FracG V 0.1

FractureGraph - Fault and fracture analysis and meshing software

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What is FracG?

FRACG is a command line based software that performs analysis of discontinuity data (e.g. faults and fractures). The input need to be provided as line-vector data in shape-file format. The software performs a statistical analysis on the data that can be extended by providing relevant raster data in GeoTIFF format. The main analysis derives metrics from a graph representation. The line data can also be converted into 2D and 3D finite element meshes. Note that the 3D meshing is considered experimental at the moment and only a vertical extrusion to a specified depth can be performed at the moment.

Installation

FracG is designed for Debian GNU/Linux and requires third-party libraries as outlined below. First, get the latest GDAL/OGR version, add the PPA to your sources:

```
sudo add-apt-repository ppa:ubuntugis/ppa && sudo apt-get update
```

Now the necessary libraries can be installed from the terminal:

```
sudo apt-get install build-essential \
libgdal-dev \
libboost-all-dev \
liblapack-dev \
libblas-dev \
libgsl-dev \
libgmsh-dev \
libgrmadillo-dev
```

Export the environmental variables for gdal

```
export CPLUS_INCLUDE_PATH=/usr/include/gdal
export C_INCLUDE_PATH=/usr/include/gdal
```

To obtain *gmsh* visit http://gmsh.info/. Note that open cascade is used for creating the mesh. Three options are available for compilation:

xutils

This options uses a shell script and and Imake file.

```
sudo apt-get install xutils-dev
```

If the libraries are all installed correctly the permission for the file "install.sh" need to be changed.

```
cd ...FracG
sudo chmod -x install.sh
```

Now the installation script can be executed:

```
sudo ./install.sh
```

FracG can now be executed from the command line.

cmake

This option is recommended

```
sudo apt-get install cmake
```

In the FracG directory, type;

```
mkdir build \
cd build \
cmake .. \
sudo make install
```

FracG can now be executed from the command line.

Manual compilation

You can compile FracG without the install script or cmake.

```
cd ...FracG/src
g++ -o FracG main.cpp graph.cpp GeoRef.cpp geometrie.cpp stats.cpp model.cpp
-larmadillo -lgsl -lgdal -lgmsh
```

Executing FracG

After installation with cmak FracG will be set as a global executable in your environment. In a terminal choose your current directory in which files you wish to analyse are located. The first argument is a shape file containing the fault or fracture traces. This is always required. The second and thirds optional arguments are the raster file in GeoTiff format and a point shape file containing source and target node for shortest path.

Options

FracG has several parameters that can defined by the user. To see what options are available type:

The optional parameters can be set after after the name of input file. They do not have to be in a specific order and you can add as many of them as you want:

```
FracG <.shp> --option1 --option2 ...
```

Input files

The input files can be defined on the command line. This might be necessary if several input parameters should be user-defined. In the following, we list the optional parameters defined by a keyword, their data-type and their default value.

```
shapefile | < std :: string >
```

default: na

Path/name of the line-shape-file including extension.

```
raster\_file \mid < std :: string >
```

default: na

Path/name of the raster-file in GeoTIFF format including extension. This file has to be in the same reference system as the line-shape file.

```
\mathbf{source\_file} \mid < std :: string >
```

default: na

Path/name of the point-shape-file including extension. This file needs to contain two points that will be used for computing the shortest path and maximum flow between them. If not source file is given the shortest path will not be computed. This file has to be in the same reference system as the line-shape file.

Output

```
\mathbf{out\_dir} \mid < std :: string >
```

default: fracg_output_ + <name-of-shapefile>

The main directory in which all results will be written.

${\bf graph_results_file} \mid < std :: string >$

default: graph_vertices & graph_branches

Filename to save graph analysis results to.

$graph_results_folder \mid < std :: string >$

default: graph

Foldername to save graph analysis results to.

Correction parameters and distances

 $dist_thresh \mid < double >$

default: 1

Distances under this distance threshold will be considered the same location. Used for merging line segments and as distance in the point index map of the graph. The units are meters.

$angl_threshold \mid < double >$

default: 25

Maximum orientation difference in degrees for merging two segments whose tips are with the critical distance.

$dfd_{-}threshold \mid < double >$

default: 1 Threshold of cumulative discrete frechet distance of two line-strings. If the cumulative distance is below this threshold the lines will be considered duplicates and one of them will be removed.

$\mathbf{split_dist_thresh} \mid < double >$

 $default: = dist_thresh$

Distance threshold to use in splitting faults into segments, for considering nearby but separate faults to actually overlap. Used for fixing flaws in digitisation leading to false intersection classification. The units are in meters.

$spur_dist_thresh \mid < double >$

 $default: = dist_thresh$

Distance threshold to use in removing spurs, remove spurs which are shorter than this distance. Used to correct for false intersection classification. The units are in meters.

${\bf classify_lineaments_dist} \mid < double >$

 $default: = dist_thresh$

Distance used in to classify lineaments in terms of intersection number along their trace. This distance represents the buffer width around the line and intersections within this distance are counted for the classification. The units are in meters.

$raster_stats_dist \mid < double >$

default: = 1.25

Distance used for analysing raster data for the line-strings. The distance is the buffer width around the traces for computing mean values, the length of the profile lines for computing cross gradient, parallel gradients and cumulative cross-gradients along the line-strings. The unit is the number of pixels of the raster of the input raster. Note that the distance in meters is derived as the mean of the x- and y-cellsize multiplied by this factor.

$raster_spacing \mid < double >$

default: 1000

Pixel size of output density/intensity maps. The units are in meters.

$raster_spacing2 \mid < double >$

default: 500

Pixel size of output distance maps. The units are in meters.

$isect_search_size \mid < double >$

 $default: = raster_spacing = 1000$

Search for intersections within this distance. Using a circular sampling window this is the radius of the window. The units are in meters.

resample |< bool >

default: false

Resampling all created raster files to a 10^{th} of the initial cell size using cubic spline interpolation.

Statistical parameters

$angle_param_penalty \mid < double >$

default: 2

Penalty per parameter, when fitting Gaussian distributions to the angle distribution.

$scanline_count \mid < int >$

default: 100

Number scanlines to construct.

scanline_spaceing |< double>

default: 10

Minimum spacing of scanlines in meters.

$graph_min_branches \mid < int >$

default: 100

Number of branches that build a connected component. Connected components with branch numbers above this threshold will be analyses separately in addition to the analysis of the entire dataset.

$component \mid < int >$

default: -1

If greater than zero extract this connected component from the graph and build a line shape-file from it.

Maximum flow options

$max_flow_cap_type \mid < std :: string >$

default: 1

Type of capacity to use in maximum flow calculations, I for length, o for orientation, lo for both.

$max_flow_gradient_flow_direction \mid < std :: string >$

default: right

Target direction of the gradient-based maximum flow (towards left, right, top, or bottom).

$max_flow_gradient_pressure_direction \mid < std :: string >$

default: right

Target direction of the gradient-based maximum flow pressure (towards left, right, top, or bottom).

$max_flow_gradient_border_amount \mid < double >$

default: 0.05; For gradient-based maximum flow, the border features are those that intersect with the bounding box that is reduced by this amount (0 to 1).

Gmsh options

$gmsh_cell_count \mid < int >$

default: 10

Target element count in x and y direction. For rectangular domains this will be the target mean element size along x and y. This will yield the usual characteristic length (cl) of the model

$gmsh_show_output \mid < bool >$

default: false

Show out put of gmsh while meshing and the final mesh in the gmsh GUI.

$gmsh_min_cl \mid < double >$

default: cl/10

Minimum characteristic length. By default this is will be a 10^{th} of the usual characteristic length (cl).

$gmsh_max_dist \mid < double >$

default: cl/2

Maximum distance for refinement around side-sets in 2D.

$gmsh_min_dist \mid < double >$

default: cl/4

Minimum distance for refinement around side-sets in 2D.

default: 100

$gmsh_ext_depth \mid < double >$

Extrusion depth for 3D mesh in meters.

$gmsh_name_ss \mid < bool >$

default: false

Name sideset individually.

$gmsh_sample_cell_count \mid < int >$

default: 2

Number of sampling windows from which 2D-meshes should be generated.

$gmsh_sample_count \mid < int >$

default: 10

Target element count in x and y direction for the sampling windows. For rectangular domains this will be the target mean element size along x and y. This will yield the usual characteristic length (cl) of the model.

$gmsh_sample_show_output \mid < bool >$

default: false

Show out put of gmsh while meshing and the final mesh in the gmsh GUI for every sampling window.

$gmsh_in_show_meters | < bool >$

default: false

Convert coordinates into meters. This can be necessary fro small scale models.

Trouble shooting

FracG comes without any warranty. If you encounter any bugs please let us know and we will do our best to fix them as soon as possible. In some cases it might be necessary to install Aramdillo manually. In that case please follow the instruction on the web page: ARMADILLO

Please send any inquires or bug-reports to us via email.