Package 'soilhypfit'

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Title Modelling of Soil Water Retention and Hydraulic Conductivity Data

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Depends nloptr(>= 1.2.1), snowfall

Imports graphics, mgcv, quadprog(>= 1.5-7), parallel, Rmpfr(>= 0.7-2), stats, SoilHyP(>= 0.1.3), utils

Suggests lattice

SystemRequirements gmp (>= 4.2.3), mpfr (>= 3.0.0)

SystemRequirementsNote 'MPFR' (MP Floating-Point Reliable Library, http://mpfr.org/) and 'GMP' (GNU Multiple Precision library, http://gmplib.org/) are required for Rmpfr

Description Provides functions for efficiently estimating properties of the Van Genuchten-Mualem model for soil hydraulic parameters from possibly sparse soil water retention and hydraulic conductivity data by multi-response parameter estimation methods (Stewart, W.E., Caracotsios, M. Soerensen, J.P. (1992) ``Parameter estimation from multi-response data" <doi:10.1002/aic.690380502>). Parameter estimation is simplified by exploiting the fact that residual and saturated water contents and saturated conductivity are conditionally linear parameters (Bates, D. M. and Watts, D. G. (1988) ``Nonlinear Regression Analysis and Its Applications" <doi:10.1002/9780470316757>). Estimated parameters are optionally constrained by the evaporation characteristic length (Lehmann, P., Bickel, S., Wei, Z. and Or, D. (2020) ``Physical Constraints for Improved Soil Hydraulic Parameter Estimation by Pedotransfer Functions" <doi:10.1029/2019WR025963>) to ensure that the estimated parameters are physically valid. Common S3 methods and further utility functions allow to process, explore and visualise estimation results.

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Contents

`soilhypfit-package` *The soilhypfit Package*

Description

This is a summary of the features and functionality of **soilhypfit**, a package in R for parameteric modelling of soil water retention and hydraulic conductivity data.

Details

Estimation approach:

The main function, `fit_wrc_hcc`, estimates parameters of models for soil water retention and hydraulic conductivity by *maximum likelihood* (ml, default), *maximum posterior density* (mpd) *estimation* (*Stewart et al., 1992*) or nonlinear *weighted least squares* (wls) from data on volumetric *soil water content*, $\boldsymbol{\theta}^{\mathrm{T}} = (\theta_1, \theta_2, ..., \theta_{n_\theta})$, and/or *soil hydraulic conductivity*, $\boldsymbol{K}^{\mathrm{T}} = (K_1, K_2, ..., K_{n_K})$, both measured at given capillary pressure head, $\boldsymbol{h}^{\mathrm{T}} = (h_1, h_2, ...)$.

For mpd and ml estimation, the models for the measurements are

$$\theta_i = \theta(h_i; \boldsymbol{\mu}, \boldsymbol{\nu}) + \varepsilon_{\theta,i}, \quad i = 1, 2, ..., n_\theta,$$

$$\log(K_j) = \log(K(h_j; \boldsymbol{\mu}, \boldsymbol{\nu})) + \varepsilon_{K,j}, \quad j = 1, 2, ..., n_K,$$

where $\theta(h_i; \boldsymbol{\mu}, \boldsymbol{\nu})$ and $K(h_j; \boldsymbol{\mu}, \boldsymbol{\nu})$ denote modelled water content and conductivity, $\boldsymbol{\mu}$ and $\boldsymbol{\nu}$ are the *conditionally linear* and *nonlinear* model parameters (see below and *Bates and Watts, 1988*,

sec. 3.3.5), and $\varepsilon_{\theta,i}$ and $\varepsilon_{K,j}$ are independent, normally distributed errors with zero means and variances σ_θ^2 and σ_K^2, respectively.

Let

$$Q_\theta(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{\theta},\boldsymbol{h}) = (\boldsymbol{\theta} - \boldsymbol{\theta}(\boldsymbol{h};\boldsymbol{\mu},\boldsymbol{\nu}))^{\mathrm{T}}\, \boldsymbol{W}_\theta\, (\boldsymbol{\theta} - \boldsymbol{\theta}(\boldsymbol{h};\boldsymbol{\mu},\boldsymbol{\nu}))\,,$$

and

$$Q_K(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{K},\boldsymbol{h}) = (\log(\boldsymbol{K}) - \log(\boldsymbol{K}(\boldsymbol{h};\boldsymbol{\mu},\boldsymbol{\nu})))^{\mathrm{T}}\, \boldsymbol{W}_K\, (\log(\boldsymbol{K}) - \log(\boldsymbol{K}(\boldsymbol{h};\boldsymbol{\mu},\boldsymbol{\nu})))\,,$$

denote the (possibly weighted) residual sums of squares between measurements and modelled values. $\boldsymbol{\theta}(\boldsymbol{h};\boldsymbol{\mu},\boldsymbol{\nu}))$ and $\boldsymbol{K}(\boldsymbol{h};\boldsymbol{\mu},\boldsymbol{\nu})$ are vectors with modelled values of water content and hydraulic conductivity, and $\boldsymbol{W}_\theta$ and $\boldsymbol{W}_K$ are optional diagonal matrices with weights $w_{\theta,i}$ and $w_{K,j}$, respectively. The weights are products of *case weights* $w'_{\theta,i}$ and $w'_{K,j}$ and *variable weights* w_θ, w_K, hence $w_{\theta,i} = w_\theta\, w'_{\theta,i}$ and $w_{K,j} = w_K\, w'_{K,j}$.

The objective function for ml and mpd estimation is equal to (*Stewart et al., 1992*, eqs 14 and 15, respectively)

$$Q(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{\theta},\boldsymbol{K},\boldsymbol{h}) = \frac{\kappa_\theta}{2}\log(Q_\theta(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{\theta},\boldsymbol{h})) + \frac{\kappa_K}{2}\log(Q_K(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{K},\boldsymbol{h})),$$

where $\kappa_v = n_v + 2$ for mpd and $\kappa_v = n_v$ for ml estimation, $v \in (\theta, K)$, and weights $w_{\theta,i} = w_{K,j} = 1$. Note that $Q(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{\theta},\boldsymbol{K},\boldsymbol{h})$ is equal to the negative logarithm of the concentrated loglikelihood or the concentrated posterior density, obtained by replacing σ_θ^2 and σ_K^2 by their conditional maximum likelihood or maximum density estimates $\widehat{\sigma}_\theta^2(\boldsymbol{\mu}),\boldsymbol{\nu})$ and $\widehat{\sigma}_K^2(\boldsymbol{\mu})$ respectively (*Stewart et al., 1992*, p. 642).

For wls the objective function is equal to

$$Q(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{\theta},\boldsymbol{K},\boldsymbol{h}) = Q_\theta(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{\theta},\boldsymbol{h}) + Q_K(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{K},\boldsymbol{h}).$$

If either water content or conductivity data are not available, then the respective terms are omitted from $Q(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{\theta},\boldsymbol{K},\boldsymbol{h})$.

The function `fit_wrc_hcc` does not attempt to estimate the parameters by minimising $Q(\boldsymbol{\mu},\boldsymbol{\nu};\boldsymbol{\theta},\boldsymbol{K},\boldsymbol{h})$ directly with respect to $\boldsymbol{\mu}$ and $\boldsymbol{\nu}$. Rather, it exploits the fact that for *given nonlinear parameters* $\boldsymbol{\nu}$, the *conditionally linear parameters* $\boldsymbol{\mu}^T = (\theta_r, \theta_s, \log(K_0))$ can be estimated straightforwardly by minimising the *conditional residual sums of squares*

$$Q^*_\theta(\theta_r,\theta_s;\boldsymbol{\theta},\boldsymbol{h},\boldsymbol{\nu}) = \left(\boldsymbol{\theta} - [\mathbf{1},\boldsymbol{S}(\boldsymbol{h};\boldsymbol{\nu})]\begin{bmatrix}\theta_r\\ \theta_s-\theta_r\end{bmatrix}\right)^{\mathrm{T}}\boldsymbol{W}_\theta\left(\boldsymbol{\theta} - [\mathbf{1},\boldsymbol{S}(\boldsymbol{h};\boldsymbol{\nu})]\begin{bmatrix}\theta_r\\ \theta_s-\theta_r\end{bmatrix}\right)$$

with respect to θ_r and $\theta_s - \theta_r$ and/or

$$Q^*_K(K_0;\boldsymbol{K},\boldsymbol{h},\boldsymbol{\nu}) = (\log(\boldsymbol{K}) - \log(K_0\,\boldsymbol{k}(\boldsymbol{h};\boldsymbol{\nu})))^{\mathrm{T}}\,\boldsymbol{W}_K\,(\log(\boldsymbol{K}) - \log(K_0\,\boldsymbol{k}(\boldsymbol{h};\boldsymbol{\nu})))\,,$$

with respect to $\log(K_0)$, where $\mathbf{1}$ is a vector of ones, $\boldsymbol{S}(\boldsymbol{h};\boldsymbol{\nu})^{\mathrm{T}} = (S(h_1;\boldsymbol{\nu}),...,S(h_{n_\theta};\boldsymbol{\nu}))$ and $\boldsymbol{k}(\boldsymbol{h};\boldsymbol{\nu})^{\mathrm{T}} = (k(h_1;\boldsymbol{\nu}),...,k(h_{n_K};\boldsymbol{\nu}))$ are vectors of modelled *water saturation* and modelled *relative conductivity* values, θ_r and θ_s are the *residual* and *saturated water content*, and K_0 is the *saturated hydraulic conductivity*.

Unconstrained conditional estimates, say $\widehat{\theta}_r(\boldsymbol{\nu})$, $\widehat{\theta}_s(\boldsymbol{\nu}) - \widehat{\theta}_r(\boldsymbol{\nu})$ and $\widehat{\log(K_0)}(\boldsymbol{\nu})$ can be easily obtained from the normal equations of the respective (weighted) least squares problems, and

quadratic programming yields conditional (weighted) least squares estimates that honour the inequality constraints $0 \leq \theta_r \leq \theta_s \leq 1$.

Let $\widehat{\boldsymbol{\mu}}(\boldsymbol{\nu})^{\mathrm{T}} = (\widehat{\theta}_r(\boldsymbol{\nu}), \widehat{\theta}_s(\boldsymbol{\nu}), \widehat{\log(K_0)}(\boldsymbol{\nu}))$ be the conditional estimates of the linear parameters obtained by minimising $Q^*_\theta(\theta_r, \theta_s; \boldsymbol{\theta}, \boldsymbol{h}, \boldsymbol{\nu})$, and $Q^*_K(K_0; \boldsymbol{K}, \boldsymbol{h}, \boldsymbol{\nu})$, respectively. `fit_wrc_hcc` then estimates the nonlinear parameters by minimising $Q(\widehat{\boldsymbol{\mu}}(\boldsymbol{\nu}), \boldsymbol{\nu}; \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h})$ with respect to $\boldsymbol{\nu}$ by a nonlinear optimisation algorithm.

For mpd and ml estimation the variances of the model errors are estimated by (*Stewart et al., 1992*, eq. 16)

$$\widehat{\sigma}^2_\theta = \frac{Q_\theta(\widehat{\boldsymbol{\mu}}(\widehat{\boldsymbol{\nu}}), \widehat{\boldsymbol{\nu}}; \boldsymbol{\theta}, \boldsymbol{h})}{\kappa_\theta},$$

and

$$\widehat{\sigma}^2_K = \frac{Q_K(\widehat{\boldsymbol{\mu}}(\widehat{\boldsymbol{\nu}}), \widehat{\boldsymbol{\nu}}; \boldsymbol{K}, \boldsymbol{h})}{\kappa_K}.$$

Furthermore, for mpd and ml estimation, the covariance matrix of the estimated nonlinear parameters may be approximated by the inverse Hessian matrix of $Q(\widehat{\boldsymbol{\mu}}(\boldsymbol{\nu}), \boldsymbol{\nu}; \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h})$ at the solution $\widehat{\boldsymbol{\nu}}$ (*Stewart and Sørensen, 1981*), i.e.

$$\mathrm{Cov}[\widehat{\boldsymbol{\nu}}, \widehat{\boldsymbol{\nu}}^T] \approx \boldsymbol{A}^{-1},$$

where

$$[\boldsymbol{A}]_{kl} = \frac{\partial^2}{\partial \nu_k \, \partial \nu_l} \, Q(\widehat{\boldsymbol{\mu}}(\boldsymbol{\nu}), \boldsymbol{\nu}; \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h})|_{\nu=\hat{\nu}} \, .$$

Details on parameter estimation:

Models for water retention curves and hydraulic conductivity functions:
Currently, `fit_wrc_hcc` allows to estimate the parameters of the simplified form of the *Van Genuchten-Mualem* (VGM) model (Van Genuchten, 1980) with the restriction $m = 1 - \frac{1}{n}$, see `wc_model` and `hc_model`. This model has the following parameters:

- $\boldsymbol{\mu}^{\mathrm{T}} = (\theta_r, \theta_s, K_0)$ (see above) and
- $\boldsymbol{\nu}^{\mathrm{T}} = (\alpha, n, \tau)$ where α is the inverse air entry pressure, n the shape and τ the tortuosity parameter.

Any of these parameters can be either estimated from data or kept fixed at the specified initial values (see arguments `param` and `fit_param` of `fit_wrc_hcc`).

Imposing physical constraints on the estimated parameters:
Parameters of models for the water retention curve and the hydraulic conductivity function may vary only within certain bounds (see `wc_model`, `hc_model` and `param_boundf` for allowed ranges). `fit_wrc_hcc` either estimates *transformed parameters* that vary over the whole real line and can therefore be estimated without constraints (see `param_transf`), or it uses algorithms (quadratic programming for estimating $\boldsymbol{\mu}$, nonlinear optimisation algorithms with box constraints for estimating $\boldsymbol{\nu}$) that restrict estimates to permissible ranges, see *Details* section of `control_fit_wrc_hcc`.
In addition, for natural soils, the parameters of the VGM model cannot vary independently from each other over the allowed ranges. Sets of fitted parameters should always result in soil hydraulic quantities that are physically meaningful. One of these quantities is the *characteristic length* L_c of *stage-I* evaporation from a soil (Lehmann et al., 2008). L_c can be related to the parameters of the VGM model, see Lehmann et al. (2008, 2020) and `evaporative-length`.

Using several soil hydrological databases, Lehmann et al. (2020) analysed the mutual dependence of VGM parameters and proposed regression equations to relate the inverse air entry pressure α and the saturated hydraulic K_0 to the shape parameter n, which characterises the width of the pore size distribution of a soil. Using these relations, Lehmann et al. (2020) then computed the expected value ("target") L_t of L_c for given n and tortuosity parameter τ, see `evaporative-length`. `fit_wrc_hcc` allows to constrain estimates of the nonlinear parameters $\boldsymbol{\nu}$ by defining permissible lower and upper bounds for the ratio $L_\mathrm{c}/L_\mathrm{t}$, see arguments `ratio_lc_lt_bound` of `fit_wrc_hcc` and `settings` of `control_fit_wrc_hcc`.

Choice of optimisation algorithm for estimating the nonlinear parameters:
To estimate $\boldsymbol{\nu}$, `fit_wrc_hcc` minimises $Q(\widehat{\boldsymbol{\mu}}(\boldsymbol{\nu}), \boldsymbol{\nu}; \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h})$ either by a nonlinear optimisation algorithm available in the library *NLopt* (Johnson, see `nloptr`) or by the Shuffled Complex Evolution (SCE) optimisation algorithm (Duan et al., 1994, see `SCEoptim`). The choice of the algorithm is controlled by the argument `settings` of the function `control_fit_wrc_hcc`:

1. global optimisation without constraints for the ratio $L_\mathrm{c}/L_\mathrm{t}$ (`settings = "uglobal"` or `settings = "sce"`),
2. global optimisation with inequality constraints for the ratio $L_\mathrm{c}/L_\mathrm{t}$ (`settings = "cglobal"`),
3. local optimisation without constraints for the ratio $L_\mathrm{c}/L_\mathrm{t}$ (`settings = "ulocal"`),
4. local optimisation with inequality constraints for the ratio $L_\mathrm{c}/L_\mathrm{t}$ (`settings = "clocal"`).

The `settings` argument also sets reasonable default values for the termination (= convergence) criteria for the various algorithms, see NLopt documentation, section Termination conditions. The NLopt documentation contains a very useful discussion of (constrained) optimisation problems in general, global vs. local optimisation and gradient-based vs. derivative-free algorithms. Note that besides the `settings` argument of `control_fit_wrc_hcc`, the arguments `nloptr` and `sce` along with the functions `control_nloptr` and `control_sce` allow to fully control the nonlinear optimisation algorithms, see `control_fit_wrc_hcc` for details.

Computing initial values of parameters:
For local optimisation algorithms "good" initial values of $\boldsymbol{\nu}$ are indispensable for successful estimation. `fit_wrc_hcc` allows to compute initial values of α and n for the Van Genuchten model from water retention data by the following procedure:

1. Smooth the water retention data, $(\theta_i, y_i = \log(h_i)), i = 1, 2, ...n_\theta$, by an additive model.
2. Determine the saturation, S^*, and the logarithm of capillary pressure head, $y^* = \log(h^*)$, at the inflection point of the additive model fit.
3. Find the root, say $\widehat{m}$, of $S^* = (1+1/m)^{-m}$. One obtains the right-hand side of this equation by solving $\frac{\partial^2}{\partial y^2}\left[S_\mathrm{VG}(\exp(y); \boldsymbol{\nu})\right] = 0$ for y and plugging the result into the expression for $S_\mathrm{VG}(\exp(y); \boldsymbol{\nu})$, see `wc_model`.
4. Compute $\widehat{n} = 1/(1-\widehat{m})$ and $\widehat{\alpha} = 1/\exp(y^*)\,(1/\widehat{m})^{1-\widehat{m}}$. The second expression is again a result of solving $\frac{\partial^2}{\partial y^2}\left[S_\mathrm{VG}(\exp(y); \boldsymbol{\nu})\right] = 0$.

Initial values for local optimisation algorithms can of course also be obtained by first estimating the parameters by a global algorithm. These estimates can be "refined" in a second step by a local unconstrained algorithm, possibly followed by a third call of `fit_wrc_hcc` to constrain the estimated parameters by the ratio $L_\mathrm{c}/L_\mathrm{t}$. The method `coef.fit_wrc_hcc` can be used to extract the estimated parameters from an object of class `fit_wrc_hcc` and to pass them as initial values to `fit_wrc_hcc`, see `fit_wrc_hcc` for examples.

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Van Genuchten, M. Th. (1980) A closed-form equation for predicting the hydraulic conductivity of unsaturated soils. *Soil Science Society of America Journal*, **44**, 892–898, doi:10.2136/sssaj1980.03615995004400050002x.

See Also

fit_wrc_hcc for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

control_fit_wrc_hcc for options to control fit_wrc_hcc;

soilhypfitmethods for common S3 methods for class fit_wrc_hcc;

vcov for computing (co-)variances of the estimated nonlinear parameters;

prfloglik_sample for profile loglikelihood computations;

wc_model and hc_model for currently implemented models for soil water retention curves and hydraulic conductivity functions;

evaporative-length for physically constraining parameter estimates of soil hydraulic material functions.

control_fit_wrc_hcc *Controlling fit_wrc_hcc*

Description

This page documents options to control fit_wrc_hcc. It describes the arguments of the functions control_fit_wrc_hcc, param_boundf, param_transf, fwd_transf, dfwd_transf, bwd_transf, control_nloptr, control_sce and control_pcmp, which all serve to steer fit_wrc_hcc.

Usage

```
control_fit_wrc_hcc(
    settings = c("uglobal", "ulocal", "clocal", "cglobal", "sce"),
    method = c("ml", "mpd", "wls"), hessian,
    nloptr = control_nloptr(), sce = control_sce(),
    wrc_model = "vg", hcc_model = "vgm",
    initial_param = c(alpha = 2., n = 1.5, tau = 0.5),
    approximation_alpha_k0 =
        c(c0 = 1, c1 = 5.55, c2 = 1.204, c3 = 2.11, c4 = 1.71),
    variable_weight = c(wrc = 1, hcc = 1),
    gam_k = 6, gam_n_newdata = 101, precBits = 256,
    min_nobs_wc = 5, min_nobs_hc = 5,
    keep_empty_fits = FALSE,
    param_bound = param_boundf(), param_tf = param_transf(),
    fwd_tf = fwd_transf(), deriv_fwd_tfd = dfwd_transf(), bwd_tf = bwd_transf(),
    pcmp = control_pcmp())

param_boundf(alpha = c(0 + 10 * sqrt(.Machine$double.eps), 500),
    n = c(1 + 10 * sqrt(.Machine$double.eps), 20), tau = c(-2, 20),
    thetar = c(0, 1), thetas = c(0, 1), k0 = c(0, Inf))

param_transf(alpha = c("log", "identity"), n = c("log1", "identity"),
    tau = c("logitlu", "identity"), k0 = c("log", "identity"))

fwd_transf(...)

dfwd_transf(...)

bwd_transf(...)

control_nloptr(...)

control_sce(reltol = 1.e-8, maxtime = 20, ...)

control_pcmp(ncores = detectCores() - 1L,
    fork = !identical(.Platform[["OS.type"]], "windows"))
```

Arguments

settings	a keyword with possible values "uglobal" (default), "ulocal", "clocal", "cglobal" or "sce" to choose the approach for the nonlinear optimisation, see *Details* and soilhypfitIntro.
method	a keyword with possible values "ml" (maximum likelihood, default), "mpd" (maximum posterior density) or "wls" (weighted least squares) to choose the method for estimating the nonlinear parameters $\boldsymbol{\nu}$, see soilhypfitIntro.
hessian	a logical scalar controlling whether the Hessian matrix of the objective function should be computed with respect to the possibly transformed nonlinear parameters $\boldsymbol{\nu}$ at the solution (default: TRUE if settings %in% c("uglobal", "ulocal", "sce") && method %in% c("mpd", "ml") and FALSE otherwise).
nloptr	a list of arguments passed to the optimiser nloptr by its argument opts, or a function such as control_nloptr that generates such a list. Note that control_fit_wrc_hcc chooses sensible default values for the various components of the list in dependence of the value chosen for settings, but these defaults can be overridden by the arguments of control_nloptr, see *Details*.
sce	a list of arguments passed to the optimiser SCEoptim by its argument control, or a function such as control_sce that generates such a list. Note that control_fit_wrc_hcc chooses sensible default values for the various components of the list, but these defaults can be overridden by the arguments of control_sce, see *Details*.
wrc_model	a keyword choosing the parametrical model for the water retention curve. Currently, only the *Van Genuchten* model ("vg") is implemented, see wc_model and soilhypfitIntro.
hcc_model	a keyword choosing the parametrical model for the hydraulic conductivity function. Currently, only the *Van Genuchten-Mualem* model ("vgm") is implemented, see hc_model and soilhypfitIntro.
initial_param	a named numeric vector with default initial values for the *nonlinear* parameters $\boldsymbol{\nu}$ of the models for the water retention curve and/or hydraulic conductivity function. Currently, initial values must be defined for the parameters α (default 1.5 [m^{-1}]), n (default 2 [-]) and τ (default 0.5 [-]), hence the elements of initial_param must be named "alpha", "n" and "tau".
approximation_alpha_k0	a named numeric vector with constants to approximate the parameter α and the saturated hydraulic conductivity K_0 when constraining the estimated *nonlinear* parameters $\boldsymbol{\nu}$ of the Van Genuchten-Mualem model by the *characteristic evaporative length* (*Lehmann et al., 2020*), see evaporative-length and soilhypfitIntro. For consistency with other quantities, the following units should be used for the constants: • c1: m^{-1}, • c3: $\mathrm{m\,d}^{-1}$. The remaining constants are dimensionless.
variable_weight	a named numeric vector of length 2 that defines the weights of water content and hydraulic conductivity measurements in the objective function for method

= "wls". If equal to 1 (default) the weights of the variables are equal to the inverse variances of the water content and (possibly log-transformed) hydraulic conductivity measurements. If different from 1, then the inverse variances are multiplied by variable_weight to get the variable weights w_θ and w_K, see fit_wrc_hcc and soilhypfitIntro. Note that for estimation methods mpd and ml the variable weights are equal to 1 but the case weights $w'_{\theta,i}$ and $w'_{K,j}$ may differ from 1.

gam_k — the dimension of the basis of the additive model for the water retention data when computing initial values of the parameters α and n, see s and soilhypfitIntro.

gam_n_newdata — the number of evaluation points of the additive model for the water retention data when computing initial values of the parameters α and n, see soilhypfitIntro.

precBits — an integer scalar defining the default precision (in bits) to be used in high-precision computations by mpfr, see *Details*.

min_nobs_wc — an integer scalar defining the minimum number of water content measurements per sample required for fitting a model to the water content data.

min_nobs_hc — an integer scalar defining the minimum number of hydraulic conductivity measurements per sample required for fitting a model to hydraulic conductivity data.

keep_empty_fits — a logical scalar controlling whether missing fitting results should be dropped for samples without any data or failed fits (FALSE, default) or kept (TRUE).

param_bound — a named list of numeric vectors of length 2 that define the allowed lower and upper bounds (box constraints) for the parameters of the models or a function such as param_boundf which generates this list, see *Details*.

param_tf — a named vector of keywords that define the transformations to be applied to the model parameters before estimation or a function such as param_transf, which generates this vector, see *Details*.

fwd_tf — a named list of invertible functions to be used to transform model parameters or a function such as fwd_transf, which generates this list, see *Details*.

deriv_fwd_tfd — a named list of functions corresponding to the first derivatives of the functions defined in fwd_tf or a function such as dfwd_transf, which generates this list, see *Details*.

bwd_tf — a named list of inverse functions corresponding the functions defined in fwd_tf or a function such as bwd_transf, which generates this list, see *Details*.

pcmp — a list to control parallel computations or a function such as control_pcmp that generates this list, see control_pcmp for allowed arguments.

alpha — either a numeric vector of length 2 defining the allowed lower and upper bounds for the nonlinear parameter α (param_boundf), or a keyword defining the transformation to be applied to the parameter α before estimation (param_transf), see *Details*.

n — either a numeric vector of length 2 defining the allowed lower and upper bounds for the nonlinear parameter n (param_boundf), or a keyword defining the transformation to be applied to the parameter n before estimation (param_transf), see *Details*.

tau	either a numeric vector of length 2 defining the allowed lower and upper bounds for the nonlinear parameter τ (param_boundf), or a keyword defining the transformation to be applied to the parameter τ before estimation (param_transf), see *Details*.
thetar	a numeric vector of length 2 defining the allowed lower and upper bounds for the linear parameter θ_r (param_boundf), see *Details*.
thetas	a numeric vector of length 2 defining the allowed lower and upper bounds for the linear parameter θ_s (param_boundf), see *Details*.
k0	either a numeric vector of length 2 defining the allowed lower and upper bounds for the linear parameter K_0 (param_boundf), or a keyword defining the transformation to be applied to the parameter K_0 before estimation (param_transf), see *Details*.
reltol	a numeric scalar defining (one possible) convergence criterion for the optimiser SCEoptim, see argument control of SCEoptim for details.
maxtime	a numeric scalar defining the maximum duration of optimisation in seconds by the optimiser SCEoptim, see see argument control of SCEoptim for details.
ncores	an integer defining the number of cores for parallel computations. Defaults to the number of available cores minus one. ncores = 1 suppresses parallel computations.
fork	a logical scalar controlling whether forking should be used for parallel computations (default: TRUE on Unix and MacOS and FALSE on Windows operating systems). Note that stetting fork = TRUE on Windows suppresses parallel computations.
...	further arguments, such as details on parameter transformations (fwd_transf, dfwd_transf, bwd_transf) or control options passed to the optimisers nloptr and SCEoptim, see *Details*.

Details

Enforcing bounds on the estimated parameters:

Parameters of models for the water retention curve and the hydraulic conductivity function may vary only within certain bounds (see param_boundf for allowed ranges). fit_wrc_hcc uses two mechanisms to constrain parameter estimates to permissible ranges:

1. *Parameter transformations*
   If a local algorithm is used for nonlinear optimisation (settings = "ulocal" or settings = "clocal") and a transformation not equal to "identity" is specified in param_tf for any of the *nonlinear* parameters $\boldsymbol{\nu}$, then the elements of $\boldsymbol{\nu}$ are transformed by the functions given in param_tf. The values of the transformed parameters vary then over the whole real line, and an unconstrained algorithm can be used for nonlinear optimisation.
   Note that the *linear* parameters θ_r (residual) and θ_s (saturated water content) are never transformed and for the saturated hydraulic conductivity, K_0, only "log" (default) or "identity" can be chosen. Quadratic programming (see solve.QP) is employed to enforce the box constraints specified in the argument param_bound for θ_r and θ_s. Quadratic programming is also used to enforce the positivity constraint on K_0 if K_0 is not log-transformed ("identity"). Otherwise, the logarithm of K_0 is estimated unconstrainedly, see soilhypfitIntro for further details.

2. *Box constraints*
   If a global algorithm is used for the optimisation (`settings` equal to `"uglobal"` `"cglobal"` or `"sce"`) or if `"identity"` transformations are specified for all elements of $\boldsymbol{\nu}$, then an optimisation algorithm is deployed that respects the box constraints given in `param_bound`. If parameters are estimated for several soil samples in a single call of `fit_wrc_hcc` and if sample-specific box constraints should be used then the lower and upper bounds of the box-constraints must be specified by the arguments `lower_param` and `upper_param` of the function `fit_wrc_hcc`, see explanations there.
   Further note that the transformations specified by `param_tf` for the nonlinear parameters $\boldsymbol{\nu}$ are ignored when a global optimisation algorithm is used.

Parameter transformations:

The arguments `param_tf`, `fwd_tf`, `deriv_fwd_tfd`, `bwd_tf` define how the model parameters are transformed for estimation by local optimisation algortihms (see above and `soilhypfitIntro`). The following transformations are currently available:

`"log"`: $\log(x)$,

`"log1"`: $\log(x-1)$,

`"logitlu"`: $\log((x-l)/(u-x))$ with l and u the allowed lower and upper bounds for a parameter, see `param_boundf`,

`"identity"`: no transformation.

These are the possible values that the various arguments of the function `param_transf` accept (as quoted character strings), and these are the names of the list components returned by `fwd_transf`, `dfwd_transf` and `bwd_transf`.

Additional transformations can be implemented by:

1. Extending the function definitions by arguments like
   `fwd_tf = fwd_transf(my_fun = function(x) your transformation)`,
   `deriv_fwd_tfd = dfwd_transf(my_fun = function(x) your derivative)`,
   `bwd_tf = bwd_transf(my_fun = function(x) your back-transformation)`,
2. Assigning to a given argument of `param_transf` the name of the new function, e.g. `alpha = "my_fun"`.

Note that the values given to the arguments of `param_transf` must match the names of the functions returned by `fwd_transf`, `dfwd_transf` and `bwd_transf`.

High-precision numerical computations:

Estimation of $\log(K_0)$ is somewhat delicate for large values of the shape parameter n and/or small values of α. The water saturation and the relative conductivity are then close to zero for capillary pressure head exceeding $1/\alpha$. To avoid numerical problems caused by limited accuracy of double precision numbers, `fit_wrc_hcc` uses the function `mpfr` of the package **Rmpfr** for high-accuracy computations. The argument `precBits` of `control_fit_wrc_hcc` controls the accuracy. Increase its value if computation fails with a respective diagnostic message.

Options to choose the approach for nonlinear optimisation:

The argument `settings` defines sets of default options to control the optimisers. The following `settings` are currently defined:

`"uglobal"`: unconstrained optimisation by any of the global algorithms (named `"NLOPT_G..."`) of the NLopt library.

"cglobal": constrained optimisation by the global algorithm "NLOPT_GN_ISRES" of NLopt that allows for inequality constraints.

"ulocal": unconstrained optimisation by any of the local algorithms (named "NLOPT_L...") of NLopt.

"clocal": constrained optimisation by any of the local algorithms ("NLOPT_LN_COBYLA", "NLOPT_LN_AUGLAG", "NLOPT_LD_AUGLAG", "NLOPT_LD_SLSQP", "NLOPT_LD_MMA"), "NLOPT_LD_CCSAQ") of NLopt that allow for inequality constraints.

"sce": unconstrained optimisation by the global algorithm implemented in SCEoptim.

The functions control_nloptr and control_sce allow finer control of the optimisers. control_nloptr and control_sce take any argument available to pass controlling options to the optimisers nloptr (by its argument opts) and SCEoptim (by its argument control), respectively.

Controlling nloptr:
The function nloptr.print.options prints all options available to control nloptr by its argument opts. Detailed information on the options can be found in the NLopt documentation.

The function control_fit_wrc_hcc sets meaningful defaults for opts in dependence of the chosen optimisation approach as specified by the argument settings, and it checks the consistency of the arguments of control_nloptr if they are explicitly passed to fit_wrc_hcc.

The following defaults are set by control_fit_wrc_hcc for the argument opts of nloptr (:

1. Unconstrained, global optimisation (settings = "uglobal"):

```
nloptr = control_nloptr(
  algorithm = "NLOPT_GN_MLSL_LDS",
  local_opts = list(
    algorithm = "NLOPT_LN_BOBYQA",
    xtol_rel = -1.,
    ftol_rel = 1.e-6
  ),
  xtol_rel = -1,
  ftol_rel = -1,
  maxeval = 125,
  maxtime = -1)
```

   In addition, any parameter transformations specified by param_tf are overridden and the untransformed parameters ("identity") are estimated when settings = "uglobal" is chosen.

2. Constrained, global optimisation (settings = "cglobal"):

```
nloptr = control_nloptr(
  algorithm = "NLOPT_GN_ISRES",
  xtol_rel = -1,
  ftol_rel = -1,
  maxeval = 1000,
  maxtime = -1)
```

   In addition, any parameter transformations specified by param_tf are overridden and the untransformed parameters ("identity") are estimated when settings = "cglobal" is chosen.

3. Unconstrained, local optimisation (`settings = "ulocal"`):

```
nloptr = control_nloptr(
  algorithm = "NLOPT_LN_BOBYQA",
  xtol_rel = -1,
  ftol_rel = 1.e-8,
  maxeval = 250,
  maxtime = -1)
```

4. Constrained, local optimisation (`settings = "clocal"`):

```
nloptr = control_nloptr(
  algorithm = "NLOPT_LD_CCSAQ",
  xtol_rel = -1,
  ftol_rel = 1.e-8,
  maxeval = 1000,
  maxtime = -1)
```

If the algorithm `"NLOPT_LD_AUGLAG"` is used for constrained, local optimisation then

```
nloptr = control_nloptr(
  algorithm = "NLOPT_LD_AUGLAG",
  local_opts = list(
    algorithm = "NLOPT_LD_LBFGS",
    xtol_rel = -1.,
    ftol_rel = 1.e-6
  ),
  xtol_rel = -1,
  ftol_rel = 1.e-8,
  maxeval = 1000,
  maxtime = -1)
```

For other, unspecified elements of `opts` default values as listed by `nloptr.print.options` are used.

Controlling SCEoptim:
The function `control_sce` sets meaningful defaults for the argument `control` of `SCEoptim`. Currently, the following defaults are defined:

```
sce = control_sce(
  reltol = 1e-08,
  maxtime = 20)
```

In addition, any parameter transformations specified by `param_tf` are overridden and the untransformed parameters (`"identity"`) are estimated when `settings = "sce"` is chosen.

Value

`control_fit_wrc_hcc` creates a list with components `settings`, `hessian`, `method`, `nloptr`, `sce`, `wrc_model`, `hcc_model`, `initial_param`, `approximation_alpha_k0`, `variable_weight`, `gam_k`, `gam_n_newdata`, `precBits`, `min_nobs_wc`, `min_nobs_hc`, `keep_empty_fits`, `param_bound`, `param_tf`,

fwd_tf, deriv_fwd_tfd, bwd_tf, pcmp corresponding to its arguments and some further components (delta_sat_0, grad_eps, use_derivative) that cannot be changed by the user.

control_nloptr and control_sce create lists with control parameters passed to nloptr and SCEoptim, respectively, see *Details*.

param_boundf generates a list with allowed lower and upper bounds of the model parameters.

param_transf generates a list with keywords that define what transformations are used for estimating the model parameters, and fwd_transf, bwd_transf and dfwd_transf return lists of functions with forward and backward transformations and the first derivatives of the forward transformations, see *Details*.

control_pcmp generates a list with control parameters for parallel computations.

Author(s)

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References

Johnson, S.G. The NLopt nonlinear-optimisation package. https://github.com/stevengj/nlopt.

Lehmann, P., Assouline, S., Or, D. (2008) Characteristic lengths affecting evaporative drying of porous media. *Physical Review E*, **77**, 056309, doi:10.1103/PhysRevE.77.056309.

Lehmann, P., Bickel, S., Wei, Z., Or, D. (2020) Physical Constraints for Improved Soil Hydraulic Parameter Estimation by Pedotransfer Functions. *Water Resources Research* **56**, e2019WR025963, doi:10.1029/2019WR025963.

See Also

soilhypfitIntro for a description of the models and a brief summary of the parameter estimation approach;

fit_wrc_hcc for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

soilhypfitmethods for common S3 methods for class fit_wrc_hcc;

vcov for computing (co-)variances of the estimated nonlinear parameters;

prfloglik_sample for profile loglikelihood computations;

wc_model and hc_model for currently implemented models for soil water retention curves and hydraulic conductivity functions;

evaporative-length for physically constraining parameter estimates of soil hydraulic material functions.

Examples

```
# use of \donttest{} because execution time exceeds 5 seconds
data(sim_wrc_hcc)

# estimate parameters for a single soil sample by maximizing loglikelihood ...

# ... with unconstrained, global optimisation algorithm NLOPT_GN_MLSL
```

```
coef(
  fit1 <- fit_wrc_hcc(
    wrc_formula = wc ~ head, hcc_formula = hc ~ head,
    data = sim_wrc_hcc, subset = id == 2
  ), gof = TRUE)

# ... as fit1 but fitting parameter tau as well
coef(
  fit2 <- update(fit1,
    fit_param = default_fit_param(tau = TRUE)
  ), gof = TRUE)

plot(fit1, y = fit2)

# ... with unconstrained, local optimisation algorithm NLOPT_LN_BOBYQA,
#     initial values for alpha and n are computed from data and
#     transformed nonlinear parameters are estimated without box-constraints
coef(
  fit3 <- update(
    fit2,
    control = control_fit_wrc_hcc(settings = "ulocal"),
    verbose = 2), gof = TRUE)

# estimate parameters by unconstrained, weighted least squares minimisation with
#     algorithm NLOPT_LD_LBFGS, giving larger weight to conductivity data,
#     using specified initial values for alpha and n and
#     fitting untransformed nonlinear parameters with default box constraints
#     defined by param_boundf()
#     diagnostic output directly from nloptr
coef(
  fit4 <- update(
    fit2,
    param = c(alpha = 1.7, n = 2),
    control = control_fit_wrc_hcc(
      settings = "ulocal", method = "wls",
      variable_weight = c(wrc = 1, hcc = 2),
      nloptr = control_nloptr(algorithm = "NLOPT_LD_LBFGS", print_level = 3),
      param_tf = param_transf(alpha = "identity", n = "identity", tau = "identity")
    ), verbose = 0), gof = TRUE)

# ... as fit4 but giving even larger weight to conductivity data
coef(
  fit5 <- update(
    fit4,
    control = control_fit_wrc_hcc(
      settings = "ulocal", method = "wls",
      variable_weight = c(wrc = 1, hcc = 5),
      nloptr = control_nloptr(algorithm = "NLOPT_LD_LBFGS", print_level = 3),
      param_tf = param_transf(alpha = "identity", n = "identity", tau = "identity")
    ), verbose = 0), gof = TRUE)

plot(fit4, y = fit5)
```

`evaporative-length` *Evaporative Characteristic Length*

Description

The functions `lc` and `lt` compute the *characteristic length* L_c of *stage-I* evaporation from a soil and its "target" (expected) value L_t, used to constrain estimates of nonlinear parameters of the Van Genuchten-Mualem (VGM) model for water retention curves and hydraulic conductivity functions.

Usage

```
lc(alpha, n, tau, k0, e0, c0 = NULL, c3 = NULL, c4 = NULL)

lt(n, tau, e0, c0, c1, c2, c3, c4)
```

Arguments

`alpha`	parameter α (inverse air entry pressure) of the VGM model, see `wc_model` and `hc_model`. For consistency with other quantities, the unit of α should be **1/meter** $[\mathrm{m}^{-1}]$.
`n`	parameter n (shape parameter) of the VGM model, see `wc_model` and `hc_model`.
`tau`	parameter τ (tortuosity parameter) of the VGM model, see `hc_model`.
`k0`	saturated hydraulic conductivity K_0, see `hc_model`. If `k0` is missing or equal to `NA` in calls of `lc` then `k0` is approximated by the same relation as used for `lt`, see *Details*. For consistency with other quantities, the unit of K_0 should be **meter/day** $[\mathrm{m\,d}^{-1}]$.
`e0`	a numeric scalar with the *stage-I* rate of evaporation E_0 from a soil, see *Details* and `soilhypfitIntro`. For consistency with other quantities, the unit of E_0 should be **meter/day** $[\mathrm{m\,d}^{-1}]$.
`c0`, `c1`, `c2`, `c3`, `c4`	numeric constants to approximate the parameter α and the saturated hydraulic conductivity K_0 when computing L_t, see *Details* and `control_fit_wrc_hcc`. For consistency with other quantities, the following units should be used for the constants: • `c1`: m^{-1}, • `c3`: $\mathrm{m\,d}^{-1}$. The remaining constants are dimensionless.

Details

The *characteristic length* of *stage-I evaporation* L_c (*Lehmann et al., 2008, 2020*) is defined by

$$L_c(\boldsymbol{\nu}, K_0, E_0) = \frac{\frac{1}{\alpha\, n}\left(\frac{2n-1}{n-1}\right)^{\frac{2n-1}{n}}}{1 + \frac{E_0}{K_{\mathrm{eff}}}}$$

where $\boldsymbol{\nu}^{\mathrm{T}} = (\alpha, n, \tau)$ are the nonlinear parameters of the VGM model, K_0 is the saturated hydraulic conductivity, E_0 the stage-I evaporation rate and $K_{\mathrm{eff}} = 4\,K_{\mathrm{VGM}}(h_{\mathrm{crit}}; K_0, \boldsymbol{\nu})$ is the effective hydraulic conductivity at the critical pressure

$$h_{\mathrm{crit}} = \frac{1}{\alpha}\left(\frac{n-1}{n}\right)^{\frac{1-2n}{n}},$$

see `hc_model` for the definition of $K_{\mathrm{VGM}}(h; K_0, \boldsymbol{\nu})$.

The quantity L_{t} is the expected value ("target") of L_{c} for given shape (n) and tortuosity (τ) parameters. To evaluate L_{t}, the parameters α and K_0 are approximated by the following relations

$$\widehat{\alpha} = g_{\alpha}(n; c_0, c_1, c_2) = c_1\,\frac{n - c_0}{1 + c_2\,(n - c_0)},$$

$$\widehat{K}_0 = g_{K_0}(n; c_0, c_3, c_4) = c_3\,(n - c_0)^{c_4}.$$

The default values for c_0 to c_4 (see argument `approximation_alpha_k0` of `control_fit_wrc_hcc`) were estimated with data on African desert regions of the database *ROSETTA3* (*Zhang and Schaap, 2017*), see *Lehmann et al. (2020)* for details.

Value

A numeric scalar with the characteristic evaporative length (`lc`) or its expected value (`lt`).

Author(s)

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References

Lehmann, P., Assouline, S., Or, D. (2008) Characteristic lengths affecting evaporative drying of porous media. *Physical Review E*, **77**, 056309, doi:10.1103/PhysRevE.77.056309.

Lehmann, P., Bickel, S., Wei, Z., Or, D. (2020) Physical Constraints for Improved Soil Hydraulic Parameter Estimation by Pedotransfer Functions. *Water Resources Research* **56**, e2019WR025963, doi:10.1029/2019WR025963.

Zhang, Y. , Schaap, M. G. 2017. Weighted recalibration of the Rosetta pedotransfer model with improved estimates of hydraulic parameter distributions and summary statistics (Rosetta3). *Journal of Hydrology*, **547**, 39-53, doi:10.1016/j.jhydrol.2017.01.004.

See Also

`soilhypfitIntro` for a description of the models and a brief summary of the parameter estimation approach;

`fit_wrc_hcc` for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

`control_fit_wrc_hcc` for options to control `fit_wrc_hcc`;

`soilhypfitmethods` for common S3 methods for class `fit_wrc_hcc`;

`vcov` for computing (co-)variances of the estimated nonlinear parameters;

prfloglik_sample for profile loglikelihood computations;

wc_model and hc_model for currently implemented models for soil water retention curves and hydraulic conductivity functions;

Examples

```
# use of \donttest{} because execution time exceeds 5 seconds

# estimate parameters of 4 samples of the Swiss forest soil dataset
# that have water retention (theta, all samples), saturated hydraulic conductivity
# (ksat) and optionally unsaturated hydraulic conductivity data
# (ku, samples "CH2_4" and "CH3_1")

data(swissforestsoils)

# select subset of data
sfs_subset <- droplevels(
  subset(
    swissforestsoils,
    layer_id %in% c("CH2_3", "CH2_4", "CH2_6", "CH3_1")
  ))

# extract ksat measurements
ksat <- sfs_subset[!duplicated(sfs_subset$layer_id), "ksat", drop = FALSE]
rownames(ksat) <- levels(sfs_subset$layer_id)
colnames(ksat) <- "k0"

# define number of cores for parallel computations
if(interactive()) ncpu <- parallel::detectCores() - 1L else ncpu <- 1L


# unconstrained estimation (global optimisation algorithm NLOPT_GN_MLSL)
# k0 fixed at measured ksat values
rsfs_uglob <- fit_wrc_hcc(
  wrc_formula = theta ~ head | layer_id,
  hcc_formula = ku ~ head | layer_id,
  data = sfs_subset,
  param = ksat,
  fit_param = default_fit_param(k0 = FALSE),
  control = control_fit_wrc_hcc(
    settings = "uglobal", pcmp = control_pcmp(ncores = ncpu)))
summary(rsfs_uglob)
coef(rsfs_uglob, lc = TRUE, gof = TRUE)


# constrained estimation by restricting ratio Lc/Lt to [0.5, 2]
# (global constrained optimisation algorithm NLOPT_GN_MLSL)
# k0 fixed at measured ksat values
rsfs_cglob <- update(
  rsfs_uglob,
  control = control_fit_wrc_hcc(
    settings = "cglobal", nloptr = control_nloptr(ranseed = 1),
```

```
    pcmp = control_pcmp(ncores = ncpu)))
summary(rsfs_cglob)
coef(rsfs_cglob, lc = TRUE, gof = TRUE)

# get initial parameter values from rsfs_cglob
ini_param <- cbind(
  coef(rsfs_cglob)[, c("alpha", "n")],
  ksat
)
# constrained estimation by restricting ratio Lc/Lt to [0.5, 2]
# (local constrained optimisation algorithm NLOPT_LD_CCSAQ)
# k0 fixed at measured ksat values
rsfs_cloc <- update(
  rsfs_uglob,
  param = ini_param,
  control = control_fit_wrc_hcc(
    settings = "clocal", nloptr = control_nloptr(ranseed = 1),
    pcmp = control_pcmp(ncores = ncpu)))
summary(rsfs_cloc)
coef(rsfs_cloc, lc = TRUE, gof = TRUE)

op <- par(mfrow = c(4, 2))
plot(rsfs_uglob, y = rsfs_cglob)
on.exit(par(op))

op <- par(mfrow = c(4, 2))
plot(rsfs_uglob, y = rsfs_cloc)
on.exit(par(op))
```

`fit_wrc_hcc` *Parametric Modelling of Soil Hydraulic Properties*

Description

The function `fit_wrc_hcc` estimates parameters of models for the soil water retention curve and/or soil hydraulic conductivity function from respective measurements by nonlinear regression methods, optionally subject to physical constraints on the estimated parameters. `fit_wrc_hcc` uses optimisation algorithms of the NLopt library (Johnson, see `nloptr-package`) and the Shuffled Complex Evolution (SCE) algorithm (Duan et al., 1994) implemented in the function `SCEoptim`.

Usage

```
fit_wrc_hcc(
    wrc_formula, hcc_formula, data,
    subset = NULL, wrc_subset = subset, hcc_subset = subset,
    weights = NULL, wrc_weights = weights, hcc_weights = weights,
    na.action, param = NULL, lower_param = NULL, upper_param = NULL,
    fit_param = default_fit_param(),
```

```
    e0 = 2.5e-3, ratio_lc_lt_bound = c(lower = 0.5, upper = 2),
    control = control_fit_wrc_hcc(), verbose = 0)

default_fit_param(
    alpha = TRUE, n = TRUE, tau = FALSE,
    thetar = TRUE, thetas = TRUE, k0 = TRUE)
```

Arguments

wrc_formula	an optional two-sided formula such as wc ~ head or wc ~ head | id, specifying the variables for the water content (wc), the capillary pressure head and, optionally, for sample ids when model parameters are estimated for several soil samples at the same time, see formula and *Details*.
hcc_formula	an optional two-sided formula such as hc ~ head or hc ~ head | id, specifying the variables for the hydraulic conductivity (hc), the capillary pressure head and, optionally, for sample ids when model parameters are estimated for several soil samples at the same time. See formula and *Details*.
data	a mandatory data frame containing the variables specified in the formula, the subset and weights arguments.
subset	an optional expression generating a vector to choose a subset of water content and hydraulic conductivity data. The expression is evaluated in the environment generated by model.frame(wrc_formula, data) and model.frame(hcc_formula, data), respectively.
wrc_subset	an optional expression generating a vector to choose a subset of water content data. The expression is evaluated in the environment generated by model.frame(wrc_formula, data). Defaults to subset.
hcc_subset	an optional expression generating a vector to choose a subset of hydraulic conductivity data. The expression is evaluated in the environment generated by model.frame(hcc_formula, data). Defaults to subset.
weights	an optional expression generating a numeric vector of case weights $w'_{\theta,i}$ and $w'_{K,i}$ (default: 1, see soilhypfitIntro) for water content and hydraulic conductivity data. The expression is evaluated in the environment generated by model.frame(wrc_formula, data) and model.frame(hcc_formula, data), respectively.
wrc_weights	an optional expression generating a numeric vector of case weights $w'_{\theta,i}$ (see soilhypfitIntro) for water content data. The expression is evaluated in the environment generated by model.frame(wrc_formula, data). Defaults to weights.
hcc_weights	an optional expression generating a numeric vector of case weights $w'_{K,i}$ (see soilhypfitIntro) for hydraulic conductivity data. The expression is evaluated in the environment generated by model.frame(hcc_formula, data). Defaults to weights.
na.action	a function which indicates what should happen when the data contain NAs. The default is set by the na.action argument of options, and is na.fail if that is unset. The "factory-fresh" default is na.omit. Another possible value is NULL, no action. Value na.exclude can be useful.

`param` an optional named numeric vector (or a numeric matrix or a dataframe with specified row and column names, see *Details*) with initial values for the model parameters. Currently, `param` may have elements (or columns) named `"alpha"`, `"n"`, `"tau"`, `"thetar"`, `"thetas"`, `"k0"`, see `wc_model` and `hc_model` and *Details*. For consistency with other quantities, the unit of α should be **1/meter** [m^{-1}] and the unit of K_0 **meter/day** [$\mathrm{m\,d}^{-1}$].

`lower_param` an optional named numeric vector (or a numeric matrix or a dataframe with specified row and column names, see *Details*) with lower bounds for the parameters of the models. Currently, `lower_param` may have elements (or columns) named `"alpha"`, `"n"`, `"tau"`, `"thetar"`, `"thetas"`, `"k0"`, see `wc_model`, `hc_model` and `param_boundf`. For consistency with other quantities, the unit of α should be **1/meter** [m^{-1}] and the unit of K_0 **meter/day** [$\mathrm{m\,d}^{-1}$]. If lower bounds are specified for θ_r but not for θ_s then the lower bounds specified for θ_r will also be used for θ_s.

`upper_param` an optional named numeric vector (or a numeric matrix or a dataframe with specified row and column names, see *Details*) with upper bounds for the parameters of the models. Currently, `upper_param` may have elements (or columns) named `"alpha"`, `"n"`, `"tau"`, `"thetar"`, `"thetas"`, `"k0"`, see `wc_model`, `hc_model` and `param_boundf`. For consistency with other quantities, the unit of α should be **1/meter** [m^{-1}] and the unit of K_0 **meter/day** [$\mathrm{m\,d}^{-1}$]. If upper bounds are specified for θ_s but not for θ_r then the upper bounds specified for θ_s will also be used for θ_r.

`fit_param` a named logical vector (or a logical matrix or a dataframe with specified row and column names, see *Details*) containing flags that control whether model parameters are estimated (`TRUE`) or kept fixed at the initial values (`FALSE`) as specified in `param`. This vector can be generated easily by the function `default_fit_param`. Currently, `fit_param` may have elements (or columns) named `"alpha"`, `"n"`, `"tau"`,
`"thetar"`, `"thetas"`, `"k0"`, see *Details*, `wc_model` and `hc_model`.

`e0` a numeric scalar (or named vector, see *Details*) with the stage-I rate of evaporation E_0 from a soil (default $2.5 \cdot 10^{-3}$ $\mathrm{m\,d}^{-1}$) required to evaluate the *characteristic evaporative length*, see `evaporative-length` and `soilhypfitIntro`. For consistency with other quantities, the unit of E_0 should **meter/day** [$\mathrm{m\,d}^{-1}$]. Note that `e0` is ignored when an unconstrained nonlinear optimisation algorithm is chosen (argument `settings` of `control_fit_wrc_hcc` equal to `"uglobal"`, `"sce"` or `"ulocal"`).

`ratio_lc_lt_bound`
a named numeric vector of length 2 (or a matrix with specified rownames and two columns, see *Details*) defining the default `lower` and `upper` bounds of the ratio $L_\mathrm{c}/L_\mathrm{t}$ (*Lehmann et al., 2008, 2020*) for constrained estimation of the *nonlinear* parameters $\boldsymbol{\nu}^\mathrm{T} = (\alpha, n, \tau)$ (default values 0.5 (`lower`) and 2 (`upper`)), see `evaporative-length` and `soilhypfitIntro`. Note that `ratio_lc_lt_bound` is ignored when an unconstrained nonlinear optimisation algorithm is chosen (argument `settings` of `control_fit_wrc_hcc` equal to `"uglobal"`, `"sce"` or `"ulocal"`).

`control` a list with options to control `fit_wrc_hcc` or a function such as `control_fit_wrc_hcc` that generates such a list. The main argument `settings`

of control_fit_wrc_hcc selects a nonlinear optimisation approach, see soilhypfitIntro and control_fit_wrc_hcc for full details.

verbose — positive integer controlling logging of diagnostic messages to the console and plotting of data and model curves during fitting.

verbose < 0 suppresses all output,

verbose >= 0 suppresses all output except warning messages,

verbose >= 1 displays at the end the measurements along with the fitted model curves,

verbose >= 2 prints for each iteration the parameters values, the value of the objective function and the ratio of L_c/L_t (see evaporative-length),

verbose >= 3 plots for each iteration the measurements along with the fitted model curves.

Logging of further diagnostics during fitting is controlled in addition by the arguments print_level, check_derivatives, check_derivatives_print when nloptr is used (see nloptr.print.options and control_nloptr) and by the argument trace of SCEoptim (see control_sce).

alpha — a logical scalar controlling whether the inverse air entry pressure α should be estimated (TRUE, default) or kept fixed at the initial value (FALSE) specified by param, see wc_model and hc_model.

n — a logical scalar controlling whether the shape parameter n should be estimated (TRUE, default) or kept fixed at the initial value (FALSE) specified by param, see wc_model and hc_model.

tau — a logical scalar controlling whether the tortuosity parameter τ should be estimated (TRUE) or kept fixed at the initial value (FALSE, default) specified by param, see hc_model.

thetar — a logical scalar controlling whether the residual water content θ_r should be estimated (TRUE, default) or kept fixed at the initial value (FALSE) specified by param, see wc_model.

thetas — a logical scalar controlling whether the saturated water content θ_s should be estimated (TRUE, default) or kept fixed at the initial value (FALSE) specified by param, see wc_model.

k0 — a logical scalar controlling whether the saturated hydraulic conductivity K_0 should be estimated (TRUE, default) or kept fixed at the initial value (FALSE) specified by param, see hc_model.

Details

Estimating model parameters of water retention curves and/or hydraulic conductivity functions:

The function fit_wrc_hcc estimates model parameters of water retention curves and/or hydraulic conductivity functions by *maximum likelihood* (ml, default) *maximum posterior density* (mpd) or *nonlinear least squares* (wls), see argument method of control_fit_wrc_hcc. Measurements must be available for at least one of the two material functions. If one type of data is missing then the respective formula, subset and weights arguments should be omitted.

If both types of data are available then measurements are weighed when computing the residual sum of squares for wls, but unit weights are used by default for mpd and ml estimation, see soilhypfitIntro. The wls weights are the product of the *case weights* $w'_{\theta,i}$ and $w'_{K,i}$ given by weights, wrc_weights and hcc_weights, respectively, and the *variable weights* as specified by the argument variable_weight of control_fit_wrc_hcc. Note that the variable_weights are not used "as is", but they are multiplied by the inverse variances of the water content or the (log-transformed) conductivity data per sample to obtain the variable weights w_θ, w_K mentioned in soilhypfitIntro.

Estimating model parameters for a single or multiple soil samples:

When parameters are estimated for a single soil sample only, then model formulae are of the form wc ~ head and/or hc ~head, and there is no need to specify a sample id. In this case param, lower_param, upper_param, fit_param and ratio_lc_lt_bound are vectors and e0 is a scalar.

fit_wrc_hcc allows to fit models to datasets containing data for multiple soil samples. The model formula must then be of the form wc ~ head|id and/or hc ~ head|id where the factor id identifies the soil samples. If param, lower_param, upper_param, fit_param and ratio_lc_lt_bound are vectors and e0 is a scalar then this information is used for processing all the soil samples. If specific information per sample should be used then param, lower_param, upper_param, fit_param and ratio_lc_lt_bound must be matrices (or data frames) with as many rows as there are soil samples. The matrices (or data.frames) must have rownames matching the levels of the factor id defining the soil samples. e0 must be a named vector with as many elements as there are soil samples and the names must again match the levels of id.

By default, fit_wrc_hcc processes data of multiple soil samples in parallel, see control_pcmp for options to control parallel computing. Note that diagnostic output (except warning messages) is suppressed in parallel computations. If computations fail for a particular soil sample, the id of the sample with the failed fit can be extracted by the utility function select_failed_fits and the respective error messages by extract_error_messages.

Controlling fit_wrc_hcc:

The argument control is used to pass a list of options to fit_wrc_hcc that steer the function, see soilhypfit-package and control_fit_wrc_hcc for full details.

Value

fit_wrc_hcc returns an object of class fit_wrc_hcc, which is a list with the following components:

fit — a list with as many components as there are soils samples. Each component is in turn a list that contains the estimated parameters and accompanying information:

- converged: an integer code issued by SCEoptim or nloptr on (non-) convergence, see convergence_message and NLopt return values.
- message: a character string issued by SCEoptim or nloptr on (non-)convergence.
- evaluation: the number of evaluations of the objective function and of the gradient.
- objective: the value of the objective function evaluated at the solution. The contributions of the water retention curve and the hydraulic conductivity function to objective are reported as attributes obj_wc and obj_hc. The attributes ssq_wc and ssq_hc are the respective residual sums of squares Q_θ and Q_K, see soilhypfitIntro.

- `gradient`: the gradient of the objective function at the solution with respect to the (possibly transformed) nonlinear parameters.
- `lp`: the estimated values for the linear parameters $\boldsymbol{\mu}^{\mathrm{T}} = (\theta_r, \theta_s, K_0)$, see `wc_model` and `hc_model`.
- `nlp`: the estimated values for the nonlinear parameters $\boldsymbol{\nu}^{\mathrm{T}} = (\alpha, n, \tau)$, see `wc_model` and `hc_model`.
- `inequality_constraint`: a list with the values of the inequality constraints evaluated at the solutions. Currently, values are reported for the expressions.

$$-(\frac{L_\mathrm{c}}{L_\mathrm{t}} - l)$$
$$\frac{L_\mathrm{c}}{L_\mathrm{t}} - u$$

  in the only component `lc`, where l and u are the lower and upper bounds of the ratio, see argument `ratio_lc_lt_bound` and `evaporative-length`. The values of L_c, L_t, the imposed bounds on their ratio and `e0` are returned as attributes of `lc`.
- `hessian`: optionally the Hessian matrix of the objective function with respect to the possibly transformed nonlinear parameters at the solution.
- `variable_weight`: a named vector with the variable weights w_θ and w_K, see *Details*.
- `weight`: a list with one or two components named `weights_wc` and or `weights_hc` with the case weights $w_{\theta,i}$ and $w_{K,i}$ used in the objective function, see *Details*.
- `fitted`: a list with one or two components named `wrc` and/or `hcc` with the fitted values for the water retention and the (possibly log-transformed) hydraulic conductivity data.
- `residuals`: a list with one or two components named `wrc` and/or `hcc` with residuals for the water retention and the (possibly log-transformed) hydraulic conductivity data.
- `model`: a list with one or two components named `wrc` and/or `hcc` with the `model.frame`s for the water retention and hydraulic conductivity data.
- `initial_objects`: a list with the values for `param`, `fit_param`, `variable_weight`, `param_bound`, `e0` and `ratio_lc_lt_bound` as taken from the (processed) arguments of the call of `fit_wrc_hcc`.

`sample_id_variable`
: the name of the variable defining the samples.

`wrc`
: logical scalar signalling whether water retention data were used to estimate the parameters.

`wrc_formula`
: formula defining the variables for water content and hydraulic head of the water retention data (`NULL` if not `wrc` equal to `FALSE`).

`wrc_mf`
: unevaluated call of `model.frame` to build the modelframe for the water retention data (`NULL` if `wrc` equal to `FALSE`).

`wrc`
: logical scalar signalling whether water retention data were used to estimate the parameters.

hcc_formula	formula defining the variables for hydraulic conductivity and hydraulic head of the hydraulic conductivity data (NULL if not hcc equal to FALSE).
hcc_mf	unevaluated call of model.frame to build the modelframe for the water retention data (NULL if hcc equal to FALSE).
control	a list with the options used to control fit_wrc_hcc, see control_fit_wrc_hcc.
call	the matched call.
na.action	information returned by model.frame on the special handling of NAs.

Author(s)

Andreas Papritz <papritz@retired.ethz.ch>.

References

Duan, Q., Sorooshian, S., and Gupta, V. K. (1994) Optimal use of the SCE-UA global optimisation method for calibrating watershed models, *Journal of Hydrology* **158**, 265–284, doi:10.1016/0022-1694(94)900574.

Johnson, S.G. The NLopt nonlinear-optimisation package. https://github.com/stevengj/nlopt.

Lehmann, P., Assouline, S., Or, D. (2008) Characteristic lengths affecting evaporative drying of porous media. *Physical Review E*, **77**, 056309, doi:10.1103/PhysRevE.77.056309.

Lehmann, P., Bickel, S., Wei, Z., Or, D. (2020) Physical Constraints for Improved Soil Hydraulic Parameter Estimation by Pedotransfer Functions. *Water Resources Research* **56**, e2019WR025963, doi:10.1029/2019WR025963.

See Also

soilhypfitIntro for a description of the models and a brief summary of the parameter estimation approach;

control_fit_wrc_hcc for options to control fit_wrc_hcc;

soilhypfitmethods for common S3 methods for class fit_wrc_hcc;

vcov for computing (co-)variances of the estimated nonlinear parameters;

prfloglik_sample for profile loglikelihood computations;

wc_model and hc_model for currently implemented models for soil water retention curves and hydraulic conductivity functions;

evaporative-length for physically constraining parameter estimates of soil hydraulic material functions.

Examples

```
# use of \donttest{} because execution time exceeds 5 seconds

data(sim_wrc_hcc)

# define number of cores for parallel computations
if(interactive()) ncpu <- parallel::detectCores() - 1L else ncpu <- 1L
```

```
# estimate parameters for single sample ...

#  ... from wrc and hcc data
plot(rfit_wrc_hcc <- fit_wrc_hcc(
  wrc_formula = wc ~ head, hcc_formula = hc ~ head,
  data = sim_wrc_hcc, subset = id == 1))
coef(rfit_wrc_hcc, gof = TRUE)

#  ... from wrc data
plot(rfit_wrc <- fit_wrc_hcc(
  wrc_formula = wc ~ head,
  data = sim_wrc_hcc, subset = id == 1))
coef(rfit_wrc, gof = TRUE)

#  ... from  hcc data
plot(rfit_hcc <- fit_wrc_hcc(
  hcc_formula = hc ~ head,
  data = sim_wrc_hcc, subset = id == 1))
coef(rfit_hcc, gof = TRUE)

# ... from wrc and hcc data
#     keeping some parameters fixed
plot(rfit_wrc_hcc_fixed <- fit_wrc_hcc(
  wrc_formula = wc ~ head, hcc_formula = hc ~ head,
  data = sim_wrc_hcc, subset = id == 1,
  param = c(alpha = 2.1, thetar = 0.1),
  fit_param = default_fit_param(alpha = FALSE, thetar = FALSE)),
  y = rfit_wrc_hcc)
coef(rfit_wrc_hcc, gof = TRUE)
coef(rfit_wrc_hcc_fixed, gof = TRUE)


# estimate parameters for 3 samples simultaneously by ...

# ... unconstrained, global optimisation algorithm NLOPT_GN_MLSL (default)
rfit_uglob <- fit_wrc_hcc(
  wrc_formula = wc ~ head | id, hcc_formula = hc ~ head | id,
  data = sim_wrc_hcc,
  control = control_fit_wrc_hcc(pcmp = control_pcmp(ncores = ncpu)))
summary(rfit_uglob)
op <- par(mfrow = c(3, 2))
plot(rfit_uglob)
on.exit(par(op))

# ... unconstrained, global optimisation algorithm SCEoptim
rfit_sce <- update(
  rfit_uglob,
  control = control_fit_wrc_hcc(
    settings = "sce", pcmp = control_pcmp(ncores = ncpu)))
coef(rfit_sce, gof = TRUE, lc = TRUE)
convergence_message(2, sce = TRUE)
op <- par(mfrow = c(3, 2))
```

```
plot(rfit_sce, y = rfit_uglob)
on.exit(par(op))

# ... unconstrained, local optimisation algorithm NLOPT_LN_BOBYQA,
#     logging iteration results to console
rfit_uloc <- update(
  rfit_uglob,
  param = as.matrix(coef(rfit_uglob, what = "nonlinear")),
  control = control_fit_wrc_hcc(
    settings = "ulocal", pcmp = control_pcmp(ncores = 1L)),
  verbose = 2)
coef(rfit_uloc, gof = TRUE, lc = TRUE)

# ... constrained, global optimisation algorithm NLOPT_GN_ISRES
rfit_cglob <- update(
  rfit_uglob,
  ratio_lc_lt_bound = c(lower = 0.8, upper = 1.2),
  control = control_fit_wrc_hcc(
    settings = "cglobal", nloptr = control_nloptr(ranseed = 1),
    pcmp = control_pcmp(ncores = ncpu)))
coef(rfit_cglob, gof = TRUE, lc = TRUE)

# ... constrained, local optimisation algorithm NLOPT_LD_CCSAQ
#     starting from unconstrained, locally fitted initial values
rfit_cloc_1 <- update(
  rfit_uglob,
  param = coef(rfit_uloc, what = "nonlinear"),
  ratio_lc_lt_bound = c(lower = 0.8, upper = 1.2),
  control = control_fit_wrc_hcc(
    settings = "clocal", pcmp = control_pcmp(ncores = ncpu)))
coef(rfit_cloc_1, gof = TRUE, lc = TRUE)
op <- par(mfrow = c(3, 2))
plot(x = rfit_uloc, y = rfit_cloc_1)
on.exit(par(op))

# ... constrained, local optimisation algorithm NLOPT_LD_CCSAQ
#     starting from constrained, globally fitted initial values,
#     using sample-specific evaporation rates and limits for ratio lc/lt
rfit_cloc_2 <- update(
  rfit_uglob,
  param = as.matrix(coef(rfit_cglob, what = "nonlinear")),
  e0 = c("1" = 0.002, "2" = 0.0025, "3" = 0.003),
  ratio_lc_lt_bound = rbind(
    "1" = c(lower = 0.7, upper = 2),
    "2" = c(lower = 0.8, upper = 1.4),
    "3" = c(lower = 0.8, upper = 2)
  ),
  control = control_fit_wrc_hcc(
    settings = "clocal", pcmp = control_pcmp(ncores = ncpu)))
coef(rfit_cloc_2, gof = TRUE, lc = TRUE, e0 = TRUE)

# ... global optimisation algorithm NLOPT_GN_MLSL
#     with sample-specific box-constraints
```

```
rfit_uglob_bc <- update(
  rfit_uglob,
  lower_param = rbind(
    "1" = c(alpha = 2.4, n = 1.11, thetar = 0.2, thetas = 0.6),
    "2" = c(alpha = 1.2, n = 1.12, thetar = 0.2, thetas = 0.6),
    "3" = c(alpha = 1.2, n = 1.13, thetar = 0.2, thetas = 0.6)
  ),
  upper_param = rbind(
    "1" = c(alpha = 20.1, n = 2.51, thetar = 0.6, thetas = 0.6),
    "2" = c(alpha = 20.2, n = 1.5,  thetar = 0.6, thetas = 0.6),
    "3" = c(alpha = 1.3, n = 2.53,  thetar = 0.6, thetas = 0.6)
  )
)
coef(rfit_uglob, gof = TRUE)
coef(rfit_uglob_bc, gof = TRUE)
```

`hc_model` *Models for Soil Hydraulic Conductivity Functions*

Description

The functions `hc_model` and `hcrel_model` compute, for given capillary pressure head h, the *hydraulic conductivity* $K(h)$ and the *relative hydraulic conductivity* $k(h)$ respectively, of a soil by parametrical models.

Usage

```
hcrel_model(h, nlp, precBits = NULL, hcc_model = "vgm")

hc_model(h, nlp, lp, precBits = NULL, hcc_model = "vgm")
```

Arguments

`h`	a mandatory numeric vector with values of capillary pressure head for which to compute the hydraulic conductivity. For consistency with other quantities, the unit of head should be **meter** [m].
`nlp`	a mandatory named numeric vector, currently with elements named `"alpha"`, `"n"` and `"tau"`, which are the *nonlinear* parameters $\boldsymbol{\nu}^{\mathrm{T}} = (\alpha, n, \tau)$, where α, n and τ are the inverse air entry pressure, the shape and the tortuosity parameters, see *Details*. For consistency with other quantities, the unit of α should be **1/meter** $[\mathrm{m}^{-1}]$.
`lp`	a mandatory named numeric vector, currently with a single element named `"k0"`, which is the saturated hydraulic conductivity K_0, the only *linear* parameter of the model, see *Details*. For consistency with other quantities, the unit of K_0 should be **meter/day** $[\mathrm{m\,d}^{-1}]$.

`precBits` an optional integer scalar defining the maximal precision (in bits) to be used in high-precision computations by `mpfr`. If equal to `NULL` (default) then `mpfr` is not used and the result of the function call is of storage mode `double`, see `soilhypfitIntro`.

`hcc_model` a keyword denoting the parametrical model for the hydraulic conductivity function. Currently, only the *Van Genuchten-Mualem* model (`wrc_model = "vgm"`) is implemented, see *Details*.

Details

The functions `hcrel_model` and `hc_model` currently model soil hydraulic conductivity functions only by the simplified form of the Van Genuchten-Mualem model (*Van Genuchten, 1980*) with the restriction $m = 1 - \frac{1}{n}$, i.e. by

$$k_{\mathrm{VGM}}(h;\boldsymbol{\nu}) = S_{\mathrm{VG}}(h;\boldsymbol{\nu})^{\tau}\left[1-\left(1-S_{\mathrm{VG}}(h;\boldsymbol{\nu})^{\frac{n}{n-1}}\right)^{\frac{n-1}{n}}\right]^2,$$

$$K_{\mathrm{VGM}}(h;\mu,\boldsymbol{\nu}) = K_0\, k_{\mathrm{VGM}}(h;\boldsymbol{\nu}),$$

where $\mu = K_0$ is the saturated hydraulic conductivity ($K_0 > 0$), $\boldsymbol{\nu}^{\mathrm{T}} = (\alpha, n, \tau)$ are the inverse air entry pressure ($\alpha > 0$), the shape ($n > 1$) and tortuosity parameter ($\tau > -2$), and $S_{\mathrm{VG}}(h;\boldsymbol{\nu})$ is the the volumetric water saturation, see `sat_model` for an expression.

Note that μ and $\boldsymbol{\nu}$ are passed to the functions by the arguments `lp` and `nlp`, respectively.

Value

A numeric vector with values of (relative) hydraulic conductivity.

Author(s)

Andreas Papritz <papritz@retired.ethz.ch>.

References

Van Genuchten, M. Th. (1980) A closed-form equation for predicting the hydraulic conductivity of unsaturated soils. *Soil Science Society of America Journal*, **44**, 892–898, doi:10.2136/sssaj1980.03615995004400050002x.

See Also

`soilhypfitIntro` for a description of the models and a brief summary of the parameter estimation approach;

`fit_wrc_hcc` for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

`control_fit_wrc_hcc` for options to control `fit_wrc_hcc`;

`soilhypfitmethods` for common S3 methods for class `fit_wrc_hcc`;

`vcov` for computing (co-)variances of the estimated nonlinear parameters;

`evaporative-length` for physically constraining parameter estimates of soil hydraulic material functions.

Examples

```
## define capillary pressure head (unit meters)
h <- c(0.01, 0.1, 0.2, 0.3, 0.5, 1., 2., 5.,10.)

## compute (relative) hydraulic conductivity
hcrel <- hcrel_model(h, nlp = c(alpha = 1.5, n = 2, tau = 0.5))
hc <- hc_model(h, nlp = c(alpha = 1.5, n = 2, tau = 0.5), lp = c(k0 = 5))

## display hydraulic conductivity function
op <- par(mfrow = c(1, 2))
plot(hcrel ~ h, log = "xy", type = "l")
plot(hc ~ h, log = "xy", type = "l")
on.exit(par(op))
```

---

profile_loglikelihood *Profile Loglikelihoods and Confidence Intervals for Parametric Modelling of Soil Hydraulic Properties*

Description

The function `prfloglik_sample` computes for a single soil sample a loglikelihood profile as a function of the specified values for subsets of model parameters of soil water retention and/or soil hydraulic conductivity functions. The function `confint_prfloglik_sample` computes a confidence interval of one model parameter based on the likelihood ratio test for a single soil sample, and the S3 method `confint` computes confidence intervals of *nonlinear* model parameters for multiple soil samples.

Usage

```
prfloglik_sample(object, values, soil_sample,
    ncores = min(detectCores() - 1L, NROW(values)), verbose = 0)

confint_prfloglik_sample(object, parm = names(default_fit_param()),
    soil_sample, level = 0.95, test = c("F", "Chisq"),
    denominator_df = c("nonlinear", "all"), param_bound = NULL,
    root_tol = .Machine$double.eps^0.25, froot_tol = sqrt(root_tol),
    verbose = 0)

## S3 method for class 'fit_wrc_hcc'
confint(object,
    parm = names(object[["control"]][["initial_param"]]), level = 0.95,
    subset = NULL, type = c("loglik", "normal"), test = c("F", "Chisq"),
    denominator_df = c("nonlinear", "all"),
    root_tol = .Machine$double.eps^0.25, froot_tol = sqrt(root_tol),
    ncores = detectCores() - 1L, verbose = 0, ...)
```

Arguments

object	an object of class fit_wrc_hcc, see fit_wrc_hcc.
values	a data.frame or a matrix with the values of the conditionally linear ($\boldsymbol{\mu}$) and nonlinear parameters ($\boldsymbol{\nu}$) that should be kept fixed to compute the likelihood profile (mandatory argument, see soilhypfitIntro, wc_model and hc_model for information about the parametrization of models for soil water retention curves and/or soil hydraulic conductivity functions.). The names of the columns of values must match the names of model parameters.
soil_sample	a character scalar with the label of the soil sample for which the loglikelihood profile or the confidence interval should be computed. If object contains parameter estimates for a single soil sample then soil_sample is ignored, otherwise soil_sample is a mandatory argument.
ncores	an integer defining the number of cores for parallel computations. ncores = 1 suppresses parallel computations.
verbose	positive integer controlling logging of diagnostic messages to the console and plotting of data and model curves during fitting, see fit_wrc_hcc.
parm	character scalar (confint_prfloglik_sample) or vector (confint) with name(s) of parameter(s) for which to compute the confidence interval(s). Note that confint_prfloglik_sample allows to compute a confidence interval for *all* parameters (including the linear ones), whereas the confint method computes confidence intervals for the *nonlinear* parameters ($\boldsymbol{\nu}$) only, see soilhypfitIntro for information about the parametrization of models for soil water retention curves and/or soil hydraulic conductivity functions.
level	numeric scalar with the confidence level required to compute the confidence interval.
test	character keyword specifying whether to use the asymptotic χ^2-distribution or the finite sample approximate F-distribution for the likelihood ratio test statistic when computing the confidence interval, see *Details*.
denominator_df	character keyword specifying whether the denominator degrees of freedom for the F-distribution of the test statistic is equal to the number of estimated *nonlinear* parameters ("nonlinear", default) or equal to the total number of estimated parameters ("all").
param_bound	a numeric vector of length 2 with the allowed range of the parameter for which the confidence interval is computed. The limits of the confidence interval are searched within this range, see *Details*. When equal to NULL (default) param_bound is taken from to component initial_objects of the selected component fit of object, see *Details*.
root_tol	a numeric scalar defining the desired accuracy (convergence tolerance) for root finding by uniroot when computing the confidence interval.
froot_tol	a numeric scalar defining the desired accuracy (function value tolerance) for deciding whether a root has been found.
subset	an integer, character or logical vector to the choose the soil samples for which confidence intervals should be computed. Defaults to NULL which computes the intervals for all samples contained in object.

type	character keyword specifying whether to compute confidence intervals based on the likelihood ratio test ("logik", default) by confint_prfloglik_sample or based on the asymptotic normal distribution of maximum likelihood estimates ("normal"), see *Details*.
...	additional arguments passed to methods, currently not used.

Details

Computing likelihood profiles:

The function prfloglik_sample computes the loglikelihood profile for a subset of the nonlinear ($\boldsymbol{\nu}$) (and linear $\boldsymbol{\mu}$) model parameters. We denote the profiling parameters of interest by $\boldsymbol{\phi}$ and the nuisance parameters that are estimated when computing the profile by $\boldsymbol{\psi}$, so that $(\boldsymbol{\phi}^{\mathrm{T}}, \boldsymbol{\psi}^{\mathrm{T}})^{\mathrm{T}}$ is a rearranged version of the parameter vector $(\boldsymbol{\mu}^{\mathrm{T}}, \boldsymbol{\nu}^{\mathrm{T}})^{\mathrm{T}}$, see soilhypfitIntro. prfloglik_sample computes the estimates $\widehat{\boldsymbol{\psi}}(\boldsymbol{\phi})$ of $\boldsymbol{\psi}$ and the profile loglikelihood $Q(\boldsymbol{\phi}, \widehat{\boldsymbol{\psi}}(\boldsymbol{\phi}); \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h})$ as a function of $\boldsymbol{\phi}$.

To compute the estimates, $\boldsymbol{\psi}$ is partitioned into the nonlinear and conditionally linear parameters, see soilhypfitIntro for details. prfloglik_sample uses the packages **parallel** and **snowfall** for parallelized computation of loglikelihood profiles.

Computing likelihood based confidence intervals:

The function confint_prfloglik_sample computes the confidence interval of a single model parameter $\phi \in (\boldsymbol{\mu}^{\mathrm{T}}, \boldsymbol{\nu}^{\mathrm{T}})^{\mathrm{T}}$ based on the likelihood ratio test. The interval is computed by assuming either

- that the test statistic $T(\phi) = \Delta Q(\phi) = 2(Q(\widehat{\phi}, \widehat{\boldsymbol{\psi}}; \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h}) - Q(\phi, \widehat{\boldsymbol{\psi}}(\phi); \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h}))$ follows a χ^2_1-distribution with 1 degrees of freedom (possible choice when both water retention and/or hydraulic conductivity data were used to estimate the parameters), or
- that the transformed test statistic $T(\phi) = (\exp(\Delta Q(\phi)/n) - 1)\,(n-p)$ follows an $F(1, n-p)$-distribution where n is the number of water content or hydraulic conductivity measurements, and p is the number of estimated parameters (see Uusipaikka, 2008, p. 115, for details). denominator_df allows to control how p is chosen. Note that this test distribution can only be chosen (and is then the default) when only water retention or only hydraulic conductivty data were used to estimate the parameters.

confint_prfloglik_sample computes profile loglikelihoods $Q(\phi, \widehat{\boldsymbol{\psi}}(\phi); \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h})$ by prfloglik_sample and then uses uniroot to search the roots of the equation

$$f(\phi) = q_T(\gamma) - T(\phi)$$

in the interval defined by param_bound. $q_T(\gamma)$ is the γ-quantile of the chosen test distribution for T.

Computing confidence intervals for several soil samples:

The confint method computes 2 types of confidence intervals (in dependence of type) of only the nonlinear parameters $\boldsymbol{\nu}$, possibly for multiple soil samples at the same time:

- intervals based on the asymptotic normal distribution of maximum likelihood estimates with standard errors computed by the vcov method for class fit_wrc_hcc (see vcov.fit_wrc_hcc,
- intervals based on the likelihood ratio test by confint_prfloglik_sample. The intervals for several soil samples are computed in parallel.

Requirements for computing loglikelihood profiles:

The parameters contained in `object` must be estimated by maximum likelihood (`method = "ml"`, see `soilhypfitIntro` and `control_fit_wrc_hcc`. Use of other estimation methods results an error.

Preferably an unconstrained local algorithm (`settings = "ulocal"`, see `soilhypfitIntro` and `control_fit_wrc_hcc`)) is used for minimizing the negative loglikelihood when generating `object`. Use of other algorithms generates a warning message.

Value

The function `prfloglik_sample` returns a `data.frame` with the columns of `values` (parameters of interest ϕ), a column `loglik` with the maximized profile loglikelihood $Q(\phi, \widehat{\psi}(\phi); \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h})$, columns with the estimated parameters $\widehat{\psi}(\phi)$, columns with the gradient of the loglikelihood with respect to the estimated nonlinear parameters (missing if all nonlinear parameters were fixed) and a column `converged`, indicating whether convergence has occurred (see `convergence_message`) when estimating the nonlinear parameters (equal to `NA` when all nonlinear parameters are fixed).

The function `confint_prfloglik_sample` returns a numeric vector of length 2 with the lower and upper limits of the confidence interval. If no roots were found then `NA` is returned. The returned result further contains as attribute list `prfloglik` the parameter estimate $\widehat{\phi}$ (`param_estimate`), the maximized loglikelihood $Q(\widehat{\phi}, \widehat{\psi}; \boldsymbol{\theta}, \boldsymbol{K}, \boldsymbol{h})$ (`loglik`), the quantile of the test distribution $q_{\text{test}}(\gamma)$ (`qtest`), the type of test distribution used (`test`), the significance `level`, the number of water content (`nobs_wrc`) and conductivity measurements (`nobs_hcc`) and the function values of $f(\phi)$ evaluated at the roots returned by `uniroot`.

The method `confint` returns a dataframe with the lower and upper limits of the confidence intervals for the estimated nonlinear parameters.

Author(s)

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References

Uusipaikka, E. (2008) Confidence Intervals in Generalized Regression Models. Chapman & Hall/CRC Press, Boca Raton doi:10.1201/9781420060386.

See Also

`soilhypfitIntro` for a description of the models and a brief summary of the parameter estimation approach;

`fit_wrc_hcc` for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

`control_fit_wrc_hcc` for options to control `fit_wrc_hcc`;

`soilhypfitmethods` for common S3 methods for class `fit_wrc_hcc`;

`vcov` for computing (co-)variances of the estimated nonlinear parameters;

`wc_model` and `hc_model` for currently implemented models for soil water retention curves and hydraulic conductivity functions;

evaporative-length for physically constraining parameter estimates of soil hydraulic material functions.

Examples

```
# use of \donttest{} because execution time exceeds 5 seconds

if(!requireNamespace("lattice", quietly = TRUE)){
  install.packages("lattice", repos = "https://cloud.r-project.org/")
  requireNamespace("lattice", quietly = TRUE)
}

data(sim_wrc_hcc)

# define number of cores for parallel computations
if(interactive()) ncpu <- parallel::detectCores() - 1L else ncpu <- 1L

# estimate parameters for 3 samples simultaneously by ...

# ... unconstrained, global optimisation algorithm NLOPT_GN_MLSL (default)
rfit_uglob <- fit_wrc_hcc(
  wrc_formula = wc ~ head | id, hcc_formula = hc ~ head | id,
  data = sim_wrc_hcc,
  control = control_fit_wrc_hcc(param_bound = param_boundf(
      alpha = c(0.00001, 50), n = c(1.0001, 7), tau = c(-1, 5)
    ), pcmp = control_pcmp(ncores = ncpu)))

# ... unconstrained, local optimisation algorithm NLOPT_LN_BOBYQA,
rfit_uloc <- update(
  rfit_uglob,
  param = as.matrix(coef(rfit_uglob, what = "nonlinear")),
  control = control_fit_wrc_hcc(
    settings = "ulocal", param_tf = param_transf(
      alpha = "identity", n = "identity", tau = "identity"
    ), param_bound = param_boundf(
      alpha = c(0.00001, 50), n = c(1.0001, 7), tau = c(-1, 5)
    ), pcmp = control_pcmp(ncores = ncpu)))

# extract estimated parameters for sample id == "1"
coef_id_1 <- unlist(coef(rfit_uloc, gof = TRUE, se = TRUE, subset = "1"))

# compute loglikelihood profile of parameter alpha for sample id == "1"
rprfloglik_alpha <- prfloglik_sample(
  rfit_uloc, values = data.frame(alpha = seq(1.5, 3.0, length = 40L)),
  soil_sample = "1", ncores = ncpu)
# plot loglikelihood profile along with 95% confidence intervals
plot(loglik ~ alpha, rprfloglik_alpha, type = "l")
abline(v = coef_id_1["alpha"])
# 95% confidence intervall based on likelihood ratio test
abline(h = -coef_id_1["obj"] - 0.5 * qchisq(0.95, 1), lty = "dashed")
# 95% confidence intervall based on asymptotic normal distribution
segments(
  x0 = coef_id_1["alpha"] + qnorm(0.025) * coef_id_1["se.alpha"],
```

```
  x1 = coef_id_1["alpha"] + qnorm(0.975) * coef_id_1["se.alpha"],
  y0 = min(rprfloglik_alpha$loglik)
)

# compute loglikelihood profile of parameter n for sample id == "1"
rprfloglik_n <- prfloglik_sample(
  rfit_uloc, values = data.frame(n = seq(1.7, 2.25, length = 40L)),
  soil_sample = "1", ncores = ncpu
)
# plot loglikelihood profile along with 95% confidence intervals
plot(loglik ~ n, rprfloglik_n, type = "l")
abline(v = coef_id_1["n"])
# 95% confidence intervall based on likelihood ratio test
abline(h = -coef_id_1["obj"] - 0.5 * qchisq(0.95, 1), lty = "dashed")
# 95% confidence intervall based on asymptotic normal distribution
segments(
  x0 = coef_id_1["n"] + qnorm(0.025) * coef_id_1["se.n"],
  x1 = coef_id_1["n"] + qnorm(0.975) * coef_id_1["se.n"],
  y0 = min(rprfloglik_n$loglik)
)

# compute loglikelihood profile of parameters alpha and n for sample id == "1"
rprfloglik_alpha_n <- prfloglik_sample(
  rfit_uloc, values = expand.grid(
    alpha = seq(1.5, 3.0, length = 40L), n = seq(1.7, 2.25, length = 40L)),
  soil_sample = "1", ncores = ncpu
)

# joint confidence region of alpha and n based on likelihood ratio test
lattice::levelplot(loglik ~ alpha + n, rprfloglik_alpha_n,
  panel = function(x, y, z, subscripts, ...){
    lattice::panel.levelplot(x, y, z, subscripts, ...)
    lattice::panel.levelplot(x, y, z, subscripts, region = FALSE, contour = TRUE,
      at = -coef_id_1["obj"] - 0.5 * qchisq(0.95, 2),
      lty = "solid"
    )
    lattice::panel.levelplot(x, y, z, subscripts, region = FALSE, contour = TRUE,
      at = -coef_id_1["obj"] - 0.5 * qchisq(0.9, 2),
      lty = "dashed"
    )
    lattice::panel.levelplot(x, y, z, subscripts, region = FALSE, contour = TRUE,
      at = -coef_id_1["obj"] - 0.5 * qchisq(0.8, 2),
      lty = "dotted"
    )
    lattice::panel.points(coef_id_1["alpha"], coef_id_1["n"], pch = 16)
    lattice::panel.lines(
      x = rprfloglik_alpha[, c("alpha", "n")], col = "blue"
    )
    lattice::panel.lines(
      x = rprfloglik_n[, c("alpha", "n")], col = "magenta"
    )
  }, key = list(
    corner = c(1, 0.97),
```

```
    lines = list(
      lty = c(rep("solid", 3), "dashed", "dotted"),
      col = c("blue", "magenta", rep("black", 3))
    ),
    text = list(c(
        "estimate of n as a function of fixed alpha",
        "estimate of alpha as a function of fixed n",
        "95% joint confidence region",
        "90% joint confidence region",
        "80% joint confidence region"
      ))
  ))

# compute 95%-confidence interval
(ci.alpha <- confint_prfloglik_sample(
  rfit_uloc, parm = "alpha", soil_sample = "1"
))
# use limits to draw loglikelihood profile for alpha
rprfloglik_alpha <- prfloglik_sample(
  rfit_uloc, values = data.frame(
    alpha = seq(0.9 * ci.alpha[1], 1.1 * ci.alpha[2], length = 40L)),
  soil_sample = "1", ncores = ncpu)
plot(loglik ~ alpha, rprfloglik_alpha, type = "l")
lines(
  ci.alpha,
  rep(diff(unlist(attr(ci.alpha, "prfloglik")[c("qtest", "loglik")])) , 2)
)
abline(v = attr(ci.alpha, "prfloglik")["estimate"], lty = "dashed")

# 95%-confidence intervals of all nonlinear parameters based for all
# soil samples asymptotic normal distribution of maximum likelihood estimates
confint(rfit_uloc, type = "normal")

# 95%-confidence intervals of all nonlinear parameters based for all
# soil samples based on likelihood ratio test
confint(rfit_uloc, ncores = ncpu)
```

`sim_wrc_hcc` *Simulated Soil Water Retention and Hydraulic Conductivity Data*

Description

The data give simulated values of water content and hydraulic conductivity at given capillary pressure head for 3 soil samples.

Usage

```
data(sim_wrc_hcc)
```

Format

A data frame with 28 observations on the following 4 variables.

`id` a factor with levels `1`, `2`, `3` coding the soil samples.

`head` a numeric vector with values of capillary pressure head (unit m).

`wc` a numeric vector with values of simulated volumetric water content (unit -)

`hc` a numeric vector with values of simulated hydraulic conductivity (unit $\mathrm{m\,d^{-1}}$).

Details

The values of `wc` and `hc` were simulated by the model of *Van Genuchten Mualem* (Van Genuchten, 1980, see `wc_model` and `hc_model`) with the following parameters:

sample id	θ_r	θ_s	K_0 [$\mathrm{m\,d^{-1}}$)]	α [$\mathrm{m^{-1}}$]	n	τ
1	0.05	0.45	0.1	2	2	0.5
2	0.1	0.5	5	1.5	1.5	0.5
3	0.05	0.45	2	1.4	1.3	0.5

Normally distributed errors were added to the model values (`wc`: sd: 0.05; `log(hc)`: sd 0.5).

References

Van Genuchten, M. Th. (1980) A closed-form equation for predicting the hydraulic conductivity of unsaturated soils. *Soil Science Society of America Journal*, **44**, 892–898, doi:10.2136/sssaj1980.03615995004400050002x.

Examples

```
data(sim_wrc_hcc)

if(!requireNamespace("lattice", quietly = TRUE)){
  install.packages("lattice", repos = "https://cloud.r-project.org/")
  requireNamespace("lattice", quietly = TRUE)
}

lattice::xyplot(wc ~ head|id, type = "l", sim_wrc_hcc,
  scales = list( x = list(log=TRUE)))
lattice::xyplot(hc ~ head|id, type = "l", sim_wrc_hcc,
  scales = list( log = TRUE))
```

soilhypfitS3methods *Common S3 Methods for Class* fit_wrc_hcc

Description

This page documents the methods coef, summary, print, plot and lines for the class fit_wrc_hcc.

Usage

```
## S3 method for class 'fit_wrc_hcc'
coef(object, what = c("all", "nonlinear", "linear"),
    subset = NULL, residual_se = FALSE, se = FALSE, gof = FALSE, lc = FALSE,
    e0 = FALSE, bound = lc, ...)

## S3 method for class 'fit_wrc_hcc'
summary(object, what = c("all", "nonlinear", "linear"),
    subset = NULL, gof = TRUE, lc = TRUE, ...)

## S3 method for class 'fit_wrc_hcc'
print(x, ...)

## S3 method for class 'fit_wrc_hcc'
plot(x, what = c("wrc", "hcc"), y = NULL,
    subset = NULL, ylim_wc = NULL, ylim_hc = NULL,
    head_saturation = 0.01,
    beside = identical(sum(par("mfrow")), 2L), pch = 1, col_points = "black",
    col_line_x = "blue", lty_x = "solid",
    col_line_y = "orange", lty_y = "dashed",
    xlab_wc = "head [m]", ylab_wc = "water content [-]",
    xlab_hc = "head [m]", ylab_hc = "hyd. conductivity [m/d]",
    draw_legend = TRUE, draw_parameter = FALSE, cex_legend = 0.7, ...)

## S3 method for class 'fit_wrc_hcc'
lines(x, what = c("wrc", "hcc"), id = 1,
    head_saturation = 0.01, ...)
```

Arguments

object, x, y	an object of class fit_wrc_hcc, see fit_wrc_hcc.
what	character keyword indicating the type of parameters to return (coef) or the type of data to plot.
subset	an integer, character or logical vector to the choose the soil samples for which data and model curves are displayed or extracted. Defaults to NULL which displays results for all soil samples.

residual_se	a logical scalar to control whether residual standard errors (= standard deviations of residuals) should be returned, see *Details*.
se	a logical scalar to control whether standard errors of the nonlinear parameters $\boldsymbol{\nu}$ should be returned, see *Details* and vcov.fit_wrc_hcc.
gof	a logical scalar to control whether goodness-of-fit statistic should be returned.
lc	a logical scalar to control whether the characteristic evaporative length should be returned, see evaporative-length.
e0	a logical scalar to control whether the evaporation rate should be returned. This is only effective for constrained estimation, see evaporative-length.
bound	a logical scalar to control whether the lower and upper bounds of the ratio L_c/L_t should be returned. This is only effective for constrained estimation, see evaporative-length.
ylim_wc	optional numeric vector of length 2 to set the range of water content values displayed on the y-axis (default NULL for automatic axis scales).
ylim_hc	optional numeric vector of length 2 to set the range of hydraulic conductivity values displayed on the y-axis (default NULL for automatic axis scales).
head_saturation	head value (unit m) assigned to zero head values in plots with logarithmic head scale.
beside	a logical scalar controlling whether water retention curves and hydraulic conductivity functions of a sample should be plotted side by side.
pch	plotting 'character', i.e., symbol to use for the measurements, see points.
col_points	color code or name for symbol colors for the measurements, see par.
col_line_x	color code or name for the line representing the fitted model x, see par.
lty_x	type of line representing the fitted model x, see par.
col_line_y	color code or name for the line representing the fitted model y, see par.
lty_y	type of line representing the fitted model y, see par.
xlab_wc	a character string with the annotation for the *x*-axis of a water retention curve.
ylab_wc	a character string with the annotation for the *y*-axis of a water retention curve.
xlab_hc	a character string with the annotation for the *x*-axis of a hydraulic conductivity function.
ylab_hc	a character string with the annotation for the *y*-axis of a hydraulic conductivity function.
draw_legend	a logical scalar controlling whether a legend with the values of the arguments x and y and the residual sums of squares is drawn if y is non-NULL.
draw_parameter	a logical scalar controlling whether the parameters are drawn (default FALSE).
cex_legend	a character expansion factor for annotations by legend.
id	a character string or integer scalar to select the sample for which to plot the modelled water retention curve or hydraulic conductivity function.
...	additional arguments passed to methods.

Details

Residual standard errors, standard errors of the nonlinear parameters and confidence intervals based on the asymptotic normal distribution are computed only for mpd and ml estimates, see soilhypfitIntro, control_fit_wrc_hcc and vcov.

The plot method for class fit_wrc_hcc displays for each sample the measurements of the water retention curve and/or the hydraulic conductivity function, along with the fitted model curve(s). Optionally, the curves of a second model fit (specified by the argument y) can be plotted for each sample.

The lines method adds the curve of a fitted model to an existing plot.

Value

The method coef returns a dataframe with the estimated parameters (and optionally standard errors), optionally the value of the objective function along with convergence code and/or information on the characteristic evaporative length.

The method summary generates a list (of class summary.fit_wrc_hcc) with the following components:

data a named integer vector with the total number of samples (nsamp) and the number of samples with only water retention (nwrc), only hydraulic conductivity (nhcc) and both type of measurements (nwrchcc).

control a list with a subset (settings, nloptr, sce, approximation_alpha_k0, param_bound, param_tf) of the components of object[["control"]], see fit_wrc_hcc and control_fit_wrc_hcc.

result a dataframe with the estimated parameters and optionally the residual sum of squares along with convergence code and/or information on the characteristic evaporative length.

call the call component of object.

Note that only a print method is available for class summary.fit_wrc_hcc.

Author(s)

Andreas Papritz <papritz@retired.ethz.ch>.

See Also

soilhypfitIntro for a description of the models and a brief summary of the parameter estimation approach;

fit_wrc_hcc for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

control_fit_wrc_hcc for options to control fit_wrc_hcc;

vcov for computing (co-)variances of the estimated nonlinear parameters;

prfloglik_sample for profile loglikelihood computations;

wc_model and hc_model for currently implemented models for soil water retention curves and hydraulic conductivity functions;

evaporative-length for physically constraining parameter estimates of soil hydraulic material functions.

Examples

```
# use of \donttest{} because execution time exceeds 5 seconds

data(sim_wrc_hcc)

# define number of cores for parallel computations
if(interactive()) ncpu <- parallel::detectCores() - 1L else ncpu <- 1L

# estimate parameters for 3 samples by unconstrained, global optimisation
# algorithm NLOPT_GN_MLSL
# sample 1: use only conductivity data
# sample 2: use only water content data
# sample 3: use both types of data
rfit_uglob <- fit_wrc_hcc(
  wrc_formula = wc ~ head | id, hcc_formula = hc ~ head | id,
  wrc_subset = id != 1, hcc_subset = id != 2,
  data = sim_wrc_hcc, fit_param = default_fit_param(tau = TRUE),
  control = control_fit_wrc_hcc(param_bound = param_boundf(
      alpha = c(0.00001, 50), n = c(1.0001, 7), tau = c(-1, 5)
    ), pcmp = control_pcmp(ncores = ncpu)))
print(rfit_uglob)
summary(rfit_uglob)
coef(rfit_uglob, what = "nonlinear")
coef(rfit_uglob, what = "linear", gof = TRUE)
coef(vcov(rfit_uglob), status = TRUE, se = FALSE)
op <- par(mfrow = c(3, 2))
plot(rfit_uglob)
on.exit(par(op))
```

---

swissforestsoils *Physical properties of Swiss forest soils*

Description

The data give basic physical properties, water content and hydraulic conductivity measurements (at given capillary pressure head) for 128 soil layers (horizons) measured at 23 forest sites in Switzerland.

Usage

```
data(swissforestsoils)
```

Format

A data frame with 1373 observations on the following 21 variables.

`profile_id` a factor with short labels for the 23 sites.

`profile` a factor with with long labels for the 23 sites.

`longitude` a numeric vector with the latitude of the site in degree.

`latitude` a numeric vector with the latitude of the site in degree.

`layer_id` a factor with labels for the 128 soil layer.

`layer_ub`, `layer_lb` numeric vectors with the upper and lower depth (unit cm) of the soil layer for the measurements of the basic physical properties (`particle_density`, ..., `ksat`).

`particle_density` a numeric vector with the density of the solid soil material (unit $\mathrm{g\,cm^{-3}}$).

`bulk_density` a numeric vector with soil (bulk) density (unit $\mathrm{g\,cm^{-3}}$).

`porosity` a numeric vector with the soil porosity (unit volume percentage).

`clay` a numeric vector with the clay content (unit mass percentage).

`silt` a numeric vector with the silt content (unit mass percentage).

`sand` a numeric vector with the sand content (unit mass percentage).

`ksat` a numeric vector with the saturated hydraulic conductivity (unit $\mathrm{m\,d^{-1}}$).

`head` a numeric vector with capillary pressure head at which `theta` (water retention curve) and/or `ku` (hydraulic conductivity function) were measured (unit m).

`layer_ub_theta`, `layer_lb_theta` numeric vectors with the upper and lower depth (unit cm) of the soil layer for which the water retention curve was measured.

`theta` a numeric vector with volumetric water content measurements (dimensionless) of the water retention curve.

`layer_ub_ku`, `layer_lb_ku` a numeric vector with the upper and lower depth (unit cm) of the soil layer for which the water retention curve was measured.

`ku` a numeric vector with hydraulic conductivity measurements (unit $\mathrm{m\,d^{-1}}$) of the hydraulic conductivity function.

Details

`clay`, `silt` and `sand` refer to soil particles with diameter less than 2, between 2 and 50 and larger than 50 μm.

Source

Richard, F. & Lüscher, P. 1978 – 1987. Physikalische Eigenschaften von Böden der Schweiz. Lokalformen Bände 1 – 4. Eidgenössische Anstalt für das forstliche Versuchswesen, Birmensdorf.

Examples

```
# use of \donttest{} because execution time exceeds 5 seconds

# estimate parameters using all samples (samples with water retention,
# hydraulic conductivity, or with both type of measurements)

data(swissforestsoils)

# define number of cores for parallel computations
if(interactive()) ncpu <- parallel::detectCores() - 1L else ncpu <- 1L
```

```
# unconstrained estimation (global optimisation algorithm NLOPT_GN_MLSL)
r_uglob <- fit_wrc_hcc(
  wrc_formula = theta ~ head | layer_id,
  hcc_formula = ku ~ head | layer_id,
  data = swissforestsoils,
  control = control_fit_wrc_hcc(
    settings = "uglobal", pcmp = control_pcmp(ncores = ncpu)))
summary(r_uglob)
coef(r_uglob)
```

`utility-functions` *Utility functions for package soilhypfit*

Description

This page documents the functions `convergence_message`, `extract_error_messages`, `select_failed_fits` and `check_param_boundf`.

Usage

```
convergence_message(x, sce = FALSE)

extract_error_messages(object, start = 1, stop = 80, prt = TRUE)

select_failed_fits(object)

check_param_boundf(y, compare_thetar_thetas = TRUE)
```

Arguments

`x`	an integer scalar issued by the optimisers of `nloptr` or `SCEoptim` on (non-)convergence
`sce`	a logical scalar to select the optimiser `nloptr` (`FALSE`, default) or `SCEoptim` (`TRUE`).
`object`	an object of class `fit_wrc_hcc`, see `fit_wrc_hcc`.
`prt`	a logical scalar controlling whether the error messages should be printed.
`start, stop`	integer scalar with the first and last character to print.
`y`	a named list of numeric vectors of length 2 that define the allowed lower and upper bounds (box constraints) for the parameters of the models, see argument `param_bound` of `control_fit_wrc_hcc`.
`compare_thetar_thetas`	logical scalar to control cross-comparison of valid ranges of parameters `thetar` and `thetas`.

Details

The function convergence_message prints a message that explains the convergence codes, for nloptr, see NLopt return values. The function extract_error_messages extract the error messages of estimations that failed and optionally prints sub-strings of them.
The function select_failed_fits returns the ids of the soil samples for which parameter estimation failed. The function check_param_boundf checks the validity and consistecy of bounds of box constraints of model parameters.

Value

The function convergence_message and extract_error_messages return invisibly the convergence code or the error messages.

Author(s)

Andreas Papritz <papritz@retired.ethz.ch>.

References

Duan, Q., Sorooshian, S., and Gupta, V. K. (1994) Optimal use of the SCE-UA global optimisation method for calibrating watershed models, *Journal of Hydrology* **158**, 265–284, doi:10.1016/0022-1694(94)900574.

Johnson, S.G. The NLopt nonlinear-optimisation package. https://github.com/stevengj/nlopt.

See Also

soilhypfitIntro for a description of the models and a brief summary of the parameter estimation approach;

fit_wrc_hcc for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

control_fit_wrc_hcc for options to control fit_wrc_hcc;

soilhypfitmethods for common S3 methods for class fit_wrc_hcc;

vcov for computing (co-)variances of the estimated nonlinear parameters;

prfloglik_sample for profile loglikelihood computations;

wc_model and hc_model for currently implemented models for soil water retention curves and hydraulic conductivity functions;

evaporative-length for physically constraining parameter estimates of soil hydraulic material functions.

Examples

```
convergence_message(3)
convergence_message(2, sce = TRUE)
```

vcov vcov *Method for Class* fit_wrc_hcc

Description

This page documents the method vcov for the class fit_wrc_hcc and its coef method. vcov extracts the covariance matrices of the nonlinear parameters $\widehat{\boldsymbol{\nu}}$ estimated by maximum likelihood or maximum posterior density.

Usage

```
## S3 method for class 'fit_wrc_hcc'
vcov(object, subset = NULL, grad_eps,
    bound_eps = sqrt(.Machine$double.eps), ...)

## S3 method for class 'vcov_fit_wrc_hcc'
coef(object, se = TRUE, correlation = se,
    status = FALSE, ...)
```

Arguments

object	either an object of class fit_wrc_hcc for vcov or an object of class vcov_fit_wrc_hcc for coef.
subset	an integer, character or logical vector to the choose the soil samples for which covariance matrices should be extracted. Defaults to NULL, which extracts the covariances for all soil samples.
grad_eps	a numeric scalar defining a critical magnitude of the moduli of scaled gradient components so that they are considered to be approximately equal to zero, see *Details*.
bound_eps	a numeric scalar defining the critical difference between parameter estimates and the boundaries of the parameter space so that the estimates are considered to be identical to the boundary values, see *Details*.
se	a logical scalar to control whether standard errors of the estimated nonlinear parameters $\widehat{\boldsymbol{\nu}}$ should be returned (TRUE, default) or variances (FALSE).
correlation	a logical scalar to control whether correlations (TRUE, default) or covariances (FALSE) of the esitmated nonlinear parameters $\widehat{\boldsymbol{\nu}}$ should be returned.
status	a logical scalar to control whether diagnostics should be returned along with the results.
...	additional arguments passed to methods, currently not used.

Details

The function vcov extracts (co-)variances of the nonlinear parameters from the inverse Hessian matrix of the objective function at the solution $\widehat{\nu}$ for mpd and ml estimates, see soilhypfitIntro and *Stewart and Sørensen (1981)*.

vcov checks whether the gradient at the solution is approximately equal to zero and issues a warning if this is not the case. This is controlled by the argument grad_eps which is the tolerable largest modulus of the scaled gradient (= gradient divided by the absolute value of objective function) at the solution. The function control_fit_wrc_hcc selects a default value for grad_eps in the dependence of the chosen optimisation approach (argument settings of control_fit_wrc_hcc).

vcov sets covariances equal to NA if the parameter estimates differ less than bound_eps from the boundaries of the parameter space as defined by param_boundf.

Value

The method vcov returns an object of of class vcov_fit_wrc_hcc, which is a list of covariance matrices of the estimated nonlinear parameters for the soil samples. The attribute status of the matrices qualifies the covariances.

The coef method for class vcov_fit_wrc_hcc extracts the entries of the covariances matrices, optionally computes standard errors and correlation coefficients and returns the results in a dataframe.

Author(s)

Andreas Papritz <papritz@retired.ethz.ch>.

References

Stewart, W.E. and Sørensen, J.P. (1981) Bayesian estimation of common parameters from multiresponse data with missing observations. *Technometrics*, **23**, 131–141, doi:10.1080/00401706.1981.10486255.

See Also

soilhypfitIntro for a description of the models and a brief summary of the parameter estimation approach;

fit_wrc_hcc for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

control_fit_wrc_hcc for options to control fit_wrc_hcc;

soilhypfitmethods for common S3 methods for class fit_wrc_hcc;

prfloglik_sample for profile loglikelihood computations;

wc_model and hc_model for currently implemented models for soil water retention curves and hydraulic conductivity functions;

evaporative-length for physically constraining parameter estimates of soil hydraulic material functions.

Examples

```
# use of \donttest{} because execution time exceeds 5 seconds

data(sim_wrc_hcc)

# define number of cores for parallel computations
if(interactive()) ncpu <- parallel::detectCores() - 1L else ncpu <- 1L

# estimate parameters for 3 samples by unconstrained, global optimisation
# algorithm NLOPT_GN_MLSL
# sample 1: use only conductivity data
# sample 2: use only water content data
# sample 3: use both types of data
rfit_uglob <- fit_wrc_hcc(
  wrc_formula = wc ~ head | id,
  hcc_formula = hc ~ head | id,
  wrc_subset = id != 1,
  hcc_subset = id != 2,
  data = sim_wrc_hcc,
  control = control_fit_wrc_hcc(pcmp = control_pcmp(ncores = ncpu)))
print(rfit_uglob)
summary(rfit_uglob)
coef(rfit_uglob, what = "nonlinear")
coef(rfit_uglob, what = "linear", gof = TRUE)
coef(vcov(rfit_uglob), status = TRUE, se = FALSE)
op <- par(mfrow = c(3, 2))
plot(rfit_uglob)
on.exit(par(op))
```

---

wc_model *Models for Soil Water Retention Curves*

Description

The functions `sat_model` and `wc_model` compute, for given capillary pressure head h, the volumetric water saturation $S(h)$ and the volumetric water content $\theta(h)$, respectively, of a soil by parametrical models.

Usage

```
sat_model(h, nlp, precBits = NULL, wrc_model = "vg")

wc_model(h, nlp, lp, precBits = NULL, wrc_model = "vg")
```

Arguments

h	a mandatory numeric vector with values of capillary pressure head for which to compute the volumetric water saturation or content. For consistency with other quantities, the unit of pressure head should be **meter** [m].
nlp	a mandatory named numeric vector, currently with elements named "alpha" and "n", which are the *nonlinear* parameters $\boldsymbol{\nu}^{\mathrm{T}} = (\alpha, n)$, where α and n are the inverse air entry pressure and the shape parameter, see *Details*. For consistency with other quantities, the unit of α should be **1/meter** [m^{-1}].
lp	a mandatory named numeric vector, currently with elements named "thetar" and "thetas", which are the *linear* parameters $\boldsymbol{\mu}^{\mathrm{T}} = (\theta_r, \theta_s)$ where θ_r and θ_s are the residual and saturated water content, respectively, see *Details*.
precBits	an optional integer scalar defining the maximal precision (in bits) to be used in high-precision computations by mpfr. If equal to NULL (default) then mpfr is not used and the result of the function call is of storage mode double, see soilhypfitIntro.
wrc_model	a keyword denoting the parametrical model for the water retention curve. Currently, only the *Van Genuchten model* (wrc_model = "vg") is implemented, see *Details*.

Details

The functions sat_model and wc_model currently model soil water retention curves only by the simplified form of the model by *Van Genuchten (1980)* with the restriction $m = 1 - \frac{1}{n}$, i.e.

$$S_{\mathrm{VG}}(h; \boldsymbol{\nu}) = (1 + (\alpha\, h)^n)^{\frac{1-n}{n}}$$

$$\theta_{\mathrm{VG}}(h; \boldsymbol{\mu}, \boldsymbol{\nu}) = \theta_r + (\theta_s - \theta_r)\, S_{\mathrm{VG}}(h; \boldsymbol{\nu})$$

where $\boldsymbol{\mu}^{\mathrm{T}} = (\theta_r, \theta_s)$ are the residual and saturated water content ($0 \le \theta_r \le \theta_s \le 1$), respectively, and $\boldsymbol{\nu}^{\mathrm{T}} = (\alpha, n)$ are the inverse air entry pressure ($\alpha > 0$) and the shape parameter ($n > 1$).

Note that $\boldsymbol{\mu}$ and $\boldsymbol{\nu}$ are passed to the functions by the arguments lp and nlp, respectively.

Value

A numeric vector with values of volumetric water saturation (sat_model) or water content (wc_model).

Author(s)

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References

Van Genuchten, M. Th. (1980) A closed-form equation for predicting the hydraulic conductivity of unsaturated soils. *Soil Science Society of America Journal*, **44**, 892–898, doi:10.2136/sssaj1980.03615995004400050002x.

See Also

soilhypfitIntro for a description of the models and a brief summary of the parameter estimation approach;

fit_wrc_hcc for (constrained) estimation of parameters of models for soil water retention and hydraulic conductivity data;

control_fit_wrc_hcc for options to control fit_wrc_hcc;

soilhypfitmethods for common S3 methods for class fit_wrc_hcc;

vcov for computing (co-)variances of the estimated nonlinear parameters;

prfloglik_sample for profile loglikelihood computations;

evaporative-length for physically constraining parameter estimates of soil hydraulic material functions.

Examples

```
## define capillary pressure head (unit meters)
h <- c(0.01, 0.1, 0.2, 0.3, 0.5, 1., 2., 5.,10.)

## compute water saturation and water content
sat <- sat_model(h, nlp = c(alpha = 1.5, n = 2))
theta <- wc_model(
  h ,nlp = c(alpha = 1.5, n = 2), lp = c(thetar = 0.1, thetas = 0.5))

## display water retention curve
op <- par(mfrow = c(1, 2))
plot(sat ~ h, log = "x", type = "l")
plot(theta ~ h, log = "x", type = "l")
on.exit(par(op))
```

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