

Flow prediction in vegetative channel using hybrid artificial neural network approach

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ABSTRACT

Channel flow–vegetation interaction has been extensively studied in the past few decades and many equations have been developed which essentially differ from each other in derivation and form. As the process is extremely complex, getting deterministic or analytical forms of process phenomena are too difficult. A hybrid neural network model (combining genetic algorithm with neural network), which is particularly useful in modeling processes about which adequate knowledge of the physics is limited, is presented here as a complementary tool to model channel flow–vegetation interactions in submerged vegetation conditions. The prediction capability of the model has been found to be satisfactory. The input significance of the different parameters has been analyzed in the present work in order to find out the influence of these parameters on channel flow velocity.

Key words | flow velocity, genetic algorithm, input significance, sensitivity analysis, vegetation density, vegetation height

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INTRODUCTION

Estimating the flow resistance of vegetation is of great importance in river management, since it may have a significant effect on the conveyance of the channel (Järvelä 2002). The presence of vegetation in a channel also leads to a change in the hydrodynamic characteristics of the channel (Nepf 1999; Shields & Cooper 2000; Millar 2002; Montgomery & Piegay 2003). Interactions between flow and vegetation are complex and depend on environmental factors and plant characteristics such as mean flow velocity, turbulence, channel morphology, water temperature, plant morphology, age and size, and the spatial distribution of plant patches (Yen 2002). Knowledge of flow–vegetation interactions is important for both engineering and ecological applications (Nikora *et al.* 2008). The literature is replete with experimental and numerical works on flow–vegetation interactions. Experiments in laboratory flumes have been carried out to quantify flow–vegetation interactions by various researchers (Wilson *et al.* 2003; Armanini *et al.* 2005). Green (2006) has done field measurements in natural vegetated fields. Also, detailed numerical simulations of flow through vegetation have been performed (Neary 2003; Choi & Kang 2004).

Several empirical (Kouwen & Fathi-Moghadam 2000) and theoretical relations (Stone & Shen 2002; Stephan & Gutknecht 2002) have been proposed to describe flow–vegetation interactions. Huthoff *et al.* (2007) have ascertained that due to their simplicity, empirical equations have better field applicability than theoretical ones. However, empirical relations have the drawback that their applicability is limited to the range of conditions for which they were derived. On the other hand, theoretical descriptions are often complex. Besides, they may require poorly understood closure parameters and sometimes pose practical difficulties when gathering required input data (Huthoff *et al.* 2007). Galema (2009), based on data available in the literature, has compared different predictors of flow characteristics and concluded that no simple predictor exists for both conditions.

Data mining techniques can help in modeling such processes about which the level of available knowledge is too limited to put the relevant information in a mathematical framework (Bhattacharya *et al.* 2004). Data mining is presently being utilized in almost all branches of science as an alternative and complementary technique to the more

traditional physically based modeling systems. Use of artificial neural networks (ANNs) remains in the forefront of this complementary modeling practice. Neural network techniques have been used to study several hydrologic and hydraulic phenomena (Dolling & Varas 2002; Srinivasulu & Jain 2006; Kumar 2011; Wei 2013).

Although conventional ANN setups have been employed in several instances, there are certain difficulties that ANNs are prone to when used with the back propagation training algorithm. The selection of network architecture and its efficient training procedure are major obstacles for any ANN-based model. Due to the profile of the error hyper-surface, ANN generally misses out the global optima. More recent algorithms (genetic algorithm (GA), particle swarm optimization, etc.) from the paradigm of evolutionary computing have made it possible to overcome such limitations of training algorithms (Koza 1992; Giustolisi & Savic 2006). A GA could represent a valid alternative to ANNs to establish a relationship among the parameters (vegetation characteristics, channel characteristics, etc.) influencing the flow. GA, based on natural selection and genetic mechanism, is a global search that grabbles from one population of points to another. As the algorithm continuously samples the parameter space, the search is directed toward the area of the best solution (Goldberg 1989). GA has been applied to many problems in hydrology and water resources (Cheng *et al.* 2002; Wang *et al.* 2007; Nasseri *et al.* 2008).

The primary objective of the present work is to develop the flow prediction model of a submerged vegetative channel. Recent studies have provided a variety of methods for quantifying and interpreting the contributions of the variables in neural networks (Duh *et al.* 1998; Olden & Jackson 2002). Thus the methods developed for input significance testing through neural networks have been applied in the present work to find out variability of flow velocity with other different physical parameters.

FUNCTIONAL ANALYSIS OF THE FLOW-VEGETATION INTERACTION

The use of neural networks allows developing a model incorporating many complex variables of the systems to be built without requiring the explicit formulation of the

possible relationships that may exist among variables. Neural network does function mapping of the input vectors of real values to the real output values. Thus, it needs a functional form of a physical system describing its dependent and independent variables. Yen (2002) has analyzed several equations in terms of their dependent and independent variables. The physical modeling of channel flow through vegetation is very complex. According to Lopez & Garcia (1998), it is a function of many variables: the fluid properties, flow properties, vegetation characteristics, and channel characteristics. A closer look (Table 1) at the existing equations shows that flow-vegetation interactions can be described by the following function:

$$f(u, h, k, i, m, C_d, D) = 0 \quad (1)$$

where u is the mean velocity, h is the flow depth, k is the height of the vegetation, i is the channel slope. D is the diameter of cylindrical vegetation and m is the number of cylinders per m^2 horizontal area. C_d is the non-dimensional drag coefficient. Equation (1) is for homogeneous vegetation of fixed height and diameter. Flow is assumed to be steady and uniform. Assuming a wide channel, sidewall effects can be neglected (Borovkov & Yurchuk 1994). The condition of fully submerged vegetation is $h \gg 5k$. In this case, the vegetation does not block the velocity at the upper part of the water column (Galema 2009). The submerged condition also follows $5k > h > k$. In these conditions, the vegetation is relatively high in relation to the flow depth; as a

Table 1 | Available descriptors for flow velocity in vegetative channel (submerged vegetation)

Author(s)	Equation(s)
Stone & Shen (2002)	$u = \sqrt{\frac{2g}{C_d m D}} \sqrt{i} (1 - D\sqrt{m}) \sqrt{\left(\frac{h}{k} - \frac{1}{4} \pi m D^2\right) \frac{h}{k}}$
Van Velzen <i>et al.</i> (2003)	$u = \sqrt{\frac{2g}{C_d m D}} \sqrt{i} + 18(h - k)^{3/2} \frac{\sqrt{i}}{k} \log \frac{12(h - k)}{1.6k^{0.7}}$
Baptist <i>et al.</i> (2006)	$u = \left(\sqrt{\frac{2g}{C_d m D}} \sqrt{i} \sqrt{k} + \frac{\sqrt{g}}{0.4} \ln \left(\frac{h}{k} \right) \sqrt{i} \right) \sqrt{h}$
Huthoff (2007)	$u = \sqrt{\frac{2g}{C_d m D}} \sqrt{i} \left(\sqrt{\frac{k}{h}} + \frac{h - k}{k} \left(\frac{(h - k)\sqrt{m}}{1 - D\sqrt{m}} \right)^{2/3} \right)$

consequence, the velocity profile changes a great deal over depth. In channels with sufficient density of vegetation, the effect of bottom roughness can also be neglected (Stone & Shen 2002). The present work aims at predicting the flow velocity of a submerged vegetated channel, so u can be made as dependent variable which depends on the various other factors as shown in Equation (1). Keeping this in view, Equation (1) can be written as:

$$u = f(h, k, i, m, C_d, D) \quad (2)$$

The objective here is to construct a neural network model that approximates an unknown input–output mapping of Equation (2) on the basis of given real observations. The goal, however, is not to provide an exact fit to the data but to develop a model that captures the underlying relationship so that it can be used to predict the output at some future observation of the input.

NEURAL NETWORK MODELING

In this paper, a hybrid algorithm using GA technique is proposed to optimize parameters of the ANN. This is done by considering the following steps: first, define the structure of the ANN; second, define the encoding of the interconnection weights; and third, optimize the interconnection weights of the ANN by GA. In this paper, the most popular models of neural networks (NNs) such as feedforward networks are considered. Supervised learning is one of the most effective weight training algorithms, whereby efforts are made to find an optimal set of connective weights for a NN according to some optimality criteria. One of the most popular supervised learning training algorithms for feedforward NNs is back propagation (BP). The BP is a gradient descent search algorithm. It is based on minimization of the total mean square error between the actual output and the desired output (Irie & Miyanki 1988). This error is used to guide the search of the BP algorithm in the weight space. However, the problem of the BP algorithm is that it is very often trapped in local minima and the learning and adaptation speed are very slow in searching for the global minimum of the search space. Some developed evolutionary algorithms, notably GA, have attracted great attention in the NN from many communities (Chen *et al.* 1999; Salajegheh &

Gholizadeh 2005; Giustolisi & Simeone 2006; Singh *et al.* 2007). GAs are parallel stochastic optimization procedures which are good at exploring a large and complex space in an intelligent way to find values close to the global optimum (Goldberg 1989). As compared with BP, GA is more qualified for neural networks if only the requirement of a global searching is considered. A simple chromosome representation is used, which contains information about connections, weights, and biases of the multiple layer perceptron neural network (MPLNN). The parameter learning process, based on GA technique and BP algorithm, is a two-step learning process. In the first step, the initial parameters, such as weights of the NN are tuned by the GA. In the second step, the BP algorithm and the Levenberg–Marquardt method is introduced to train the NN to yield optimal values of weights of the NN. The block diagram of the proposed hybrid algorithm is depicted by Figure 1.

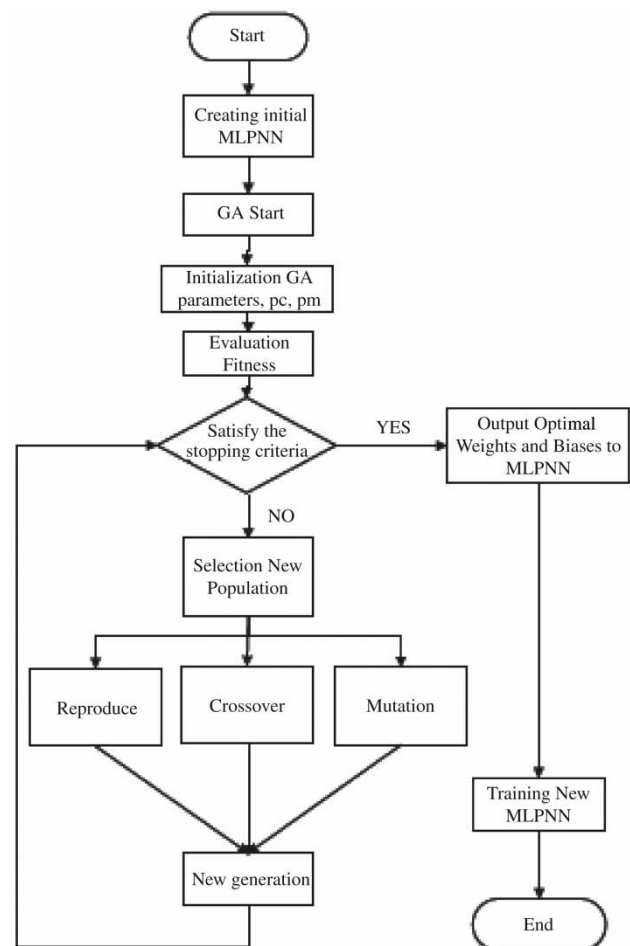


Figure 1 | Flowchart of the learning algorithm.

The procedure of the hybrid algorithm is presented as follows in the sections below.

Chromosome representation

An MLPNN can be represented by a directed graph, encoded on a chromosome with each parameter (weights and biases). All these parameters are memorized by a row vector $C = (c_i)$, $i = 1, 2 \dots M$, where M is the number of all NN parameters. The chromosome can be written as:

$$C = [W_1, W_2] \quad (3)$$

where W_1 denotes the connective weight of link between the input layer and the first hidden layer and W_2 is the connective weight of link between the hidden layer and output layer.

Fitness function

The fitness function is dependent on problem and is used to evaluate the performance of each individual. The error signal of the output neuron j at iteration n (i.e., presentation of the n th training example) is defined by:

$$e_j(n) = d_j(n) - y_j(n), \quad \text{neuron } j \text{ is an output neuron} \quad (4)$$

The instantaneous value of the error energy for neuron j can also be defined as: $(1/2)e_j^2(n)$. Correspondingly, the value of $\xi(n)$ is obtained by summing $(1/2)e_j^2(n)$ over all neurons in the output layer (Haykin 1999).

$$\xi(n) = \frac{1}{2} \sum_{j=C} e_j^2(n) \quad (5)$$

where the set C includes all the neurons in the output layer of the network. For MLPNN it is the sum squared error. The fitness is defined as by summing $\xi(n)$ over all n with respect to the set size N , as shown by:

$$F = \sum_{n=1}^N \xi(n) \quad (6)$$

Here, the objective is to minimize $F(\cdot)$ subject to weights and biases.

Selection

The selection operator is to select individuals from the population for reproduction based on the relative fitness value of each individual. The extraction can be carried out in several ways. One of the most commonly used selection methods is the roulette wheel selection (Goldberg 1989), where individuals are extracted in probability following a Monte Carlo procedure. The extraction probability $p_r(X_i)$ of each individual X_i is proportional to its fitness $F(X_i)$ as a ratio to the average fitness of all the individuals. The offsprings are produced based on this selection.

Crossover

To apply the standard crossover operator the individuals of the population are randomly paired. Crossover takes two parents and performs an interpolation of the two parents. Each pair is then recombined, and the new individuals (offsprings) are formed by the interpolation of parents.

Mutation

After crossover, the new individuals are subjected to mutation. Mutation prevents the algorithm being trapped in a local minimum. A variable is selected with a certain probability and its value is modified by a random value. Here, we choose non-uniform mutation method. Non-uniform mutation changes one of the genes of the parent based on a non-uniform probability distribution.

Levenberg–Marquardt method

The equations for changing the weights (ΔW) during training in the Levenberg–Marquardt method are given as follows:

$$\Delta W = (J^T J + \mu I)^{-1} J^T e \quad (7)$$

where J is the Jacobian matrix of the derivative of each error to each weight, μ is a scalar, and e is an error vector. The Levenberg–Marquardt algorithm performs very well and its efficiency is found to be of several orders above the

conventional back propagation with learning rate and momentum factor.

RESULTS AND DISCUSSION

The entire modeling and analysis has been done by the use of the neural network toolbox of MATLAB[®] software. Galema (2009) documented a comprehensive database of different types of vegetations with hydraulic properties from different sources. Table 2 shows the ranges of the observations used in the present work. The total number of observations used in the modeling is 449. The sample size of the data set is an important issue in neural networks. Neural networks typically require larger sample sizes than conventional statistical procedures for model building and validation. In general, the larger the sample is, the better is the chance for a neural network to adequately approximate the underlying complex patterns without suffering from the problem of overfitting/underfitting. For development of the ANN model, 60% of observations (independent of the source, data have been randomized before selecting training, test, and validation sets) have been assigned as training sets and 20% for testing and 20% for validation. All data sources are not cited in the references of this paper as they can be found in Galema (2009). It should be noted that, like all empirical models, ANNs perform best in interpolation rather than extrapolation (Masters 1993); consequently, the extreme values of the available data are included in the training set. Once data have been divided into their subsets, the input and output variables are preprocessed by scaling them between 0.0 and 1.0 to eliminate their dimension and to ensure that all variables receive equal attention during training. The architecture of the initial neural network (Figure 2) has been chosen based on maximum R^2 over complete observations with increasing number of nodes in the hidden layers (Appendix 1, available online at <http://www.iwaponline.com/jh/016/055.pdf>). The transfer function used in the model is log sigmoid and tan sigmoid at input to hidden and hidden to output layers, respectively. The number of iterations has been fixed at 1,000.

The results of neural modeling are shown in Figure 3. It can be clearly seen from Figure 3 that the linear coefficient of correlation is very high between observed data and values

predicted through neural nets and it is 0.993, 0.984, and 0.988 in training, validation, and testing, respectively. Overall, the linear coefficient of correlation is 0.982 as shown in Figure 3. This shows the learning and generalization performance of the network is good. The non-linear and non-parametric nature of the neural network model is desirable for field applications, it also brings about more opportunities to go wrong in the modeling and application process. Thus, it is worthwhile noting that predictions from ANN models are better when used for ranges of input variables similar to those utilized in model training. Two standard training algorithms – the Levenberg–Marquardt (LM) optimization and the BFGS (Broyden, Fletcher, Goldfarb, and Shanno) quasi-Newton method along with multiple regression analysis have been used to compare the GA-ANN results. The performance of gradient-based methods (ANN-LM and ANN-BFGS) was found to be comparable with those of the GA-ANN algorithm. The model equation generated by multiple regression analysis is given in Appendix 2 (available online at <http://www.iwaponline.com/jh/016/055.pdf>). However, the results of training the algorithm with GA-ANN gave superior predictive models, as reflected in Table 3.

In the present work, formulas listed in Table 1 have also been tested to quantify their prediction capability. Performance analysis of these formulas is shown in Figure 4. As shown in Figure 4, Huthoff's (2007) formula predicts better than other formulas with R^2 around 0.67. The functional form of the empirical formulas does not hold true for all the ranges, for example, Huthoff's (2007) formula works only when h is greater than k . ANN is free from such functional form and will work on all the ranges used in deriving the model.

In the neural network, the connection weights between neurons are the linkages between the input and the output of the network, and therefore are the link between the problem and the solution (Olden & Jackson 2002). Garson's algorithm or 'weights' method includes partitioning the connection weights to determine the relative importance of the various inputs. In the present work, connection weights (Olden & Jackson 2002), Garson's algorithm (Garson 1991) and partial derivative method (Ng *et al.* 2004) have been implemented to know the importance of the input variables on output. Connection weights calculate the product of the raw input–hidden and hidden–output connection weights between each input neuron and output neuron and sum

Table 2 | Summary of hydraulic and vegetation characteristics of the data

Source	No. of observations	D (m)		m (m ⁻²)		k (m)		C _d		h (m)		i		u	
		Min value	Max value	Min value	Max value	Min value	Max value	Min value	Max value	Min value	Max value	Min value	Max value	Min value	Max value
Einstein & Banks (1950)	20	0.0064	0.0064	3	108		0.038		1.4	0.073	0.108	0.00774	0.01024	0.8	1.2
Tsujimoto & Kitamura (1990)	6	0.0015	0.0015	2.500	2.500		0.046		1.46	0.073	0.095	0.001	0.00703	0.117	0.331
Shimizu & Tsujimoto (1994)	13	0.001	0.0015	2.500	10.000	0.041	0.046		1	0.0631	0.1052	0.001	0.00886	0.096	0.331
Dunn <i>et al.</i> (1996)	12	0.0064	0.0064	42	384		0.12	1.13	1.13	0.164	0.335	0.00359	0.01607	0.308	0.854
Meijer (1998)	48	0.008	0.008	64	256	0.45	1.5	0.96	1	0.99	2.5	0.00055	0.00205	0.175	1.242
Stone & Shen (2002)	92	0.00318	0.0127	173	696	0.124	0.124	0.96	1.11	0.151	0.314	0.00009	0.04402	0.017	0.502
Poggi <i>et al.</i> (2004)	5	0.004	0.004	67	1.072		0.12	1.5	1.5	0.6	0.6	0.00004	0.00032	0.3	0.313
Murphy <i>et al.</i> (2007)	24	0.006	0.0064	250	800		0.0139	0.61	1	0.088	0.467	0.00000055	0.000596	0.013	0.197
Fenzl (1962)	26	0.00238	0.00238	11	1.808	0.051	0.152	1.01	1.17	0.058	0.181	0.00163	0.00285	0.052	0.289
Kouwen <i>et al.</i> (1969)	27	0.005	0.005	5.000	5.000	0.05	0.1		3	0.149	0.4	0.0005	0.01001	0.03	0.609
Ree & Crow (1977)	30	0.005	0.005	1.464	1.076	0.203	0.305		1	0.242	0.751	0.00042	0.0021	0.046	0.427
Murota <i>et al.</i> (1984)	8	0.00024	0.00024	4.000	4.000	0.048	0.06	2.75	2.75	0.092	0.116	0.0005	0.00383	0.085	0.258
Tsujimoto <i>et al.</i> (1991)	12	0.0015	0.0015	2.500	2.500	0.0238	0.0419	3.14	3.14	0.07	0.11	0.001	0.007	0.0945	0.4836
Tsujimoto <i>et al.</i> (1993)	12	0.00062	0.00062	10.000	10.000	0.061	0.065		2	0.1	0.16	0.00051	0.01076	0.078	0.385
Ikeda & Kanazawa (1996)	7	0.00024	0.00024	20.000	20.000	0.04	0.045		1	0.142	0.19	0.00247	0.00641	0.349	0.606
Meijer (1998)	7	0.0057	0.0057	254	254	1.55	0.165	1.81	1.81	1.75	2.5	0.00109	0.00208	0.142	0.393
Rowinski & Kubrak (2002)	8	0.000825	0.000825	2.500	10.000	0.165	0.165	1.22	1.35	0.1962	0.2475	0.0087	0.0174	0.1587	0.2948
Järvelä (2003)	12	0.0028	0.003	12.000	512	0.155	0.295		1	0.306	0.7065	0.0002	0.0051	0.072	0.33
Carollo <i>et al.</i> (2005)	80	0.0045	0.0045	28.000	44.000	0.044	0.082		1	0.061	0.272	0.001	0.05	0.211	1.047

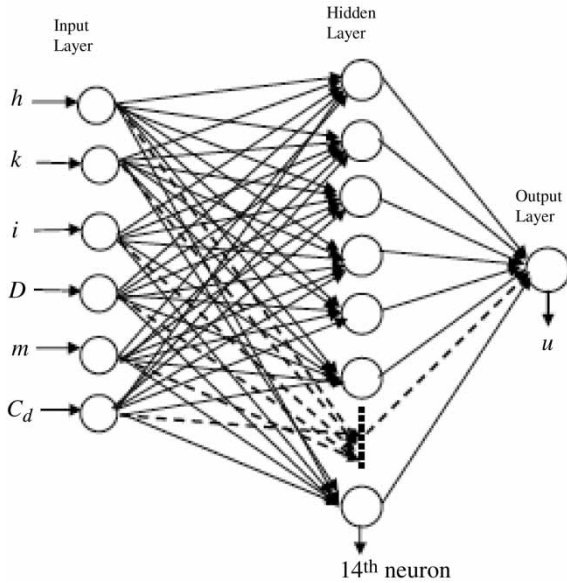


Figure 2 | Optimal artificial neural network.

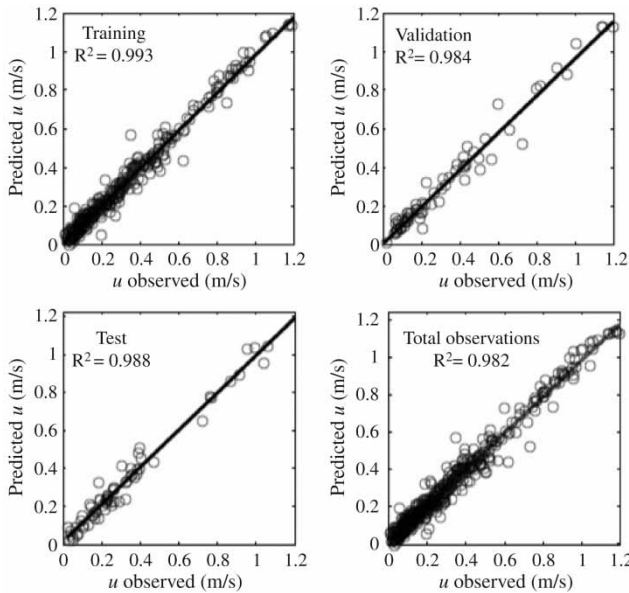


Figure 3 | Results of neural network model.

the products across all hidden neurons. The relative importance of input parameter j is determined through the following formula:

$$Imp(j) = \sum_{x=1}^n CW_{jh(x)} CW_{ho(x)} \tag{8}$$

Table 3 | Comparative analysis of different intelligent predictors

Methodology	R^2	
	Training	Testing
GA-ANN	0.993	0.988
ANN-LM	0.965	0.92
ANN-BFGS	0.94	0.88
Multiple regression	0.88	0.81

where $Imp(j)$ is the relative importance of parameter j , n the total number of hidden nodes, x the index number of hidden node, $CW_{jh(x)}$ the connectivity weight between input parameter j and hidden node x , $CW_{ho(x)}$ is the connectivity weight between hidden node x and the output node. The methodology for Garson’s algorithm is as follows:

- (a) For each hidden neuron H , divide the absolute value of the input–hidden layer connection weight by the sum of the absolute value of the input–hidden layer connection weight of all input neurons.

For $H = 1$ to nH , for $j = 1$ to n_j ,

$$A_{jH} = \frac{|W_{jH}|}{\sum_{i=1}^{n_j} |W_{iH}|} \tag{9}$$

- (b) For each input neuron j , divide the sum of the A_{jH} for each hidden neuron by the sum for each hidden neuron of the sum for each input neuron of A_{jH} , multiply by 100. The relative importance of all output weights attributable to the given input variable is then obtained.
- $j: 1$ to n_j

$$RI(\%)_j = \frac{\sum_{n=1}^{nH} A_{jH}}{\sum_{n=1}^{nH} \sum_{j=1}^{n_j} A_{jH}} \times 100 \tag{10}$$

Connection weight and Garson’s algorithm are based on the weight matrices of the neural network. Partial derivative method computes the partial derivatives of the ANN output with respect to the input neurons by making use of the Taylor expansion to approximate the output perturbations with respect to either the input or weight perturbations. Calculations of all approaches for input significance testing are tabulated in Table 4. The values depicted in Table 4 have been derived from both initial and final weight values of the ANN model. As can be seen from Table 4, flow depth

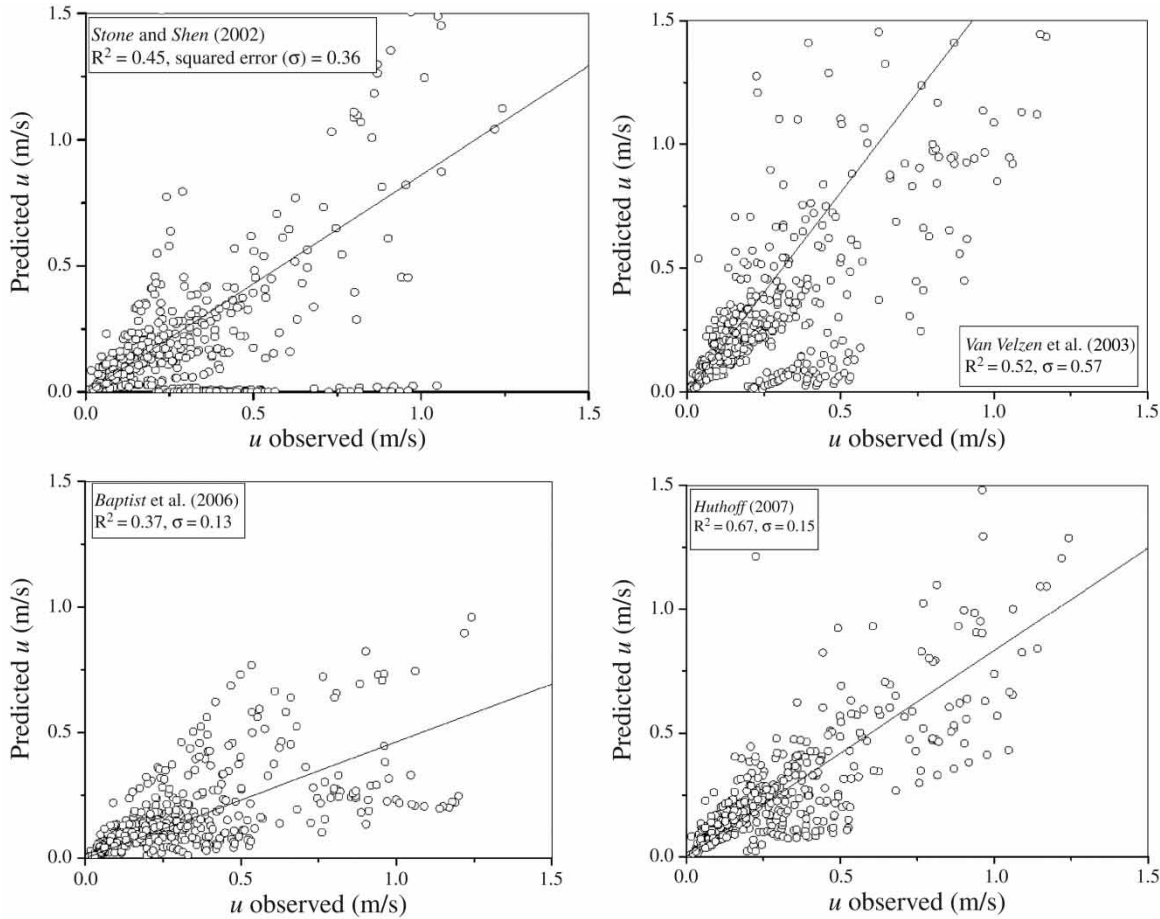


Figure 4 | Performance analysis of flow velocity predictors in vegetative channel.

Table 4 | Input significance and ranking of the variables

Variables	Connection weights		Garson's algorithm		Partial derivatives	
	Initial weights	Final weights	Initial weights	Final weights	Initial weights	Final weights
i	0.08 (6)	0.01 (6)	7.05 (5)	9.12 (6)	0.03 (6)	0.02 (6)
m	0.165 (4)	0.13 (5)	19.1 (3)	15.33 (4)	0.165 (4)	0.12 (5)
k	0.163 (5)	0.15 (4)	4.4 (6)	11.58 (5)	0.09 (5)	0.14 (4)
D	0.164 (3)	0.19 (3)	9.45 (4)	19.86 (3)	0.17 (3)	0.27 (2)
h	0.171 (2)	0.23 (2)	30.5 (1)	23.02 (1)	0.28 (1)	0.15 (3)
C_d	0.25 (1)	0.28 (1)	29.5 (2)	21.03 (2)	0.25 (2)	0.29 (1)

The value in brackets shows rank importance.

and drag coefficient have much more influence on the mean velocity of a vegetative channel. For the channel with densely distributed vegetation, the drag of vegetation becomes the major parameter to characterize the channel

hydrodynamics. Turbulent vortices are created in the wakes downstream of the protruding stems in a vegetative channel (Akilli & Rockwell 2002). The associated energy losses of the mean flow field cause the flow to slow down.

These drag effects can become very important, which is also shown in Table 4. Calculations shown in Table 4 indicate that any error in determining flow depth and drag coefficient may result in under- or over-estimation of mean velocity. Table 4 also highlights the importance of D on mean velocity prediction.

Contribution plots for each of the predictor variables have been constructed in the present work, as shown in Figure 5. Figure 5(a) shows the variation of u with h and i by keeping the other variables at mean value. The range of i has been divided into 10 parts, and plots of h and u have been drawn at each value of i . It can be seen from Figure 5(a): (i) mean velocity increases with an increase in h and (ii) it also increases with an increment in the channel slope (i). Mean velocity is directly proportional to flow depth and channel slope as per any flow equations existing in the literature. It is probably more relevant to analyze the trends shown in Figure 5(a) through Manning's formulation as it is most often employed in practice. The trend shown in Figure 5(a) matches with the most used Manning's formula for mean velocity. Mean velocity will increase with flow depth and channel slope as per Manning's formula. Manning's formula, however, has Manning's coefficient as

controlling factor, thereby limiting the increasing trend of mean velocity with flow depth and channel slope. The same can be seen in Figure 5(a). When flow depth and channel slope have high values, increment in mean velocity is slow and it tends towards attaining an asymptotic nature. Figure 5(b) shows the combined variation of D and k with u . D and k are descriptors of vegetation properties. Except for Stone and Shen's equation (Stone & Shen 2002), most of the equations given in Table 1 show that u will decrease non-linearly with an increase in D . Equations given in Table 1 do not show a general behavior of k with u ; however, Figure 5(b) shows that u will increase with k . The trend that u will decrease non-linearly with an increase in D is also shown in Figure 5(c). m represents the density of vegetation. When m is very high, flow will be very low because of high density of vegetation, which is shown in Figure 5(c). Drag opposes the fluid motion; this general behavior is shown in Figure 5(d). u will decrease with an increase in C_d , which may be due to increasing moment absorbing area. The non-linearity, shown in Figures 5(a) to 5(d), probably originates from attenuation effects of vegetation that reduce bed surface contribution to the overall resistance (Armanini *et al.* 2005).

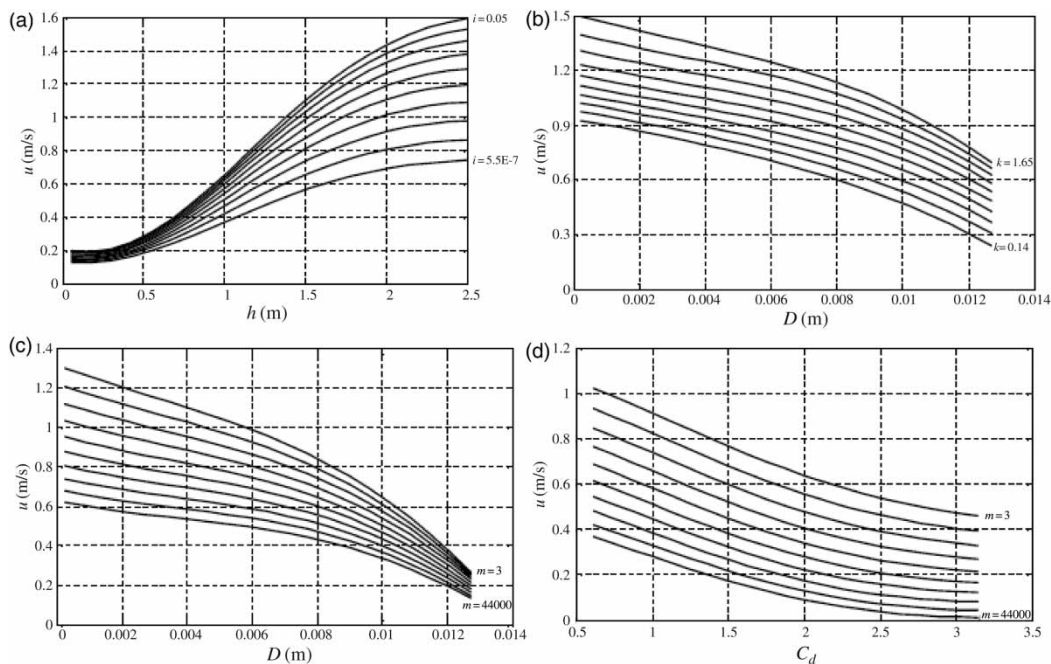


Figure 5 | Variability of u with input parameters: (a) relationship between u , h , and i ; (b) variation of D and k with u ; (c) variation of D and k with u ; (d) relationship between u , C_d , and m .

CONCLUSION

Vegetation–flow interactions are central to many problems of practical interest to hydraulic engineers, including flood risk studies, sediment transport studies, and the analysis of the hydraulic performance of river restoration schemes. Existing predictors are too complex and limited in terms of their applicability. Based on a large database of submerged vegetation conditions (rigid and flexible), the present work formulates a neural network model of mean velocity in a vegetative channel comprising all the relevant parameters. In this way, a qualitative effect of all types of parameters can be incorporated in the prediction of flow velocity in a vegetative channel. The generalization capacity of the model is very good considering there are so many different parameters affecting flow–vegetation interactions. In engineering applications, errors in determining the governing parameters affect the accuracy of the prediction. The present analysis based on the input significant test shows that in order to get the accurate flow velocity prediction, there should be more accurate determination of flow depth and drag coefficient due to vegetation. The variability of flow velocity with different parameters is also important to understanding the process. The present work shows the mean velocity decreases with an increase in vegetation diameter, vegetation density, and drag coefficient. The variability of mean velocity with vegetation height is an increasing one.

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