

Multistep-Ahead River Flow Prediction Using LS-SVR at Daily Scale

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Received April 5, 2012; revised May 5, 2012; accepted June 1, 2012

ABSTRACT

In this study, potential of Least Square-Support Vector Regression (LS-SVR) approach is utilized to model the daily variation of river flow. Inherent complexity, unavailability of reasonably long data set and heterogeneous catchment response are the couple of issues that hinder the generalization of relationship between previous and forthcoming river flow magnitudes. The problem complexity may get enhanced with the influence of upstream dam releases. These issues are investigated by exploiting the capability of LS-SVR—an approach that considers Structural Risk Minimization (SRM) against the Empirical Risk Minimization (ERM)—used by other learning approaches, such as, Artificial Neural Network (ANN). This study is conducted in upper Narmada river basin in India having Bargi dam in its catchment, constructed in 1989. The river gauging station—Sandia is located few hundred kilometer downstream of Bargi dam. The model development is carried out with pre-construction flow regime and its performance is checked for both pre- and post-construction of the dam for any perceivable difference. It is found that the performances are similar for both the flow regimes, which indicates that the releases from the dam at daily scale for this gauging site may be ignored. In order to investigate the temporal horizon over which the prediction performance may be relied upon, a multistep-ahead prediction is carried out and the model performance is found to be reasonably good up to 5-day-ahead predictions though the performance is decreasing with the increase in lead-time. Skills of both LS-SVR and ANN are reported and it is found that the former performs better than the latter for all the lead-times in general, and shorter lead times in particular.

Keywords: Multistep-Ahead Prediction; Kernel-Based Learning; Least Square-Support Vector Regression (LS-SVR); Daily River Flow; Narmada River

1. Introduction

River flow is an important component in hydrological cycle, which is directly available to the community. In hydrology, river flow plays an important role in establishing some of the critical interactions that occur between physical, ecological, social or economic processes. Accurate or at least reasonably reliable prediction of river flow is an important foundation for preventing flood, reducing natural disasters, and the optimum management of water resource. Constructions of major and minor dams are essential in order to effective use of available water resources. This is more crucial for the monsoon dominated countries, where most of the annual rainfall occurs during couple of months, and rest of the months are mostly no-rainfall months. Existence of dams adds to the complexity in the river flow modelling, which is influenced by the releases from upstream reservoirs as well as the influence of the inputs from the catchment area between immediate downstream of the dam and the gauging site.

Depending on the natural variability and complexity, traditional statistical methods, such as, transfer function model, Box-Jenkins approach etc. may be inadequate due to their underlying assumptions. As a consequence many new methodologies have been introduced to understand the variations of hydrologic variables and to predict for the future time steps.

Development of different techniques to predict river flow is having a long history. Among the parametric linear approaches, Auto-Regressive Integrated Moving Average (ARIMA) model is one of the most popular approaches. Since last decade, machine learning techniques are being applied in this field. For instance, Artificial Neural Networks (ANN), fuzzy logic, genetic programming, etc., have been widely used in the modelling and prediction of hydrologic variables. More recently, kernel-based learning approaches have gained wide popularity.

The statistical learning framework proposed by Vapnik [1] led to the introduction of the Support Vector Machine (SVM), which has been successfully applied for

nonlinear classification and regression in learning problems. Kernel based approaches are expected to perform better due its nonlinear, even smooth enough, feature space development based on the available historical record. One such kernel based machines learning approach is the Least Squares-Support Vector Machine (LS-SVM) [2].

1.1. Support Vector Machines (SVMs) and Kernel Based Learning

SVMs are a kind of supervised machine learning system that use a linear high dimensional hypothesis space called feature space. These systems are trained using a learning algorithm, which is based on optimization theory. SVMs belong to a family of generalized linear classifier [3]. The SVMs can be applied to both classification and regression problems. Application of SVMs to regression problem was made in late nineties [1,4]. Popularity of SVM increases very rapidly since then in fields of text classification, pattern recognition, remote sensing and so on.

The basic idea of working principle of SVMs is provided by the use of kernel functions that implicitly map the data to a higher dimensional space. According to Cover's theorem, linear solution in the higher dimensional feature space corresponds to a non-linear solution in the original lower dimensional input space. This makes SVM a powerful tool for solving a variety of hydrologic problems, which are non-linear in nature. There are methods available which use nonlinear kernels in their approach towards regression problems while applying SVMs.

1.2. Application of SVMs in Hydrologic Problems

Application of SVMs in the field of hydrology is gaining wide popularity and the results are found to be encouraging. Such applications range from remotely sensed image classification [5], statistical downscaling [6], soil water forecasting [7], stream flow forecasting [8] and so on. Liong and Sivapragasam [9] compared SVM performance with other machine learning model, such as, ANN in forecasting flood stage and reported a superior performance of SVM. Bray and Han [10] used SVM for rainfall runoff modelling and that model was compared with a transfer function model. The study outlined a promising area of research for further application of SVMs in unexplored areas. Samui [11] used LS-SVM to determine evaporation loss of reservoir and it is established to be a powerful approach for the determination of evaporation loss. She and Basketfield [12] forecasted spring and fall season stream flows in Pacific Northwest region of US using SVM and reported superior results in fore-

casting. Zhang *et al.* [13] studied two machine learning approaches—ANN and SVM and compared for approximating the Soil and Water Assessment Tool (SWAT) model. The results showed that SVM in general exhibited better generalization ability than ANN. Khadam and Kaluarachchi [14] discussed the impact of accuracy and reliability of hydrological data on model calibration. This, coupled with application of SVMs, was used to identify the faulty model calibration, which would have been undetected otherwise. Applicability of SVMs was also demonstrated in downscaling Global Circulation Models (GCMs), which are among the most advanced tools for estimating future climate change scenarios. The results presented SVMs as a compelling alternative to traditional Artificial Neural Networks (ANN) to conduct climate impact studies [10,11] downscaled monthly precipitation to basin scale using SVMs and reported the results to be encouraging in their accuracy while showing large promise for further applications.

1.3. Advancement of SVM

Apart from the general benefit of SVM pointed out in the aforementioned studies, SVMs are sometimes criticized by its large number of parameters and high level of computational effort, particularly in case of large dataset. Chunking is one of the proposed remedies to the latter problem. However, according to Suyken *et al.* [15], it is worthwhile to investigate the possibility of simplifying the approach to the extent possible without losing the any of its advantages. Thus, they proposed a modification over the SVM approach, which essentially leads to Least Square-Support Vector Machines (LS-SVM).

The main advantage of LS-SVM is in its higher computational efficiency than that of standard SVM method, since the training of LS-SVM requires only the solution of a set of linear equations instead of the long and computationally demanding quadratic programming problem involved in the standard SVM [2]. Qin *et al.* [16] investigated the application of LS-SVM for the modelling of water vapor and carbon dioxide fluxes and they found the excellent generalization property of LS-SVM and its potential for further applications in area of general hydrology. Maity *et al.* [13] investigated the potential of support vector regression, which is also based on LS-SVM principle, for prediction of streamflow using endogenous property of the monthly time series. In this study, potential of LS-SVM for Regression (LS-SVR) is investigated for the objective outlined as follows.

1.4. Objective of This Study

Potential of LS-SVM for Regression (LS-SVR) is exploited for multistep-ahead river flow prediction at daily scale, to assess its performance with the increasing time

horizon. Most of the major river basins are being populated with major and minor dams. River flow modelling is supposed to be influenced by the effect of the release from these dams if the site location is on the downstream side of the dam. However, the effect of its existence will gradually reduce with increase of the distance from the dam location. The study is carried out with observed daily river flow in the upper Narmada River basin with Sandia gauging station at the outlet. Bargi dam exists few hundred km upstream from the outlet of the watershed. Details of this dam are provided in the ‘‘Study Area’’ section later. Investigation is carried out to assess the necessity of consideration of daily releases from upstream dam on the daily river flow variation at the outlet of the study area. Also, a multistep-ahead prediction is carried out to assess maximum temporal horizon over which the prediction results may be relied upon. The results are compared with the performance of neural networks approach that uses Empirical Risk Minimization (ERM).

2. Methodology

2.1. Data Normalization

The observed river flow data, as commonly used in data-driven models, is normalized to prevent the model from being dominated by the large values. The performance of LS-SVM with normalized input data has shown to outperform the same with non-normalized input data [16]. Therefore, the data is normalized and finally the model outputs are back transformed to their original form by denormalisation. The normalization (also back transformation) is carried out using

$$y_i = 0.1 + \frac{S_i}{1.2 \max(S_i)} \quad (1)$$

where y_i is the normalized data for i^{th} day and S_i is the observed value for i^{th} day.

2.2. Least Square-Support Vector Regression (LS-SVR)

Let us consider a given training set $\{\mathbf{x}_i, y_i\}_{i=1}^N$, where

$$\left[y_{i-n}, y_{i-(n-1)}, \dots, y_{i-1} \right]^T \in \mathfrak{R}^n$$

represents a n -dimensional input vector and $y_i \in \mathfrak{R}$ is a scalar measured output, which represents system output. Subscript i indicates the i^{th} training pattern. In the context of multistep-ahead river flow prediction, \mathbf{x}_i is the vector comprising of n -previous days normalized river flow values, y_i is the target river flow with certain lead-time in day(s) and subscript i indicates the reference time from which n -previous days and the lead-time are counted. The goal is

to construct a function $y = f(\mathbf{x})$, which represents the dependence of the output y on the input \mathbf{x} . The form of this function is as follows:

$$y = \mathbf{w}^T \phi(\mathbf{x}) + b \quad (2)$$

where \mathbf{w} is known as weight vector and b as bias.

This regression model can be constructed using a nonlinear mapping function $\phi(\cdot)$. The function $\phi(\cdot): \mathfrak{R}^n \rightarrow \mathfrak{R}^h$ is a mostly nonlinear function, which maps the data into a higher, possibly infinite, dimensional feature space. The main difference from the standard SVM is that the LS-SVR involves equality constraints instead of inequality constraints, and works with a least square cost function. The optimization problem and the equality constraints are defined by the following equations:

$$\left. \begin{aligned} \text{Minimize } J(\mathbf{w}, e) &= \frac{1}{2} \mathbf{w}^T \mathbf{w} + \gamma \frac{1}{2} \sum_{i=1}^N e_i^2 \\ \text{such that } y_i &= \mathbf{w}^T \phi(\mathbf{x}_i) + b + e_i \quad i = 1, \dots, N \end{aligned} \right\} \quad (3)$$

where e_i is the random error and $\gamma \in \mathfrak{R}^+$ is a regularization parameter in optimizing the trade-off between minimizing the training errors and minimizing the model’s complexity. The objective is now to find the optimal parameters that minimize the prediction error of the regression model. The optimal model will be chosen by minimizing the cost function where the errors, e_i , are minimized. This formulation corresponds to the regression in the feature space and, since the dimension of the feature space is high, possibly infinite, this problem is difficult to solve. Therefore, to solve this optimization problem, the following Lagrange function is given,

$$L(w, b, e; \alpha) = J(w, e) - \sum_{i=1}^N \alpha_i \{ \mathbf{w}^T \phi(\mathbf{x}_i) + b + e_i - y_i \} \quad (4)$$

The solution of above can be obtained by partially differentiating with respect to w , b , e_i and α_i , *i.e.*

$$\frac{\partial L}{\partial w} = 0 \rightarrow \mathbf{w} = \sum_{i=1}^N \alpha_i \phi(\mathbf{x}_i) \quad (5)$$

$$\frac{\partial L}{\partial b} = 0 \rightarrow \sum_{i=1}^N \alpha_i = 0 \quad (6)$$

$$\frac{\partial L}{\partial e_i} = 0 \rightarrow \alpha_i = \gamma e_i, \quad i = 1, \dots, N \quad (7)$$

$$\frac{\partial L}{\partial x_i} = 0 \rightarrow \mathbf{w}^T \phi(\mathbf{x}_i) + b + e_i - y_i = 0, \quad i = 1, \dots, N \quad (8)$$

From the set of Equations (5)-(8), \mathbf{w} and e can be eliminated and finally, the estimated values of b and α_i , *i.e.* b and $\hat{\alpha}_i$, can be obtained by solving the linear system. Replacing \mathbf{w} in Equation (2) from Equation (5),

the kernel trick may be applied as follows:

$$K(\mathbf{x}, \mathbf{x}_i) = \varphi(\mathbf{x}_i)^T \varphi(\mathbf{x}) \quad (9)$$

Here, the kernel trick means a way to map the observations to an inner product space, without actually computing the mapping and it is expected that the observations will have meaningful linear structure in that inner product space.

Thus, the resulting LS-SVR model can be expressed as

$$\hat{y} = \sum_{i=1}^N \hat{\alpha}_i K(\mathbf{x}, \mathbf{x}_i) + \hat{b} \quad (10)$$

where $K(\mathbf{x}, \mathbf{x}_i)$ is a kernel function.

In comparison with some other feasible kernel functions, the RBF is a more compact and able to shorten the computational training process and improve the generalization performance of LS-SVR (LS-SVM, in general), a feature of great importance in designing a model [13]. Aksornsingchai and Srinilta [17] studied support vector machine with polynomial kernel (SVM-POL), and support vector machine with Radial Basis Function kernel (SVM-RBF) and found SVM-RBF is the accurate model for statistical downscaling methods. Also, many works have demonstrated the favorable performance of the radial basis function [9,15]. Therefore, the radial basis function is adopted in this study. The nonlinear radial basis function (RBF) kernel is defined as:

$$K(\mathbf{x}, \mathbf{x}_i) = \exp\left(-\frac{1}{\sigma^2} \|\mathbf{x} - \mathbf{x}_i\|^2\right) \quad (11)$$

where σ is the kernel function parameter of the RBF kernel. The symbol $\|\cdot\|$ is the norm of the vector \cdot and thus, $\|\mathbf{x} - \mathbf{x}_i\|^2$ is basically the Euclidean distance between the vectors \mathbf{x} and \mathbf{x}_i . In the context of river flow prediction, \mathbf{x}_i is the new vector of previous river flow, based on which multi-step ahead prediction (\hat{y}_i) is made with certain lead-time. y_i (observed) and \hat{y}_i are compared to assess the model performances.

2.3. Model Calibration and Parameter Estimation

The regularization parameter γ determines the trade-off between the fitting error minimization and smoothness of the estimated function. It is not known beforehand which γ and σ^2 are the best for a particular problem to achieve maximum performance with LS-SVR models. Thus, the regularization parameter γ and the RBF kernel parameter σ^2 have to be calibrated during model development period. These parameters are interdependent, and their (near) optimal values are often obtained by a trial-and-error method. Interrelationship is also coupled with the number of previous river flow values to be considered, which is denoted as n . In order

to find all these parameters (γ , σ^2 and n) grid search method is employed in parameter space. Once the parameters are estimated from the training dataset, the obtained LS-SVR model is complete and ready to use for modelling new river flow data period. Performance of the developed model is then assessed with the data set during testing period. Different models (parameter sets) are developed for different prediction lead-times in case of multi-step ahead prediction.

2.4. Comparison with Artificial Neural Networks (ANN)

The flexible computing based ANN models have been extensively studied and used for time series forecasting in many hydrologic applications since late 1990s. This model has the capability to execute complex mapping between input and output and to form a network that approximates non-linear functions. A single hidden layer feed forward network is the most widely used model form for time series modeling and forecasting [18]. This model usually consists of three layers: the first layer is the input layer where the data are introduced to the network followed by the hidden layer where data are processed and the final or output layer is where the results of the given input are produced.

The number of input nodes and output nodes in an ANN are dependent on the problem to which the network is being applied. However, there is no fixed method to find out the number of hidden layer nodes. If there are too few nodes in the hidden layer, the network may have difficulty in generalizing the problems. On the other hand, if there are too many nodes in the hidden layer, the network may take an unacceptably long time to learn any thing from the training set [19]. Increase in the number of parameters may slow the calibration process [20]. In a study by Zealand *et al.* [21], networks were initially configured with both one and two hidden layers. However, the improvement in forecasting results was only marginal for the two hidden layer cases. Therefore, it is decided to use a single hidden layer in this study. In most of the cases, suitable number of neurons in the hidden layer is obtained based on the trial-and-error method [22]. Maity and Nagesh Kumar, [23] proposed a GA based evolutionary approach to decide the complete network structure. The output X_i of an ANN, assuming a linear output neuron having a single hidden layer with h sigmoid hidden nodes, is given by:

$$X_i = g\left(\sum_{j=1}^h w_j f(s_j) + b_k\right) \quad (12)$$

where $g(\cdot)$ and b_k are the linear transfer function and bias respectively of the output neuron k , w_j is the connection weights between j^{th} neuron of hidden la-

yers and output units, $f = (\cdot)$ is the transfer function of the hidden layer [24]. The transfer functions can take several forms and the most widely used transfer functions are Log-sigmoid, Linear, Hyperbolic tangent sigmoid etc. In this study, hyperbolic tangent sigmoid is used:

$$f(s_i) = \frac{2}{1 + \exp(-2s_i)} - 1 \quad (13)$$

where $s_i = \sum_{i=0}^n w_i x_i$ is the input signal referred to as the weighted sum of incoming information. Several optimization algorithms can be used to train the ANN. Among the various training algorithms available, the back-propagation is most popular and widely used algorithm [25]. Details of this techniques is well established in the literature and can be found elsewhere (ASCE 2000 and references therein) [26].

3. Study Area and Data Sets

Narmada River is the largest west flowing river of Indian peninsula. It is the fifth largest river in India. The study area is up to Sandia gauging station, which is in the upstream part of Narmada river basin as shown in **Figure 1**. The upstream part Narmada river basin is in the state of Madhya Pradesh, India. The river originates from Maikala ranges at Amarkantak and flows westwards over a length of 334 km up to Sandia. The elevation difference

in study area is from 289 to 1134 m. The basin lies between east longitudes 78°30' and 81°45', and north latitudes 21°20' and 23°45'. Bargi dam (later renamed as Rani Avanti Bai Sagar Project) is a major structure in the basin up to Sandia, which is located few hundred km upstream of Sandia. The latitude and longitude of the dam are 22°56'30"N and 79°55'30"E, respectively. It was constructed in late eighties and being operated from early nineties. Thus, pre-construction period (1978-1986) is considered for training. For testing, two sets of data are used—pre-construction data set (1986-1990) and post-construction period (1990-2000). However, out of this entire range four years data is missing (1, 1981 - 31 May 1982; June 1, 1987 - May 31, 1988; June 1, 1993 - May 31, 1994; June 1, 1998 - May 31, 1999). These periods are ignored from the analysis. River flow data from Sandia station, operated by the Water Resources Agency, is obtained from Central Water Commission, Govt. of India. Among these records, a daily data (June 1, 1978 to May 31, 1986) is used for training and (June 1, 1986 to May 31, 2000) is used to test the model performance.

4. Results and Discussions: Performance of LS-SVR for River Flow Prediction

4.1. Data Pre-Processing and Parameter Estimation

Observed river flow data at Sandia is normalized as explained in the methodology before proceeding to parameter

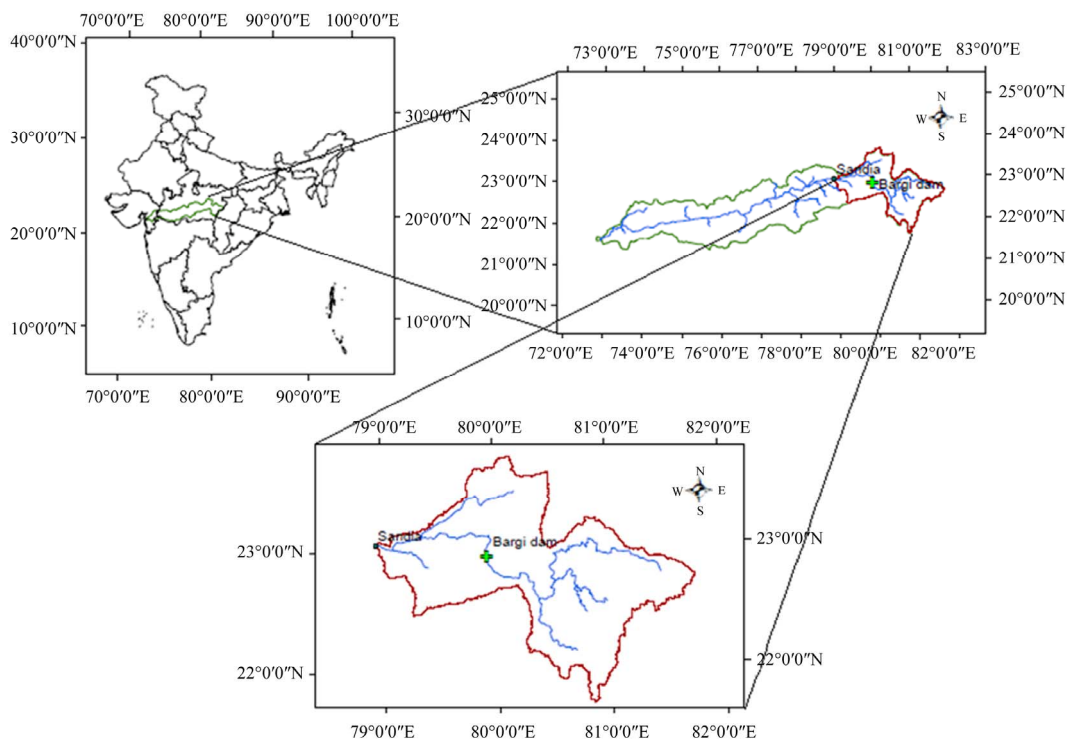


Figure 1. Narmada river basin with study area up to Sandia station.

meter estimation. Optimum number of previous river flow values to be considered is denoted as n . This parameter along with the regularization parameter γ and the RBF kernel parameter σ^2 is calibrated during model development period (training period). To select best combination of n , γ and σ^2 , grid search method is used. Model performances for different combinations of these parameters are assessed based on statistical measures, such as, Correlation Coefficient (CC), Root Mean Square Error (RMSE) and Nash-Sutcliffe Efficiency (NSE). Ten different values of n (1 through 10) are tested to decide the optimum number of previous daily river flow to be considered for the best possible results. Range of γ is considered to be 25 to 1000 with a resolution of 25 and σ^2 in the range of 0.01 to 1 with a resolution of 0.01. Approximately (because of different lag and lead-times considered) 2556 data points are used for training purpose. Performance of each model is assessed with the remaining testing data points. Model performances statistic is obtained between observed and modelled river flow values during training and testing period. The combination that yields comparable performance during training and testing period is selected, which ensures the best parameter values without the fear of overfitting. Results are shown in **Table 1** for prediction lead-time of 1 day. This is achieved in the following way: for each value of n , two (training period and testing period) 2-D surfaces are obtained for each of the performance measures (CC, RMSE and NSE). The values of γ and σ^2 are identified for which the training and testing period performances are “closest”. Priority is given to CC and corresponding RMSE and NSE are reported for same values of γ and σ^2 . These values are named as “best γ ” and “best σ^2 ” in **Table 1** for a particular n . It is found from that the best combination of n , γ and σ^2 is 5 days, 175 and 0.21 respectively. However, these parameters are for prediction lead-time of 1 day. Opti-

num combinations of parameters are computed separately for different prediction lead-times and reported later.

4.2. Model Performance during Pre- and Post-Construction Testing Period

Model performances are tested during both pre- and post-construction period. For this purpose, the model is trained with river flow data during pre-construction period and the developed model is used to assess the performance during both pre-construction and post construction period. It is found that the performance remains almost similar in both pre- and post-construction period (**Table 2**). Flow regimes are supposed to be modified on the immediate downstream of a newly constructed dam due to the effect of human controlled releases from the dam. However, such effect is supposed to get diminished with increase in distance from the dam location towards downstream. This is being reflected in case of river flow at Sandia gauging station indicating that the station is sufficiently away towards downstream. Model performance during training period is shown in **Figure 2** and the model performances during testing period—both pre-construction and post-construction period are shown in **Figures 3** and **4** respectively. The immediate next question would be to find the temporal horizon up to which the prediction would be reliable.

4.3. Performance of River Flow Prediction for Different Lead-Times

The multistep-ahead river flow prediction is carried out for time steps T , $T + 1$, $T + 2$, $T + 3$ and $T + 4$. In other words, five different prediction lead-times are tested for prediction performance, *i.e.*, 1 day, 2 days, 3 days, 4 days and 5 days in advance. As stated earlier, optimum combinations of parameters are computed separately for different prediction lead-times. Results are shown in **Table 3**.

Table 1. Model performances during training (testing) period for different numbers of previous daily river flows considered as inputs (Lead-time = 1 day) with corresponding “best γ ” and “best σ^2 ” values.

Number of previous river flow values input (n)	Best γ	Best σ^2	CC	RMSE	NSE
1	265	0.01	0.810 (0.810)	0.025 (0.030)	0.657 (0.655)
2	960	0.03	0.829 (0.829)	0.024 (0.029)	0.687 (0.679)
3	865	0.11	0.827 (0.827)	0.024 (0.029)	0.683 (0.675)
4	845	0.31	0.825 (0.825)	0.024 (0.030)	0.681 (0.662)
5	175	0.21	0.827 (0.827)	0.0243 (0.029)	0.684 (0.665)
6	795	0.34	0.829 (0.829)	0.024 (0.029)	0.688 (0.674)
7	710	0.32	0.831 (0.831)	0.024 (0.029)	0.690 (0.681)
8	235	0.22	0.833 (0.833)	0.024 (0.029)	0.694 (0.681)
9	530	0.28	0.838 (0.838)	0.024 (0.028)	0.702 (0.688)
10	750	0.45	0.839 (0.839)	0.024 (0.028)	0.703 (0.696)

Table 2. Comparison of performance during pre- and post-construction of Bargi dam.

Statistical Measures	Training Period (5-Jun-78 to 31-May-86)	Testing performance	
		Before dam construction (5-Jun-86 to 31-May-90)	After dam construction (5-Jun-90 to 31-May-97)
Correlation Coefficient	0.83	0.84	0.83
RMSE	0.024	0.017	0.033
NSE	0.68	0.71	0.66

Table 3. Optimum river flow lags and parameter estimation for different lead-times.

Lead-time (in days)	Number of previous river flow values input (n)	Gamma (γ)	Sigma (σ)
1	5	175	0.21
2	2	75	0.14
3	1	150	0.40
4	1	50	0.38
5	1	75	0.37

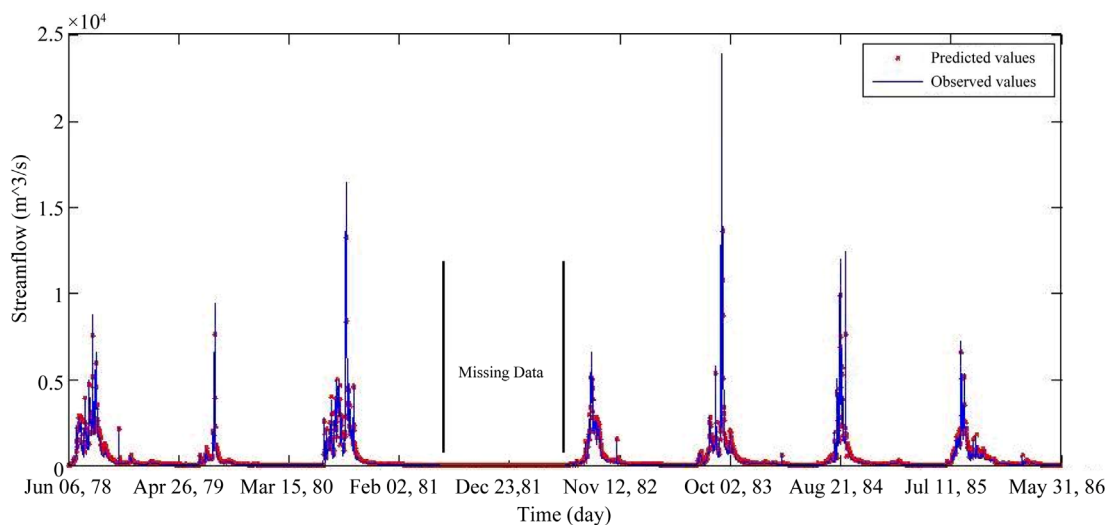


Figure 2. Comparison between observed and predicted river flow for training period (1-day-ahead).

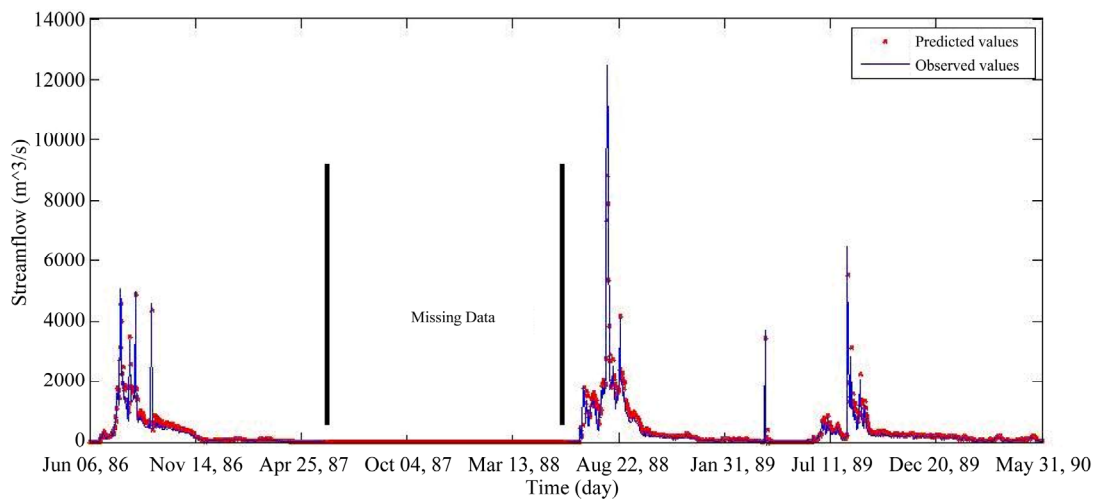


Figure 3. A plot between observed and predicted river flow (1-day-ahead) during testing period before construction of dam.

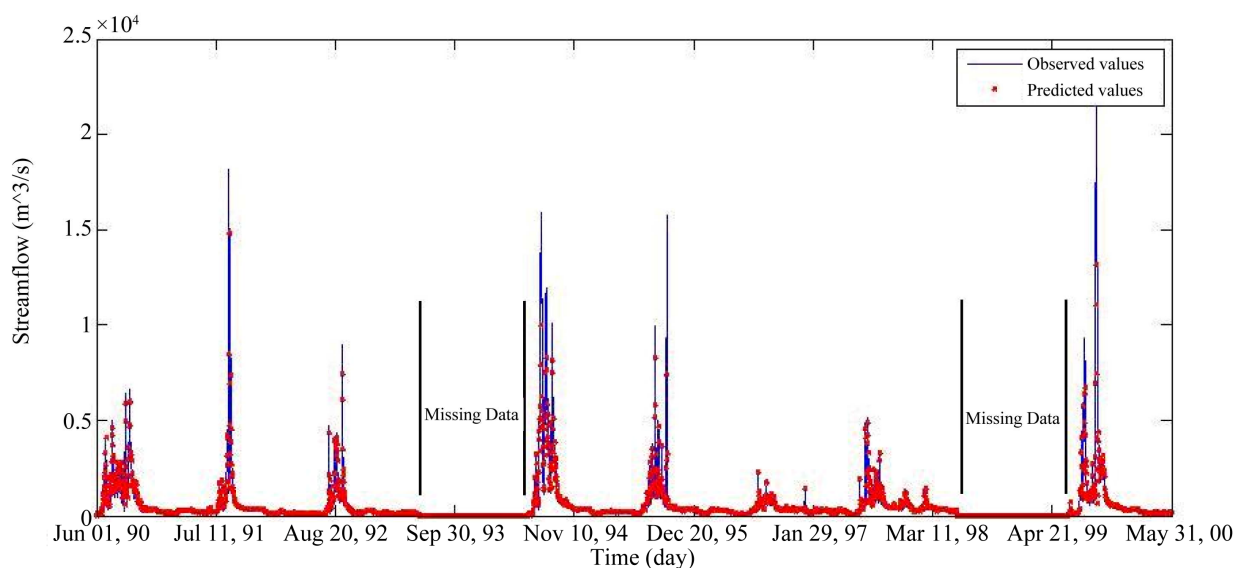


Figure 4. A plot between observed and predicted river flow (1-day-ahead) during testing period after construction of dam.

From these results, it is observed that for different prediction lead-times the best combination of model parameters varies. The optimum numbers of previous river values to be considered differ for different lead-times. This is decreasing with the increase of lead-time. For instance, five previous days input is best for 1-day ahead prediction; similarly for 2-day lead-time best is two previous days inputs and for further lead-times one previous day input is resulting the best river flow prediction.

Model performance is investigated to determine ability of model to predict multistep-ahead river flow through CC, RMSE and NSE. Results are shown in Table 4. It is found that the prediction performance decreases with the increase in lead-time. For instance, 69% (square of correlation coefficient) of daily variation is explained in case of 1-day-ahead prediction, whereas 50%, 41%, 36% and 34% of daily variation is explained in case of 2-day, 3-day, 4-day and 5-day ahead predictions respectively. RMSE and NSE values are also increasing and decreasing respectively with the increase in lead-times, indicating that the performance is better for the shorter prediction lead-times.

Plots between observed and predicted river flow for Training and Testing periods for 1-day lead-times are shown in Figures 2-4 sequentially.

It is observed that predicted river flow values well correspond to the observed one at low and medium range of river flows. However, the highest peak is not captured properly. As these types of peaks are very short-lived (at daily scale), the reason could be the effect of some other factor, such as, sudden flash rain, which is shorter than the daily temporal resolution and thus, not available from the information contained in the previous days river flows.

Table 4. Performance of LS-SVR for multistep-ahead daily river flow prediction for different lead-times.

Lead-time (in days)	CC	RMSE	NSE
1	0.83	0.029	0.67
2	0.71	0.036	0.49
3	0.64	0.040	0.39
4	0.60	0.041	0.35
5	0.58	0.042	0.32

4.4. Comparison of LS-SVR and ANN Model Performances

The relative skill of LS-SVR is compared with the performance of ANN. Basics of ANN approaches are discussed in methodology section. The architecture of ANN consists of different layers and combinations of neurons as discussed earlier. Architecture of ANN networks adopted in this study is having n neurons in the input layer as river flow values from n previous days are used as input. As explained earlier, n is the optimum number of previous river flow values to be considered. Output layer consists of one neuron as the target is to predict the daily river flow with a particular lead-time. The adopted networks for different lead-times also differ from each other with respect to the number of neurons in the hidden layers. The trial-and-error method is used to find out best combination of neuron for hidden layer. In trial-and-error method, different combinations of hidden layer neurons are tried and performance is checked. Based upon the performance, optimum number of hidden layer neurons is selected. It is found that 5, 3, 3, 3 and 3 neurons in hidden layer provide the best performance for lead-times of 1, 2, 3, 4 and 5 days respectively. It might be noted

here that back propagation algorithm (as explained earlier) is used to train these networks using 2556 to 2552 (from lead-time 1-day through 5-day) training data set. The largest network (5-5-1) is having $25 + 5 + 5 + 1 = 31$ parameters, which is much less than number of training patterns.

Different networks for different lead-times are trained separately. Statistical measures showing the prediction performances are obtained during training and testing period for different lead-time cases and the results are shown in **Table 5**. After comparison between LS-SVR and ANN, it is found that, in general, LS-SVR performs better than ANN for all the lead-times. For instance, at 1-day lead-time, 69% (CC = 0.83) and 53.2% (CC = 0.73) of daily river flow variation are captured by LS-SVR and ANN respectively during the entire testing period (**Table 5**). Also, for this case, the NSE values are found to be 0.67 and 0.52 in case of LS-SVR and ANN respectively. The performance of ANN also decreases with the increase in lead-time. It is also found that the difference in prediction performance obtained from LS-SVR and ANN is more prominent for shorter lead times. As the lead time increases, the performance of LS-SVR and ANN becomes comparable. Different error measures, *i.e.*, Mean Absolute Error (MAE), Mean Absolute Relative Error (MARE) and Root Mean Square Error (RMSE) are also determined for LS-SVR and ANN based predictions. Results are shown in **Table 6**. Based on these measures also, it is found that error measures for LS-SVR are bet-

ter compared to that of ANN. These observations indicate the higher capability of LS-SVR approach to capture the river flow variation using the previous information.

For a visual comparison of prediction performances by LS-SVR and ANN simultaneously, both LS-SVR and ANN model predicted river flow values are plotted with the observed river flow during the model testing period (**Figure 5**). This is shown in case of 1-day-ahead prediction—best for both LS-SVR and ANN. It is observed that predicted river flow values are very well corresponds to the low and medium range of river flow in case of LS-SVR and ANN. Relatively better performance of LS-SVR is discussed with respect to the statistical measures as shown in **Table 5**. However, as discussed in case of LS-SVR earlier, the peak river flow values are not captured with reasonable accuracy in case of ANN as well. Apart from the fact that these peaks are very short-lived (at daily scale), this is a shortcoming for both LS-SVR and ANN based approaches. The reason could be the same as discussed before, *i.e.*, these peaks might be the effect of some factor, such as, sudden flash rain, which is shorter than the daily temporal resolution and thus, not available from the information contained in the previous days river flows alone. Consideration and incorporation of such additional inputs could be future scope of study. However, the better performance of LS-SVR, particularly for shorter prediction lead-times, as compared to that of ANN, is noticed in this study.

The better performance of LS-SVR over ANN may be

Table 5. Performance measures for LS-SVR and ANN models for training and testing period (1978-2000). Testing period values are shown within parentheses.

Lead-time (in days)	Optimum number of previous daily flow(s) to be considered	LS-SVR		ANN	
		CC	NSE	CC	NSE
1	5	0.83 (0.83)	0.68 (0.67)	0.76 (0.73)	0.58 (0.52)
2	2	0.71 (0.71)	0.51 (0.49)	0.71 (0.71)	0.50 (0.49)
3	1	0.64 (0.64)	0.40 (0.39)	0.63 (0.59)	0.39 (0.33)
4	1	0.61 (0.60)	0.36 (0.35)	0.66 (0.55)	0.43 (0.30)
5	1	0.58 (0.58)	0.33 (0.32)	0.62 (0.56)	0.39 (0.30)

CC: Correlation Coefficient; NSE: Nash-Sutcliffe Efficiency (NSE).

Table 6. Error measures for LS-SVR and ANN models for training and testing period. Testing period values are shown within parentheses.

Lead-time (in days)	LS-SVR			ANN		
	MAE	MARE	RMSE	MAE	MARE	RMSE
1	0.006 (0.008)	0.031 (0.041)	0.024 (0.029)	0.008 (0.009)	0.055 (0.049)	0.028 (0.035)
2	0.008 (0.010)	0.044 (0.059)	0.030 (0.036)	0.008 (0.010)	0.047 (0.056)	0.031 (0.036)
3	0.009 (0.012)	0.056 (0.065)	0.033 (0.040)	0.009 (0.011)	0.053 (0.058)	0.034 (0.042)
4	0.011 (0.012)	0.066 (0.070)	0.034 (0.041)	0.010 (0.012)	0.061 (0.061)	0.033 (0.043)
5	0.011 (0.013)	0.069 (0.073)	0.035 (0.042)	0.010 (0.012)	0.060 (0.067)	0.034 (0.042)

MAE: Mean Absolute Error; MARE: Mean Absolute Relative Error; RMSE: Root Mean Square Error.

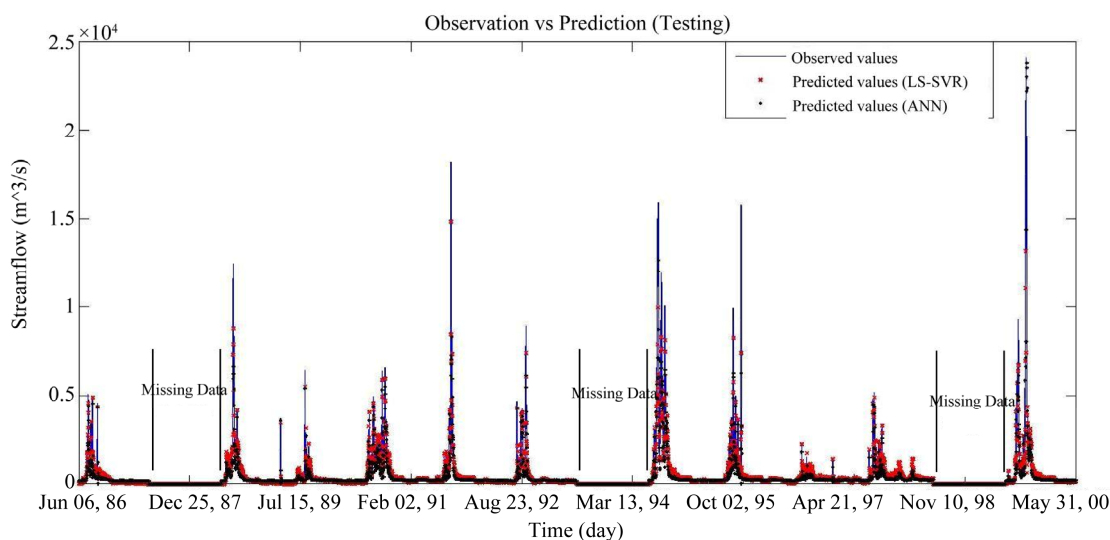


Figure 5. Comparison between observed and predicted river flow from LS-SVR and ANN (1-day-ahead).

attributed to its fundamental approach towards error minimization. Fundamental difference in the working principles of ANN and LS-SVR (SVM, in general) lies in their approaches of risk minimization. ANN follows an Empirical Risk Minimization (ERM) approach, whereas Structural Risk Minimization (SRM) principle is followed in LS-SVR. The SRM minimizes an upper bound on the generalization error, as opposed to ERM which minimizes the error on the training data. Thus, the solutions may fall in to local optima in case of ANN [27]. It is this difference which equips LS-SVR with a greater potential to generalize the regression surface without the danger of overfitting the training data set and global optimum solution is guaranteed in LS-SVR [28].

With respect to the number of parameters, LS-SVR is less complex than ANN. As discussed earlier, there are only three parameters in LS-SVR based approach, namely—regularization parameter (γ) RBF kernel parameter (σ) and number of previous daily river flow values to be considered (n). However, in ANN, the number of hidden layers, number of hidden nodes, transfer functions and so on must be determined, which are comparatively more complex. On the other hand, LS-SVR is able to provide better prediction with small sample size. This is because the decision function of LS-SVR is only determined by supporting vectors. In general, the supporting vectors are only a part of training pattern (from available river flow data) and the remaining pattern are not used in constructing the LS-SVR model. Therefore the performance of LS-SVR may still be acceptable, even if the sample size is small. In contrast, the decision function in ANN is determined by all training data sets [29]. Thus, generalization of relationship between past and future river flow values is more likely in case of LS-SVR as found in this study and thus, more suitable for mul-

tistep-ahead prediction.

5. Conclusions

In this study, daily variation of river flow is modelled and potential Least Square-Support Vector Regression (LS-SVR) is used for multistep-ahead river flow prediction. Daily river flow values from Sandia station at upper Narmada river basin in India are used for illustration. Bargi dam is located few hundred km upstream of Sandia. It is investigated whether it is required to consider the releases from the upstream dam to model the daily variation at Sandia. Parameters of LS-SVR— γ (regularization parameter) and σ (RBF kernel parameter) and optimum numbers of previous river flow values to be considered (n) are estimated based on the model performance during model development period (June 1, 1978 to May 31, 1986 excluding the missing data). The model is tested for the period June 1, 1986 to May 31, 2000 and different statistical performance measures (CC, NSE, RMSE, MAE and MARE) are obtained. These statistical measures confirm the well correspondence between observed and predicted river flow values. This correspondence is found to be better for low and medium range of flow values. However, the peak river flow values are not captured with reasonable accuracy in case of both LS-SVR and ANN.

While comparing the performance during pre- and post-construction of dam, it is found that the prediction performances are similar for both the flow regimes, which indicates that we may ignore the releases from the dam at daily scale for this gauging site. In order to investigate the temporal horizon over which this may be relied upon, a multistep-ahead prediction is carried out and the model performance is investigated up to 5-day-ahead

predictions. The performance is found to be decrease with the increase in lead-time. In other words, the performance is better for shorter lead-times.

General comparison between LS-SVR and ANN model reveals that the performance of LS-SVR is better than that of ANN for all the lead-times—1-day-ahead through 5-day-ahead prediction. Better performance of LS-SVR, in comparison with that of ANN, becomes more prominent for shorter prediction lead-times. Thus, the better performance of LS-SVR, as compared to that of ANN, is noticed for multistep-ahead prediction. The superior performance of LS-SVR over ANN may be attributed to its fundamental approach towards error minimization, ensured global optimum solution and capability to generalize the relationship between past and future river flow values, even with short length of available data. Thus, it may be inferred that Structural Risk Minimization (SRM) is better approach as compared to Empirical Risk Minimization (ERM) and LS-SVR, being a SRM based approach, may be used for multistep-ahead prediction to obtain better performance. Use of exogenous inputs may be of further research interests.

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