Modelling the effect of land use change on hydrological model parameters via linearized calibration method in the upstream of Huaihe River Basin, China

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ABSTRACT

Conceptual rainfall–runoff models have become a basic tool for evaluating effects of land use/cover changes on the hydrologic processes in small-scale as well as large watersheds. The runoff-producing mechanism is influenced by land use/ cover changes. In this study, we analysed the effect of land use change on hydrological model parameters by calibrating the model parameters of different time periods with different land use via a linearized calibration method. The parameter calibration of a conceptual model usually involves the construction of objective function and optimization methods for good performance of observed data. However, the objective function of the minimum-sum-squared error will introduce an unrelated optimum solution for the parameter calibration problem of a conceptual model, which belongs to a highly complex nonlinear system. Thus, a linearized parameter calibration method, which searches for the optimal value on a parameter surface, is presented, based on the analysis of the problems of the objective function of the minimum-sum-squared error. Firstly, an ideal model is shown that illustrates the efficiency and applicability of this method. Secondly, the novel method is demonstrated for solving the Xinanjiang daily model parameter calibration. Finally, 50 years of data are divided into 4 different periods for parameter comparison, through which the effects of land use/cover changes on runoff in Dapoling watershed are evaluated. The results show that the linearized parameter calibration method is convergent, reasonable and effective. For example, the model parameter of evapotranspiration coefficient KC varied considerably, from 0.658 to 0.922, in response to land use/cover change within the watershed.

Keywords: land use/cover change; parameter calibration; linearized; upper Huaihe River Basin

INTRODUCTION

Conceptual rainfall-runoff models have been widely used for quantification of the effect of land use and land cover change on runoff dynamics (Burns et al., 2005; Jiang et al., 2004, Wan and Yang, 2004; Wang et al., 2007; Yuan and Shi, 2001; Zhang et al., 2004). The runoff-producing mechanism and hydrological response will be influenced by land use/cover changes, and this influence can be reflected by the calibrated parameter values of different time periods (Tao et al., 2015; Zhang et al., 2016; Zhong-Bo et al., 2014). A good model performance is critical to evaluate the effect of land use/cover change on hydrological processes. The major difficulty is that these models generally have a large number of parameters that cannot be obtained directly from field measurements; hence, model calibration is needed (Chen and Chau, 2006; Wang et al., 2009; Wu et al., 2009). The process of model calibration is normally done either manually or by using computer-based, automated procedures. Because of the time-consuming nature of manual calibration and the difficulty in explicitly assessing the confidence of model simulations, automation of the calibration process has been explored (Gupta et al., 1998; Gupta et al., 1998; Hendrickson et al., 1988; Johnston and Pilgrim, 1976; Yapo et al., 1996). Research into automated optimization methods has led to the development of stochastic global optimization searching

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http://dx.doi.org/10.4314/wsa.v43i2.11 Available on website http://www.wrc.org.za ISSN 1816-7950 (Online) = Water SA Vol. 43 No. 2 April 2017 Published under a Creative Commons Attribution Licence methods based on extending the search area or increasing the number of initial optima (Cheng et al., 2002; Cheng et al., 2008; Duan et al., 1992; Franchini, 1996; Franchini and Galeati, 1997; Ritzel et al., 1994; Savic et al., 1999; Vasquez et al., 2000; Wang, 1991).

Rainfall-runoff model calibration is a highly complex nonlinear problem (Chau, 2007; Cheng et al., 2005; Muttil and Chau, 2006; Taormina et al., 2012). A successful calibration depends not only on effective optimization methods, but also on the calibration objective. Minimum-sum-squared error (MSSE) has been widely used as an objective function (Bárdossy and Singh, 2008; Bao, 1991; Bao et al., 2004; Rusli et al., 2015; Zádor et al., 2006). This objective function has been shown to work well in linear parameter calibration, which has been proved to attain the unique optimal value. However, it will introduce an unrelated local optimal into parameter calibration of the nonlinear function.

Therefore, a more convenient and efficient method for parameter estimation of hydrological models is to study the hydrological model parameters, model structure and hydrological simulation. In this paper, a methodology for system decomposition and degeneration of dimensionality has been developed in the calibration of the Xinanjiang model. The objective of degeneration is to reduce the number of parameters being optimized simultaneously to at least 3 or fewer. An ideal model was used to verify the efficiency and applicability of this method. Finally, this linearized calibration method was applied for Xinanjiang model parameter calibration to evaluate the effect of land use/cover change on the runoff of the upper Huaihe River Basin. We only compare the best calibrated parameters of different periods to assess the influence of land use change on runoff generation via this fast and convenient parameter calibration method. Model parameters might be different when the model is calibrated to different time periods. The change of land use and the impact on hydrological response processes are analysed by the model parameters at different stages.

METHODS

The Xinanjiang model

The Xinanjiang model (Zhao, 1992) was chosen for the study. The primary feature of the Xinanjiang model is the concept of runoff formation on repletion of storage, which means that runoff is not produced until the soil moisture content of the aeration zone reaches field capacity. The hypothesis was first proposed in China in the 1960s, and much subsequent experience supports its validity for humid and semi-humid regions. According to the original formulation, runoff generated was separated into 2 components using Horton's concept of a final, constant, infiltration rate. Infiltrated water was assumed to go to the groundwater storage and the remainder to surface or storm runoff. However, evidence of variability in final infiltration rate, and the unit hydrograph to connect the storm runoff to the discharge from each sub-basin, suggested the necessity of a third component. Guided by the work of Kirkby (1978), an additional component, interflow, was provided in the model in 1980 (Zhao, 1992). The modified model is now successfully and widely used in China (Guo et al., 2016; Ju et al., 2009; Yao et al., 2012).

In the Xinanjiang model, the basin is divided into a set of sub-basins. The outflow from each sub-basin is first simulated and then routed down the channels to the main basin outlet. The flow chart is shown in Fig.1. Based on the concept of runoff formation on repletion of storage, the simulation of outflow from each sub-basin is computed in 4 major steps:

- 1. Evapotranspiration and runoff production: the evapotranspiration that generates the deficit of the soil storage is divided into 3 layers: upper, lower, and deep; the runoff production that produces the runoff according to the rainfall and soil storage deficit.
- Separation of runoff components: the runoff separation that divides the previously determined total runoff into 3 components: surface runoff, subsurface runoff, and groundwater runoff.
- 3. Overland flow concentration within each sub-basin: this is represented by the convolution with an empirical unit hydrograph or by linear reservoir to produce *Q*, the sub-basin outflow. In this paper, we calculate the surface flow, interflow and groundwater of sub-basin outflow. Respectively, by linear reservoir (parameters *CS*, *CI* and *CG* in Fig. 1).
- 4. River network flow concentration: flood routing from the sub-basin outlets to the total basin outlet is achieved by applying the Muskingum method to successive sub-reaches (parameters *KE* and *XE* in Fig. 1).

The flow chart for the Xinanjiang model is given in Fig. 1. The inputs of the model were rainfall (P) and measured pan evaporation (EM). The outputs were the discharge (TQ) from the whole basin and the actual evapotranspiration (E), which includes 3 components, EU, EL, and ED.

The state variables were the area mean tension water storage (W) and the area mean free water storage (S). The area mean tension water (W) had 3 components (WU, WL, and WD) in the upper, lower, and deep layer. *FR* is the runoff-contributing area factor that was related to *W*. The rest of the symbols inside the

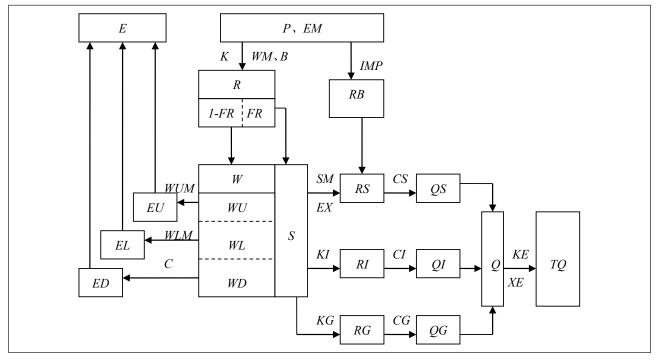


Figure 1

Flow chart for the Xinanjiang model: all symbols inside the blocks are variables, including inputs, outputs, state variables, and internal variables, whereas those outside the blocks are parameters.

blocks were all internal variables. RB was the runoff directly from the small portion of impervious area. R was the runoff produced from the pervious area and divided into 3 components (RS, RI, and RG), referred to as surface runoff, subsurface runoff, and groundwater runoff, respectively. The three components were further transferred into QS, QI, and QG, and together form the total inflow to the channel network of the sub-basin. The outflow of the sub-basin was Q. Then, the Muskingum method was used to calculate the discharge (TQ).

The model had 16 parameters: evapotranspiration parameters *K*, *WUM*, *WLM* and *C*; runoff production parameters *WM*, *B* and *IMP*; parameters of runoff separation *SM*, *EX*, *KI* and *KG*; runoff concentration parameters *CS*, *CI* and *CG*; Muskingum parameters *KE* and *XE*. The details of the Xinanjiang model and the parameter calibration refer to Qu et al. (2012). There are 14 parameters used in the Xinanjiang model. Additionally, another 2 parameters are used for flow routing along the main rivers (see Table 1 for a list of the 16 model parameters). Generally, the output is more sensitive to 9 parameters (Zhao 1992): *KC*, *SM*, *KI*, *KG*, *CS*, *CI*, *CG*, *KE* and *XE*. For the daily model, there are only 4 sensitive parameters, that is, *KC*, *SM*, *KI* and *KG* (Song et al., 2013; Zhijia et al., 2011).

Change detection (parameters comparison)

Qu et al. (2012) used 3 different methods to evaluate the land use/cover change effects on hydrological processes in Dapoling basin. In this study, we only compare the best calibrated parameters of different periods to assess the influence of land use change on runoff generation.

Model parameters might be different when the model is calibrated to different time periods. The Xinanjiang model consists of 16 parameters, and 9 of these (*KC*, *SM*, *KI*, *KG*, *CS*, *CI*, *CG*, *KE*, *XE*) (Table 1) are very sensitive to the calibration process (Song et al., 2013; Zhao, 1992; Zhijia et al., 2011). As a result, the calibration process is time consuming. To reduce this complexity, a methodology for system decomposition and degeneration of dimensionality has been developed in the calibration of the Xinanjiang model. By decomposition, the model structure is analysed hierarchically. The calibration is carried out from a lower to higher hierarchy; to this end, different objective functions are used for different levels. The objective of degeneration is to reduce the number of parameters being optimized simultaneously to at least 3 or fewer. The reduction is based on the sensitivity analysis and structure constraint. For example, there is a daily model and an hourly model for the Xinanjiang model. The objective function for the daily model is to keep the mean annual water balance, which means KC is the most important parameter that needs to be calibrated. Those insensitive parameter values (e.g. WUM, WLM, C and so on) can be chosen by experience. In this study, the daily model and linearized parameter calibration method are used for automatic parameter calibration of the Xinanjiang model.

Linearized parameter calibration method

The linearized calibration method is a new optimization algorithm. It was developed to solve the theoretical problem of unrelated local optima produced in the nonlinear model parameter calibration by using the objective function based on error sum of squares. The calibrated parameter values are stable when using different initial parameter values via the linearized calibration method, which can find the true parameter values without producing unrelated local optima. Furthermore, compared with the SCE-UA method and the simplex method, the linearized calibration method has higher calculation accuracy and convergence rate and the parameter calibration results are also very stable, i.e., not influenced by the different initial parameter values. In the real model study, the method can also find the unique global optima and has good performance in parameter calibration. The linearized calibration method is an efficient, effective, and robust calibration method (Bao and

	TABLE 1			
Parameter	Model parameters and variation range Explanation	unit	Lower bound	Upper bound
Evapotranspi	ration calculation			
KC	Ratio of potential evapotranspiration to the pan evaporationww		0.6	1.2
WUM	Tension water capacity of upper layer	mm	5	20
WLM	Tension water capacity of lower layer	mm	60	90
С	Deeper evapotranspiration coefficient		0.08	0.18
Runoff genera	ation calculation			
WM	Areal mean tension water capacity	mm	120	220
В	Exponential of the distribution of tension water capacity		0.1	0.4
IMP	Ratio of impervious area to the total area of the basin		0.01	0.02
Water source	separation			
SM	Free water storage capacity	mm	10	50
EX	Exponential of distribution water capacity		1	1.5
KG	Groundwater outflow coefficient of free water storage		0.2	0.6
KI	Interflow outflow coefficient of free water storage		0.2	0.6
Concentratio	n calculation			
CS	Recession constant of surface water storage		0.4	0.7
CI	Recession constant of interflow storage		0.5	0.9
CG	Recession constant of groundwater storage		0.990	0.998
KE	Residence time of water	h	0.5	1.5
XE	Muskingum coefficient		0	0.5

Zhao, 2014; Bao et al., 2013). The basic principles and derivation process of the linearized calibration method are as follows:

The traditional parameter estimation method is based on the objective function of the minimum-sum-squared error. For any function *f*:

$$f = f(\theta, X) \tag{1}$$

where: θ is parameter vector, *X* is independent variable vector. The objective function of minimum-sum-squared error is constructed as:

$$\min_{\theta \in \Re} \left\{ F(\theta, X) = \sum_{i=1}^{L} \left(f(\theta, X_i) - f_i \right)^2 \right\}$$
(2)

where: $\{X_i, f_i; i = 1, 2, \dots, L\}$ is the sample series for parameter calibration. It can be observed that Eq. 2 increases the power of linear parameter from 1 to 2 and for a linear function this is clearly correct. However, the sum-squared causes the increase in unrelated optimal values when solving nonlinear function parameter optimization problems. For example, if the power of parameters is equal to or greater than 2, sum-squared will double the power of parameters, which will be reduced only by 1 through first-order derivation – thus, it will increase the number of parameter solutions. A 1-parameter function with 1 sample is employed to demonstrate the problem. It can be seen from the analysis that the objective function of the minimum-sumsquared error works well in parameter optimization for a linear function; however, this is not the case for a nonlinear function.

Based on this analysis, we developed a linearized parameter calibration method. Firstly, we linearize the nonlinear function using parameters as independent variables. Secondly, the linearized parameters can be calibrated with the objective function of minimum-sum-squared error. Finally, iteration is adopted to gradually approach the optimal value of nonlinear function parameters.

The forward difference of the partial derivative of *f* as in Eq. 1 with respect to parameter θ_i :

$$\frac{\partial f}{\partial \theta_i} = \frac{f([\theta_1, \cdots \theta_i + \Delta \theta_i, \cdots \theta_n]^T, X)}{\Delta \theta_i} - \frac{f([\theta_1, \cdots \theta_i, \cdots \theta_n]^T, X)}{\Delta \theta_i}$$
(3)

In the calibration process, $\theta_{i,j}$ represents estimation of θ_i in Step j and $\theta_{i,j+1}$ is j+1 step estimation of θ_i . Substituting $\theta_{i,j}$ and $\theta_{i,j+1}$ into Eq. 3, we obtain:

$$\frac{\partial f}{\partial \theta_i}\Big|_{\theta_i = \theta_{i,j}} = \frac{f([\theta_1, \cdots, \theta_{i,j+1}, \cdots, \theta_n]^T, X)}{\theta_{i,j+1} - \theta_{i,j}} - \frac{f([\theta_1, \cdots, \theta_{i,j}, \cdots, \theta_n]^T, X)}{\theta_{i,j+1} - \theta_{i,j}}$$
(4)

Set $f_{i,j+1} = f([\theta_1, \dots, \theta_{i,j+1}, \dots, \theta_n]^T, X),$ $f_{i,j} = f([\theta_1, \dots, \theta_{i,j}, \dots, \theta_n]^T, X)$ Eq. 4 can be rewritten as:

$$f_{i,j+1} = f_{i,j} + \frac{\partial f}{\partial \theta_i}\Big|_{\theta_i = \theta_{i,j}} \left(\theta_{i,j+1} - \theta_{i,j}\right)$$
(5)

Eq. 5 shows the relation of the parameters, function values and partial derivation with respect to parameters in the adjacent 2 steps in the parameter searching process. The partial derivation with respect to a parameter exhibits the degree of function variation resulting from the variation of parameters. Usually, the greater the degree of function change, the more sensitive the parameter is in the function, which is called parameter sensitivity. Considering the variation of all parameters in the consecutive 2 steps, Eq. 5 can be transformed into:

$$f_{j+1} = f_j + \frac{\partial f}{\partial \theta_1} \bigg|_{\theta_1 = \theta_{1,j}} (\theta_{1,j+1} - \theta_{1,j}) + \frac{\partial f}{\partial \theta_2} \bigg|_{\theta_2 = \theta_{2,j}} (\theta_{2,j+1} - \theta_{2,j}) + \cdots + \frac{\partial f}{\partial \theta_n} \bigg|_{\theta_n = \theta_{n,j}} (\theta_{n,j+1} - \theta_{n,j})$$

$$(6)$$

Eq. 6 is the linear approximate of nonlinear function f with respect to parameter θ . In the process of parameter calibration, we assume L groups of observation samples:

$$(X_1, f_1) (X_2, f_2) \cdots (X_L, f_L)$$
 (7)

Substitute it into Eq. 6:

$$\begin{cases} f(\theta_{j+1}, X_1) = f(\theta_j, X_1) \\ + \frac{\partial f(\theta_j, X_1)}{\partial \theta_1} (\theta_{1,j+1} - \theta_{1,j}) \\ + \dots + \frac{\partial f(\theta_j, X_1)}{\partial \theta_n} (\theta_{n,j+1} - \theta_{n,j}) \\ \vdots \\ f(\theta_{j+1}, X_L) = f(\theta_j, X_L) \\ + \frac{\partial f(\theta_j, X_L)}{\partial \theta_1} (\theta_{1,j+1} - \theta_{1,j}) \\ + \dots + \frac{\partial f(\theta_j, X_L)}{\partial \theta_n} (\theta_{n,j+1} - \theta_{n,j}) \end{cases}$$
(8)

The vector matrix of Eq. 8 is:

$$f_{j+1} = f_j + S(\boldsymbol{\theta}_{j+1} - \boldsymbol{\theta}_j) \tag{9}$$

where:

$$\begin{split} f_{j+1} &= \left[f(\theta_{j+1}, X_1), f(\theta_{j+1}, X_2), \cdots, f(\theta_{j+1}, X_L) \right]^{t} \\ f_j &= \left[f(\theta_j, X_1), f(\theta_j, X_2), \cdots, f(\theta_j, X_L) \right]^{t} , \\ \theta_j &= \left[\theta_{1,j}, \theta_{2,j}, \cdots, \theta_{n,j} \right]^{t}, \theta_{j+1} = \left[\theta_{1,j+1}, \theta_{2,j+1}, \cdots, \theta_{n,j+1} \right]^{t} \end{split}$$

$$S = \begin{bmatrix} \frac{\partial f(\theta_j, X_1)}{\partial \theta_1} & \cdots & \frac{\partial f(\theta_j, X_1)}{\partial \theta_n} \\ \vdots \\ \frac{\partial f(\theta_j, X_L)}{\partial \theta_1} & \vdots \\ \frac{\partial f(\theta_j, X_L)}{\partial \theta_1} & \frac{\partial f(\theta_j, X_L)}{\partial \theta_n} \end{bmatrix}$$
 is called parameter sensitivity matrix. Eq. 9 is the linear approximation of actual function *f*. Substitute the observation sample into Eq. 9 and it can be written as:

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$$f = f_j + S(\theta_{j+1} - \theta_j) + E$$
⁽¹⁰⁾

where $E = [e_1, e_2, ..., e_L]^T$ is the difference between f and f_{i+1} . As the iterative process continues, the difference between f and f_{i+1} approaches a minimum. We obtain:

$$\min_{\theta \in \Re} \left\{ F_{j+1} = (f - f_{j+1})^T (f - f_{j+1}) \right\}$$
(11)

where $f = [f_1, f_2, \dots, f_i$ s observation vector. The optimal estimation value of $\theta_{_{i+1}}$ can be inferred from Eq. 10 in the linear condition:

$$\theta_{j+1} = \theta_j + (S^T S)^{-1} S^T (f - f_j)$$
(12)

Considering Eq. 9 is the linear approximation of actual function $f(\theta, X)$, which cannot guarantee that θ_{i+1} is the optimum in the searching direction, a scale coefficient b needs to be multiplied to make the actual error function the optimum in the searching direction.

$$\boldsymbol{\theta}_{j+1} - \boldsymbol{\theta}_j = b(S^T S)^{-1} S^T (f - f_j)$$
(13)

The variation range of b is (0, 1), which can be determined by minimum-sum-squared error of actual function:

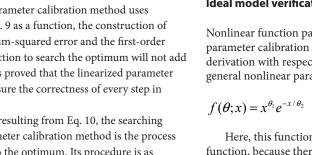
$$\min_{b} \left\{ e^{T} e = (f - f(\theta_{j+1}, b, X))^{T} (f - f(\theta_{j+1}, b, X)) \right\}$$
(14)

Eq. 12 determines the searching direction, and Eqs 13 and 14 determine the step length of parameter, which not only ensures the correct searching direction, but also satisfies the minimum error in the searching direction. Eqs 12, 13 and 14 comprise the linearized parameter calibration method, which has a similar form to the Gauss-Newton iteration method, but a different theoretical basis.

Since the linearized parameter calibration method uses linear approximation of Eq. 9 as a function, the construction of the objective function of sum-squared error and the first-order derivation of objective function to search the optimum will not add unrelated local optima. It is proved that the linearized parameter calibration method can ensure the correctness of every step in searching direction.

Considering the error resulting from Eq. 10, the searching process of linearized parameter calibration method is the process of step-by-step approach to the optimum. Its procedure is as follows:

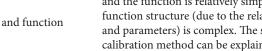
- 1. Set an initial value of parameter θ_0
- 2. Calculate the parameter sensitivity matrix S and function vector f_i according to the parameter vector
- 3. Determine new parameter vector θ_{i+1} and search direction $\Delta \theta$ $= \theta_{i+1} - \theta_i$ with Eq. 12
- 4. Determine scale coefficient b with Eq. 14
- 5. The search will be stopped if θ_{i+1} arrives at the optimum individual. If not, return to Step (2) and go on to the next iterative operation. The framework flow chart for the linearized parameter calibration method is shown in Fig. 2.

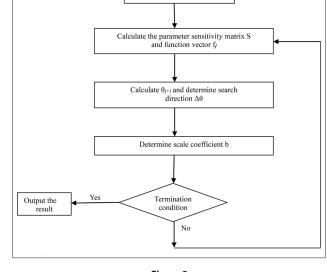


Here, this function is adopted as the nonlinear parameter function, because there is only one independent variable xin this function and 2 parameters θ_1 , θ_2 . In addition, these 2 parameters for the independent variables are highly nonlinear, and the function is relatively simple in the expression, but the function structure (due to the relationship between variables and parameters) is complex. The superiority of the parameter calibration method can be explained more clearly by using this function.

The derivation function of a parameter is:

$$\begin{cases} \frac{\partial f(\theta, x)}{\partial \theta_1} = \ln(x) x^{\theta_1} e^{-x/\theta_2} \\ \frac{\partial f(\theta, x)}{\partial \theta_2} = \frac{x^{\theta_1 + 1}}{\theta_2^2} e^{-x/\theta_2} \end{cases}$$
(16)





Set initial parameters:00

Figure 2 The flow chart of the linearized calibration method

It can be seen that the linearized parameter calibration method is simpler than the conventional optimization method from Step (3) of the calculation process. The optimal parameter vector $\theta_{i,i}$ can be determined through θ_{i} and linear condition using Eq. 12 instead of the trial-and-error method.

To prove the validity of Step (3) that ensures the correct searching direction $\Delta \theta = \theta_{j+1} - \theta_j$, we only need to verify the error F_{i+1} corresponding to new parameter vector θ_{i+1} of any step calculated by Eq. 12 is less than F_i corresponding to the above step parameter vector θ_i . The proof is shown in Appendix 1.

APPLICATION EXAMPLE

Ideal model verification

Nonlinear function parameter calibration refers to nonlinear parameter calibration of explicit function, which can obtain derivation with respect to parameters directly. The following general nonlinear parameter function is used in this paper:

$$f(\theta; x) = x^{\theta_1} e^{-x/\theta_2} \tag{15}$$

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Ideal model refers to the model with no error in the independent variable, dependent variable and parameters. We assume $\theta_1 = 2$ and $\theta_2 = 10$ in the ideal model, as Eq. 15 and *x* varies in the range of [1, 100] with equal intervals.

Table 2 lists the process of searching parameters in every step, starting from a group of initial parameters ($\theta_{1,0} = 1.2427$; $\theta_{2,0} = 49.4716$). It can be seen from Table 2 that the parameters sought from every step follow a loop approach; the true value and the objective function of sum-squared error gradually decrease in the process of searching, as does the searching step length in every loop. In the beginning, sum-squared error and searching step length are large and they decrease sharply after several iterations, which demonstrates the significant convergence and effectiveness of the linearized parameter calibration method.

Table 3, on the next page, summarizes the calibration results starting from different initial parameter values. In the table, *J* is an iterative step for searching for optimum. According to the results given in Table 3:

- 1. The influence of initial parameter value is minor for linearized parameter calibration. The square deviation of a parameter decreased from the initial value of 37.2895 and 1 804.38 to the final optima of 7.2×10^{-7} and 7.8×10^{-5} , respectively.
- 2. The linearized parameter calibration method is objective with a high degree of accuracy. The relative error between the results and true value of every trial are 9.5×10^{-5} and 16.3×10^{-5} .
- 3. The iteration number of the linearized parameter calibration is low, and the effect is good. According to the results from Table 3, the minimum iteration number is 11, the maximum number is 35 and the average loop number is 20.6.

Table 4 demonstrates the calibration results with consideration of error in the dependent variable. The error was generated using a normal distribution model. The error proportion is the average per cent relative to the true value of the function $f(\theta; x)$. *D* is the distance between the evaluated parameter and the true value, calculated by Eq. 17:

$$D = \sqrt{(2 - \theta_1)^2 + (\mathbf{0} - \theta_2)^2}$$
(17)

According to the results from Table 4, the error of dependent variable has certain but little influence on the parameter estimation. The average of relative error distance is only 1.21%.

Case study

Dapoling Station was used as a case study application, situated in the upper reaches of Xixian basin with a drainage basin of 1 640 km²; the length of its main channel is 73 km (Fig. 3).

The river flows through mountainous terrain with many tributaries and a steep slope; runoff is intermittent and vulnerable to zero flow in the dry season. There are not many water conservancy projects in the basin and rice is the dominant crop grown there, followed by wheat.

TABLE 2							
The values of linearizing search in each step							
Cycle number J	$\theta_{_{1,j}}$	θ _{2, j}	F _j				
0	1.2427	49.4716	104 003.162				
1	1.3404	22.0497	21 896.339				
2	1.5067	16.0065	16 093.434				
3	1.7169	12.0299	10 299.993				
4	1.9406	10.0383	1 967.064				
5	1.9783	10.0294	258.892				
6	1.9916	10.0126	39.079				
7	1.9967	10.0051	6.1208				
8	1.9987	10.0021	0.971				
9	1.9995	10.0008	0.155				
10	1.9998	10.0003	0.025				
11	2.0000	10.0001	0.004				

TABLE 4							
The results of the error ideal model parameter calibrated by the linearized method							
Proportion of error	F ₀	θ_{1}	θ_{2}	D	F		
0.05	104 732.68	1.9999	9.9531	0.0469	36.7		
0.10	102 795.88	1.9956	10.053	0.0532	137.8		
0.15	105 853.17	2.0093	9.9851	0.0176	301.77		
0.20	104 049.51	2.0036	9.9513	0.0488	569.63		
0.25	104 492.94	2.0027	9.9525	0.0476	944.49		
0.30	106 575.46	2.0089	9.8865	0.1138	1 132.8		
0.35	104 433.01	1.9909	10.165	0.1652	1 579.4		
0.40	105 554.93	1.9833	10.182	0.1828	2 205.9		
0.45	103 573.51	2.0243	9.7111	0.2899	2 481.5		
0.50	110 931.04	2.0149	9.697	0.3034	2 990.5		
0.60	105 255.62	1.9983	10.086	0.0860	4 230.7		
Average	105 295.25	2.0029	9.9657	0.1232	1 510.1		

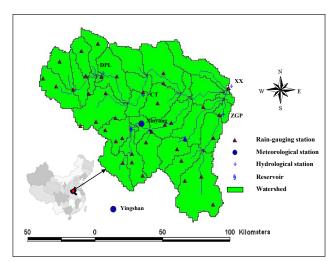


Figure 3 Location of the study watershed (DPL = Dapoling, CTG = Changtaiguan, ZGP = Zhuganpu, XX = Xixian)

	TABLE 3						
	The results of the no error ideal model parameter calibrated by the linearized method						
$ heta_{_{1,0}}$	$\theta_{_{2,0}}$	F ₀	$\theta_{_1}$	θ_{2}	F	J	
3.4177	24.3693	7.0×10^{11}	1.999832308	10.00088456	6.9 x 10 ⁻³	28	
1.7915	17.4012	9.0 x 10 ⁴	1.999728607	10.00317219	4.4 x 10 ⁻³	12	
3.0856	21.2353	$1.8 \ge 10^{10}$	1.999762626	10.00167424	8.3 x 10 ⁻³	24	
1.9876	17.6943	7.6 x 10 ⁵	1.999751796	10.00313901	4.4 x 10 ⁻³	14	
2.7204	6.7910	3.1 x 10 ⁵	2.000286033	9.997430342	6.7 x 10 ⁻³	13	
1.5509	19.6627	$1.8 \ge 10^4$	1.999782815	10.00131889	9.3 x 10 ⁻³	11	
4.2809	28.0269	2.9 x 10 ¹⁵	1.999841901	10.00051831	9.9 x 10 ⁻³	35	
0.9825	8.8556	$7.0 \ge 10^4$	1.999798566	10.00125761	7.6 x 10 ⁻³	13	
2.9236	9.0509	1.5 x 10 ⁷	2.000027461	10.00087218	7.8 x 10 ⁻³	17	
1.7553	29.0297	1.4 x 10 ⁶	1.999787343	10.00176653	4.5 x 10 ⁻³	15	
2.5507	9.7885	1.6 x 10 ⁶	2.000032618	10.00069936	6.4 x 10 ⁻³	15	
2.4268	26.9291	3.0 x 10 ⁸	1.999773288	10.00160568	7.5 x 10 ⁻³	20	
4.1490	19.0226	6.6 x 10 ¹³	1.999890526	10.00056829	3.0 x 10 ⁻³	33	
1.9032	6.1336	5.0 x 10 ⁴	1.999809336	10.00160602	3.5 x 10 ⁻³	12	
2.5548	8.0481	3.3 x 10 ⁵	2.000108207	9.999409939	2.7 x 10 ⁻³	14	
2.6308	24.5795	9.9 x 10 ⁸	1.999735991	10.00190456	9.8 x 10 ⁻³	21	
3.9853	20.5735	2.9 x 10 ¹³	1.999864084	10.00072722	4.4 x 10 ⁻³	32	
2.4599	30.8056	7.2 x 10 ⁸	1.999802061	10.00118197	7.9 x 10 ⁻³	21	
4.3424	19.9380	5.0 x 10 ¹⁴	1.999905642	10.00037873	3.0 x 10 ⁻³	35	
3.9952	18.9385	1.7 x 10 ¹³	1.999818607	10.00101991	7.4 x 10 ⁻³	31	
2.6486	18.8275	2.5 x 10 ⁸	1.999765416	10.00234184	3.5 x 10 ⁻³	20	
3.5827	21.0177	1.1 x 10 ¹²	1.999888487	10.00063858	2.7 x 10 ⁻³	29	
3.6108	15.0113	1.2 x 10 ¹¹	1.999755857	10.00230066	4.3 x 10 ⁻³	26	
3.2511	21.4369	7.8 x 10 ¹⁰	1.999842998	10.00103336	4.2 x 10 ⁻³	26	
0.8960	27.3718	$5.3 \ge 10^4$	2.000038402	9.99825104	9.9 x 10 ⁻³	11	
2.5298	25.1485	4.8 x 10 ⁸	1.999835158	10.00121636	3.6 x 10 ⁻³	21	
4.2369	25.2893	1.1 x 10 ¹⁵	1.999860661	10.00054026	6.7 x 10 ⁻³	35	
1.9197	17.7015	4.0 x 10 ⁵	1.999671436	10.00411829	7.5 x 10 ⁻³	13	
37.2895	1 804.38		7.2 x 10 ⁻⁷	7.8 x 10 ⁻⁵	5.7 x 10 ⁻³	20.6	

Xixian basin is situated in the transition zone between the northern subtropical region and the warm temperate zone. Rainfall for the flood season is mainly affected by monsoon; and the long-term mean annual precipitation is 1 145 mm (calculated by the data from 1954 to 2000). Half of the precipitation (~50%) falls between June and September. For further details of the watershed refer to Qu et al. (2012).

The Dapoling basin is serviced by 13 precipitation stages and one discharge stage (Dapoling stage). There is no evaporation stage in this catchment; because of this, the Nanwanevaporation stage data were used for the evaporation calculation. Series of observed data included daily precipitation of 13 rainfall stages, daily evaporation of Nanwan stage and daily discharge of Dapoling stage from 1966 to 2005. The total series were divided into 10-year periods for parameter calibration, 4 periods in all; that is, from 1966–1975, 1976–1985, 1986–1995 and 1996 to 2005. The initial parameter values are presented in Table 5.

Also, Table 5 shows the results of parameter calibration in different periods. Table 6 lists the performance of the calibrated

0%greatest relative error is 0.13% and all errors are less than 0.15%,
which means that the proposed method is reliable in terms of
water budgets. From Table 5, it can be seen that the change of
the ratio of potential evapotranspiration to pan evaporation, KC,
is highly significant, increasing from 0.658 to 0.922. A higher
ratio means higher evapotranspiration from the basin, which
causes a decrease in runoff. Additionally, the increase in free
water storage capacity, SM, reflects the decrease in surface runoff.
Although the sum of the outflow coefficient KG (free water
storage to the groundwater flow) and KI (free water storage to
the interflow) is kept constant, the single value of each parameter
has changed. KI changed towards a larger proportion (increase
of interflow), while KG decreased from 0.334 to 0.252 (Table 5),
causing higher peak flows and a quicker recession.

parameters. From Table 6, the calibrated parameters using

the linearized parameter calibration method can give good

simulation results: relative errors of 4 different periods from 1964

to 2005 are -0.1%, 0.05%, 0.13% and -0.05%, respectively. The

CONCLUSIONS

Streamflow at the outlet is the comprehensive response of the whole basin system to the input, precipitation, and thus, invariably, to the change of underlying surface in the basin. In appreciation of this fact, a number of studies have focused on using conceptual rainfall–runoff models to quantify the effect of land use and land cover change on runoff dynamics. However, because of the large number of parameters in the hydrological model, an effective automatic calibration method becomes important. Thus, a linearized parameter calibration method is presented for solving this problem. Unlike the previous study by Qu et al. (2012), we only use parameter comparison to assess the influence of land use/cover change on runoff of the basin. Compared to the method based on the error square sum, the linearized parameter calibration method has the following advantages:

- It can solve the theoretical problem of added unrelated local optimums
- The convergence of the method can be strictly proven in theory
- The determination of seeking direction and step length is objective and effective

In addition, this method was applied for parameter calibration of the Xinanjiang model in the upper Huaihe River Basin to evaluate the effect of land use change on hydrological processes through the change of parameters, for example, *KC*, *SM*, *KI* and *KG*. An increase in *KC* means higher evapotranspiration in the upper Huaihe River Basin, which causes a decrease in runoff. An increase in *SM* means a decrease in surface runoff. The increase in *KI* and decrease in *KG* in Dapoling basin reveal that the present land use pattern causes high peak flows and quicker recession. Because of limited data availability, only the daily model was applied for parameter calibration and evaporation, and runoff generation parameters were compared, while water source separation and concentration parameters were not.

The linearized parameter calibration method is a new calibration method. Future studies will involve tested the method in more basins and using more hydrological model parameter calibrations.

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APPENDIX 1

The validity of parameter search is justified as follow. From the text we can get the sum of squared errors F_j corresponding to the above step parameter vector θ_i .

$$F_{j} = (f - f_{j})^{T} (f - f_{j})$$
(A-1)

TABLE 5 Comparison of parameters calibrated for different periods						
Parameter	Initial value	1966–1975	1976–1985	1986–1995	1996-2005	
KC	0.741	0.658	0.675	0.710	0.922	
WUM	20	20	20	20	20	
WLM	80	80	80	80	80	
С	0.16	0.16	0.16	0.16	0.16	
WM	150	150	150	150	150	
В	0.3	0.3	0.3	0.3	0.3	
SM	16.813	5.431	10.809	15.849	19.94	
EX	1.5	1.5	1.5	1.5	1.5	
KG	0.398	0.334	0.294	0.294	0.252	
KI	0.363	0.366	0.406	0.406	0.448	
CS	0.45	0.45	0.45	0.45	0.45	
CI	0.88	0.88	0.88	0.88	0.88	
CG	0.995	0.995	0.995	0.995	0.995	
KE	1	1	1	1	1	
XE	0.38	0.38	0.38	0.38	0.38	

TABLE 6 Simulation results from different periods					
Period	R _{obs} (mm)	R _{sim} (mm)	ER (%)	DC _{ave} *	
1966-1975	4 877	4 882	-0.1	0.46	
1976-1985	3 847	3 845	0.05	0.51	
1986-1995	2 979	2 975	0.13	0.395	
1996-2005	4 037	4 039	-0.05	0.470	

The sum of squared errors F_{i+1} can be expressed as:

$$F_{j+1} = (f - f_{j+1})^{T} (f - f_{j+1})$$
(A-2)

Substitute Eq. 9 into Eq. A-2:

$$F_{j+1} = (f - f_j - S(\theta_{j+1} - \theta_j))^T \cdot (f - f_j - S(\theta_{j+1} - \theta_j))$$
(A-3)

then:

$$\begin{split} F_{j+1} &= (f - f_j)^T \cdot (f - f_j) \\ &- (\theta_{j+1} - \theta_j)^T S^T (f - f_j) \\ &- (f - f_j)^T S(\theta_{j+1} - \theta_j) \\ &+ (\theta_{j+1} - \theta_j)^T S^T S(\theta_{j+1} - \theta_j) \end{split} \tag{A-4}$$

$$F_{j+1} = F_j - (\theta_{j+1} - \theta_j)^T S^T (f - f_j)$$

- $(f - f_j)^T S(\theta_{j+1} - \theta_j)$
+ $(\theta_{j+1} - \theta_j)^T S^T S(\theta_{j+1} - \theta_j)$ (A-5)

Substitute Eq. 12 into Eq. A-5:

$$\begin{split} F_{j+1} &= F_j - (S^T S)^{-1} S^T (f - f_j)^T S^T (f - f_j) \\ &- (f - f_j)^T S (S^T S)^{-1} S^T (f - f_j) \\ &+ (S^T S)^{-1} S^T (f - f_j)^T S^T (f - f_j) \\ &= F_j - (f - f_j)^T S (S^T S)^{-1} S^T (f - f_j) \\ &< F_j \end{split}$$
(A-6)

It can be observed that the new parameter vector θ_{j+1} of any step determined by Eq. 12 satisfying the relation:

$$F_{j+1} < F_j \tag{A-7}$$

According to Eq. A-7, every searching step will get less sumsquared. As the iterative process continues, the sum-squared error approaches the minimum value and the optimal parameter will be obtained. So the searching direction determined by Eq. 12 is correct.

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