Pitfalls and a feasible solution for using KGE as an informal likelihood function in MCMC methods: DREAM_(ZS) as an example

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Abstract. The Kling-Gupta Efficiency (KGE) is a widely used performance measure because of its advantages in orthogonally considering bias, correlation and variability. However, in most Markov chain Monte Carlo (MCMC) algorithms, error-based formal likelihood functions are commonly applied. Due to its statistically informal characteristics, using the original KGE in MCMC methods leads to problems in posterior density ratios due to negative KGE values and high proposal acceptance rates resulting in less identifiable parameters. In this study we propose adapting the original KGE using a gamma distribution to solve these problems and to apply KGE as an informal likelihood function in the DiffeRential Evolution Adaptive Metropolis DREAM(ZS), which is an advanced MCMC algorithm. We compare our results with the formal likelihood function to show whether our approach is robust and plausible to explore posterior distributions of model parameters and to reproduce the system behaviors. For that, we set three case studies that contain different uncertainties and different types of observation data. Our results show that model parameters cannot be identified and the uncertainty of discharge simulations is large when directly using the original KGE. The adapted KGE finds similar posterior distributions of model parameters derived from the formal likelihood function. Even though the acceptance rate of the adapted KGE is lower than the formal likelihood function for some systems, the convergence rate (efficiency) is similar between the formal and the adapted KGE approaches for the calibration of real hydrological systems showing generally acceptable performances. We also show that both the adapted KGE and the formal likelihood function provide low performances for low flows, while the adapted KGE has a balanced performance for both low and high flows. Furthermore, the adapted KGE shows a general better performance for calibrations of solute concentrations. Thus, our study provides a feasible way to use KGE as an informal likelihood in the MCMC algorithm and provides possibilities to combine multiple data for better and more realistic model calibrations.

1 Introduction

Markov chain Monte Carlo (MCMC) techniques are extremely useful in uncertainty assessments and parameter estimations of hydrological models (Smith & Marshall, 2008). Among those MCMC methods, Vrugt et al. (2008, 2009) developed a

DiffeRential Evolution Adaptive Metropolis (DREAM) algorithm, which has found numerous applications in various fields (Vrugt, 2016). It is an adaptation of the SCEM-UA algorithm (Vrugt et al., 2003a) that can efficiently estimate the posterior probability distribution of model parameters in the presence of high-dimensional and complex response surfaces with multiple local optima.

The formal likelihood function, e.g. mean squared error (MSE) or root mean squared error (RMSE), obtained from first-order statistical principles based on error series derived from simulations and observations, is commonly used in the DREAM algorithm. The formal likelihood function strongly relies on error assumptions, which can highly influence the shape of parameter posterior distributions (Beven et al., 2008). The informal likelihood functions such as Nash-Sutcliffe efficiency (NSE) and the alternative Kling-Gupta Efficiency (KGE) are often used in hydrological studies to indicate the general performance of model simulations (Gupta et al., 2009). These metrics represent an important measure of model performance, so-called goodness-of-fit (Pool et al., 2018). These likelihood functions are not directly derived from stochastic error series, but can be easily used to combine different types of data.

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There are studies that discussed how to adjust the calculation of NSE in order to overcome the problems using NSE in MCMC methods. For example, McMillan & Clark (2009) introduced a constant K into the adapted NSE calculation. By adjusting this constant, they can mimic the weight such that small improvements in NSE can also be distinctly identified leading to the chain evolution. They even found that the informal likelihoods can provide a more complete exploration of the behavioral regions of the response space and hence more accurate estimation of total uncertainty (McMillan & Clark, 2009). Freer et al. (1996) introduced a parameter as an exponent symbolled with N. They argued that higher N values have the effect of accentuating the weight given to the better simulations.

However, how to properly use KGE in the MCMC methods has not been studied. Directly using KGE in MCMC methods, e.g. DREAM algorithm, may raise difficulties such as incorrect posterior ratios due to negative KGE values, and nonlinearity between model performance and KGE values. These difficulties essentially affect chain evolutions such as the acceptance rate, indicating how easy a proposal is accepted, and the convergence rate, denoting how fast a chain converges to a stationary distribution. As a consequence, considering the computational cost with a limited number of realizations in practice the informal character of KGE and its use in MCMC methods influences the exploration of posterior parameter distribution and model uncertainty, such as density of identifiable parameters. Studies showed that using informal likelihood functions in generalized likelihood uncertainty estimation (GLUE) may lead to unsatisfactory posterior distributions of model parameters (Mantovan & Todini, 2006; Stedinger et al., 2008). Using NSE as the likelihood function, the number of measurements cannot be considered. Therefore, with increasing number of measurements, the information added to the performance measure is little, thus preventing the improvement of chain evolution (Mantovan & Todini, 2006). Therefore, to feasibly use KGE in MCMC methods requires solving problems in drawing better proposals to avoid a very flat posterior

distribution, to account for the influence of observation size (the amount of information included in calibration) on parameter estimations and to achieve reasonable acceptance and convergence rates.

In this study we propose adapting the gamma distribution and KGE to find a feasible solution for properly using KGE as an informal likelihood function in DREAM_(ZS). We test the robustness of this approach with three case studies representing known and unknown systems with varying amount of observations and also different types of data using two hydrological models (a lumped and a semi-distributed). The aim of our study is to attempt to form the probability calculation based on KGE in a pseudo formal way. The derivation of probability based on KGE in a statistically sound manner is beyond the scope of this study and will need future work. We compare the performance between the original KGE, GLUE (generalized likelihood uncertainty estimation), the formal likelihood function, the log-transformation and the adapted KGE. We aim to show whether the adapted KGE is robust and plausible to explore posterior distributions of model parameters and to reproduce the hydrological behaviors. Thus, we will compare performance (1) regarding acceptance, convergence, and uncertainty between the adapted KGE, the original KGE, a formal likelihood function and a log-transformation; (2) regarding discharge simulations in terms of general performance, variability, bias and correlation for total, low and high flows; and (3) model performance combining discharge and solutes. These comparisons allow us to provide recommendations for more reliable applications of KGE in MCMC methods in different research areas.

2 Material and Methods

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2.1 Kling-Gupta Efficiency

Kling-Gupta Efficiency (KGE) takes account of variability (α), non-scaled bias (β) and correlation (r) by computing the Euclidian distance (ED) of the three components from the ideal point, which avoids the underestimation of variability and enables the comparison of the bias term between catchments (Gupta et al., 2009).

$$KGE = 1 - ED \tag{1}$$

$$ED = \sqrt{(r-1)^2 + (\alpha - 1)^2 + (\beta - 1)^2}$$
with $\alpha = \frac{\sigma_s}{\sigma_o}$ and $\beta = \frac{\mu_s}{\mu_o}$ (2)

where (μ_s, σ_s) and (μ_o, σ_o) are the mean and standard deviation of simulations and observations, respectively. KGE ranges from $-\infty$ to 1 with the optimal value at unity. A value larger than -0.41 indicates that a model improves upon using the mean (Knoben et al., 2019).

2.2 Adapting KGE in DREAM(ZS)

2.2.1 Basics of DREAM(ZS)

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DREAM_(ZS) is one type of DREAM algorithms, which uses sampling from an archive of past states to generate candidate points in each individual chain. It automatically tunes the scale and orientation of the proposal distribution towards the target distribution and maintains a detailed balance and ergodicity (Vrugt et al., 2008, 2009). We take DREAM_(ZS) as an example and investigate the appropriate way to use KGE within DREAM_(ZS). Since our goal is to adapt KGE as an informal likelihood function, we focus on the Metropolis probability, $p_{acc}(X^i \to X_p^i)$, which calculates the probability to accept a proposal.

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$$p_{acc}(\mathbf{X}^i \to \mathbf{X}^i_p) = min[1, p(\mathbf{X}^i_p)/p(\mathbf{X}^i)]$$
 (3)

where $p(X_p^i)$ and $p(X^i)$ denote the probability density of the proposal and the present location of the *i*-th chain. If $p_{acc}(X^i \to X_p^i)$ is larger than a random value drawn from the uniform distribution U(0, 1), the proposal will be accepted, otherwise the chain remains in the present location. After chain evolutions, the Gelman-Rubin \widehat{R}_j convergence diagnostic is computed for each parameter $j=\{1,\ldots,d\}$. If $\widehat{R}_j\leq 1.2$, the convergence can be declared.

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$$\widehat{R}_{j} = \sqrt{\frac{N+1}{N} \cdot \frac{\widehat{\sigma}_{+}^{2(j)}}{W_{j}} - \frac{T-2}{N \cdot T}}$$
 (4)

where N and T signify, respectively, the number of chains and the number of samples in each chain. W_j is the within-chain variance, and $\hat{\sigma}_+^{2(j)}$ is an estimate of the variance of the j-th parameter of the target distribution.

We choose the easily applied formal likelihood function (lik=11 in Table B1, Vrugt, 2016) to calculate the log-likelihood (LogL), which assumes the error residuals to be normally distributed. It is described as:

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$$\operatorname{LogL} = -\frac{n}{2} \log \{ \sum_{t=1}^{n} e_t(\mathbf{X})^2 \}$$
 (5)

where $e_t(X)$ denotes the t-th error residuals. n is the total number of observations.

2.2.2 Deriving pseudo probability density based on KGE

Based on above-mentioned basics of DREAM_(ZS), using KGE as a likelihood function becomes the proper calculation of pseudo probability density $p(X_p^i)$. From the straightforward and easily applied way, one would directly use KGE as $p(X_p^i)$ by setting the negative KGE values to zeros. However, it results in two problems (Fig. 1):

• Problem 1: the probability density $p(X_p^i)$ needs to be positive in order to get a positive ratio $p(X_p^i)/p(X^i)$ to determine the right orientation to accept a proposal. However, KGE ranges from $-\infty$ to 1. Setting negative KGE

values to zero can work, but then we lose the orientation of proposals with negative KGE values. Thus, it reduces the efficiency of chain evolution.

- Problem 2: the model performance does not linearly increase with the linear increase of KGE. Therefore, directly using positive KGE as the pseudo probability density $p(X_p^i)$ will lead to a high possibility to accept poor proposals. For instance, the probability is 0.75 (0.6/0.8) to accept a proposal with KGE=0.6 under the present sample with KGE=0.8. However, the performance of a simulation with KGE=0.8 is much better compared to one with KGE=0.6.
- We propose solving these two problems with adapting the gamma distribution and KGE to derive a proper probability density (Fig. 1). The gamma distribution has two parameters, the shape parameter (*k*) and the scale parameter (*θ*), and is evaluated for variables with positive values. When the shape parameter *k*=1, the probability distribution is one-side with nonlinear decreasing probability for increasing variable values. Therefore, we can use 1-KGE (ranges from 0 to ∞) as the variable for gamma distribution and get higher probability for larger KGE values. When choosing the scale parameter *θ*=0.5, the increasing rate of probability becomes faster when KGE>0.5, especially when KGE>0.7 the probability increases much faster. This helps chain evolution to find proposals which leads to high model performance. The non-linear increase of probability with increasing KGE is shown as the red line in the top right box of Fig. 1. The log-likelihood is used in DREAM_(ZS), thus we derive the pseudo log-likelihood (LogL) using gamma distribution and KGE. To include the influence of observation size on the parameter estimations, analogy to the formal likelihood function we take account of the number of observations (the term n/2 in the formal log-likelihood) in deriving the pseudo log-likelihood. Finally, we have the pseudo log-likelihood using KGE and gamma distribution as:

$$LogL = \frac{n}{2}log(f(1 - KGE|\mathbf{X}; k, \theta)), \text{ with } k=1 \text{ and } \theta=0.5$$
 (6)

where, f() is the gamma probability density function; k and θ are the shape and scale parameters, respectively. n is the number of observations. K represents the parameter vector of the calibrated model. The purpose of this approach is to provide a feasible way to incorporate KGE as the informal likelihood function for MCMC methods, in our case DREAM_(ZS). Another goal is to achieve similar performance as using the formal likelihood function such as RMSE so that we can compare the model simulations and predictions between formal and informal (our approach, KGE) likelihood functions.

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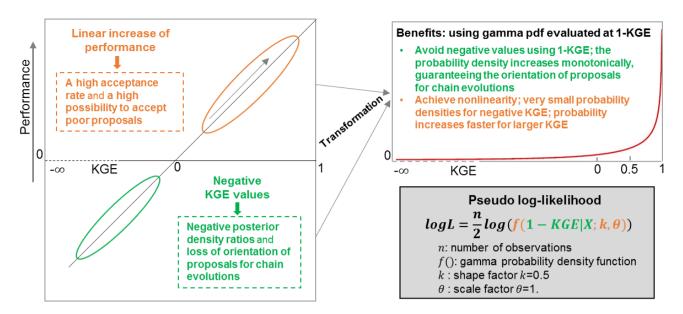


Figure 1 Concept of adapting KGE in DREAM_(ZS). It shows the problems of using the original KGE as the likelihood function in DREAM_(ZS) and how to adapt gamma distribution and KGE to get a proper informal likelihood.

2.3 Case studies

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To test the robustness of our new approach, we define three case studies: (1) true and pseudo-analytical posterior distributions of model parameters are known by a virtual experiment, and uncertainties in model structures and input data are absent; (2) calibrations and evaluations with a long observation time-series using a rainfall-runoff model, which allows comparing the performance between three approaches by varying the amount of data in calibration. The parameterization of the system is unknown and there are uncertainties in model structure, input data and observations; and (3) a model calibration combining hydrodynamics and simple solute transport for a more complex karst system with a large subsurface heterogeneity and processes for fast recharge and groundwater discharge from conduit networks. The observation period is short and uncertainties exist in model structure, input and observation data, and model parameter estimations.

150 2.3.1 Case study 1: virtual experiment

We generate a virtual experiment using a rainfall-runoff model (the HBV model). We obtain the forcing data, daily mean temperature and daily precipitation (2001–2008), from the German site in Liu et al. (2021). The HBV model represents typical catchment rainfall-runoff processes considering one soil water storage and two groundwater storages (Lindström et al., 1997). In this virtual experiment, we use the model version without snow processes, which contains nine parameters. Since our goal is not the model itself but the calibration of model parameters, we only provide descriptions of model parameters (Table 1). For the model structure and equations, refer to Liu et al. (2021).

Table 1 Names, description, ranges and virtual true values of the HBV model parameters for the virtual experiment

Parameter	Description	Unit	Parameter ranges		True
			lower	upper	value
BETA	Shape coefficient of recharge function	[-]	1	6	4.5
FC	Maximum water storage in the unsaturated-zone store	[mm]	50	700	600
<i>K0</i>	Additional recession coefficient of upper groundwater store	$[d^{-1}]$	0.05	0.99	0.5
K1	Recession coefficient of upper groundwater store	[d ⁻¹]	0.01	0.8	0.25
K2	Recession coefficient of lower groundwater store	$[d^{-1}]$	0.001	0.15	0.07
LP	Soil moisture value above which actual evaporation reaches potential evaporation	[-]	0.3	1	0.55
PERC	Maximum percolation to lower zone	[mmd ⁻¹]	0	6	3
UZL	Threshold parameter for extra outflow from upper zone	[mm]	0	100	60
MAXBAS	Length of equilateral triangular weighting function	[d]	1	3	2

note: K0, UZL and MAXBAS (shaded ones) are insensitive parameters in this case study, thus are fixed to the true values

Since the analytical posterior distributions of model parameters of a hydrological model is hardly achievable, we use the following procedure to generate the pseudo-analytical posterior distribution. Firstly, we set the catchment area of 100 km² and run the model with "true" parameters for 2004–2008 to obtain the simulated discharge. Secondly, assuming a normal distribution for error residuals (a common assumption for hydrological modeling), we generate random values from a normal distribution (mean=0, standard deviation=5% of the mean simulated discharge) and add these random values as measurement errors to the simulated discharge to form the observations. Finally, due to no uncertainty in input data and model structure, using this setting for measurement errors we can use the formal likelihood function (equation 5) to derive the pseudo-analytical posterior distribution of model parameters. We performed a local parameter sensitivity analysis before model calibrations to find insensitive parameters (*KO*, *UZL* and *MAXBAS* in Table 1, changing these parameters only affects model performance KGE by 0.001). Therefore, we fixed the three parameters in calibration, resulting in six parameters to be calibrated.

170 2.3.2 Case study 2: long observations for rainfall-runoff modeling

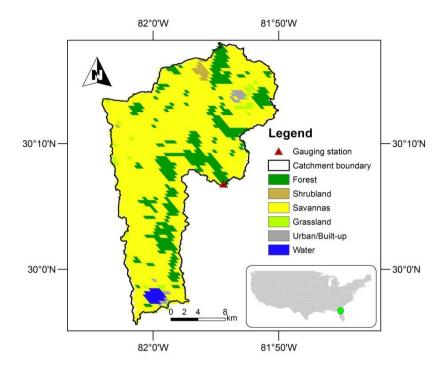
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In order to test the capability of our approach for a real system with uncertainties in forcing, observations, model structure and model parameters, we select a catchment from the CAMELS-US dataset (Newman et al., 2014, 2015) and simulate the rainfall-runoff processes with the HBV model as case study 1. We also compare the performance of our transformed KGE with the GLUE method. We have the following criteria to select this catchment: 1) catchment area is between 100 and 500 km² to avoid the influence of channel routing; 2) the snow fraction is zero to avoid the snow processes to reduce the HBV model parameters and the number of parameters remain nine (table 1); and 3) the carbonate rock fraction is zero since we

will test a karst catchment as a separate case study. We then choose the first catchment (smallest number in catchment ID) that fulfills our criteria, the catchment 02246000 (gauging station at North Fork black creek NR Middleburg, FLA, USA). It has a catchment area of 451 km² with the mean annual precipitation of 1352 mm. The main land cover is savannas, accounting for 77% of the total area. Details of the catchment properties can be found in CAMELS-US dataset (Newman et al., 2014, 2015).



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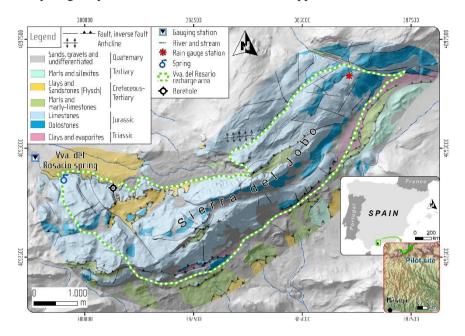
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Figure 2 Study area of the catchment with the gauging station at North Fork black creek NR Middleburg, FLA, USA for the case study 2.

2.3.3 Case study 3: short observations for a heterogeneous karst system

In order to test the capability of our new approach for a complex system, we set case study 3 in a karst system that has conduit systems resulting in fast recharge and discharge. It has uncertainties from the forcing data, the model structure and observation errors. Daily discharge time-series and weekly solute (Cl⁻, NO₃⁻ and SO₄²-) concentrations of the hydrological years 10.1.2006–9.30.2009 are combined for model calibrations. The study site is located in Southern Spain with a recharge area of 13.85 km² (the study site in Hartmann et al., 2014). In this case, we use the same model, the VarKarst model with the solute transport routine, for spring discharge and solute simulations since it was successfully applied to this site before (Hartmann et al., 2014). The VarKarst model is a semi-distributed hydrological model, which considers subsurface heterogeneity, soil and epikarst storage dynamics and groundwater hydrodynamics. It uses a mixing routine to simply reproduce the solute transport. These processes are represented by 10 parameters (Table 2). Details of the VarKarst model

processes and assumptions can be found in Hartmann et al. (2014). Discussion of the transport processes is beyond the scope of this study, we, therefore, choose the same processes as published in Hartmann et al., (2014). Our study then focuses on comparing the performance of different calibration approaches.



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Figure 3 Study area of the Rosario spring for case study 3. This map is an updated version of the map in Hartmann et al. (2014)

Table 2 Names, descriptions and ranges of the VarKarst model parameters

Parameter	Description	Unit	Parameter ranges	
1 arameter	Description	Omt	lower	upper
$V_{mean,S}$	Mean soil storage capacity	[mm]	0	500
$V_{mean,E}$	Mean epikarst storage capacity	[mm]	0	500
a_{SE}	Soil/epikarst depth variability constant	[-]	0.1	2
$K_{mean,E}$	Epikarst mean storage coefficient	[d]	1	50
a_{fsep}	Recharge separation variability constant	[-]	0.1	2
K_C	Conduit storage coefficient	[d]	1	20
a_{GW}	Groundwater variability constant	[-]	0.1	2
a_{Geo}	Equilibrium concentration variability constant	[-]	0	2
$log_{10}~Geo_{Cl}$	Equilibrium concentration of Cl in matrix	$[\log_{10} (mg L^{-1})]$	0	5
$log_{10}~Geo_{SO4}$	Equilibrium concentration of SO ₄ in matrix	$[log_{10} (mg L^{-1})]$	0	5

2.3.4 Calibration and evaluation

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For calibration of the three case studies, we have the GLUE approach and DREAM_(ZS) with different likelihood functions: (1) using the original KGE as the likelihood function ("KGE_{ori}" thereafter) – here negative KGE is set to zero to avoid negative posterior density ratios; (2) using the traditional formal likelihood assuming error is normally distributed("formal" thereafter); (3) using the log-transformation ("formal_{log}" thereafter), which is suggested good for low flows (McInerney et al., 2017) and (4) using our new approach adapting gamma distribution and KGE to derive the pseudo log-likelihood ("KGE_{gamma}" thereafter). We use three parallel Markov chains (default setting in DREAM_(ZS)), and set 20,000 realizations for case studies 1 and 2, while 30,000 realizations for case study 3 (due to more processes and parameters). The last 25% of realizations are used to approximate the posterior distributions and the corresponding parameter sets are used for the discharge simulations in the evaluation period. They are also used to derive the parameter uncertainty and the total simulation uncertainty (parameter uncertainty + randomly sampled error from a normal distribution with mean=0 and standard deviation=root mean squared error of the simulation with the maximum a posteriori parameter) in DREAM_(ZS).

For case study 1, following a standard calibration procedure we use 2001-2003 as the warm-up period and 2004-2008 as the calibration period. The posterior distribution of model parameters derived from DREAM_(ZS) using the formal likelihood function will be deemed as the pseudo-analytical posterior distribution. By comparing to it, we can investigate whether calibrations using KGE_{ori} and KGE_{gamma} can explore the right posterior distribution and true model parameters.

For case study 2, the true model parameters are unknown. We use 25 hydrological years (10.1.1980–9.30.2005) to perform the calibration and evaluation. We choose the Daymet forcing data to drive our hydrological model since this meteorological

data has potential evapotranspiration (PET) estimates and was used to calculate the catchment climatic properties (Addor et al., 2017). We use the first 5 years as the warm-up period, then the following 10 years for calibration and the last 10 years for evaluation. We test the performance of four approaches (GLUE, formal, formal_{log} and KGE_{gamma}) using 1, 3, 5, 8 and 10 years of observations for calibrations, respectively to check the capability of the four approaches for calibration of a real system with varying amount of available observations. For the GLUE method, we use Nash-Schetcliff efficiency (NSE) as the objective function. We set 20,000 realizations and choose the top 25% in performance (keep the same as DREAM_(ZS)) as the behavioral parameter sets to explore the posterior distribution.

For case study 3, the true model parameters are unknown too. We calibrate the hydrodynamics and solute transport simultaneously. For calibrations using the formal likelihood function, we normalize each observation variable by its mean to exclude the influence of units and magnitudes of discharge and solute concentrations. We compare the performance of three approaches: (i) We use the formal likelihood function and use the normalized daily discharge and normalized weekly concentrations of three solutes as the combined observations ("formal_{norm}" thereafter). Here we do not consider the difference in the total number of observations between discharge and three solutes (the total number of discharge observations is ten times of each solute); (ii) We also use the formal likelihood function. But we replicate each normalized solute observations ten times to have the same weight for discharge and each solute ("formal_{norm},"). Issues regarding different weights for discharge and solutes and different ways to obtain weights are out of the scope of our study; and (iii) We use the KGE_{gamma} approach. Firstly, we calculate KGE for discharge and each solute using their observations and simulations (totally four KGE values, for discharge, Cl⁻, NO₃⁻ and SO₄²-). Then we calculate the mean of the four KGE values (equal weight for the four variables, same as (ii)) and use it in the KGE_{gamma} approach. We use three hydrological years (10.1.2003–9.30.2006) for warm-up of the simulations and the three following hydrological years (10.1.2006–9.30.2009) for calibration. It is the same as the calibration setting in Hartmann et al. (2014) considering the short observations of each solute.

The model performance for calibration and evaluation is examined using KGE and its three components representing variability (α) , bias (β) and correlation (r). The calculations of KGE, α , β and r refer to equations 1 and 2. We evaluate the model performance using the four metrics for total flow, low flow (smaller than 10^{th} percentile of observed discharge) and high flow (larger than 90^{th} percentile of observed discharge) and also for three solutes.

3 Results

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3.1 Case study 1: posterior parameter exploration

When using the original KGE (set negative KGE values to zeros) as the likelihood function, the posterior parameter range is only slightly reduced for all sensitive parameters compared to the prior uniform distribution. In addition, the density around the true values of model parameters are still very flat, indicating true model parameters are barely identified (Fig. 4). When

applying the adapted KGE (adapting the gamma distribution and KGE to derive probability density), the posterior parameter range is much more reduced and the reduced range is more or less centered at the true values as shown in Fig. 4. Compared to the pseudo-analytical posterior distributions of all model parameters derived from the formal likelihood function using the special virtual setting, the adapted KGE approach (KGE_{gamma}) shows similar magnitude and shape regarding the posterior distributions. It suggests that our adapted KGE approach performs similarly as using the traditional formal likelihood function and can explore the right parameter posterior distributions.

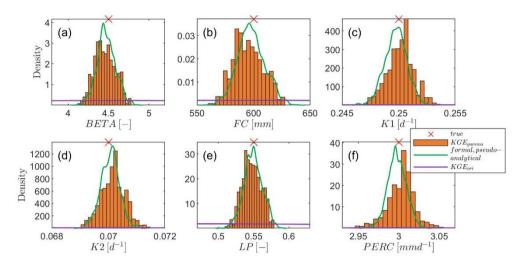


Figure 4 Posterior distributions of sensitive model parameters for the virtual experiment. The red cross symbol denotes the true model parameter value. KGE_{ori} indicates using the original KGE (set negative KGE to zero) as the likelihood function, while KGE_{gamma} represents our new approach using gamma distribution and KGE to derive the probability density, and the case *formal*, *pseudo-analytical* means pseudo-analytical posterior distribution derived from the formal likelihood function. Note that the densities of using KGE_{ori} is very low and close to the x axis.

As expected, using the original KGE, we have a very high acceptance rate (ca. 60-80%, Fig. 5a), leading to a very fast convergence (Fig. 5b). This results in a large uncertainty bound in the discharge simulations (Fig. 5c), and the uncertainty of peak discharges is particularly large. With the adapted KGE, we see that the acceptance rate becomes smaller and the convergence gets slower. This can be explained by introducing the nonlinearity of the adapted KGE: probability densities for large and small KGE values are more distinct compared to the original KGE. Fig. 5a also shows that the acceptance rate of our approach is 5-10%, which is lower than ca. 20% of the formal likelihood function. Similarly, the convergence rate of our approach is slower than the formal likelihood function (Fig. 5b). It suggests that using the formal likelihood function has a higher efficiency than the approach adapting KGE for calibrations of a system that only contains little uncertainty (only small observation errors in our case). However, when more uncertainties appear, e.g. uncertainties in forcing data and model structures, the convergence rates (efficiency) become similar between the adapted KGE and the formal likelihood function (refer to the following subsections, Fig. 6b). Compared to the width of the discharge uncertainty bound using the original KGE (Fig. 5c), calibration using the formal likelihood function and the adapted KGE both reduce the average width of total

discharge uncertainty bound by ca 85%. Since this virtual experiment does not assume uncertainty in the input data and the model structure, the adapted KGE shows a similar performance in the uncertainty estimation as using the formal likelihood function; both can closely reproduce observations.

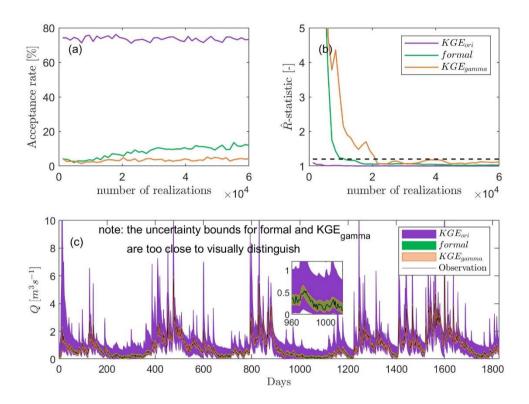


Figure 5 Acceptance rate (a), convergence rate shown with \widehat{R} -statistic (b), and uncertainty of discharge simulations (total uncertainty) for the virtual experiment (c). KGE_{ori} indicates using the original KGE as the likelihood function, while KGE_{gamma} represents our new approach using gamma distribution and KGE to derive the probability density, and the case *formal* means using the traditional formal likelihood function. Note the uncertainty bounds for *formal* and KGE_{eamma} are too small to be visually seen.

3.2 Case study 2: model parameter calibrations with long observations

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Calibration for a real system (with uncertainties in forcing, observations, and model structure and parameters), the acceptance rate of the adapted KGE is lower than that of the formal likelihood function, but higher than that of the log-transformation (Fig. 6a) for calibrations using both short and long observations., The convergence rate is almost identical between the formal likelihood function and the adapted KGE (higher than the log-transformation, Fig. 6b) This indicates that our approach has a same efficiency as the formal likelihood function and a higher efficiency than the log-transformation for calibrations of a system with more uncertainties. With more observations in calibrations, the unidentified parameters *KO* and *UZL* (Fig. 6c and 6d) using the adapted KGE and the formal likelihood function become identified (Fig. 6g and 6h). The

identified parameter values for *K0* (Fig. 6g) show a similar distribution that is different from the log-transformation, while the identified parameter values for *UZL* (Fig. 6h) differ between the three approaches. The identified parameter *K1* with 1-year observations in calibration (Fig. 6e) shows a similar distribution as using 10-year observations (Fig. 6i) between the adapted KGE and the formal likelihood function, which is different from the log-transformation. The density is higher at the peak when using more observations (Fig. 6i). For the identified parameter *MAXBAS* between the three approaches is similar when using 1-year observations for calibration (Fig. 6f), but it changes after adding more observations into calibration (Fig. 6g), where the adapted KGE approach shows a similar distribution as the log-transformation. This suggests that using different likelihood functions may lead to different identified model parameters for a system with various uncertainties due to parameter interactions. More information may be needed to confine the model parameters.

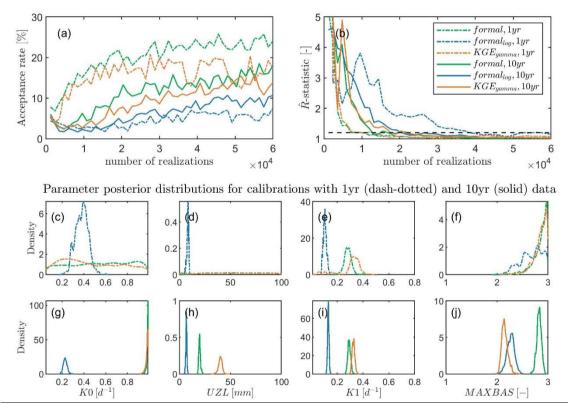


Figure 6 Acceptance rate (a), convergence rate shown with \widehat{R} -statistic (b), and posterior distributions of selected parameters for calibrations with 1-year (c-f) and 10-year (g-j) observations, respectively. KGE_{gamma} represents our new approach using gamma distribution and KGE to derive the probability density, the case *formal* means using the traditional formal likelihood function, and *formal*_{log} means using the log transformation. The subscripts lyr and l0yr denote calibrations with 1-year and 10-year observations, respectively. The four parameters (nine parameters in total) are selected to represent different cases from unidentified (less observations for calibrations) to identified (more observations for calibrations) parameter and show how the identified parameters change when using different amount of observations for calibrations.

In this section, we focus on analyzing the performance in the evaluation period to show the prediction ability of the four approaches. Generally, the uncertainty of the model performance (represented by the interquartile of KGE, α , β and r) of the GLUE approach is much larger than the other three approaches (Fig. 7) regardless of total flow, low flow or high flow. With the increasing amount of observations added to calibrations, the performance of GLUE does not change significantly for all four metrics and for all flow conditions, while using the adapted KGE or the formal likelihood function we can see an increasing trend of the model performance. The log-transformation only has an improved performance regarding low flow with increasing observation data. In the following, we focus on comparing the performance between the adapted KGE, the formal likelihood function and the log-transformation. The log-transformation has a better performance for low flow as expected, but a lower performance for high flow. The formal likelihood function without transformation has a better performance for high flow but a lower performance for low flow. The adapted KGE combines these advantages, leading to a good and balanced performance for low and high flows. For the total flow, the general performance KGE of our approach is higher than using the other two formal likelihood functions. The three approaches perform similarly for the variability (α) for calibrations with 3-5 years of data, while the adapted KGE tends to overestimate and the other two formal likelihood functions underestimate variability when more data is added to calibrations. The adapted KGE has a smaller overestimation of bias (β) than the formal likelihood functions. They have similar performance in terms of the correlation (r) for the total flow. For the low flow, the performances (all metrics) of all three approaches are poor. The adapted KGE and the logtransformation have a similar general performance in KGE, which is better than the formal likelihood function without logtransformation. The adapted KGE has a lower overestimation of variability and a better simulation of bias, while the two formal likelihood functions has a better performance in correlation. For the high flow, the adapted KGE and the formal likelihood function perform similarly in terms of KGE, bias (β) and correlation (r), which are both better than the logtransformation. But the adapted KGE has a better representation of variability (a) than the two formal likelihood functions.

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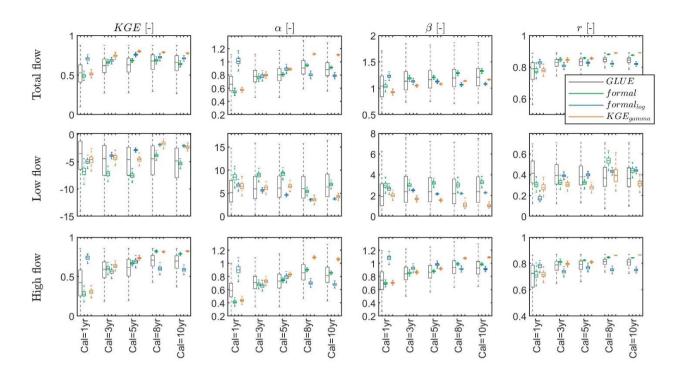


Figure 7 General performance (KGE), variability (α), non-scaled bias (β) and correlation (r) for total flow, low flow (smaller than 10th percentile of observed discharge) and high flow (larger than 90th percentile of observed discharge) during the evaluation period using the GLUE approach (*GLUE*), the formal likelihood function (*formal*), the log-transformation (*formallog*) and our approach using KGE and gamma distribution to derive probability density (*KGEgamma*) with varying amount of observations (1-year to 10-year) in calibration, for instance, calibration with 1-year observations is shown as Cal=1yr. The boxplot shows the performance of the last 25% of all simulations
 (top 25% in performance for GLUE), which is used to approximate the "true" system behavior in DREAM_(ZS). The performance is only shown for the evaluation period to avoid too much information in the boxplot and to represent the prediction ability of different approaches. The performance of the calibration period is provided in Fig. S1 in the supplement. The optimal value for KGE, α, β and r is one, and the closer to one the better the performance.

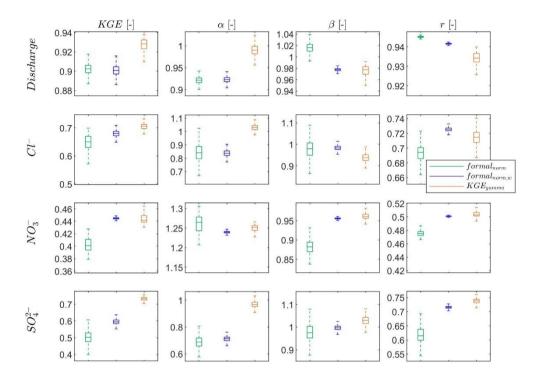
3.3 Case study 3: model parameter calibrations for a heterogeneous karst system

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For calibration combining discharge and solute concentrations at this heterogeneous karst system with short observation records, the adapted KGE is superior than the formal likelihood function regardless of the weight given to discharge and solutes (Fig. 8). For the general performance measured by KGE, the adapted KGE approach performs best, followed by the formal likelihood function with same weights in discharge and each solute, and then the calibration with different weights (the number of discharge data is 10 times of each solute). The performance regarding discharge is similar between the three approaches (the mean KGE is around 0.9) with a slight higher performance for the adapted KGE approach. However, the adapted KGE approach improves the mean performance regarding Cl^- , NO_3^- and SO_4^{2-} by 7%, 10% and 44% compared to the formal likelihood function using discharge and solutes. The adapted KGE approach has a very good representation of variability (α) compared to the other two approaches, especially for discharge, Cl^- and SO_4^{2-} where the variability metric α is

centered around 1. The performance in terms of bias (β) is in the range of 0.9 and 1.1 for all three approaches. The correlation of simulated and observed discharge is all larger than 0.9 for the three approaches. But the adapted KGE and the formal likelihood function using the same weight in discharge and solutes have a higher performance (improvement is up to 17%) regarding correlation for the three solutes compared with the formal likelihood function using different amount of observation data.



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Figure 8 General performance (KGE), variability (α), non-scaled bias (β) and correlation (r) for discharge and solutes (Cl⁻, NO₃⁻ and SO₄²-) for a heterogeneous karst system using the formal likelihood function with normalized observations ($formal_{norm}$), the formal likelihood function with equal weights of normalized discharge and each solute ($formal_{norm,w}$) and our approach using KGE and gamma distribution to derive probability density (KGE_{gamma}). The boxplot shows the performance of the last 25% of all simulations, which is used to approximate the "true" system behavior in DREAM(ZS). The optimal value for KGE, α , β and r is one, and the closer to one the better the performance.

Since the formal likelihood function with the same weight for discharge and solutes ($formal_{norm,w}$) has a better performance than the formal likelihood function with different amount of observation data for discharge and solutes ($formal_{norm}$), we only show the comparison regarding the total uncertainty between $formal_{norm,w}$ and KGE_{gamma} in Fig. 9. The two approaches have a similar uncertainty estimate (both total uncertainty in Fig. 9 and parameter uncertainty in Fig. S2) for discharge and NO₃⁻.

However, the *formal*_{norm,w} approach has a big underestimation for Cl⁻ and SO₄²⁻ compared to the adapted KGE approach even though the uncertainty width is similar. From Fig. 9, we can see the adapted KGE approach can cover most very high and very low concentration values in the total uncertainty band. For the parameter uncertainty (Fig. S2), the adapted KGE approach performs better for Cl⁻ and SO₄²⁻ as well. This indicates that the adapted KGE approach can better represent the uncertainty when using multiple types of data for calibration such as discharge and three solutes in this case study.

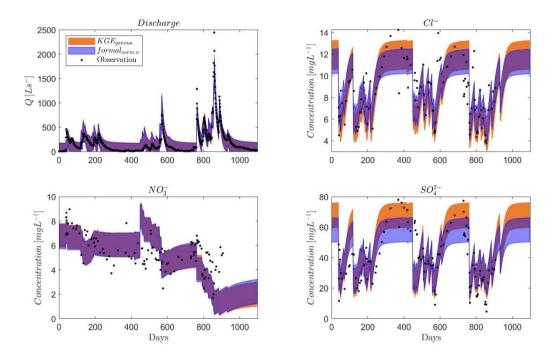


Figure 9 Total uncertainty for discharge and solutes (Cl^- , NO_3^- and SO_4^{2-}) for a heterogeneous karst system using the formal likelihood function with equal weights of normalized discharge and each solute ($formal_{norm,w}$) and our approach using KGE and gamma distribution to derive probability density (KGE_{gamma}). The total uncertainty is estimated based on the last 25% of all simulations. The parameter uncertainty is shown in Fig. S2 in the supplement.

4 Discussion

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Using the original KGE as the likelihood function, model parameters are not well identifiable, which results in a very large uncertainty in the simulation. This is because directly using the original KGE as the likelihood estimate assumes a linear increase of probability density with the linear increase of KGE. It leads to the identification of parameter proposals with good KGE performance more difficult and inefficient. The difference between large and small KGE values is not distinctly large enough that the probability to accept poor proposals is high. This is why we find a very large acceptance rate and a very fast convergence rate. Mantovan and Todini (2006) and Stedinger et al. (2008) also mentioned that using the informal likelihood function, such as Nash-Sutcliffe Efficiency (NSE), in the generalized likelihood uncertainty estimation (GLUE) as

objectives cannot find proper posterior distributions of model parameters. Therefore, directly using the original KGE should be avoided and some adaptations to solve the incapability of exploring the posterior distributions such as our approach are needed in MCMC methods.

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The adapted KGE can well estimate the pseudo-analytical posterior distributions of model parameters derived from the formal likelihood function in case study 1. This suggests that it is capable of exploring the parameter posterior distributions. The adapted KGE has a lower acceptance rate and convergence rate compared to the formal likelihood function for the virtual experiment (case study 1). The possible reason is that one KGE value can cover multiple error combinations with the same RMSE around the true optimum, which makes the RMSE a bit more efficient to draw proposals for parameters very close to the true optimum (known parameters in case study 1) for a system that only contains little uncertainty. However, calibrations of real systems usually contain more uncertainties e.g., uncertainties in forcing (including the spatial averaging), observation data (measurement errors), and uncertainty in model structures. The adapted KGE has a similar convergence rate (efficiency) as the formal likelihood for the real-world calibrations (case study 2). In particular, the acceptance rate of the adapted KGE is around 20% for a system that we have good input and observations (case study 2). This is similar to the formal likelihood function and is also close to the theoretically optimal acceptance rate (0.234) in Metropolis algorithms with random walk (Yang et al., 2020).

The uncertainty bound of discharge simulations in case study 1 is almost identical between the adapted KGE and the formal likelihood function. This indicates our approach can behave similarly concerning discharge uncertainty estimation as the formal likelihood function. For the calibration to the real system, the adapted KGE even has a higher general performance in terms of the mean KGE of the evaluation for the total, low and high flows than the formal likelihood function and the logtransformation. McMillan & Clark (2009) had a similar finding that using another informal likelihood function, NSE, in MCMC methods outperforms the formal likelihood for calibrations with high variability and multi-optimum. The formal likelihood functions go along with the strong assumption that errors are distributed normally (Vrugt et al., 2008, 2009), the informal likelihood function KGE takes into account more variability without strict assumptions on error sources (Gupta et al., 2009). The adapted KGE performs similar to the formal likelihood function regarding the correlation between simulations and observations shown in case study 2. However, they all have lower performance for the low flow simulations compared to the total and high flow. While the log-transformation works well for low flow (case study 2), the adapted KGE has a good and balanced performance for both high and low flows. It also shows a lower overestimation of bias in low flows shown as the metric β in case study 2. Jeannin et al. (2021) found that using formal likelihood functions such as RMSE as objectives has a large bias in baseflow simulations. For RMSE as the objective, each individual error has the same weight. To have a high overall performance, the optimization tends to firstly fit the high flows since its error is relatively larger and the contribution to RMSE is thus larger. The log-transformation can improve the calibration for low values but is not good for high values. The adapted KGE and the formal likelihood function both have a better representation of variability (a) with 415 more observations included in calibrations. This makes sense since more data is involved in calibration more information on variability will be captured in calibrations.

While our approach has a similar performance as the formal likelihood function for discharge simulations, we find similar posterior distributions for certain parameters but also inconsistent posterior distributions for some parameters between the formal and informal approaches. This is because some model processes interplay with other processes such that there are compensates of one parameter for another, i.e. parameter interactions. Adding additional information, e.g. solutes in case study 3, can help to further constrain model parameters (Hartmann et al., 2017) that represents the complexity of real hydrological systems. Our study show that the adapted KGE approach is superior on simultaneously calibrating model parameters with different types of data than the formal likelihood function. This can improve the model calibration using the traditional separate steps such as firstly calibrating discharge and then solute processes in Liu et al. (2020). Many studies have shown that multi-objective calibrations allow to adequately and properly estimate important characteristics of a system (Vrugt, et al., 2003b; Yapo et al., 1998). Using KGE can provide a feasible way to combine various types of observations as a measure of multi-objective performance and avoids issues regarding data units, scales and frequency.

Even though our approach adapts the gamma distribution to compute the probability density for KGE, the way we formulate the likelihood function based on KGE is still informal. It means the derivation of the likelihood is not from a strict theoretical probability framework, which is a limitation of our approach. Nevertheless, our approach provides a feasible and pragmatic way and a close solution to the formal likelihood function to avoid the pitfalls of directly using the original KGE in MCMC methods. Future work is needed to find a solution to link probability density and KGE in order to incorporate KGE in a statistical manner as much as possible.

5 Conclusions

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Our study demonstrates that using the original KGE in DREAM_(ZS) results in a very high acceptance rate and a large uncertainty bound of discharge simulations. This is due to the confusing evolution orientation for negative KGE values and the nonlinear performance of KGE. To solve these two problems, we propose adapting KGE with the gamma distribution to formulate the pseudo log-likelihood function to avoid negative posterior density ratios and to include a proper nonlinearity of performance. With three case studies we demonstrate that the adapted KGE performs as good as the formal likelihood function for the exploration of the posterior distributions of model parameters. Through the calibrations varying the amount of observations included in the calibration, we show that the adapted KGE is robust and has a good and balanced performance for both low and high flows compared to the formal likelihood function and the log-transformation. Our approach even has a higher general performance, the mean KGE of the evaluation, and a smaller bias overestimation of low flows than the formal likelihood function. Our study shows that the adapted KGE approach outperforms the formal likelihood function for calibrations using discharge and solutes. The limitation of our approach is the lack of theoretical

probability derivation. Besides that formal limitation, our approach keeps the advantages of KGE, e.g. consideration of variability and possibilities to combine multiple types of data, and performs like a pseudo formal likelihood. Thus, it provides a feasible way to use KGE as an informal likelihood function in MCMC methods.

Data and code availability

450 All data used in this study has been published in Liu, et al. (2021). Water Resources Research, 57(1). https://doi.org/10.1029/2020WR028598; Hartmann, et al. (2014). Water Resources Research, 50(8), 6507–6521. https://doi.org/10.1002/2014WR015685 and the dataset publicly available via https://ral.ucar.edu/solutions/products/camels. The Matlab code on using our approach to calculate the likelihood is provided in the supplement.

Author contribution

YL conceptualized the study, developed and applied the adapted KGE approach, and visualized the results. YL and JFO wrote the paper, and analyzed the results. MM and AH provided supervision and advice throughout developing the adapted KGE approach and supported the development of this manuscript. All authors contributed to the revision of the manuscript.

Competing interests

The authors declare that they have no conflict of interest.

460 Acknowledgement

We thank Jasper A. Vrugt for his constructive comments and suggestions. We thank the editor, Jiangjiang Zhang and another anonymous reviewer for their valuable comments during the peer-review phase. Yan Liu and Andreas Hartmann were supported by the Emmy-Noether-Programme of the German Research Foundation (DFG, grant number: HA 8113/1-1, project "Global Assessment of Water Stress in Karst Regions in a Changing World"). Jaime Fernández-Ortega and Matías Mudarra were supported by the European Project "Karst Aquifer Resources availability and quality in the Mediterranean Area (KARMA)" PRIMA, ANR-18-PRIM-0005 (PCI2019-103675), and by the project PID2019-111759RB-I00 funded by the Spanish Research Agency. Additionally, it is a contribution to the Research Group RNM-308 of Junta de Andalucía. Jaime Fernández-Ortega was also supported by the Erasmus+ Programme of the European Commission.

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