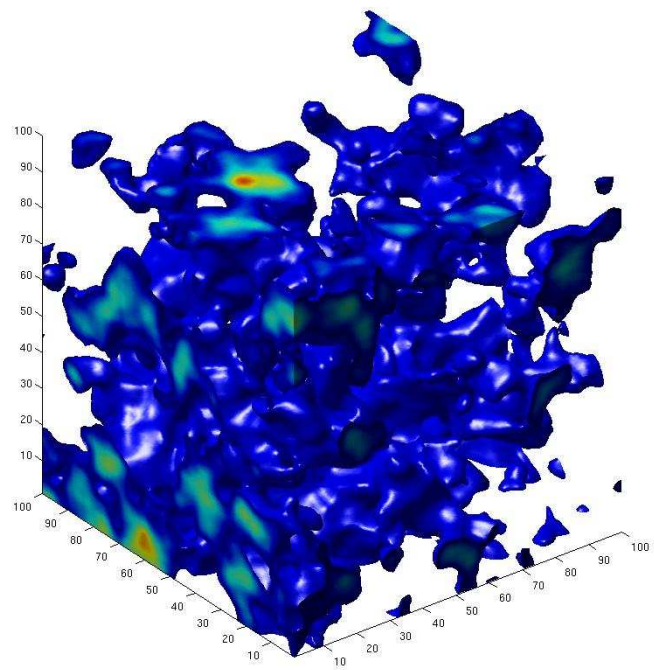


TFRACGEN
MATLAB SCRIPT
FOR
GENERATING TRUNCATED FRACTAL RANDOM
FIELDS



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Falk Heße
UFZ - Helmholtz Center for Environmental Research
Department of Computational Hydrosystems
Permoserstr. 15
D-04318 Leipzig
GERMANY

e-mail: falk.hesse@ufz.de

1 Introduction and concepts

TFRACGEN was created for generating 1-dimensional, 2-dimensional and 3-dimensional random fields with a truncated power-law variogram. When using commercial or free software, it is often hard to understand what is going on exactly when pushing buttons which is unacceptable from a scientific point of view and is why the TFRACGEN project was started. An other motivation for the ongoing development of TFRACGEN is that it can be used on any computer system with a working Matlab environment. There is no fancy GUI so you will need knowledge of Matlab.

2 Reference Manual

2.1 Basic structure

TFRACGEN is made up of several Matlab files, which in conjunction provide the functionality described above. In the following these files are listed and explained.

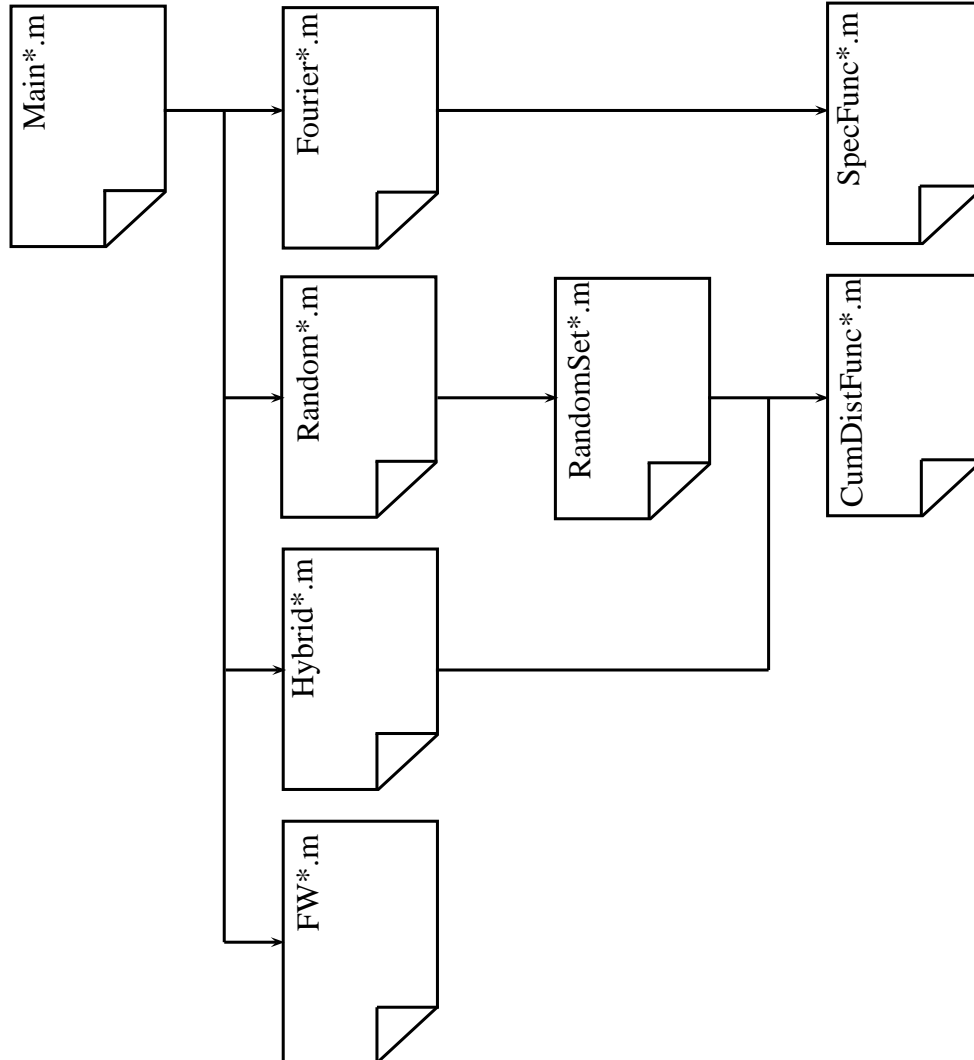


Figure 2.1: Schematic of the different files. The wildcard character is used in order to replace the dimension.

The main file is the eponymous `Main*.m`, where the different variables are declared (see schematic in Figure 2.1). Here the wildcard character is a placeholder for the dimension of the problem, i.e. 1D, 2D or 3D.

2.2 Variables

Several variables are declared in the main file, that determine the shape and characteristics of the generated random fields. In the following these variables are listed and explained.

Geostatistical parameters There are several parameters, that describe the geostatistical characteristics of the random field

- `muK`, i.e. expectation value μ of the K -field ,
- `si2K`, i.e. variance σ^2 of the K -field,
- `muY`, i.e. the expectation value μ of the Y -field,
- `si2Y`, i.e. the variance σ^2 of the Y -field,
- `L_geo`, which is the maximum geological length scale,
- `l_geo`, which is the minimum geological length scale,
- `H`, which is the Hurst coefficient as well as
- `func`, which defines the variogram model function of the random field

The first four parameters describe the one-point distribution of the hydraulic conductivity field. The algorithm itself only creates a Gaussian random field u with zero mean $\mu = 0$ and unit variance $\sigma^2 = 1$. The field of log-hydraulic conductivity Y is then computed according to $Y = \sigma_Y u + \mu_Y$. The field of hydraulic conductivity K is then computed according to $K = \exp(Y)$. Due to the fixed relationship between the Y and the K field only two of these parameters need to be specified. If the values of the K field are given, one can compute the corresponding Y values according to

$$\mu_Y = \ln(\mu_K) - \frac{1}{2} \ln \left(1 + \frac{\sigma_K^2}{\mu_K^2} \right) \quad (2.1a)$$

$$\sigma_Y^2 = \ln \left(1 + \frac{\sigma_K^2}{\mu_K^2} \right). \quad (2.1b)$$

The maximum geological length scale `L_geo` in case of a truncated fractal field is the largest length scale of heterogeneities, that is represented in the random field. In case of a non-fractal field this value is the correlation length of the field. The minimum geological length scale `l_geo` in case of a truncated fractal field is the smallest length scale of heterogeneities, that is represented in the random field. In case of a non-fractal field this value is ignored.

The Hurst coefficient `H` describes the fractality of the generated random field and is related to the fractal dimension D_f of the generated structures by the following relationship $H = 2 - D_f$.

The parameter `func` is a string variable that specifies the variogram model function. There are two fractal models, i.e. the truncated power law based on Gaussian modes `tFracGauss` and the truncated power law based on exponential modes `tFracExp`. The generator can also generate two non-fractal models, i.e. a Gaussian model `Gauss` and an exponential model `Exp`.

Geometrical parameters There are several parameters, that describe the geometrical characteristics of the random field

- L_{num} , which is the maximum numerical length scale,
- l_{num} , which is the minimum numerical length scale,
- α , which is the angle between the main axis of the random field and
- γ , which is the angle between the main axis of the random field and the numerical grid.

The numerical grid on which the random field is defined can be a regular square lattice but also any unstructured grid if the Randomization method or the Hybrid method is used for the generation of the random field.

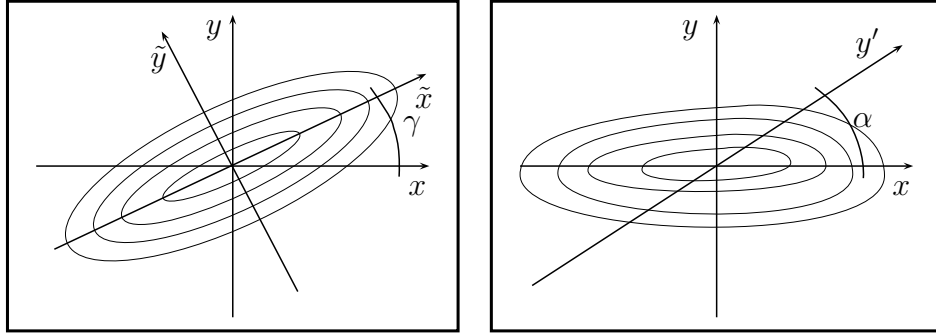


Figure 2.2: Schematic of the different angles, which are used in the generator.

In case of two or three dimensions the main axis of the random field can be at an angle, too. Here α is the angle between the main axis of the spectral density function of the random field. In two dimension α has one component being the angle between the x - and y -axis, whereas in three dimension α has two components the second being the angle between the y - and z -axis. The parameter γ is the angle between the x -axis of the random field and the numerical grid. The rotation is defined as $\tilde{\mathbf{x}} = R\mathbf{x}$ with the rotation matrix R being given according to

$$R = \begin{pmatrix} \cos(\gamma) & -\sin(\gamma) \\ \sin(\gamma) & \cos(\gamma) \end{pmatrix} \quad (2.2a)$$

$$R = \begin{pmatrix} \cos(\gamma) & -\sin(\gamma) & 0 \\ \sin(\gamma) & \cos(\gamma) & 0 \\ 0 & 0 & 1 \end{pmatrix} \quad (2.2b)$$

in two and three dimensions respectively. The inverse rotation is simply computed according to $\mathbf{x} = R^{-1}\tilde{\mathbf{x}}$. The transformation is defined as $\mathbf{x}' = T\mathbf{x}$ with the transformation matrix T being given according to

$$T = \begin{pmatrix} 1 & \tan(\alpha) \\ 0 & 1 \end{pmatrix} \quad (2.3a)$$

$$T = \begin{pmatrix} 1 & \tan(\alpha_1) & 0 \\ 0 & 1 & \tan(\alpha_2) \\ 0 & 0 & 1 \end{pmatrix} \quad (2.3b)$$

in two and three dimensions respectively. The inverse transformation is simply computed according to $\mathbf{x} = T^{-1}\mathbf{x}'$.

Numerical parameters There are several parameters, that describe the numerical characteristics of the generation of the random field

- `nMode`, which is the number of grid points in the Fourier space.

The parameter `nMode` is specifying the number of grid points in the Fourier space by which the integral over the spectral density is approximated. The more points the better the approximation and the more details can be represented in the generated numerical random field.