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# Nonlinear regression modeling and spike detection via Gaussian basis expansions

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## Abstract

We consider the problem of constructing nonlinear regression models in the case that the structure of data has abrupt change points at unknown points. We propose two stage procedure where the spikes are detected by fused lasso signal approximator at the first stage, and the smooth curve are effectively estimated along with the technique of regularization method at the second. In order to select tuning parameters in the regularization method, we derive a model selection criterion from information-theoretic viewpoints. Simulation results and real data analysis demonstrate that our methodology performs well in various situations.

*Key Words and Phrases:* Basis expansion, Information criterion, Lasso, Nonlinear regression, Regularization, Spike detection.

## 1 Introduction

For analysis of data with complex structure, flexible model is absolutely imperative. As a useful tool to analyse such diverse phenomena, nonlinear regression model based on basis expansions is widely used. The essential idea behind basis expansions is to express a regression function as a linear combination of known functions, called basis functions (Bishop, 2006; Konishi and Kitagawa, 2008; Hastie *et al.*, 2009). In constructing the model, various functions are used to represent a regression function according to the structure of data or analysis objective. For example, natural cubic splines (Green and Silverman, 1994),  $B$ -splines (Eilers and Marx, 1996; de Boor, 2001; Imoto and Konishi, 2003) and radial basis functions (Kawano and Konishi, 2007; Ando *et al.*, 2008; Hastie *et al.*, 2009) involving Gaussian basis functions, thin plate splines and so on. In particular, Gaussian basis functions have been most commonly used to construct nonlinear regression

models. Gaussian basis functions have center and width parameters that have to be determined from observed data. The width parameters adjust the amount of overlapping among basis functions and notably play the essential role to capture the structure in the data over the region of the input space. Moody and Darken (1989) used the  $k$ -means clustering algorithm to determine the width parameters. In applying nonlinear regression models based on  $k$ -means-based Gaussian basis functions, it is assumed that the structure of phenomena are smooth.

However, the underlying true structure which is generating data cannot be smooth in practice at some points where jump discontinuity may occur. Thus, the application of a usual nonlinear regression model described above will lead difficulty of obtaining effective information from the data in which the mean structure is suddenly changed.

In order to overcome this problem, we propose the method of appropriately estimating a nonlinear structure with the spiky change points by applying fused lasso signal approximator (FLSA; Friedman *et al.*, 2007) which is a special version of fused lasso (Tibshirani *et al.*, 2005) to construction of Gaussian basis functions. We present a two-stage procedure to fit spiky regression curve.

In the first stage, we apply FLSA estimation procedure to get the information of locations of spikes. FLSA estimation encourages flatness between the resembled data and the abrupt change before and behind the jump. And then, we introduce the new Gaussian basis functions based on FLSA estimate. In the second stage, we fit the nonlinear regression model with the FLSA-based Gaussian basis functions by the method of regularization. The unknown parameters in regression model are estimated by maximizing the penalized log-likelihood function.

It is a crucial issue to determine the tuning parameters, including a smoothing parameter and a hyperparameter associated with Gaussian basis functions. To choose these parameters, we derive model selection criterion from information-theoretic viewpoint. The proposed nonlinear modeling procedure is investigated through the numerical examples.

This paper is organized as follows. Section 2 describes the framework of basis expansions and Gaussian basis function models. In Section 3 we present a new Gaussian basis functions based on FLSA. Section 4 provides nonlinear regression modeling strate-

gies based on proposed asymmetric Gaussian basis functions using maximum penalized likelihood estimation procedure. Section 5 gives a model selection criterion for evaluating statistical models estimated by the regularization method. In Section 6 we investigate the performance of our nonlinear regression modeling techniques through Monte Carlo simulations and real data example. Some concluding remarks are presented in Section 7.

## 2 Nonlinear regression model with basis expansions

Suppose that we have  $n$  independent observations  $\{(y_\alpha, x_\alpha); \alpha = 1, 2, \dots, n\}$ , where  $y_\alpha$  are random response variables and  $x_\alpha$  are explanatory variables. We consider the regression model

$$y_\alpha = g(x_\alpha) + \epsilon_\alpha, \quad \alpha = 1, 2, \dots, n, \quad (1)$$

where  $g(\cdot)$  is an unknown smooth function and  $\epsilon_\alpha$  are independently, normally distributed with mean zero and variance  $\sigma^2$ . It is assumed that the function  $g(\cdot)$  can be expressed as a linear combination of basis functions  $b_j(x)$  ( $j = 1, 2, \dots, m$ ) in the form

$$g(x; \mathbf{w}) = w_0 + \sum_{j=1}^m w_j b_j(x) = \mathbf{w}^T \mathbf{b}(x), \quad (2)$$

where  $\mathbf{b}(x) = (1, b_1(x), \dots, b_m(x))^T$  is a vector of basis functions and  $\mathbf{w} = (w_0, w_1, \dots, w_m)^T$  is an unknown coefficient parameter vector. A variety of basis functions are used according to the structure of data.

One of the many basis functions is Gaussian basis function given by

$$b_j(x) = \exp \left\{ -\frac{(x - c_j)^2}{2h_j^2} \right\}, \quad j = 1, 2, \dots, m, \quad (3)$$

where  $c_j$  is the center of the basis function,  $h_j^2$  is a parameter that determines the dispersion. However, basis functions (3) often yield inadequate results because of the lack of overlapping among basis functions. In order to overcome this problem, Ando *et al.* (2008) proposed the use of Gaussian basis functions with a hyperparameter, i.e. functions of the form

$$b_j(x) = \exp \left\{ -\frac{(x - c_j)^2}{2\nu h_j^2} \right\}, \quad j = 1, 2, \dots, m, \quad (4)$$

where  $\nu$  is a hyperparameter that adjusts the dispersion of basis functions. Ando *et al.* (2008) showed that nonlinear models with these basis functions were effective in capturing the information from the data.

Unknown parameters in the Gaussian basis functions (4) include the centers  $c_j$  and dispersion parameters  $h_j^2$ . These parameters are generally determined by using unsupervised learning. For example, Moody and Darken (1989) determined the centers  $c_j$  and dispersion  $h_j^2$  using the  $k$ -means clustering algorithm to avoid local minimum and identification problems. The data set of observations of the explanatory variables  $\{x_1, \dots, x_n\}$  is divided into  $m$  clusters  $\{C_1, \dots, C_m\}$ ; centers  $c_j$  and dispersions  $h_j^2$  are determined by

$$\hat{c}_j = \frac{1}{n_j} \sum_{x_\alpha \in C_j} x_\alpha, \quad \hat{h}_j^2 = \frac{1}{n_j} \sum_{x_\alpha \in C_j} (x_\alpha - \hat{c}_j)^2, \quad (5)$$

where  $n_j$  is the number of observations included in the  $j$ th cluster  $C_j$ . Replacing  $c_j$  and  $h_j^2$  in equation (3) by  $\hat{c}_j$  and  $\hat{h}_j^2$  respectively, we obtain a set of  $m$  basis functions

$$b_j(x; \hat{c}_j, \hat{h}_j^2) = \exp \left\{ -\frac{(x - \hat{c}_j)^2}{2\nu\hat{h}_j^2} \right\}, \quad j = 1, 2, \dots, m. \quad (6)$$

However, the models with these  $k$ -means-based Gaussian basis functions will lead to smooth curve estimates, even though spikes are present. Therefore, they will underestimate spikes and sudden change points will not be visible in resulting curve. In order to overcome this problem, we construct new Gaussian basis functions using fused lasso signal approximator (FLSA; Friedman *et al.*, 2007).

### 3 Gaussian basis functions based on FLSA

For  $n$  independent observations  $\{(y_\alpha, x_\alpha); \alpha = 1, \dots, n\}$ , FLSA procedure minimizes

$$\frac{1}{2} \sum_{\alpha=1}^n (y_\alpha - \beta_\alpha)^2 + \lambda_1 \sum_{\alpha=1}^n |\beta_\alpha| + \lambda_2 \sum_{\alpha=1}^{n-1} |\beta_\alpha - \beta_{\alpha+1}|, \quad \alpha = 1, \dots, n, \quad (7)$$

where  $\beta_\alpha$  is an estimate of response variable  $y_\alpha$  taken at position  $\alpha$ , and  $(\lambda_1, \lambda_2)$  are positive smoothing parameters to be chosen appropriately. The first penalty encourages sparsity in  $\beta$ s and second penalty encourages sparsity in their neighboring differences. As the value of  $\lambda_2$  increases, the number of fused parameters increases. Applying FLSA, the data set of  $n$  observations of the explanatory variables  $\{x_1, \dots, x_n\}$  is divided into  $n_F$

sets  $\{U_1, \dots, U_{n_F}\}$  where  $U_j = \{x_\alpha | \alpha \in F_j\} = \{x_1^{F_j}, \dots, x_{n_j}^{F_j}\}$ ,  $F_j$  is the  $j$ th subset of  $\{1, \dots, n\}$  that are considered to be fused at the  $\lambda_2$ , that is,  $\hat{\beta}_{\alpha_1} = \hat{\beta}_{\alpha_2}$  ( $\forall \alpha_1, \alpha_2 \in F_j$ ) and  $n_F$  is the number of such sets.

We take the sets  $U_j$  ( $j = 1, \dots, n_F$ ) the clusters and then we construct the asymmetric Gaussian basis functions based on FLSA as follows:

$$b_j(x) = \begin{cases} \exp\left\{-\frac{(x - c_j)^2}{2\nu h_{1j}^2}\right\}, & (x < c_j) \\ 1, & (x = c_j) \\ \exp\left\{-\frac{(x - c_j)^2}{2\nu h_{2j}^2}\right\}, & (x > c_j) \end{cases} \quad j = 1, 2, \dots, m. \quad (8)$$

We take the number of clusters  $n_F$  as that of basis functions (i.e.  $m = n_F$ ) and set basis functions on each center of clusters. And then we determine dispersion parameters using gradients of FLSA estimates; for fixed  $N_{ij}$ , centers  $c_j$  and dispersions  $h_j^2$  are determined by

$$\hat{c}_j = \frac{1}{n_j} \sum_{x_\alpha \in U_j} x_\alpha, \quad \hat{h}_{ij}^2 = \frac{N_{ij}}{|G_{ij}|} \exp\left(-\frac{1}{2}N_{ij}^2\right), \quad i = 1, 2, \quad j = 1, \dots, m, \quad (9)$$

where  $n_j$  is the number of observations included in the  $j$ th clusters  $F_j$ , and we defined gradients  $G_{ij}$  as follows:

$$G_{1j} = \frac{\hat{\beta}_{F_j} - \hat{\beta}_{F_{j-1}}}{x_{n_j}^{F_j} - x_{n_{j-1}}^{F_{j-1}}}, \quad G_{2j} = \frac{\hat{\beta}_{F_{j+1}} - \hat{\beta}_{F_j}}{x_1^{F_{j+1}} - x_1^{F_j}}. \quad (10)$$

These dispersion parameters  $\hat{h}_{ij}$  are the solutions that the gradients of Gaussian basis function (3) at  $c_j - N_{ij}h_{ij}$  and  $c_j + N_{2j}h_{ij}$  equal  $G_{1j}$  and  $G_{2j}$  respectively.  $N_{ij}$  are parameters which adjust the widths of basis functions (8). As value of  $N$  increases, the width of Gaussian basis function decreases. In this study, we fixed  $N$  as follows:

$$N_{ij} = \begin{cases} 3, & (\exp(-1/2)/|G_{ij}| < \delta) \\ 1, & (\text{otherwise}) \end{cases} \quad (11)$$

where  $\delta$  is enough small.

Replacing  $c_j$  and  $h_j^2$  in equation (8) by  $\hat{c}_j$  and  $\hat{h}_j^2$  respectively, we obtain a set of  $m$  basis functions

$$b_j(x) = \begin{cases} \exp\left\{-\frac{(x - \hat{c}_j)^2}{2\nu \hat{h}_{1j}^2}\right\}, & (x < \hat{c}_j) \\ 1, & (x = \hat{c}_j) \\ \exp\left\{-\frac{(x - \hat{c}_j)^2}{2\nu \hat{h}_{2j}^2}\right\}, & (x > \hat{c}_j) \end{cases} \quad j = 1, 2, \dots, m. \quad (12)$$

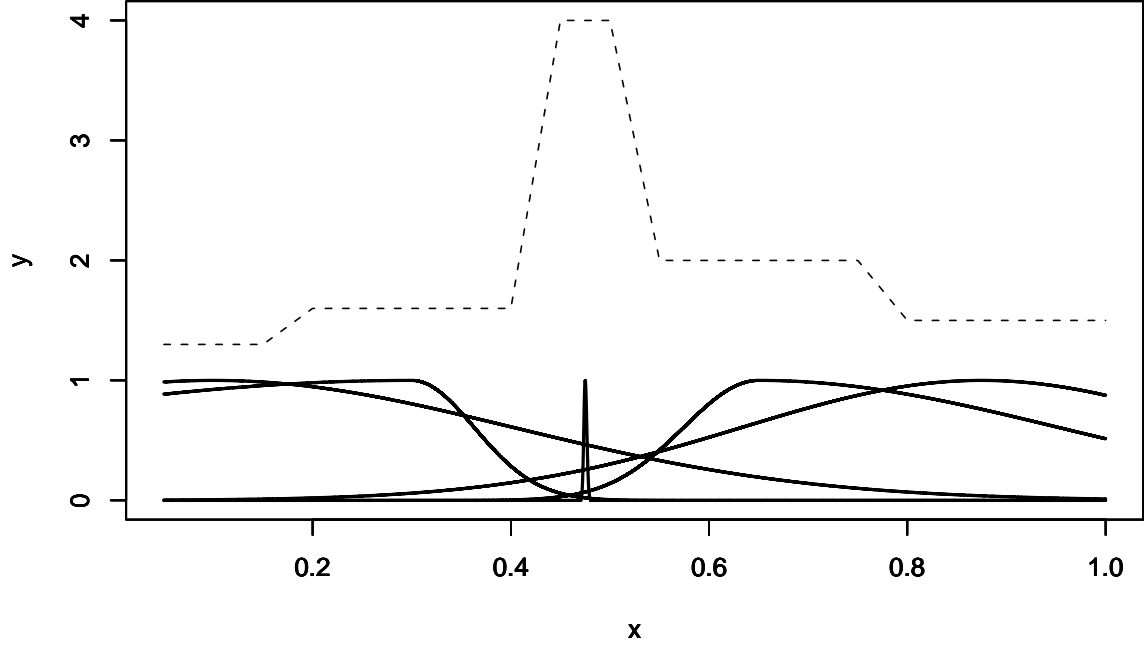


Fig. 1: Simple example of constructing our proposed Gaussian basis functions for  $\nu = 1$  (solid line) based on FLSA estimates (dashed line).

The width of Gaussian basis function  $b(x)$  will be very small if the absolute gradients of FLSA  $|G_{1j}|$  and  $|G_{2j}|$  are very large. Figure 1 shows FLSA estimates and our proposed Gaussian basis functions. Actually, in Figure 1, the central basis function become very narrow compared with the others.

## 4 Maximum penalized likelihood estimation

For  $n$  independent observations  $\{(y_\alpha, x_\alpha); \alpha = 1, \dots, n\}$ , the nonlinear regression model based on Gaussian basis functions  $b_j(x)$  ( $j = 1, \dots, n$ ) given in Section 3 is expressed as

$$y_\alpha = \mathbf{w}^T \mathbf{b}(x_\alpha) + \epsilon_\alpha, \quad \alpha = 1, \dots, n, \quad (13)$$

where  $\mathbf{b}(x_\alpha) = (1, b_1(x_\alpha), \dots, b_m(x_\alpha))^T$ ,  $\mathbf{w} = (w_0, w_1, \dots, w_m)^T$  and  $\epsilon_\alpha$  are error terms. If the error terms  $\epsilon_\alpha$  are independently and normally distributed with mean 0 and variance



$\sigma^2$ , the nonlinear regression model (13) has a probability density function

$$f(y_\alpha|\mathbf{w}, \sigma^2) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp \left[ -\frac{\{y_\alpha - \mathbf{w}^T \mathbf{b}(x_\alpha)\}^2}{2\sigma^2} \right], \quad \alpha = 1, \dots, n. \quad (14)$$

And then, the coefficient parameters  $w_j$  ( $j = 0, 1, \dots, m$ ) are estimated by the maximum penalized likelihood method.

The maximum likelihood estimates of the coefficient vectors  $\mathbf{w}$  and  $\sigma^2$  are respectively given by

$$\hat{\mathbf{w}} = (B^T B)^{-1} B^T \mathbf{y}, \quad \hat{\sigma}^2 = \frac{1}{n} (\mathbf{y} - B\hat{\mathbf{w}})^T (\mathbf{y} - B\hat{\mathbf{w}}),$$

where  $B = (\mathbf{b}(x_1)^T, \dots, \mathbf{b}(x_n)^T)^T$  and  $\mathbf{y} = (y_1, \dots, y_n)^T$ . However, when fitting a nonlinear model to data with a complex structure the maximum likelihood method often yields unstable estimates and leads to overfitting. We therefore estimate  $\mathbf{w}$  and  $\sigma^2$  by the method of regularization. Instead of using the log-likelihood function, we consider maximizing the penalized log-likelihood function imposing ridge penalty (Hoerl and Kennard, 1970)

$$l_\gamma(\boldsymbol{\theta}) = \sum_{\alpha=1}^n \log f(y_\alpha|\mathbf{w}, \sigma^2) - \frac{n\gamma}{2} \mathbf{w}^T K \mathbf{w}, \quad (15)$$

where  $\boldsymbol{\theta} = (\mathbf{w}^T, \sigma^2)^T$ ,  $\gamma (> 0)$  is a smoothing parameter that controls the smoothness of the fitted model and  $K$  is a known  $(m+1)$ th square matrix (Konishi and Kitagawa, 2008). The typical form of  $K$  is given by  $K = I_{m+1}$  for the identity matrix or  $K = D_2^T D_2$  for a second-order difference matrix. Then, the maximum penalized likelihood estimates of  $\mathbf{w}$  and  $\sigma^2$  are respectively given by

$$\hat{\mathbf{w}} = (B^T B + n\gamma \hat{\sigma}^2 K)^{-1} B^T \mathbf{y}, \quad \hat{\sigma}^2 = \frac{1}{n} (\mathbf{y} - B\hat{\mathbf{w}})^T (\mathbf{y} - B\hat{\mathbf{w}}). \quad (16)$$

Note that these estimators depend on each other. Therefore, we provide an initial value for the variance  $\sigma_{x(0)}^2$  first, then  $\hat{\mathbf{w}}$  and  $\hat{\sigma}_x^2$  are updated until convergence. The ridge estimators continuously shrink the coefficients as  $\gamma$  increases.

## 5 Model selection criterion

The statistical model estimated by the regularization method depends upon the value of the smoothing parameter  $\gamma$  and the value of the hyperparameter  $\nu$  in the Gaussian basis functions. It is a crucial issue to determine these values appropriately.

Konishi and Kitagawa (1996, 2008) introduced evaluation criteria of statistical models that can be applied to the evaluation of statistical models estimated by various types of estimation procedures such as the robust and penalized likelihood procedures. By using the result, the model selection criterion for evaluating the statistical model constructed by Gaussian basis functions is given by

$$\text{GIC} = n\{\log(2\pi) + 1\} + n \log \hat{\sigma}^2 + 2\text{tr}\{R^{-1}Q\}, \quad (17)$$

where  $R$  and  $Q$  are  $(m+2)$ th square matrices and are, respectively, given by

$$R = \frac{1}{n\hat{\sigma}^2} \begin{bmatrix} B^T B + n\gamma\hat{\sigma}^2 K & \frac{1}{\hat{\sigma}^2} B^T \Lambda \mathbf{1}_n \\ \frac{1}{\hat{\sigma}^2} \mathbf{1}_n^T \Lambda B & \frac{n}{2\hat{\sigma}^2} \end{bmatrix}, \quad (18)$$

$$Q = \frac{1}{n\hat{\sigma}^2} \begin{bmatrix} \frac{1}{\hat{\sigma}^2} B^T \Lambda^2 B - \gamma K \hat{\mathbf{w}} \mathbf{1}_n^T \Lambda B & \frac{1}{2\hat{\sigma}^4} B^T \Lambda^3 \mathbf{1}_n - \frac{1}{2\hat{\sigma}^2} B^T \Lambda \mathbf{1}_n \\ \frac{1}{2\hat{\sigma}^4} \mathbf{1}_n^T \Lambda^3 B - \frac{1}{2\hat{\sigma}^2} \mathbf{1}_n^T \Lambda B & \frac{1}{4\hat{\sigma}^6} \mathbf{1}_n^T \Lambda^4 \mathbf{1}_n - \frac{n}{4\hat{\sigma}^2} \end{bmatrix} \quad (19)$$

with  $\mathbf{1}_n = (1, \dots, 1)^T$  and  $\Lambda = \text{diag}(y_1 - \hat{\mathbf{w}}^T \mathbf{b}(x_1), \dots, y_n - \hat{\mathbf{w}}^T \mathbf{b}(x_n))$ . We obtain the optimal tuning parameters that minimize GIC.

## 6 Numerical examples

In this section, we investigate the performance of our nonlinear regression modeling techniques through Monte Carlo simulations and real data example. In FLSA procedure, we fixed  $\lambda_1 = 0$  and select the value of  $\lambda_2$  by using cross validation (CV) and we use an identity matrix as  $K$  in (15). In nonlinear regression modeling procedure with our proposed Gaussian basis functions based on FLSA (FLSA-G) and  $k$ -means-clustering (KM-G), the model selection criterion GIC was used for choosing optimal values of  $(\gamma, \nu)$  and  $(\gamma, \nu, m)$  respectively.

### 6.1 Monte Carlo simulations

We conducted Monte Carlo simulations to investigate the effectiveness of our proposed nonlinear regression modeling. For the first simulation study, repeated random samples  $\{(x_\alpha, y_\alpha); \alpha = 1, \dots, n\}$  with  $n = 100$  were generated from a true regression model  $y_\alpha = u(x_\alpha) + v_S(x_\alpha) + \epsilon_\alpha$ , where  $u(x)$  is smooth function and  $v_S(x)$  is spike function that has

peak region  $S \subset \{x_1, \dots, x_n\}$ . The  $x_\alpha = \alpha/n$  are  $n$  equally spaced design points in the interval  $[0, 1]$  and the errors  $\epsilon_\alpha$  are independently, normally distributed with mean 0 and standard deviation  $\tau = 0.2R_u$  with  $R_u$  being the range of  $u(x)$  over  $x \in [0, 1]$ . We considered the following four cases for the true regression model:

$$\begin{aligned}
\text{(a)} \quad & \begin{cases} u(x) &= 0.6 \exp\{-(x-0.3)^2/0.01\} + 1.4 \exp\{-(x-0.7)^2/0.02\}, \\ v_S(x) &= -1.0 \cdot \mathbf{I}(x \in [0.44, 0.47]) - 0.2 \cdot \mathbf{I}(x \in [0.45, 0.46]), \\ S &= \{0.44, 0.45, 0.46, 0.47\}, \end{cases} \\
\text{(b)} \quad & \begin{cases} u(x) &= \sin(4\pi x), \\ v_S(x) &= -1.5 \cdot \mathbf{I}(x \in [0.36, 0.39] \cup [0.86, 0.89]) - 0.5 \cdot \mathbf{I}(x \in [0.37, 0.38] \cup [0.87, 0.88]), \\ S &= \{0.36, 0.37, 0.38, 0.39, 0.86, 0.87, 0.88, 0.89\}, \end{cases} \\
\text{(c)} \quad & \begin{cases} u(x) &= \sin(4\pi x) + 2 \cos(5\pi x), \\ v_S(x) &= -1.5 \cdot \mathbf{I}(x \in [0.12, 0.15]) - 0.3 \cdot \mathbf{I}(x \in [0.13, 0.14]) \\ &\quad + 1.5 \cdot \mathbf{I}(x \in [0.55, 0.58] \cup [0.85, 0.88]) + 0.3 \cdot \mathbf{I}(x \in [0.56, 0.57] \cup [0.86, 0.87]), \\ S &= \{0.12, 0.13, 0.14, 0.15, 0.55, 0.56, 0.57, 0.58, 0.85, 0.86, 0.87, 0.88\}, \end{cases} \\
\text{(d)} \quad & \begin{cases} u(x) &= \exp(-x) \sin(5\pi \exp(-x)), \\ v_S(x) &= 0, \\ S &= \emptyset. \end{cases}
\end{aligned}$$

where  $\mathbf{I}(\cdot)$  is an indicator function of the event  $A$ , that is,  $\mathbf{I}(A) = 1$  if  $A$  is true and 0 otherwise. We compared the performance of FLSA-G with that of KM-G and FLSA estimation procedure (FLSA).

Figure 2 compares true curves with fitted ones. The estimated curves from the 1st row to 4th row correspond to the true curves from (a) to (d). We observe that our modeling strategy is effective in capturing the true data structures well.

We also performed 1000 repetitions, then calculated averages of mean squared errors on peak region (AMSE.p), that on smooth region (AMSE.s) and that on global region (AMSE.g) defined by  $\text{MSE.p} = \sum_{x_\alpha \in S} \{(u(x_\alpha) + v(x_\alpha)) - \hat{y}_\alpha\}^2 / n_S$ ,  $\text{MSE.s} = \sum_{x_\alpha \in U \setminus S} \{(u(x_\alpha) + v(x_\alpha)) - \hat{y}_\alpha\}^2 / (n - n_S)$  and  $\text{MSE.g} = \sum_{\alpha=1}^n \{(u(x_\alpha) + v(x_\alpha)) - \hat{y}_\alpha\}^2 / n$  respectively to assess the goodness of fit.

Table 1 shows summaries of the simulation results from (a) to (d). In all situations, our proposed modeling procedure minimized all of the AMSE.p, AMSE.s and AMSE.g, thus improving the accuracy of prediction. In the case of (d), the performance of our proposed method (FLSA-G) was superior to that of Km-G even though true model has no spiky change point.

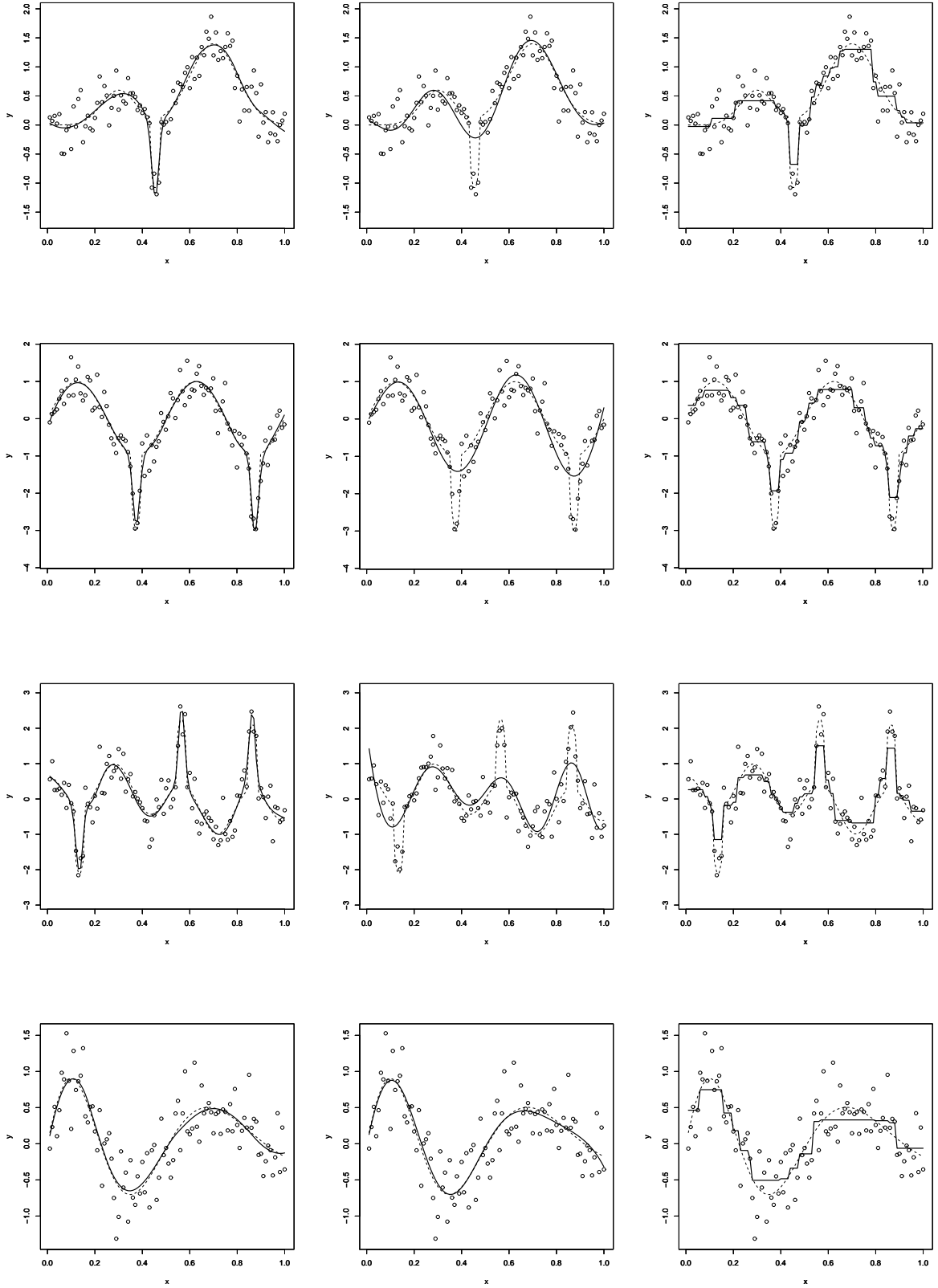


Fig. 2: These panels show true curves (dashed line) and estimated curves (solid line) obtained by FLSA-G (left), KM-G (center) and FLSA (right) for each true function.

Table 1: Comparison of results for curve fitting.

function		mean $m$	mean $\gamma$	mean $\nu$	AMSE.p	AMSE.s	AMSE.g
(a)	FLSA-G	17.50	$3.96 \times 10^{-5}$	58.26	$1.57 \times 10^{-1}$	$1.45 \times 10^{-2}$	$2.02 \times 10^{-2}$
	KM-G	9.22	$2.98 \times 10^{-2}$	4.26	$3.96 \times 10^{-1}$	$2.12 \times 10^{-2}$	$3.62 \times 10^{-2}$
	FLSA	--	--	--	$2.42 \times 10^{-1}$	$1.77 \times 10^{-2}$	$2.67 \times 10^{-2}$
(b)	FLSA-G	22.88	$4.16 \times 10^{-5}$	190.03	$1.95 \times 10^{-1}$	$2.98 \times 10^{-2}$	$4.30 \times 10^{-2}$
	KM-G	7.00	$4.07 \times 10^{-3}$	3.60	1.39	$1.08 \times 10^{-1}$	$2.10 \times 10^{-1}$
	FLSA	--	--	--	$3.12 \times 10^{-1}$	$4.65 \times 10^{-2}$	$6.78 \times 10^{-2}$
(c)	FLSA-G	21.67	$1.59 \times 10^{-5}$	189.50	$1.80 \times 10^{-1}$	$4.91 \times 10^{-2}$	$6.48 \times 10^{-2}$
	KM-G	10.52	$3.05 \times 10^{-3}$	5.62	$9.35 \times 10^{-1}$	$1.07 \times 10^{-1}$	$2.07 \times 10^{-1}$
	FLSA	--	--	--	$3.49 \times 10^{-1}$	$5.47 \times 10^{-2}$	$9.00 \times 10^{-2}$
(d)	FLSA-G	16.86	$1.04 \times 10^{-4}$	65.69	--	--	$9.92 \times 10^{-3}$
	KM-G	7.23	$9.21 \times 10^{-3}$	15.74	--	--	$1.20 \times 10^{-2}$
	FLSA	--	--	--	--	--	$1.99 \times 10^{-2}$

## 6.2 Real data example

We investigate the performance using the nursing time of beluga whale data set (Simonoff ,1996; Russell *et al.*, 1997). The data consists of 228 measurements of nursing time in seconds of a newborn male beluga whale calf named Hudson born in captivity at the New York Aquarium. This data was observed every six hours for 57 days after his birth. Russell *et al.* (1997) noted that the nursing time typically peaked at around 7-10 days postpartum from the biological view. We apply our proposed method, nonlinear regression modeling procedure with  $k$ -means-based Gaussian basis functions and FLSA estimation procedure to the Beluga data.

Figure 2 shows the result curves obtained by the three methods. Using our method, spiky change appears in around the 9th day of his life as Cheng and Raimondo (2008). We observe that our modeling procedure captures the abrupt change more remarkably than the others and lead smooth curve .

## 7 Concluding remarks

We have proposed a nonlinear regression modeling procedure along with the technique of the method of regularization. We have introduced new asymmetric Gaussian basis

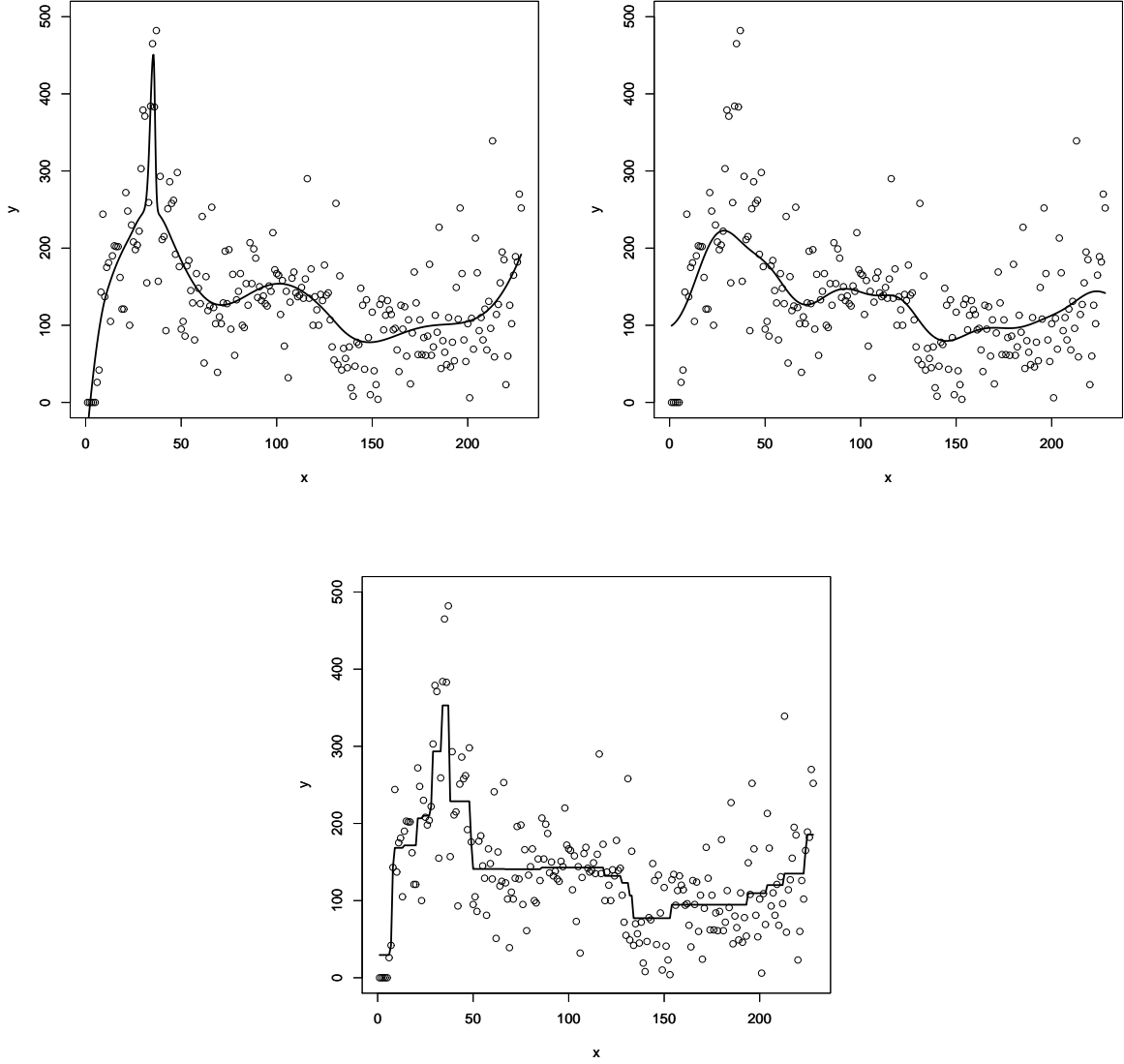


Fig. 3: Beluga data and estimated curves obtained by FLSA-G (Upper left), KM-G (Upper right) and FLSA (Lower).

functions based on FLSA estimates taking the information of spikes into account. The nonlinear regression model with these basis functions is able to capture the sudden changes in nonlinear structure of data. In order to choose tuning parameters, we presented the model selection criterion from information-theoretic approach. The simulation results reported here demonstrate the effectiveness of the proposed modeling strategy in terms of prediction accuracy and prominent visualization of spikes.

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