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Multiobjective sensitivity analysis and optimization of a distributed hydrologic model MOBIDIC

J. Yang¹, F. Castelli², and Y. Chen¹

Received: 28 February 2014 - Accepted: 16 March 2014 - Published: 26 March 2014

Correspondence to: J. Yang (yangjing@ms.xjb.ac.cn)

Published by Copernicus Publications on behalf of the European Geosciences Union.

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¹State Key Laboratory of Desert and Oasis Ecology, Xinjiang Institute of Ecology and Geography, Chinese Academy of Sciences, Xinjiang, 830011, China ²Department of Civil and Environmental Engineering, University of Florence, Italy

Calibration of distributed hydrologic models usually involves how to deal with the large number of distributed parameters and optimization problems with multiple but often conflicting objectives which arise in a natural fashion. This study presents a multiobjective sensitivity and optimization approach to handle these problems for a distributed hydrologic model MOBIDIC, which combines two sensitivity analysis techniques (Morris method and State Dependent Parameter method) with a multiobjective optimization (MOO) approach ε -NSGAII. This approach was implemented to calibrate MOBIDIC with its application to the Davidson watershed, North Carolina with three objective functions, i.e., standardized root mean square error of logarithmic transformed discharge, water balance index, and mean absolute error of logarithmic transformed flow duration curve, and its results were compared with those with a single objective optimization (SOO) with the traditional Nelder-Mead Simplex algorithm used in MOBIDIC by taking the objective function as the Euclidean norm of these three objectives. Results show: (1) the two sensitivity analysis techniques are effective and efficient to determine the sensitive processes and insensitive parameters: surface runoff and evaporation are very sensitive processes to all three objective functions, while groundwater recession and soil hydraulic conductivity are not sensitive and were excluded in the optimization; (2) both MOO and SOO lead to acceptable simulations, e.g., for MOO, average Nash-Sutcliffe is 0.75 in the calibration period and 0.70 in the validation period; (3) evaporation and surface runoff shows similar importance to watershed water balance while the contribution of baseflow can be ignored; (4) compared to SOO which was dependent of initial starting location, MOO provides more insight on parameter sensitivity and conflicting characteristics of these objective functions. Multiobjective sensitivity analysis and optimization provides an alternative way for future MOBIDIC modelling.

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With the development of information technology (e.g., high performance computing cluster and remote sensing technology), there has been a prolific development of integrated, distributed and physically-based watershed models (e.g., MIKE-SHE, Refsgaard and Storm, 1995) over the past two decades, which are increasingly being used to support decisions about alternative management strategies in the areas of land use change, climate change, water allocation, and pollution control. Though in principle parameters of distributed and physically based models should be assessable from catchment data (in traditional conceptual rainfall—runoff models, parameters are obtained through a calibration process), these models still need a parameter calibration process in practice due to scaling problems, experimental constraints, etc. (Beven and Binley, 1992; Gupta et al., 1998; Madsen, 2003). Problems, arising in calibrating distributed hydrologic models, include how to handle large number of distributed parameters and optimization problems with multiple but often conflicting objectives.

In the literature, to deal with large number of distributed model parameters, this is often done by aggregating distributed parameters (e.g., Yang et al., 2007), or screening out the unimportant parameters through a sensitivity analysis (e.g., Muleta and Nicklow, 2005; Yang, 2011). Sensitivity analysis can be used to not only screen out the most insensitive parameters, but also study the system behaviors identified by parameters and their interactions, qualitatively or quantitatively. However, most of applications in environmental modelling are based on the one-at-a-time (OAT) local sensitivity analysis, which is "predicated on assumptions of model linearity which appear unjustified in the cases reviewed" (Saltelli and Annoni, 2010), or simple linear regressions where a lot of uncertainty are not fairly accounted for. The use of global sensitivity analysis techniques is very crucial in distributed modelling. Only recently, global sensitivity analysis techniques started to appear in hydrologic modelling.

Although most hydrologic applications are based on the single objective calibration, model calibration with multiple and often conflicting objectives arises in a natural

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fashion in hydrologic modelling. This is not only due to the increasing availability of multi-variable (e.g., flow, groundwater level, etc.) or multi-site measurements, but also due to the intrinsic different system responses (e.g., peaks and baseflow in the flow series). Instead of finding a single optimal solution in the single objective optimization (SOO), the task in the multiobjective optimization (MOO) is to identify a set of optimal trade-off solutions (called a Pareto set) between conflicting objectives. In hydrology, the traditional method to solve multiobjective problems is to form a single objective, e.g., by giving different weights to these multiple objectives or applying some transfer function. Over the past decade, several MOO algorithms approaches have been applied to the conceptual rainfall-runoff models (e.g., Yapo et al., 1998; Gupta et al., 1998; Madsen, 2000; Boyle et al., 2000; Vrugt et al., 2003; Liu and Sun, 2010), and now increasing applied to distributed hydrologic models (e.g., Madsen, 2003; Bekele and Nicklow, 2007; Shafii and Smedt, 2009; MacLean et al., 2010). And there are some papers (Tang et al., 2006; Wöhling et al., 2008) to comparatively study their strengths with the application in hydrology. It is worth noting that the multiobjective calibration is different from statistical uncertainty analysis which is based on the concept (or similar concept) of "equifinality" (see discussion in Gupta et al., 1998; Boyle et al., 2000).

This paper applies two sensitive analysis techniques (Morris method and State Dependent Parameter method) and ε -NSGAII in the multiobjective sensitive analysis and calibration framework. This was implemented to calibrate a distributed hydrological model MOBIDIC with its application to the Davidson watershed, North Carolina. The purpose is to study parameter sensitivity of the hydrologic model MOBIDIC and explore the capability of MOO in calibrating the MOBIDIC compared to the traditional SOO used in MOBIDIC applications.

This paper is structured as follows: Sect. 2 gives a description of the MOBIDIC model; Sect. 3 introduces the approach in the multiobjective sensitivity analysis and optimization; Sect. 4 gives a brief introduction of the study site, model setup, objective selection, and sensitivity and calibration procedure; in Sect. 5, the results are presented

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2 Hydrologic model MOBIDIC

MOBIDIC (MOdello di Bilancio Idrologico Distribuito e Continuo; Castelli et al., 2009; Campo et al., 2006) is a distributed and raster-based hydrological balance model. MOBIDIC simulates the energy and water balances on a cell basis within the watershed. Figure 1 gives a schematic representation of MOBIDIC. The energy balance is approached by solving the heat diffusion equations in multiple layers in the soil–vegetation system, while the water balance is simulated in a series of reservoirs (i.e., boxes in Fig. 1) and fluxes between them.

For each cell, water in the soil is simulated by

$$\frac{dW_{g}}{dt} = I_{nf} - S_{per} - Q_{d} - S_{as}$$

$$\frac{dW_{c}}{dt} = S_{as} - E_{t}$$
(1)

where $W_{\rm g}$ [L] and $W_{\rm c}$ [L] are the water contents in the soil gravitational storage and capillary storage, respectively, and $I_{\rm nf}$ [LT⁻¹], $S_{\rm per}$ [LT⁻¹], $Q_{\rm d}$ [LT⁻¹], $E_{\rm t}$ [LT⁻¹], and $S_{\rm as}$ [LT⁻¹] are infiltration, percolation, interflow, evaporation, and adsorption from gravitational to capillary storage, which are modeled through following equations:

$$S_{\text{per}} = \gamma \cdot W_{\text{g}}$$

$$Q_{\text{d}} = \beta \cdot W_{\text{g}}$$

$$S_{\text{as}} = \kappa \cdot (1 - W_{\text{c}}/W_{\text{cmax}})$$

$$I_{\text{nf}} = \begin{cases} \left[P + (Q_{\text{d}} + Q_{\text{h}} + R_{\text{d}})_{\text{up}} \right] \left[1 - \exp\left(\frac{-K_{\text{s}}}{P + (Q_{\text{d}} + Q_{\text{h}} + R_{\text{d}})_{\text{up}}}\right) \right] & \text{if } W_{\text{g}} < W_{\text{gmax}} \\ 0 & \text{otherwise} \end{cases}$$

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where γ , β and κ are percolation coefficient [T⁻¹], interflow coefficient [T⁻¹], and soil adsorption coefficient [LT⁻¹], respectively, P the precipitation [LT⁻¹], Q_h and R_d Horton runoff and Dunne runoff, K_s the soil hydraulic conductivity [LT⁻¹], W_{gmax} [L] and W_{cmax} [L] the gravitational and capillary storage capacities.

Once the surface runoff ($Q_{\rm h}$ and $R_{\rm d}$) and baseflow are calculated, three different methods can be used for river routing, i.e., the lag method, the linear reservoir method, Muskingum–Cunge method (Cunge, 1969). Muskingum–Cunge method was used in this study.

MOBIDIC uses either a linear reservoir or the Dupuit approximation to simulate the groundwater balance which relates the groundwater change to the percolation, water loss in aquifers and baseflow. In this case study, the linear reservoir method was used.

Although there are many distributed parameters in MOBIDIC, normally these distributed parameters are calibrated through the "aggregate" factors (e.g., the multiplier for hydraulic conductivity) based on their initial estimations. And hereafter we use the term "factor" (instead of "model parameter") when we conduct the sensitivity analysis and optimization, to avoid the confusion with the term "model parameter" used in model description. A factor can be a model parameter or a group of model parameters, and in this paper it is a change to be applied to a group of model parameters. In MOBIDIC, normally nine factors (i.e., nine groups of parameters) need to be calibrated. These factors, their explanations, and their corresponding model parameters are listed in Table 1.

3 Methodology

The procedure applied here consists of two-step analyses, i.e., a multiobjective sensitivity analysis generally characterizing the basic hydrologic processes and single out the most insensitive parameters, and a multiobjective calibration aiming at trade-offs between different objective functions.

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Sensitivity analysis is to assess how variations in model out can be apportioned, qualitatively or quantitatively, to different sources of variations, and how the given model depends upon the information fed into it (Saltelli et al., 2008). In the literature, a lot of sensitivity analysis methods are introduced and applied, e.g., Yang (2011) applied and compared five different sensitivity analysis methods. Here we adopted an approach which combines two global sensitivity analysis techniques, i.e., the Morris method (Morris, 1991) and SDP method (Ratto et al., 2007).

3.1.1 Morris method

Morris method is based on replicated and randomized one-factor-at-a-time design (Morris, 1991). For each factor X_i , Morris method uses two statistics, μ_i and σ_i , which measure the degree of factor sensitivity, and the degree of nonlinearity or factor interaction, respectively. The higher μ_i is, the more important the factor X_i is to the model output; and the higher σ_i is, the more nonlinear the factor X_i is to the model output or more interactions with other factors (details refer to Morris, 1991; Campolongo et al., 2007). Morris method takes $m \cdot (n+1)$ model runs to estimate these two sensitivity indices for each of n factors with sample size m. The advantage is it is efficient and effective to screen out insensitive factors. Normally m takes values around 50. And according to Saltelli et al. (2008), the sensitivity measure (μ_i) is a good proxy for the total effect (i.e., S_{T_i} in Eq. (4) below), which is a robust measure in sensitivity analysis.

3.1.2 State-Dependent Parameter method (SDP)

SDP (Ratto et al., 2007) is based on the ANOVA functional decomposition, which apportions the model output uncertainty (100%, as 1 in Eq. 3) to factors and different

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$$1 = \sum_{i} S_{i} + \sum_{i} \sum_{j>i} S_{ij} + \ldots + S_{12...n}$$
 (3)

where S_i is the main effect of factor X_i representing the average output variance reduction that can be achieved when factor X_i is fixed, and S_{ij} is the first-order interaction between X_i and X_j , and so on. In ANOVA based sensitivity analysis, total effect (S_{T_i}) is frequently used, which stands for the average output variance that would remain as long as X_i stays unknown,

$$S_{Ti} = S_i + \sum_{j \neq i} S_{ij} + \ldots + S_{12...n}$$
 (4)

SDP method uses the emulation technique to approximate lower order sensitivity indices in Eq. (3) (e.g., S_i and S_{ij} in this study) by ignoring the higher order sensitivity indices. And we define $S_{Di} = S_i + \sum_j S_{ij}$ (referred to as "quasi total effect" later) as a sur-

rogate to the total effect. The advantage is that it can precisely estimate lower order sensitivity indices at a lower computational cost (normally 500 model runs, which is independent of number of factors). The disadvantage is that it cannot estimate higher order sensitivity indices.

Since these two methods are computationally efficient, these two methods are applied individually. And then, the sensitivity of each factor and its system behaviour will be discussed, qualitatively by Morris method, and quantitatively by SDP method. And then the most insensitive factors will be screened out and excluded in the calibration.

In the context of multiobjective analysis, sensitivity analysis includes: (1) to examine the sensitivity of each factor to different objective functions, qualitatively or quantitatively; (2) to single out the most sensitive factors and study the physical behaviours of the system; (3) to exclude the most insensitive factors and therefore simplify the process of calibration.

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In the literature of hydrologic modelling, most applications are single objective based, which aims at a single optimal solution. However, for example in flow calibration, there is always a case that two solutions, one solution better simulates the peaks and poorly simulates the baseflow while the other solution poorly simulates the peaks while better simulates the baseflow. These two solutions, which are called Pareto solutions, are incommensurable, i.e., better fitting of the peaks will lead to worse fitting of the baseflow, and vice versa. This belongs to the domain of MOO, aiming at finding a set of optimal solutions (Pareto solutions), instead of one single solution.

Generally a MOO problem can be formulated as follows:

$$\min F(X) = (f_1(X), f_2(X), ..., f_i(X), ..., f_k(X))$$
s.t. $G(X) = (g_1(X), g_2(X), ..., g_i(X), ..., g_i(X))$ (5)

Where X is an n-dimensional vector and in this study represents the model factors to be calibrated, $f_i(X)$ ith objective function, and $g_i(X)$ ith constraint function.

In the literature, there are many algorithms available to obtain the Pareto solutions, e.g., NSGAII (Non-dominated Sorting Genetic Algorithm-II; Deb et al., 2002), SPEA2 (Strength Pareto Evolutionary Algorithm 2; Zitzler et al., 2001), MOSCEM-UA (Multi-objective Shuffled Complex Evolution Metropolis; Vrugt et al., 2003), and ε -NSGAII (Kollat and Reed, 2006), etc. In this study, we adopt ε -NSGAII, which is efficiency, reliability, and ease-of-use. Its strengths have been comparatively studied in Kollot and Read (2006) and Tang et al. (2006).

arepsilon-NSGAII is an extension of the NSGAII (Deb et al., 2002), a second generation of multiobjective evolution algorithm. The main characteristics of arepsilon-NSGAII include: (i) selection, crossover, and mutation processes as other genetic algorithm by mimicking the process of natural evolution, (ii) an efficient non-domination sorting scheme, (iii) an elitist selection method that greatly aids in capturing Pareto front, (iv) arepsilon-dominance archiving, (v) adaptive population sizing, and (vi) automatic termination to minimize the

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As a comparison, a single objective function is defined as 2-norm of the multiple objectives F(X), which measures how close to the original point (theoretical optimum O):

$$sof = ||F(X)||_2 = \sqrt{\sum_{j=1}^k f_j(X)^2}$$
 (6)

And SOO was done with the classic Nelder–Mead algorithm (Nelder and Mead, 1965) which is widely used in MOBIDIC applications.

To analyze the Pareto solution and also compare with the solution from SOO, except for traditional methods, the "Level diagrams" proposed by Blasco et al. (2008) was also used. Compared to traditional methods, it can visualize high dimensional Pareto front and synchronizes the objective and factor diagrams. The procedure and includes two steps. In the first step, the vector of objectives (*k*-dimension) for each Pareto point is mapped to a real number (one-dimension) according to the proximity to the theoretical optimum measured with a specific norm of objectives; and in the second step, these norm values are plotted against the corresponding values of each objective or factor. 1-norm, 2-norm and ∞-norm are suggested. To compare with SOO, 2-norm was used.

4 Davidson watershed and objective selection

4.1 Davidson watershed

The Davidson watershed, located in southwest mountain area of North Carolina, drains an area of 105 km² above the station "Davidson river near Brevard" (see Fig. 2). The

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elevation ranges from 645 m to 1820 m a.s.l. Based on the NLDAS climate data, the average annual precipitation is 1900 mm and varies from 1400 mm to 2500 mm, and daily temperature changes from -19 °C to 26 °C. The average daily flow is about $3.68 \, \text{m}^3 \, \text{s}^{-1}$.

Data used in MOBIDIC model include (i) Digital Elevation Model (DEM), (ii) soil data, (iii) land cover data, (iv) climate data (precipitation, minimum and maximum temperature, solar radiation, humidity and wind speed), and (v) flow data. 9 m DEM, land cover, SSURGO soil data, one station (Davidson river near Brevard) of flow data are from US Geological Survey, and hourly NLDAS climate data from National Aeronautics and Space Administration (NASA). NLDAS integrates a large quantity of observation-based and model reanalysis data to drive offline (not coupled to the atmosphere) land-surface models (LSMs), and executes at 1/8th-degree grid spacing over central North America, enabled by the Land Information System (LIS) (Kumar et al., 2006; Peters-Lidard et al., 2007).

DEM is used to delineate the watershed and estimate the topographic parameters and river system, Land cover for evaporation parameters, soil data for soil parameters, climate data is used to drive MOBIDIC, and flow data are used to calibrate the model and assess model performance. The climate and flow data used in this study are from 1 January 1996 to 30 September 2006. As NLDAS only has hourly temperature daily instead of hourly minimum and maximum temperature needed by MOBIDIC, we compiled the hourly climate data to daily data and run the model at a daily step. After MOBIDIC setup, the initial parameter values are listed in third column of Table 1.

We split the data into a warm-up period (from 1 January 1996 to 30 September 2000), a calibration period (from 1 October 2000 to 30 September 2003), and a validation period (from 1 October 2003 to 30 September 2006).

4.2 Objective function selection

After setting up MOBIDIC in the Davidson watershed, three objective functions were used in the multiobjective sensitivity analysis and optimization:

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SRMSE =
$$\frac{\sqrt{\frac{1}{N} \sum_{j=1}^{N} \left(\log \left(Q_{j}^{\text{obs}} \right) - \log \left(Q_{j}^{\text{sim}} \right) \right)^{2}}}{\sqrt{\frac{1}{N-1} \sum_{j=1}^{N} \left(\log \left(Q_{j}^{\text{obs}} \right) - \overline{\log Q} \right)^{2}}}$$
 (7)

(2) Water Balance Index, calculated as the mean absolute error between the simulated and observed flow accumulation curves:

$$WBI = \frac{1}{N} \sum_{i=1}^{N} \left| Q_{Ci}^{obs} - Q_{Ci}^{sim} \right|$$
 (8)

(3) Mean absolute error between the logarithms of simulated and observed flow duration curves

$$MARD = \frac{1}{100} \sum_{i=1}^{N} \left| \log \left(Q_{P_i}^{obs} \right) - \log \left(Q_{P_i}^{sim} \right) \right|$$
(9)

In Eqs. (7)–(9), Q_i^{obs} and Q_i^{sim} are observed and simulated flow series at time step i, N the data length, $\overline{\log Q}$ the average of logarithmic transformed observed flows, Q_{Ci}^{obs} and $Q_{\mathrm{C}i}^{\mathrm{sim}}$ ith observed and simulated accumulated flows, and $Q_{\mathrm{P}i}^{\mathrm{obs}}$ and $Q_{\mathrm{P}i}^{\mathrm{sim}}$ ith percentiles of observed and simulated flow duration curves.

SRMSE (Eq. 7), WBI (Eq. 8), and MARD (Eq. 9) are measures of the closeness between simulated and observed flow series, water balance, and closeness between simulated and observed flow frequencies, respectively. The smaller these measures are, the better the simulation is, and the minima are (0, 0, 0) meaning a perfect match

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And accordingly, the single objective function here is the Euclidean norm (2-norm) of SRMSE, WBI, and MARD:

$$sof = \sqrt{SRMSE^2 + WBI^2 + MARD^2}$$
 (10)

5 Result and discussion

5.1 Multiobjective sensitivity analysis

Morris method and SDP method were applied individually to the initially selected factors (in Table 1).

For Morris method, its convergences for three objective functions, monitored using the method proposed in Yang (2011), were achieved around $700 \sim 800$ model simulations. Figure 3 gives the sensitivity results for objective functions SRMSE, WBI, and MARD, respectively. In each plot, the horizontal axis (μ) denotes the degree of factor sensitivity, and the vertical axis (σ) denotes the degree of factor nonlinearity or interaction with other factors.

For SRMSE, the most sensitive parameters are group $(p\alpha,p\gamma,$ and $p\kappa)$, followed by $p\beta$ and rCH, while other parameters (especially rK_s and rK_f) are not so sensitive. This applies to the degree of the factor nonlinearity or interaction. The sensitivities of $p\alpha, p\gamma$, and $p\kappa$ indicate the importance of their corresponding processes (i.e., surface runoff, percolation, and adsorption which is related to evapotranspiration) to SRMSE, while interflow $(p\beta)$ is less important and other processes/characteristics (e.g., groundwater flow, rK_f) are not important.

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For WBI, the dominating parameter is $p\kappa$, followed by group ($p\alpha$, $p\gamma$, $p\beta$ and rCH), while other parameters (especially rK_f and rW_{cmax}) are not so sensitive. WBI measures the water balance between observed and simulated flow series, and it is reasonable that $p\kappa$ which controls the water supply for evaporation is most sensitive while other factors ($p\alpha$, $p\gamma$, $p\beta$ and rCH) are sensitive mainly through interaction with this factor, as indicated by the high " σ "s of these factors.

For MARD, the results are nearly the same to SRMSE.

Figure 4 gives the sensitivity results based on SDP method for SRMSE, WBI, and MARD, from top to bottom. In each plot, the grey and black bars are S_i and S_{Di} for each factor.

For SRMSE, as indicated by R^2 in the legend, main effects (S_i) contribute to 58.7% of SRMSE uncertainty, and quasi total effects (S_{Di}) account for 83% of SRMSE uncertainty which is quite high, while other 17% due to higher interactions are not explained. Based on S_{Di} (black bar), the most sensitive factors are $p\gamma$ and $p\kappa$, followed by $p\alpha$ and rCH, and then $p\beta$ and rW_{cmax} while other factors are not sensitive. This result quantitatively corroborates the result obtain from Morris method. The main effects (S_i) of ($p\gamma$, $p\kappa$, and $p\alpha$) are high (i.e., 0.17, 0.18 and 0.14), which suggests these factors should be determined first in model calibration as they lead to the largest reduction in SRMSE uncertainty. For each factor, the difference between the black bar and grey bar shows the first order interaction with other factors. This interaction is very strong in $p\gamma$, $p\kappa$, $p\alpha$, and rCH, and very weak in other factors.

For WBI, as indicated by R^2 in the legend, the total main effects (S_i) contribute to 38.4% of the WBI uncertainty, quasi total effects (S_{Di}) only account for 57.6% of WBI uncertainty, and 40% due to higher interactions are not explained and can not be ignored. However, by analyzing the result with that from Morris method (top right in Fig. 3), we still can get some valuable results: the dominating sensitive factor is $p\kappa$ with S_{Di} 0.43 (which is same as Morris method), followed by $p\gamma$, $p\alpha$, and rCH, while other factors are not sensitive; the main effect of $p\kappa$ is as high as 0.27, and it should be fixed

Similar to the Morris results for SRMSE and MARD, the result of MARD is nearly the same as SRMSE. The similar result for SRMSE and MARD shows similar characteristic relationship between factors and the objective function. This is explainable: a good simulation measured by SRMSE will more likely result in a good measure of MARD, and vice versa.

As aforementioned, in the context of multiobjective sensitivity analysis, sensitivity analysis is to exclude the factors which are insensitive to all the objective functions considered. Based on the analysis above, four most insensitive factors are rKs, rKf, rW_{cmax},and rW_{gmax}. However, as shown in Fig. 4, rW_{cmax} is more sensitive than other three factors, and for objective function WBI, as higher order interactions are strong based on SDP (i.e., explains around 40% of model uncertainty), and evaporation is the most sensitive process to water balance (as indicated by $p\kappa$ and rCH) and rW_{cmax} is the only factor related to evaporation storage (W_c), therefore, we only exclude rK_f, rK_s, and rW_{gmax} for calibration.

5.2 Multiobjective optimization

After sensitivity analysis, only six factors were involved in the calibration. MOO converged with 482 Pareto front points after totally 22 000 model runs with the parallelized ε -NSGAII, while SOO converged after 686 model runs with the classic Nelder–Mead algorithm. Apparently, ε -NSGAII took more model simulations than the Nelder–Mead algorithm, but simulation time was compensated by the parallelized code running on high performance clusters.

Figure 5 shows optimized non-dominant sets normalized within [0, 1] and the black line is for the factor set with SOO. It is encouraging that except rW_{cmax} , factor ranges decreased a lot. This corroborates the conclusion in the sensitivity analysis: $p\gamma$, $p\kappa$, $p\beta$, $p\alpha$, and rC_H are the most sensitive and identifiable factors to these three objective functions, while rW_{cmax} is less sensitive and less identifiable. Several scattered

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values of py and dispersed rW_{omax} show that optimized factor sets are scattered in the response surface rather than concentrated in a continuous region. And the factor set with SOO is within the range of non-dominant sets.

Figure 6 shows Pareto solutions scattered in the three-dimensional space (top left), 5 and projections in two-dimensional subspaces with corresponding correlation coefficients (r) in the calibration period, with the black dot in each plot denoting the solution for SOO. Correlation coefficients are high and negative for SRMSE and WBI (-0.54), and WBI and MARD (-0.74), and this indicates strong trade-off interactions along the Pareto surface, i.e., better (lower) WBI will eventually result in worse (higher) SRMSE, and vice versa. The correlation coefficient is low (0.13) between SRMSE and MARD, and it is close to 0 when these two variables approach to their minima regions (i.e., SRMSE < 0.53 and MARD < 0.09). Table 2 lists the statistics of these three objectives associated with Pareto sets and the result of SOO. For Pareto sets, in the calibration period, the average SRMSE is 0.49 ranging from 0.47 to 0.57, which corresponds to the average NS 0.78 ranging from 0.67 to 0.78; the average WBI is 0.05 ranging from 0.02 to 0.11; and the average MEAD is 0.08 ranging from 0.03 to 0.11. In the validation period, the average SRMSE is 0.54 ranging from 0.51 to 0.62, which corresponds to the average NS 0.70 ranging from 0.61 to 0.74; the average WBI is 0.05 ranging from 0.04 to 0.09; and the average MEAD is 0.10 ranging from 0.08 to 0.13. And for SOO, SRMSE, WBI and MEAD are 0.48, 0.06 and 0.07 for the calibration period, and 0.57, 0.06 and 0.10 for the validation, and accordingly the "NS"s are 0.77 and 0.67, respectively. According to Moriasi et al. (2007) which suggests NS > 0.75 and WBI < 10 % as excellent modelling of river discharge, all Pareto solutions with MOO and the solution with SOO are close to "excellent" for both calibration and validation periods.

To better visualize Pareto sets and compare with the result of SOO, the level diagrams are plotted in Fig. 7 by applying Euclidean norm (2-norm) to evaluate the distance of each Pareto point to the ideal origin (0.0.0) (ideal values for all three normalized objectives are 0). In Fig. 7, top three plots are for three objectives and the rest for optimized factors, and the black dot in each plot is the solution for SOO. In the level

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diagrams, each objective and each factor of a point (corresponding to a Pareto solution) is represented with the same 2-norm value for all the plots. Compared with MOO, obviously, SOO was trapped in the local optima as seen in top-left plot, which means optimization with Nelder–Mead algorithm was dependent of starting point. The 2-norm has a close linear relationship with SRMSE due to values of SRMSE are 5 to 10 times of other two objective functions, and it does not have such relationship with other two objectives. The scattering of objectives and factors makes it difficult in decision making to select a single solution because there is not a clear trade-off solution (Blasco et al., 2008). However, compared to SOO, the Pareto solutions from MOO can make decision making easy as it can be converted with expert opinion or some utility function.

Figures 8 and 9 show simulated and observed flow duration curves and time series flows, respectively, with grey lines denoting the simulations with MOO and black lines with SOO. Generally, all simulations match the observation well for both the duration curve and time series flow for both calibration period and validation period. For the duration curve, simulations from MOO show a wide range in the low flows with frequencies from 0.85 to 1.0, which reflects the insensitivity of groundwater process (discussed in the sensitivity analysis, i.e., rK_f is insensitive to these three objectives). Except for this, there is a slight overestimation of flows, large flows during the calibration period with frequencies from 0.2 to 0.1, and median to large flows during the validation period with frequencies from 0.5 to 0.1. This might be due to the uncertainty in the reanalyzed climate data. And the extreme flow with frequency around 0 is underestimated, and this is because we chose the logarithm scale of the observed and simulated flows instead of normal scale when computing objectives SRMSE and MARD. With SOO, the deviation from the observed is larger. Similar conclusions can be drawn from the time series simulations in Fig. 9, i.e., the wide ranges of low flow period, and underestimation of flow peaks. Other than this, generally all simulations can mimic the observations.

Figure 10 shows the time series of watershed average storages (soil storage expressed as soil saturation, and groundwater depth), and fluxes (evaporation, surface runoff and baseflow) associated with MOO (shaded) and SOO (black line). With MOO,

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soil saturation varies from 0.2 to 1.0 and groundwater from 0 to 120 mm. The temporal fluctuation of soil moisture is higher than groundwater, but lower than fluxes in evaporation and surface runoff. And this is true for the solution with SOO except the its ranges of soil saturation and groundwater (groundwater is very close to 0 mm). For fluxes with MOO, evaporation and surface runoff have more temporal variation than baseflow, and their magnitudes are larger than baseflow. This applies to fluxes with SOO, and its baseflow is close to 0. This can be confirmed by the De Finetti diagram in Fig. 11: with MOO, the average contributions of evaporation, surface runoff, and basedflow are 49.3, 46.1, and 4.8 %, respectively while the contribution of baseflow is very insignificant. And the contribution of baseflow is almost 0 with SOO.

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This study presents a multiobjective sensitivity and optimization approach to calibrate a distributed hydrologic model MOBIDIC with its application in the Davidson watershed for three objective functions (i.e., SRMSE, WBI, and MAED). Results show:

- 1. The two sensitivity analysis techniques are effective and efficient to determine the sensitive processes and insensitive parameters: surface runoff and evaporation are very sensitive processes to all three objective functions, while groundwater recession and soil hydraulic conductivity are not sensitive and were excluded in the optimization.
- 2. For SRMSE and MAED, though all the factors have almost same sensitivities, the non-dominance of Pareto set shows the trade-off between these two objectives in the response surface.
- 3. Both MOO and SOO achieved acceptable results for both calibration period and validation period, in terms of objective functions and visual match between simulated and observed flows and flow duration curves. For example, with MOO, the

- average NS is 0.75 ranging from 0.67 to 0.78 in the calibration and 0.70 ranging from 0.61 to 0.74 in the validation period.
- 4. In the case study, evaporation and surface runoff shows similar importance to watershed water balance while the contribution of baseflow can be ignored.
- 5. Compared to MOO with ε -NSGAII, the application of SOO with the Neld–Mead algorithm was dependent of initial starting point. Furthermore, the Pareto solution provides a better understanding of these conflicting objectives and relations between objectives and parameters, and a better way in decision making.
- Acknowledgements. The research was supported by the "Thousand Youth Talents" Plan (Xinjiang Project) and the National Basic Research Program of China (973 Program: 2010CB951003). The data used in this study were acquired as part of the mission of NASA's Earth Science Division and archived and distributed by the Goddard Earth Sciences (GES) Data and Information Services Center (DISC).

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Table 1. Initial selected factors, initial estimation of the corresponding MOBIDIC parameter, and factor ranges.

Factor	Meaning of the given factor	Initial estimation of MOBIDIC parameter	Factor range
ργ	Exponential change ^a on soil percolation coefficient γ [s ⁻¹]	1.2 × 10 ⁻¹¹	[-2, 9]
рк	Exponential change on soil adsorption coefficient κ [s ⁻¹]	1.6×10^{-7}	[-6, 5]
ρβ	Exponential change on interflow coefficient β [s ⁻¹]	2.5×10^{-6}	[-7, 4]
ρα	Exponential change on surface storage decay coefficient α [s ⁻¹]	3.3×10^{-7}	[-6, 5]
rK_s	Multiplying change ^b on soil hydraulic conductivity [ms ⁻¹]	$[5.0 \times 10^{-6}, 8.9 \times 10^{-5}]$	[0.001, 100]
$\mathrm{rW}_{\mathrm{cmax}}$	Multiplying change on maximum storage of the capillary reservoir [m]	[0.017, 0.165]	[0.01, 5]
${\rm rW}_{\rm gmax}$	Multiplying change on maximum storage of the gravitational reservoir [m]	[0.107, 0.449]	[0.01, 5]
rC _H	Multiplying change on bulk turbulent exchange coefficient for heat [–]	[0.010,0.018]	[0.01, 5]
rK _f	Multiplying change on groundwater decay coefficient [s ⁻¹]	1.0 × 10 ⁻⁷	[0.001, 5]

^a Exponential change pX means the corresponding MOBIDIC parameter X will be changed according to $X = X_0 \times \exp(pX - 1)$, where X_0 is the initial estimation of X.

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^b Multiplying change rX means the corresponding MOBIDIC parameter X will be changed according to $X = X_0 \times rX$.

Table 2. Statistics of three objective functions associated with multiobjective optimization and single objective optimization.

	Multiob Calibra	-	optimiz	ation Validation		Single object	ctive optimization Validation	
	Mean	Min	Max	Mean	Min	Max		
SRMSE	0.49	0.47	0.57	0.54	0.51	0.62	0.48	0.57
WBI	0.05	0.02	0.11	0.05	0.04	0.09	0.07	0.06
MAED	0.08	0.03	0.11	0.10	0.08	0.13	0.07	0.10

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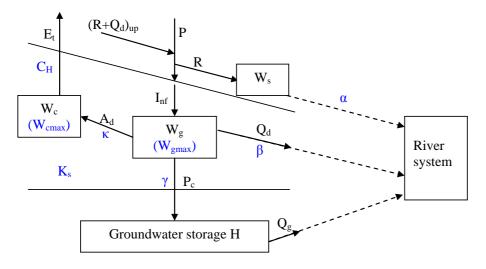


Fig. 1. A schematic representation of MOBIDIC. Boxes denote different water storages (gravitational storage $W_{\rm g}$, capillary storage $W_{\rm c}$, groundwater storage H, surface storage $W_{\rm s}$, and river system), solid arrows fluxes (evaporation $E_{\rm t}$, precipitation $P_{\rm c}$, infiltration $I_{\rm nf}$, adsorption $A_{\rm d}$, percolation $P_{\rm c}$, surface runoff R, interflow $Q_{\rm d}$, groundwater discharge $Q_{\rm g}$, and surface runoff and interflow from upper cells $(R+Q_{\rm d})_{\rm up}$), dashed arrows different routings, and blue characters major model parameters.

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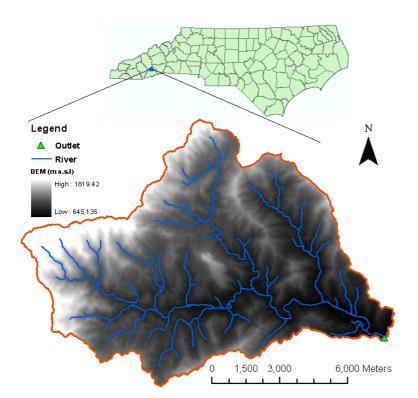


Fig. 2. The location of Davidson watershed, North Carolina, with DEM map, river system (lines), and watershed outlet (the triangle point).

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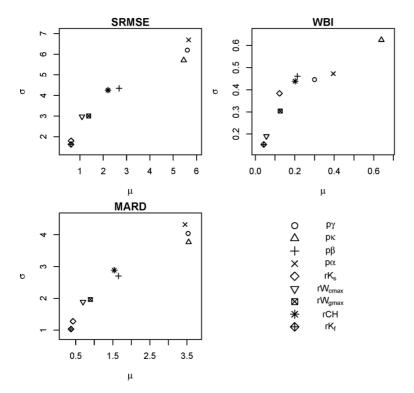


Fig. 3. Multiobjective sensitivity analysis result based on the Morris method.

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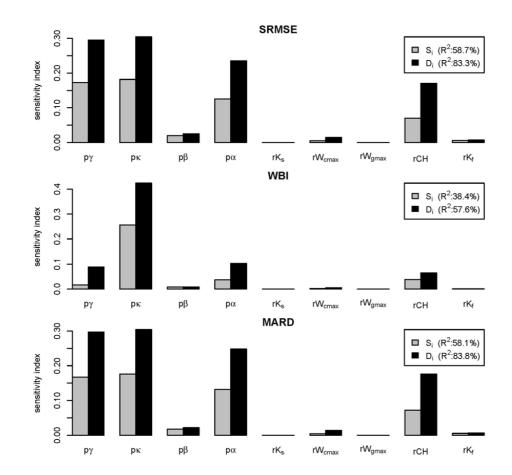


Fig. 4. Multiobjective sensitivity analysis result based on the SDP method.



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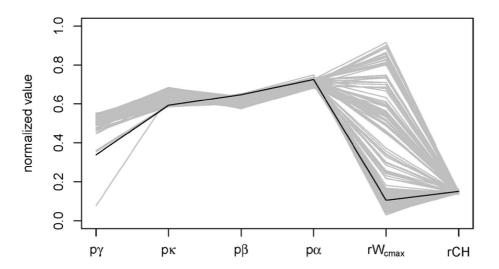


Fig. 5. The normalized factor sets associated with MOO (grey lines) and the solution with SOO (dark line).

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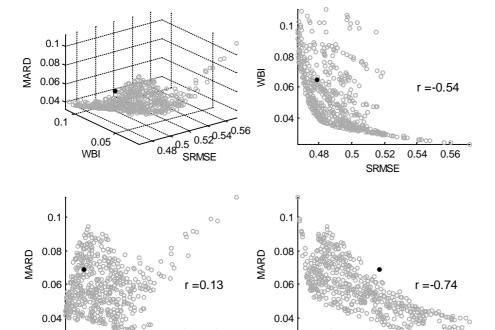


Fig. 6. The Pareto solutions in the three dimensional space (top left), and the projections in the two dimensional subspace (other plots), with MOO, and the black dot is the solution with SOO.

0.04

0.06

WBI

0.08

0.1

0.52

SRMSE

0.48

0.54

0.56

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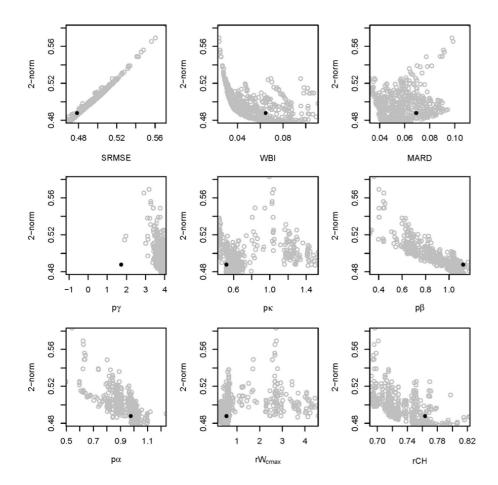


Fig. 7. 2-norm level diagrams representation of the Pareto sets with MOO, and the solution with SOO (black dot).

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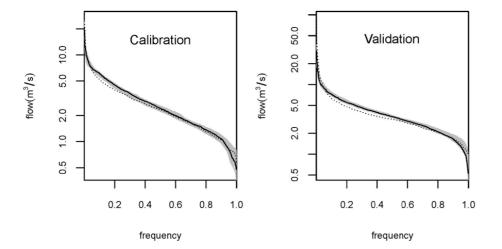


Fig. 8. Flow duration curve for observed (dotted line), and simulated with MOO (grey) and SOO (solid line).

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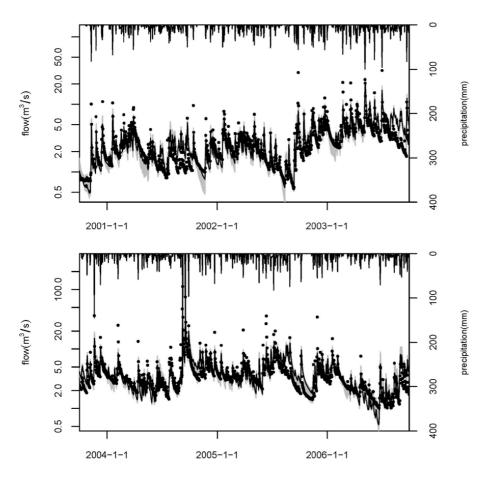


Fig. 9. Observed flows (dotted) and simulated flows with MOO (grey) and SOO (black line) for the calibration period (top) and validation period (bottom).

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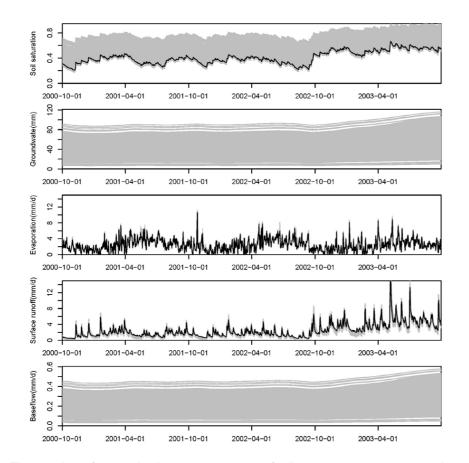


Fig. 10. Time series of watershed average storages (soil water storage expressed as soil saturation, and groundwater depth), and fluxes (evaporation, surface runoff, and baseflow) with MOO (grey) and SOO (black line). For SOO, the groundwater storage and baseflow are close to 0 and hardly seen.

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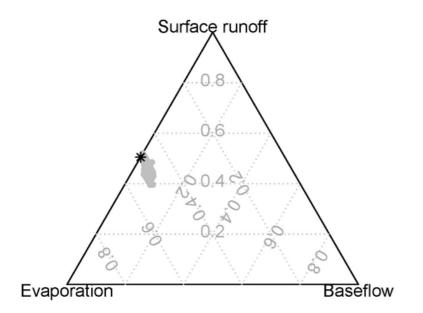


Fig. 11. De Finetti diagram (Ternary plot) of Evaporation, Surface runoff, and Baseflow with MOO (grey) and SOO (black star).