha(0, 1)$. The rate of approximation is identical with that of the Fourier series under the L_2 - or the L_∞ -metrics (e.g. Theorem 6.1 of Hesthaven *et al.* (2007)). Hence, the choice of J_n, M_n and rates are exactly the same with Example 1.

Example 4. (B-splines). If we choose the B-spline functions (Appendix) as the basis, then for $w_0 \in C^\alpha(0, 1)$, we have $e(J) \asymp J^{-\alpha}$ for either the L_2 or the L_∞ -distance. Thus, the choices of the sequences and the resulting rate ϵ_n are the same as in the case polynomial or Fourier basis. However, one distinguishing property of the B-spline basis is the non-negativity of the basis functions so positive linear combinations are positive. Further we show in Appendix B that coefficients of a B-spline basis expansion can be restricted appropriately if the target function satisfies some restrictions. The property will allow some special prior distribution on the coefficients so that posterior moments can be calculated without using MCMC techniques.

Example 5. (Wavelets). We consider a multiresolution truncated wavelet series

$$\sum_{k=1}^{2^m-1} \alpha_k \phi_k(x) + \sum_{j=0}^m \sum_{k=1}^{2^m-1} \beta_{jk} \psi_{jk}(x), \tag{2.22}$$

where the boundary corrected wavelet basis of Cohen *et al.* (1993) is used because the domain is the unit interval, which results in a finite number of terms in the earlier expansion. We put priors on m and wavelet coefficients α_k and β_{jk} for all possible values of j, k . It is well known that, for $w_0 \in C^\alpha(0, 1)$, the L_2 -approximation error is $e(m) = 2^{-m\alpha}$. Hence, we apply Theorem 2 for $J = 2^m$ and choose $\bar{J}_n = (n/\log n)^{1/(2\alpha+1)}$, $J_n = n^{1/(2\alpha+1)}(\log n)^{2\alpha/(2\alpha+1)-t_2}$, $M_n = n^{1/t_3}$ and $\bar{\epsilon}_n = (n/\log n)^{-\alpha/(2\alpha+1)}$. Doing the same calculation as in Example 1, the resulting rate ϵ_n is $n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$. This coincides with the adaptation

results for white noise models in Lian (2011) and for density estimation and regression models in Rivoirard & Rousseau (2012b).

Example 6. (Multivariate B-splines). Theorem 1 can be used in multi-dimensional situation as well. Consider the tensor-product B-splines (Schumaker, 2007) as a basis in $C^\alpha(0, 1)^s$. Then we have $e(J) \asymp J^{-\alpha/s}$ for $r = 2$ or ∞ , where $J = K^s$, and K is the number of univariate B-spline functions used in making the tensor products. Apply Theorem 2 with $\bar{J}_n = (n/\log n)^{1/(2\alpha+s)}$, $\bar{\epsilon}_n = (n/\log n)^{-\alpha/(2\alpha+s)}$, $M_n = n^{1/t_3}$ to obtain the rate ϵ_n as $n^{-\alpha/(2\alpha+s)}$ multiplied by some power of $\log n$, where the power depends on the statistical problem.

In these examples, we find that a power of $\log n$ is always present in the obtained rates. This is partly because we are dealing with a general class of problems. It is not clear whether such logarithmic terms can be removed and optimality can be established. Some negative results are given by Castillo *et al.* (2014), where a sharp rate with a precise logarithmic term is obtained under L_2 -loss. In some special situations, this logarithmic factor can be removed by using particular types of priors, such as Huang (2004), Ghosal *et al.* (2008) and Gao & Zhou (2013).

3. Density estimation

In this section, we illustrate how Theorem 2 can be used to obtain adaptive posterior contraction rate for both the univariate and the multivariate density estimation where in the latter case, the true density can be anisotropic, allowing different smoothness in different direction. We also discuss an MCMC-free method for calculating posterior moments by using a special conjugate-like prior on the coefficient vector.

3.1. Univariate density estimation

We consider estimation of a density defined on $(0, 1)$. Frequentist optimal rate of contraction $n^{-\alpha/(2\alpha+1)}$ was obtained for the maximum likelihood estimators in Hasminskii (1978). A Bayesian method using a log-spline prior was studied in Ghosal *et al.* (2000), where the optimal posterior contraction rate $n^{-\alpha/(2\alpha+1)}$ was obtained. When α is unknown, the adaptive posterior contraction rate $n^{-\alpha/(2\alpha+1)}$, possibly up to an additional logarithmic factor, was established in Ghosal *et al.* (2003, 2008).

Consider estimating a density function p on $(0, 1)$. A prior can be induced on p by using basis functions through a nonnegative, monotonic, locally Lipschitz continuous link function Ψ , that is, $p_w = \Psi(w) / \int_0^1 \Psi\{w(x)\}dx$ for $w = \theta^T \xi$, $\theta \in \mathbb{R}^J$, and J is given a prior on \mathbb{N} . If we choose Ψ as the exponential function and ξ as the B-spline, then it gives the log-spline prior. We can also choose Ψ as the identity function, use normalized B-spline basis and restrict the prior for θ on Δ_J without losing the approximation ability of original B-spline basis according to Lemma 1, part (d) in the Appendix.

Corollary 1. *Suppose that we have i.i.d observations X_1, \dots, X_n generated from a density p_0 , which satisfies $w_0 = \Psi^{-1}(p_0) \in C^\alpha(0, 1)$ and that w_0 is bounded in $[\underline{M}, \bar{M}]$ for some positive constants \underline{M} and \bar{M} . We assume that the prior satisfies Conditions (A1) and (A2), and the basis ξ satisfies (2.12) and (2.13) with $r = \infty$. If either $\log \Psi$ is Lipschitz continuous or $c(w) = \int_0^1 \Psi\{w(x)\}dx > \underline{C}$ for some constant $\underline{C} > 0$, then the posterior contraction rate is $\epsilon_n = n^{-\alpha/(2\alpha+1)} (\log n)^{\alpha/(2\alpha+1)+(1-t_1)/2}$ at p_0 with respect to the Hellinger distance.*

Proof. If w is uniformly close to w_0 , $\|\Psi(w) - \Psi(w_0)\|_\infty$ is small, and hence $c(w) = \int \Psi(w(x))dx$ is close to $\int \Psi(w_0(x))dx = \int p_0(x)dx = 1$, and hence is bounded in the

following. Thus, we have the estimate

$$\begin{aligned} \|p_w - p_0\|_\infty &\leq \left\| \frac{\Psi(w)}{c(w)} - \Psi(w) \right\|_\infty + \|\Psi(w) - \Psi(w_0)\|_\infty \\ &\leq (c(w))^{-1} |c(w) - c(w_0)| \|\Psi(w)\|_\infty + \|\Psi(w) - \Psi(w_0)\|_\infty \\ &\lesssim \|\Psi(w) - \Psi(w_0)\|_\infty \lesssim \|w - w_0\|_\infty. \end{aligned} \tag{3.1}$$

Note that because p_0 is bounded away from 0, so is p_w when $\|w - w_0\|_\infty$ is small. Now

$$h^2(p_0, p_w) = \int \frac{|p_0 - p_w|^2}{(\sqrt{p_0} + \sqrt{p_w})^2} \leq \frac{1}{M} \|p_0 - p_w\|_\infty^2 \lesssim \|w - w_0\|_\infty^2. \tag{3.2}$$

Using Lemma 8 of Ghosal & van der Vaart (2007b), we have

$$\begin{aligned} K(p_0, p_w) &\leq 2h^2(p_0, p_w) \left\| \frac{p_0}{p_w} \right\|_\infty \lesssim \|w - w_0\|_\infty^2, \\ V(p_0, p_w) &\lesssim h^2(p_0, p_w) \left(1 + \left\| \frac{p_0}{p_w} \right\|_\infty \right)^2 \lesssim \|w - w_0\|_\infty^2. \end{aligned} \tag{3.3}$$

Therefore, (2.18) holds for $r = \infty$. Next, we verify (2.17). Note that because of the i.i.d assumption, ρ_n is the Hellinger distance on p_w . As the Hellinger distance is bounded by the square root of the L_1 -distance, it suffices to bound the latter. If $\log \Psi$ is Lipschitz continuous with Lipschitz constant L , bound $\|p_{w_1} - p_{w_2}\|_1$ by

$$\begin{aligned} 2 \frac{\|\Psi(w_1) - \Psi(w_2)\|_1}{c(w_1)} &\leq 2 \|\exp[\log \Psi(w_1) - \log \Psi(w_2)] - 1\|_\infty \\ &\leq 2L \|w_1 - w_2\|_\infty e^{L\|w_1 - w_2\|_\infty}. \end{aligned}$$

On the other hand, if $c(w)$ is bounded in the following by \underline{C} , we obtain

$$\|p_{w_1} - p_{w_2}\|_1 \leq 2 \frac{\|\Psi(w_1) - \Psi(w_2)\|_1}{c(w_1)} \lesssim \|\Psi(w_1) - \Psi(w_2)\|_1 \lesssim \|w_1 - w_2\|_\infty,$$

so that the assertion holds with $a_4 = 1/2$.

Now, we apply Theorem 2 with $\bar{J}_n = (n/\log n)^{1/(2\alpha+1)}$, $J_n = n^{1/(2\alpha+1)} (\log n)^{2\alpha/(2\alpha+1)-t_2}$, $\bar{\epsilon}_n = (n/\log n)^{-\alpha/(2\alpha+1)}$, $M_n = n^{1/t_3}$ and $r = \infty$, then the posterior distribution contracts at the rate $\epsilon_n = n^{-\alpha/(2\alpha+1)} (\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$ at p_0 with respect to the Hellinger distance. \square

Remark 4. The commonly used exponential link function trivially satisfies the first requirement that $\log \Psi$ is Lipschitz continuous. The identity link function, used for MCMC-free calculation in section 3.3 satisfies $c(w) = \int w(x)dx = 1$ as w is a probability density, and hence the condition that $c(w)$ bounded away from zero trivially holds.

3.2. Anisotropic multivariate density estimation

We extend univariate density estimation to the multivariate situation by considering estimating an s -dimensional density function on $(0, 1)^s$. We induce a prior on the density using through the relation $p_w \propto \Psi(\theta^T \xi)$ with ξ chosen as the tensor-product B-spline basis of order q . The true density p_0 is allowed to have different smoothness levels at different directions. More precisely, we define an anisotropic Hölder smoothness class by

$$\mathcal{C}^\alpha(0, 1)^s = \left\{ f(x_1, \dots, x_s) : \left\| \frac{\partial^{\sum_{k=1}^s l_k} f}{\partial x_1^{l_1} \dots \partial x_s^{l_s}} \right\|_\infty < \infty, 0 \leq l_k \leq \alpha_k, k = 1, \dots, s, \sum_{k=1}^s l_k / \alpha_k < 1. \right\}$$

for some smoothness parameter $\alpha = (\alpha_1, \dots, \alpha_s)$, which are integers not greater than q . Let $J(1), \dots, J(s)$ be the number of basis functions for individual s directions and define J as their products. Given $\Psi^{-1}(p_0) \in \mathcal{C}^\alpha(0, 1)^s$, the approximation error is of the order $\sum_{k=1}^s J(k)^{-\alpha_k}$ according to Theorem 12.7 of Schumaker (2007). Hence, for the best balancing of the approximation error, we choose $J_n(k) = \bar{\epsilon}_n^{-1/\alpha_k}$ and $\bar{J}_n = \prod_{k=1}^s J_n(k) = \bar{\epsilon}_n^{-s/\alpha^*}$, where $\alpha^* = s / (\sum_{k=1}^s \alpha_k^{-1})$ is the harmonic mean of $\alpha_1, \dots, \alpha_s$, and $\bar{\epsilon}_n$ is to be chosen to match $\bar{\epsilon}_n^{-s/\alpha^*} \log n$ with $n\bar{\epsilon}_n^2$. Applying Theorem 2 with $\bar{\epsilon}_n = (n/\log n)^{-\alpha^*/(2\alpha^*+s)}$, $J_n = n^{s/(2\alpha^*+s)}(\log n)^{(2\alpha^*)/(2\alpha^*+s)-t_2}$, $M_n = n^{1/t_3}$, $a_4 = 1/2$ and $r = \infty$, the posterior distribution contracts at p_0 with respect to the Hellinger distance at the rate $\epsilon_n = n^{-\alpha^*/(2\alpha^*+s)}(\log n)^{\alpha^*/(2\alpha^*+s)+(1-t_2)/2}$. Essentially, the same rate is also obtained in Shen *et al.* (2013) (with a different logarithmic factor) using a Dirichlet mixture of normal prior.

3.3. Markov chain Monte Carlo-free computation

Next, we describe an MCMC-free calculation technique for the univariate density estimation using normalized B-splines $\{B_1^*, \dots, B_J^*\}$ as the basis (Appendix). By part (d) in Lemma 1, we can restrict the coefficients θ to a J -dimensional simplex Δ_J and maintain the same approximation rate. We put a Dirichlet prior on $\theta \sim \text{Dir}(a_1, a_2, \dots, a_J)$ for any $J \in \mathbb{N}$. Finally, we assign a prior Π on J . Thus, a prior on the density p is induced. Given the observations $X = (X_1, \dots, X_n)$ and a fixed dimension J , the posterior density of θ is a mixture of Dirichlet distribution:

$$p(\theta|X, J) \propto \prod_{k=1}^J \theta_k^{a_k-1} \prod_{i=1}^n \left\{ \sum_{k=1}^J \theta_k B_k^*(X_i) \right\} = \sum_{i_1=1}^J \dots \sum_{i_n=1}^J \prod_{k=1}^J \theta_k^{a_k-1} \prod_{s=1}^n \theta_{i_s} B_{i_s}^*(X_s).$$

Using $p(J, \theta|X) \propto p(X|J, \theta)\Pi(\theta|J)\Pi(J)$, the posterior mean of p at a point x is

$$\begin{aligned} & \frac{\sum_{j=1}^\infty \int_{\theta} p(x)p(X|J=j, \theta)\Pi(\theta|J=j)\Pi(J=j)d\theta}{\sum_{j=1}^\infty \int_{\theta} p(X|J=j, \theta)\Pi(\theta|J=j)\Pi(J=j)d\theta} \\ &= \frac{\sum_{j=1}^\infty \Pi(j) \sum_{i_0=1}^j \sum_{i_1=1}^j \dots \sum_{i_n=1}^j \int_{\theta \in \Delta_j} \prod_{k=1}^j \theta_k^{a_k-1} \prod_{s=0}^n \theta_{i_s} B_{i_s}^*(X_s)d\theta}{\sum_{j=1}^\infty \Pi(j) \sum_{i_1=1}^j \dots \sum_{i_n=1}^j \int_{\theta \in \Delta_j} \prod_{k=1}^j \theta_k^{a_k-1} \prod_{s=1}^n \theta_{i_s} B_{i_s}^*(X_s)d\theta}, \end{aligned} \tag{3.4}$$

where X_0 stands for x . Define $I_{k,j,0}^i = \sum_{s=0}^n \mathbb{1}\{i_s = k\}$ and $I_{k,j,1}^i = \sum_{s=1}^n \mathbb{1}\{i_s = k\}$. Then the expression in (3.4) can be simplified to

$$\frac{\sum_{j=1}^\infty \Pi(j) \sum_{i_0=1}^j \sum_{i_1=1}^j \dots \sum_{i_n=1}^j \prod_{k=1}^j \Gamma(a_k + I_{k,j,0}^i) \prod_{s=0}^n B_{i_s}^*(X_s) / \Gamma\left(\sum_{i=1}^j a_i + n + 1\right)}{\sum_{j=1}^\infty \Pi(j) \sum_{i_1=1}^j \dots \sum_{i_n=1}^j \prod_{k=1}^j \Gamma(a_k + I_{k,j,1}^i) \prod_{s=1}^n B_{i_s}^*(X_s) / \Gamma\left(\sum_{i=1}^j a_i + n\right)} \tag{3.5}$$

A basis function takes non-zero values only at q intervals, so the calculation involves a multiple of q^{n+1} steps. More details are given in section 5. Similar expressions can be obtained for other posterior moments, in particular, for the posterior variance.

Note that if $q = 1$, the sums over indices i_1, \dots, i_n in (3.5) will be redundant, leading to a histogram estimate whose bin length and weights are posterior averaged. The B-spline random series prior can also be viewed as a kernel mixture prior, where the kernel is a B-spline function indexed by a discrete parameter.

For multivariate situation, MCMC-free computational techniques can be developed in a similar way by using tensor products of normalized B-splines as the basis and a Dirichlet prior on the corresponding coefficients. The approximation property established in the last part of Lemma 2 justifies restricting the coefficients on the simplex.

4. Regression models

In this section, we consider several nonparametric regression problems including regression with additive Gaussian errors, binary regression, Poisson regression and functional regression. In these cases, we allow the covariates be either fixed or random and show how Theorem 2 can be used to derive contraction rates. The techniques also apply for multivariate analogues of these regression problems using the tensor-product B-spline basis as in subsection 3.2.

For fixed covariates \mathbf{Z} , define the empirical measure $\mathbb{P}_n^{\mathbf{Z}} = n^{-1} \sum_{i=1}^n \delta_{Z_i}$, and $\|\cdot\|_{2,n}$ as the norm on $L_2(\mathbb{P}_n^{\mathbf{Z}})$.

4.1. Nonparametric regression with Gaussian errors

We consider a regression model with additive error $X_i = f(Z_i) + \varepsilon_i$, where $\varepsilon_i \stackrel{\text{iid}}{\sim} N(0, \sigma^2)$, $Z_1, \dots, Z_n \in (0, 1)$. For ease of illustration, we first consider known σ and fixed covariates; the modification necessary for unknown σ and random covariates is outlined in Remark 5.

Corollary 2. *Suppose that the true regression function $f_0 \in C^\alpha(0, 1)$ and the prior satisfies Conditions (A1) and (A2). Given fixed covariates, assume that the basis ξ satisfies (2.12) and (2.13) with $r = \infty$. Then the posterior of f contracts at the rate $\epsilon_n = n^{-\alpha/(2\alpha+1)} (\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$ relative to $\|\cdot\|_{2,n}$ at f_0 .*

Proof. Let $P_{f,i}$ be the normal measure with mean $f(Z_i)$ and variance σ^2 . Then the Hellinger distance between $P_{f_1,i}$ and $P_{f_2,i}$ is of the order of the $|f_1(Z_i) - f_2(Z_i)|$ when one is, and hence both are, small. Hence, the conclusions of Lemma 2 of Ghosal & van der Vaart (2007a) hold (with different constants) for the distance $\|\cdot\|_{2,n}$. This implies that to compute entropy, we can work with $\|\cdot\|_{2,n}$ instead of ρ_n . Using the arguments in section 7.2 of Ghosal & van der Vaart (2007a), we obtain

$$\begin{aligned} & \max \left\{ n^{-1} \sum_{i=1}^n K(P_{f_0,i}, P_{f,i}), n^{-1} \sum_{i=1}^n V(P_{f_0,i}, P_{f,i}) \right\} \\ & \leq \|f_0 - f\|_{2,n}^2 / \sigma^2 \leq \|f_0 - f\|_\infty^2 / \sigma^2. \end{aligned} \tag{4.1}$$

Clearly, Condition (2.17) holds for $\|\cdot\|_{2,n}$ with $r = \infty$. Assuming that the basis ξ satisfies (2.12) and (2.13) with $r = \infty$ and choosing $\bar{J}_n = (n/\log n)^{1/(2\alpha+1)}$, $\bar{\epsilon}_n = (n/\log n)^{-\alpha/(2\alpha+1)}$, $J_n = n^{1/(2\alpha+1)} (\log n)^{2\alpha/(2\alpha+1)-t_2}$, $r = \infty$ and $M_n = n^{1/t_3}$, then we obtain the posterior contraction rate $\epsilon_n = n^{-\alpha/(2\alpha+1)} (\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$ relative to $\|\cdot\|_{2,n}$. □

Remark 5. For random covariates $Z_1, \dots, Z_n \sim G$, define $L_{2,G}$ as the L_2 -distance with respect to the probability measure G . We assume that G has a density g bounded and bounded away from zero, and ξ satisfies (2.12) and (2.13) with $r = 2$. Then $\|f_1 - f_2\|_{2,G}$ is equivalent to $\|f_1 - f_2\|_2$, and hence can be used interchangeably in entropy calculations and bounding prior concentration and posterior contraction rates. Alternatively, without any conditions on G , we can assume the basis has the L_∞ -approximation property with the same rate and bound $\|\cdot\|_{2,G}$ by $\|\cdot\|_\infty$. Hence, by applying Theorem 2 in the same way with $r = \infty$, we obtain the same rate with respect to $\|\cdot\|_{2,G}$.

When σ is unknown, we assign a prior (independent of other parameters) on it. If the prior density is positive throughout, and has exponential tail near zero and polynomial tail near infinity, then a sieve $(n^{-C_1}, \exp\{C_2 n \epsilon_n^2\})$ with sufficiently large C_1, C_2 will satisfy the conditions in Theorem 2. Note that the popular inverse gamma prior on σ^2 (or on any positive power of σ) satisfies the requirements.

4.2. Nonparametric binary regression

Assume that we have n independent observations $(Z_1, X_1), \dots, (Z_n, X_n)$ from a binary regression model $P(X = 1|Z = z) = 1 - P(X = 0|Z = z) = f_0(z)$, where X takes values in $\{0, 1\}$, and Z is either a fixed or a random covariate in some domain \mathcal{Z} . Given a link function $\Psi : \mathcal{Z} \rightarrow (0, 1)$, we can construct a random series prior on the regression function f_0 using a basis ξ as $f_\theta(z) = \Psi\{\theta^T \xi(z)\}$. Commonly, a cumulative distribution function on \mathbb{R} such as the logit or probit function is chosen as the link function and the coefficient vector θ can take any values in \mathbb{R}^J . Then any basis with approximation property for the Hölder class may be used.

Corollary 3. *Suppose that the true classification function f_0 is bounded away from 0 and 1, and satisfies $w_0 = \Psi^{-1}(f_0) \in \mathcal{C}^\alpha(0, 1)$. Given fixed covariates, and that the prior satisfies Conditions (A1) and (A2). Assume that the basis ξ satisfy (2.12) and (2.13) with $r = \infty$, and the link function Ψ is Lipschitz continuous. Then, the posterior of f contracts at the rate $\epsilon_n = n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$ relative to $\|\cdot\|_{2,n}$ at f_0 .*

Proof. Define $p_w = \Psi(w)^x(1 - \Psi(w))^{1-x}$, note that by the Lipschitz continuity of Ψ ,

$$h^2(p_{w_1}, p_{w_2}) \leq \|p_{w_1} - p_{w_2}\|_1 = 2\|\Psi(w_1) - \Psi(w_2)\|_\infty \lesssim \|w_1 - w_2\|_\infty,$$

$$\max\{K(p_{w_0}, p_w), V(p_{w_0}, p_w)\} \lesssim \|\Psi(w) - \Psi(w_0)\|_\infty^2 \lesssim \|w - w_0\|_\infty^2,$$

so the relation (2.17) holds with $a_4 = 1/2$. Now, we may apply Theorem 2 with $\bar{J}_n = (n/\log n)^{1/(2\alpha+1)}$, $M_n = n^{1/t_3}$, $J_n = n^{1/(2\alpha+1)}(\log n)^{2\alpha/(2\alpha+1)-t_1}$ and $\bar{\epsilon}_n = n^{-\alpha/(2\alpha+1)}(\log n)^{(\alpha+1)/(2\alpha+1)}$, then the posterior distribution contracts at the rate $\epsilon_n = n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$ relative to ρ_n . By Taylor's expansion of the squared Hellinger distance in a binomial model, it is easy to see that ρ_n is equivalent with the $\|\cdot\|_{2,n}$ -distance on f . □

For random covariates $Z_1, \dots, Z_n \sim G$, when G has a density g bounded and bounded away from zero, the same conclusion can be made in terms of the L_2 -distance on f , or more generally, with respect to the $L_2(G)$ -distance without any additional conditions.

When specifically the B-splines basis is used, the link function Ψ can be chosen to be the identity function in view of part (c) of Lemma 1. The expressions then simplify significantly if we use beta priors $\theta_i \stackrel{\text{ind}}{\sim} \text{Beta}(a_i, b_i)$ for some positive numbers a_i and b_i .

4.3. Nonparametric Poisson regression

Consider a Poisson regression model $X_i \stackrel{\text{ind}}{\sim} \text{Poi}\{f(Z_i)\}$, where f is an unknown monotonic function and Z is a covariate. For convenience, we assume that Z takes values in $(0, 1)$. Using a random series expansion, f can be modelled through a link function $f(z) = \Psi(\boldsymbol{\theta}^T \boldsymbol{\xi})(z)$.

Corollary 4. *Suppose that $\Psi^{-1}(f_0) \in \mathcal{C}^\alpha(0, 1)$ and f_0 is bounded away from zero and infinity. Let the prior satisfy Conditions (A1) and (A2). Assume that the basis $\boldsymbol{\xi}$ satisfies (2.12) and (2.13) with $r = \infty$, and the link function Ψ is monotonic and Lipschitz continuous on $(0, 1)$ and $\sqrt{\Psi}$ is Lipschitz continuous on $[1, \infty)$. Then the posterior of f contracts at the rate $\epsilon_n = n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$ relative to root-average squared Hellinger distance ρ_n at f_0 .*

Proof. In a Poisson model, the squared Hellinger distance is easily bounded by twice the square of the difference of the square roots of the parameters when the parameters are in $[1, \infty)$, while it is bounded by the L_1 -distance which is further bounded by the absolute difference of the parameters when they lie in $(0, 1)$. Hence, by the Lipschitz continuity assumptions on Ψ , with the choice $r = \infty$, the condition (2.17) holds with $a_3 = 0$ and $a_4 = 1/2$ when $\|w_1 - w_2\|_\infty$ is small. In fact, it is sufficient to assume that the Lipschitz continuity condition on $\sqrt{\Psi}$ holds with the Lipschitz constant growing up to polynomially in n on a sieve $\{\boldsymbol{\theta}^T \boldsymbol{\xi} \geq 1, \|\boldsymbol{\theta}\|_\infty \leq n^c \text{ and } J_n \leq n\}$. The KL divergences in Poisson model near a positive value of the parameter are bounded by a multiple of the square of the difference of parameter values, and a fixed constant can be chosen uniformly for all true parameter values lying in a compact subset of $(0, \infty)$. This leads to the verification of (2.18).

For any of the discussed basis functions, an application of Theorem 1 with the L_∞ -distance verifies the remaining conditions of Theorem 2 for $\bar{J}_n = (n/\log n)^{1/(2\alpha+1)}$, $M_n = n^{1/t_3}$, $J_n = n^{1/(2\alpha+1)}(\log n)^{(2\alpha+2)/(2\alpha+1)-t_1}$ and $\bar{\epsilon}_n = n^{-\alpha/(2\alpha+1)}(\log n)^{(\alpha+1)/(2\alpha+1)}$, then the posterior contraction rate is obtained as $\epsilon_n = n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_1)/2}$ relative to ρ_n . \square

To reinterpret this contraction rate in terms of the more desirable $\|\cdot\|_{2,n}$ -distance on f , we observe that these two rates are equivalent near the true regression function f_0 by its positivity and boundedness properties, provided that f remains in an L_∞ -bounded set with high posterior probability for most samples drawn from the true distribution. This is obviously ensured if coefficients obtain a prior confined in a bounded set, but will also hold if the posterior is consistent for the L_∞ -distance on f .

For random covariates $Z_1, \dots, Z_n \sim G$, the same contraction rate is obtained with respect to the Hellinger distance on the joint density of (X, Z) , and with respect to the $L_2(G)$ -distance on f under the aforementioned additional conditions.

If we use B-splines to form the basis, in view of Part (c) of Lemma 1, we are allowed to restrict θ_j to positive values. By choosing the identity link, then it is possible to carry out MCMC-free computation by letting $\theta_i \stackrel{\text{ind}}{\sim} \text{Gamma}(a_i, b_i)$ for some positive numbers a_i and b_i . The resulting prior satisfies all requirements for the posterior contraction rate obtained earlier.

4.4. Functional regression model

Spline functions are widely used to model functional data (Cardot *et al.*, 2003). A rate of contraction result was obtained in Hall & Horowitz (2007). A Bayesian method based on splines was given by Goldsmith *et al.* (2011). However, to the best of our knowledge, no results on posterior contraction rates for Bayesian methods are yet available. We consider two types of

functional regression model. The first one assumes only the covariates $Z(t)$ and the effects $\beta(t)$ depend on time t . The second one allows functional observations $X(t)$. We can use any basis with general approximation properties for Hölder classes under the L_2 -distance.

We first discuss the case of functional covariates with a scalar response. Suppose we observe i.i.d. copies $(Z_1, X_1), \dots, (Z_n, X_n)$ of (Z, X) , where Z is a square integrable random function defined on $(0, 1)$, and X is a scalar. A functional linear regression model can be formulated as follows:

$$X_i = \int_0^1 Z_i(t)\beta(t)dt + \varepsilon_i, \tag{4.2}$$

where $\beta(t)$ is the coefficient function we want to estimate, $\varepsilon_1, \dots, \varepsilon_n \stackrel{\text{iid}}{\sim} N(0, \sigma^2)$. We consider σ to be known; the more realistic case of unknown σ can be treated following Remark 5.

Corollary 5. *Suppose that the true regression function $\beta \in C^\alpha(0, 1)$, $EZ^2(t)$ is uniformly bounded away from 0 and ∞ for every $t \in (0, 1)$, and the basis satisfies (2.12) and (2.13) with $r = 2$. Given the prior being constructed as in (A1) and (A2), the posterior of f contracts in a rate $\epsilon_n = n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$ relative to the L_2 -distance.*

Proof. We consider a basis expansion $\beta(t) = \sum_{k=1}^J \theta_k \xi_k(t)$. Denote $W_{ik} = \int_0^1 Z_i(t)\xi_k(t)dt$, then the model can be written as $X_i = \sum_{k=1}^J \theta_k W_{ik} + \varepsilon_i$. Define $P_\beta(\cdot|Z)$ as the normal measure with mean $\int_0^1 Z(t)\beta(t)dt$ and variance σ^2 , and let E_Z be the expectation with respect to the distribution of Z . Then we can bound $K(P_{\beta_0}, P_\beta)$ and $V(P_{\beta_0}, P_\beta)$ using Cauchy–Schwarz inequality:

$$\max \{K(P_{\beta_0}, P_\beta), V(P_{\beta_0}, P_\beta)\} \lesssim \frac{1}{\sigma^2} E_Z \left(\int_0^1 Z(t)\{\beta(t) - \beta_0(t)\}dt \right)^2 \lesssim \frac{1}{\sigma^2} \|\beta - \beta_0\|_2^2.$$

For (2.17), note that the same argument used in random covariates situation in section 4.1 applies here. Hence, we can apply Theorem 2 as in section 4.1. Then the posterior contracts at the rate $\epsilon_n = n^{-\alpha/(2\alpha+1)}(\log n)^{\alpha/(2\alpha+1)+(1-t_2)/2}$ relative to the L_2 -distance. \square

Next, we consider a longitudinal type of functional model:

$$X_i(T_i) = Z_i(T_i)\beta(T_i) + \varepsilon_i. \tag{4.3}$$

For each object i , we observe its response X_i at a random time $T_i \in (0, 1)$ with a random covariate Z_i . We assume that Z_1, \dots, Z_n are i.i.d. copies of Z , and T_1, \dots, T_n are i.i.d. copies of T , $\varepsilon_i \stackrel{\text{iid}}{\sim} N(0, \sigma^2)$; they are all independent of each other, and T has a density g bounded and bounded away from zero on $(0, 1)$. Again it suffices to treat σ as known.

Suppose that the true regression function $\beta_0 \in C^\alpha(0, 1)$ and $EZ^2(t)$ are uniformly bounded away from 0 and ∞ for every $t \in (0, 1)$. Then again

$$\max \{K(P_{\beta_0}, P_\beta), V(P_{\beta_0}, P_\beta)\} \lesssim \frac{1}{\sigma^2} E \int_0^1 Z^2(t)(\beta(t) - \beta_0(t))^2 g(t)dt \lesssim \frac{1}{\sigma^2} \|\beta - \beta_0\|_2^2.$$

Hence, we obtain the same contraction rate ϵ_n if we use the same prior on β as before. This rate coincides with the optimal rate obtained in Cai & Yuan (2011) within a logarithmic factor.

5. Numerical examples

5.1. Simulation

We illustrate the use of conjugate prior structure as described in (3.4) and (3.5) on density estimation problems. We consider two examples of the true density: Beta(0.5, 0.5), and a mixture density of exponential and a normal distribution:

$$f_0(x) \propto \frac{3}{4}3e^{-3x} + \frac{1}{4} \frac{\sqrt{32}}{\sqrt{\pi}} e^{-32(x-0.75)^2}. \tag{5.1}$$

For each density, we generate $n = 20, 50, 100$ and 300 samples and then implement the random series prior for $q = 1$ and $q = 3$. When $q = 1$, the exact value of the posterior mean can be calculated. When $q = 3$, instead of evaluating all possible terms to obtain (3.5), we randomly sample $N = 3000$ of them and take the associated average values. We choose a geometric prior for J restricted between 5 and 25. The lower truncation ensures a minimum number of terms in the series expansion while an upper truncation is necessary to carry out the actual computation using a computer. For θ , we use the uniform distribution on the simplex as a default choice for the Dirichlet distribution. We obtain density estimates at 100 grid points in the unit interval.

We compare our results with that using the Gaussian process prior in Tokdar (2007) and Dirichlet mixture of normal kernels (Escobar & West, 1995). Mean absolute errors, mean squared errors (note that the theoretical results are obtained for Hellinger distance though) and computing time (in seconds) are summarized in Table 1. Standard errors are calculated based on 100 Monte Carlo replications. Comparing the performance of random series priors (RSP) using $q = 1$ with that of $q = 3$, we observe a trade-off between computation time and estimation accuracy. In terms of estimation accuracy, RSP ($q = 3$) beats Dirichlet mixture in both cases, but performs worse than Gaussian process for the mixture density estimation. Overall, RSP ($q = 1$) has the lowest computation cost because of its simple expression. It will

Table 1. *Density estimation results: mean squared error (l_2), mean absolute error (l_1) and computational time in seconds (t) using random series priors (RSP) with $q = 1$ and 3, Gaussian process (GP) and Dirichlet mixture (DM) priors*

True density		$n = 20$			$n = 50$			$n = 100$			$n = 300$		
		l_2	l_1	t	l_2	l_1	t	l_2	l_1	t	l_2	l_1	t
Mixture	RSP ($q = 1$)	.27	.40	.44	.20	.33	.58	.18	.30	.69	.17	.29	1.24
	RSP ($q = 3$)	.16	.31	255	.11	.25	317	.10	.24	320	.09	.22	425
	GP	.11	.23	53.9	.06	.17	58.8	.04	.14	61.5	.02	.10	66.9
	DM	.46	.59	9.6	.28	.44	22.0	.17	.34	33.6	.11	.28	99.5
	max s.e.	.01	.01	—	.01	.01	—	.01	.01	—	.00	.00	—
Beta (0.5, 0.5)	RSP ($q = 1$)	.35	.45	.45	.31	.42	.57	.27	.39	.67	.25	.37	1.23
	RSP ($q = 3$)	.16	.27	267	.15	.27	314	.14	.25	324	.11	.22	428
	GP	.34	.39	55.7	.27	.34	61.6	.24	.31	60.9	.19	.26	74.9
	DM	.31	.38	7.93	.32	.36	18.5	.27	.29	48.2	.25	.29	116
	max s.e.	.01	.01	—	.01	.01	—	.01	.01	—	.00	.00	—

s.e., standard errors.

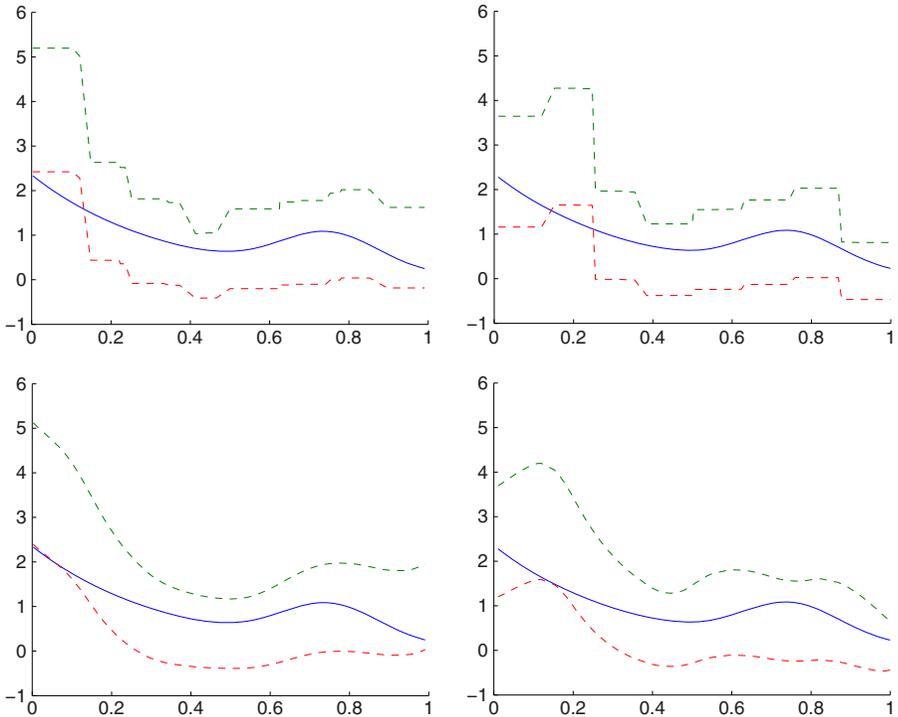


Fig. 1. Confidence bands for density estimation example (5.1). Dashed lines: confidence bands; solid line: true density. Upper left: $q = 1, n = 100$; upper right: $q = 1, n = 500$; lower left: $q = 3, n = 100$; lower right: $q = 3, n = 500$.

be interesting to consider a utility function that simultaneously evaluates the performance of estimators based on time and accuracy (Asmussen & Glynn, 2007).

Note that for RSP, the computational complexity becomes exponential in n given $q > 1$, and hence, all terms in the posterior mean cannot be computed for larger values of n . In this situation, we sample and compute a manageable number of terms and estimate the total as in sample survey for finite populations. The resulting standard error for sampling can be estimated in the usual way from the computed terms, and will be often reasonable if the terms are not very unlike each other.

We also calculate pointwise credible bands (95% nominal coverage) for the mixture true density example (5.1) based on the second moment estimation. Results are given in Fig. 1 for smoothness level $q = 1, 3$ and sample size $n = 100$ and 500. There is a significant improvement by using higher values of q . Establishing frequentist coverage properties of these intervals using Bernstein–von Mises results (e.g. (Rivoirard & Rousseau, 2012a)) will be an interesting but presumably delicate question.

5.2. Real data example

Next, we present a real data analysis of a functional linear model.

Example 7. The Tecator data (<http://lib.stat.cmu.edu/datasets/tecator>) provides an example of functional data where the spectra of meat samples are observed. The objective is to identify important chemical components and predict the fat content. The data consists of 172 training and 43 testing samples, where each sample contains 100 channel spectrum of absorbents.

We consider a functional linear model in (4.2) and use a B-spline basis expansion of $\beta(t)$. A prior is assigned by putting a Zellner's g-prior on the coefficients, a geometric distribution on J truncated between 5 and 15 and an inverse gamma distribution $IG(a, b)$ on σ^2 . We let the values of hyperparameters g, a, b range from 1 to 100, and the posterior results are quite insensitive. The MCMC-free calculation yields a root mean squared error (RMSE) of prediction 2.64 for $q = 1$ and $RMSE = 2.49$ for $q = 3$, which are generally better than the regression model results ($RMSE \geq 4$) built based on principal component analysis.

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Appendix: B-splines

Here, we provide a brief introduction to B-splines; more details are given in de Boor (2001). Let the unit interval $[0, 1]$ be divided into K equally spaced subintervals. Splines are continuous, piecewise polynomials of degree at most q , $(q - 2)$ times continuously differentiable and form a $J = q + K - 1$ dimensional linear space. B-splines provide a convenient basis for this space. B-splines are always nonnegative and add up to one, and each basis function is supported on an interval of length at most q/K .

Define the scaled B-spline basis functions $B_j^* = B_j / \int_0^1 B_j$, $j = 1, \dots, J$, so that $\int_0^1 B_j^*(z) dz = 1$, $j = 1, \dots, J$. Denote the column vector of B-spline basis functions by \mathbf{B} and that of the normalized B-spline basis functions by \mathbf{B}^* . The following results show some useful approximation properties of (tensor-product) B-splines.

Lemma 1.

- (a) For any function $f \in C^\alpha(0, 1)$, $0 < \alpha \leq q$, there exists $\boldsymbol{\theta} \in \mathbb{R}^J$ and a constant $C > 0$ that depends only on q such that $\|f - \boldsymbol{\theta}^T \mathbf{B}\|_\infty \leq C J^{-\alpha} \|f^{(\alpha)}\|_\infty$.
- (b) Further, if $f > 0$ we can choose every element of $\boldsymbol{\theta}$ to be positive.
- (c) If $0 < f < 1$, we can choose every element of $\boldsymbol{\theta}$ to be between 0 and 1.
- (d) Define $B_j^* = B_j / \int_0^1 B_j(z) dz$ for $j = 1, \dots, J$, and \mathbf{B}^* as the column vector (B_1^*, \dots, B_J^*) . If f is a density function, then there exists $\boldsymbol{\theta} \in \Delta_J$ and a constant $C > 0$ such that $\|f - \boldsymbol{\theta}^T \mathbf{B}^*\|_\infty \leq C J^{-\alpha} \|f^{(\alpha)}\|_\infty$.

Remark 6. In part (b), the condition $f > 0$ is crucial. If we approximate a nonnegative function f using nonnegative coefficients $\boldsymbol{\theta}$, then the approximation error is only $O(J^{-1})$ [cf. de Boor & Daniel (1974)], which does not adapt to smoothness levels beyond 1.

Proof of Lemma 1. The first part is a well-known spline approximation result, for example, Theorem 6.10 in Schumaker (2007).

For the second assertion, find $\epsilon > 0$ such that $f \geq \epsilon$. Using Corollaries 4 and 6 in Chapter 11 of de Boor (2001), for each θ_j , there exists a universal constant C_1 that depends only on q , such that $|\theta_j - c| \leq C_1 \sup_{x \in [t_{j+1}, t_{j+q-1}]} |f(x) - c|$ for any choice of the constant c ; here t_{j+1} and t_{j+q-1} are $(j + 1)$ th and $(j + q - 1)$ th knots. Choose $c = \inf_{z \in [t_{j+1}, t_{j+q-1}]} f(z) \geq \epsilon$ and note that the infimum is attained somewhere in $[t_{j+1}, t_{j+q-1}]$, say at t^* . By the smoothness condition on f , we have $\sup_{z \in [t_{j+1}, t_{j+q-1}]} |f(z) - c| \leq C_2 |z - t^*|^{\min(\alpha, 1)} \leq C_2 (q/J)^{\min(\alpha, 1)}$ for some constant $C_2 > 0$. Choosing $J > q(C_1 C_2 / \epsilon)^{\max(1/\alpha, 1)}$, we have $\theta_j > c - C_1 (q/J)^{\min(\alpha, 1)} \geq 0$.

Part (c) is a consequence of (b) by considering $1 - f > 0$.

For part (d), by (b), we know there exists a $\boldsymbol{\eta}_1 \in (0, \infty)^J$ such that $\|f - \boldsymbol{\eta}_1^T \mathbf{B}\| \lesssim J^{-\alpha}$. Define $\eta_{2,i} = \eta_{1,i} \int_0^1 B_j(z) dz$ for $j = 1, \dots, J$. Then $\|f - \boldsymbol{\eta}_2^T \mathbf{B}^*\|_\infty \lesssim J^{-\alpha}$, and in particular, $\|\boldsymbol{\eta}_2^T \mathbf{B}\|$ is bounded. By integration, we have $|1 - \|\boldsymbol{\eta}_2\|_1| = \left| 1 - \sum_{j=1}^J \eta_{2,j} \right| \lesssim J^{-\alpha}$. Choose $\boldsymbol{\theta} = \boldsymbol{\eta}_2 / \|\boldsymbol{\eta}_2\|_1 \in \Delta_J$. Note that $\|f - \boldsymbol{\theta}^T \mathbf{B}^*\|_\infty \leq \|f - \boldsymbol{\eta}_2^T \mathbf{B}^*\|_\infty + \|\boldsymbol{\eta}_2^T \mathbf{B}^*\|_\infty |1 - \|\boldsymbol{\eta}_2\|_1^{-1}| \lesssim J^{-\alpha}$. □

Lemma 2. Let $B_{j_1 \dots j_s}(z_1, \dots, z_s) = \prod_{k=1}^s B_{j_k}(z_k)$, $1 \leq j_k \leq J_k$, $k = 1, \dots, s$, be the tensor products of B-splines formed by univariate B-splines in J_k -dimensional space of splines, $k = 1, \dots, s$, respectively.

- (a) For any function $f \in C^\alpha(0, 1)^s$, the anisotropic Hölder class defined in section 3.2, where $\alpha_1, \dots, \alpha_s$ are positive integers less than or equal to q , there exists $\theta = (\theta_{j_1 \dots j_s} : 1 \leq j_k \leq J_k, k = 1, \dots, s) \in \mathbb{R}^{\prod_{k=1}^s J_k}$ and a constant $C > 0$ that depends only on q such that

$$\|f - \theta^T \mathbf{B}\|_\infty \leq C \sum_{k=1}^s J_k^{-\alpha_k} \left\| \frac{\partial^{\alpha_k} f}{\partial z_k^{\alpha_k}} \right\|_\infty.$$

- (b) Further, if $f > 0$ we can choose every component of θ to be positive.
 (c) If $0 < f < 1$, we can choose every element of θ to be between 0 and 1.
 (d) Define $B_{j_1 \dots j_s}^* = B_{j_1 \dots j_s} / \int_0^1 B_{j_1 \dots j_s}(z_1, \dots, z_s) dz_1 \dots dz_s$, for $j_k = 1, \dots, J_k$, $k = 1, \dots, s$, and \mathbf{B}^* as the column vector formed by the collection $B_{j_1 \dots j_s}^*$. If f is a density function, then there exists $\theta \in \Delta_{\prod_{k=1}^s J_k}$ and a constant $C > 0$ such that the same approximation order is maintained.

Proof. The first assertion is established in Theorem 12.7 in Schumaker (2007).

Proof of the second assertion proceeds as in the corresponding part of Lemma 1 using the parallel properties of tensor products of B-splines. The only relation we need to verify is $|\theta_{j_1 \dots j_s} - c| \leq C_1 \max_k \sup_{z_k \in [t_{i+1,k}, t_{i+q-1,k}]} |f(z) - c|$ for any choice of the constant c ; here $t_{i+1,k}$ and $t_{i+q-1,k}$ are $(i + 1)$ th and $(i + q - 1)$ th knots on the k th co-ordinate, $k = 1, \dots, s$. As in the univariate case, because the sum of all multivariate B-splines is one, to establish the relation, we need to bound absolute values of the coefficients using the values of the target function. Clearly, a dual basis for the multivariate B-splines is formed by tensor products of univariate dual bases, and these can be chosen to be uniformly bounded [see Theorem 4.41 of Schumaker (2007)]. Using such a dual basis, the maximum value of coefficients of spline approximations is bounded by a constant multiple of the L_∞ -norm of the target function. This gives the desired bound.

Parts (c) and (d) are established following exactly the same arguments used in the respective parts in the Proof of Lemma 1. □

Remark 7. In the isotropic case $\alpha = (\alpha, \dots, \alpha)$, the value of α need not be restricted to integers only — any $\alpha \leq q$ can be treated. This is because in this case the approximation error $\|f - \theta^T \mathbf{B}\|_\infty$ for the best multivariate spline approximation for $f \in C^\alpha(0, 1)^s$ with J terms in each direction decays at the rate $J^{-\alpha}$ for any positive $\alpha \leq q$.