



# **dfnWorks Documentation**

***Release 2.0***

**Subsurface Flow and Transport Team, LANL, LA-UR-17-22216**

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## INTRODUCTION

dfnWorks is a parallelized computational suite to generate three-dimensional discrete fracture networks (DFN) and simulate flow and transport. Developed at Los Alamos National Laboratory, it has been used to study flow and transport in fractured media at scales ranging from millimeters to kilometers. The networks are created and meshed using dfnGen, which combines FRAM (the feature rejection algorithm for meshing) methodology to stochastically generate three-dimensional DFNs with the LaGriT meshing toolbox to create a high-quality computational mesh representation. The representation produces a conforming Delaunay triangulation suitable for high performance computing finite volume solvers in an intrinsically parallel fashion. Flow through the network is simulated with dfnFlow, which utilizes the massively parallel subsurface flow and reactive transport finite volume code PFLOTTRAN. A Lagrangian approach to simulating transport through the DFN is adopted within dfnTrans to determine pathlines and solute transport through the DFN. Applications of the dfnWorks suite include nuclear waste repository science, hydraulic fracturing and CO<sub>2</sub> sequestration.

To run a workflow using the dfnWorks suite, the pydfnworks package is highly recommended. pydfnworks calls various tools in the dfnWorks suite; its aim is to provide a seamless workflow for scientific applications of dfnWorks.

### 1.1 Citing dfnWorks

Hyman, J. D., Karra, S., Makedonska, N., Gable, C. W., Painter, S. L., & Viswanathan, H. S. (2015). dfnWorks: A discrete fracture network framework for modeling subsurface flow and transport. *Computers & Geosciences*, 84, 10-19.

*BibTex:*

```
@article{hyman2015dfnworks,
  title={dfnWorks: A discrete fracture network framework
for modeling subsurface flow and transport},
  author={Hyman, Jeffrey D and Karra, Satish and Makedonska,
Nataliia and Gable, Carl W and Painter, Scott L
and Viswanathan, Hari S},
  journal={Computers \& Geosciences},
  volume={84},
  pages={10--19},
  year={2015},
  publisher={Elsevier}
}
```

### 1.2 What's new in v2.0?

- New dfnGen C++ code which is much faster than the Mathematica dfnGen. This code has successfully generated networks with 350,000+ fractures.

- Increased functionality in the pydfnworks package for more streamlined workflow from dfnGen through visualization.

## 1.3 Where can one get dfnWorks?

dfnWorks 2.0 can be downloaded from <https://github.com/dfnWorks/dfnWorks-Version2.0>

v1.0 can be downloaded from <https://github.com/dfnWorks/dfnWorks-Version1.0>

## 1.4 Installation

Tools that you will need to run the dfnWorks work flow are described in this section. VisIt and ParaView, which enable visualization of desired quantities on the DFNs, are optional, but at least one of them is highly recommended for visualization. CMake is also optional but allows faster IO processing using C++.

### 1.4.1 Python

pydfnworks is supported on Python 2.7. The software authors recommend using the Anaconda 2.7 distribution of Python, available at <https://www.continuum.io/>. pydfnworks requires the numpy and h5py modules to be installed.

### 1.4.2 pydfnworks

The source for pydfnworks can be found in the dfnWorks suite, in the folder pydfnworks.

### 1.4.3 dfnGen

dfnGen primarily involves two steps:

1. FRAM - Create DFN: Using the fractured site characterization networks are constructed using the feature rejection algorithm for meshing
2. LaGriT - Mesh DFN: The LaGriT meshing tool box is used to create a conforming Delaunay triangulation of the network.

#### FRAM

FRAM (the feature rejection algorithm for meshing) is executed using the dfnGen C++ source code, contained in the dfnGen folder of the dfnWorks repository.

#### LaGriT

The *LaGriT* (<http://lagrit.lanl.gov>) meshing toolbox is used to create a high resolution computational mesh representation of the DFN in parallel. An algorithm for conforming Delaunay triangulation is implemented so that fracture intersections are coincident with triangle edges in the mesh and Voronoi control volumes are suitable for finite volume flow solvers such as FEHM and PFLOTRAN.



### 1.4.4 dfnFlow

You will need one of either PFLOTRAN or FEHM to solve for flow using the mesh files from LaGriT.

#### PFLOTRAN

PFLOTRAN (<http://www.pflotran.org>) is a massively parallel subsurface flow and reactive transport code. PFLOTRAN solves a system of partial differential equations for multiphase, multicomponent and multiscale reactive flow and transport in porous media. The code is designed to run on leadership-class supercomputers as well as workstations and laptops.

#### FEHM

FEHM (<http://fehm.lanl.gov>) is a subsurface multiphase flow code developed at Los Alamos National Laboratory.

### 1.4.5 dfnTrans

dfnTrans is a method for resolving solute transport using control volume flow solutions obtained from dfnFlow on the unstructured mesh generated using dfnGen. We adopt a Lagrangian approach and represent a non-reactive conservative solute as a collection of indivisible passive tracer particles.

### 1.4.6 CMake

CMake (<https://cmake.org/>) is an open-source, cross-platform family of tools designed to build, test and package software. It is needed to use C++ for processing files at a bottleneck IO step of dfnWorks. Using C++ for this file processing optional but can greatly increase the speed of dfnWorks for large fracture networks. Details on how to use C++ for file processing are in the scripts section of this documentation.

### 1.4.7 VisIt

VisIt is a parallel, open-source visualisation software. PFLOTRAN can output in `.xmf` and `.vtk` format. These can be imported in VisIt for visualization.

Instructions for downloading and installing VisIt can be found at <https://wci.llnl.gov/codes/visit/download.html>

### 1.4.8 Paraview

Paraview is a parallel, open-source visualisation software. PFLOTRAN can output in `.xmf` and `.vtk` format. These can be imported in Paraview for visualization.

Instructions for downloading and installing Paraview can be found at <http://www.paraview.org>

## 1.5 Using pydfnworks in your Python scripts

To access the functionality of pydfnworks, the user must include the following line at the top of any Python script

```
import pydfnworks
```

Before doing this, one needs to ensure that the `pydfnworks` directory is in the `PYTHONPATH`. This can be done by configuring `cshrc` or `bashrc` files. Alternatively, one can add the `pydfnworks` path using `sys.path.append()` in their driver script.

## 1.6 About this manual

This manual comprises of information on setting up inputs to `dfnGen`, `dfnTrans` and `PFLOTRAN`, as well as details on the `pydfnworks` module: *pydfnworks*. Finally, the manual contains a short tutorial with prepared examples that can be found in the `tests` directory of the `dfnWorks` repository, and a description of some applications of the `dfnWorks` suite.

## 1.7 Contributors

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- Nataliia Makedonska
- Jeffrey Hyman
- Jeremy Harrod (now at Spectra Logic)
- Quan Bui (now at University of Maryland)
- Carl Gable
- Scott Painter (now at ORNL)
- Hari Viswanathan
- Nathaniel Knapp

## 1.8 Contact

For any questions about `dfnWorks`, please email [dfnworks@lanl.gov](mailto:dfnworks@lanl.gov).

## 1.9 Copyright information

Documentation:

LA-UR-17-22216

Software copyright:

LA-CC-17-027

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## 2.1 Keywords

The following is an example input file with all keywords and explanation of each keyword.

```
//=====
//General Options & Fracture Network Parameters:
/*
stopCondition: 0 /* 0: stop once nPoly fractures are accepted (Defined
below) 1: stop once all family's p32 values are equal or greater than the
families target p32 values (defined in stochastic family sections) */

nPoly: 3 /* Used when stopCondition = 0 Total number of fractures you would
like to have in the domain you defined. The program will complete once you
have nPoly number of fractures, maxPoly number of polygon/fracture
rejections, rejPoly number of rejections in a row, or reach a specified
fracture cluster size if using stoppingParameter = -largestSize */

outputAllRadii: 0 /* 0: Do not output all radii file. 1: Include file of
all raddii (accepted+rejected fractures) in output files. */

domainSize: {1,1,1} /* Mandatory Parameter. Creates a domain with dimension
x*y*z centered at the origin.*/

numOfLayers: 0 //number of layers

layers: {-500,0} {0,500}

/* Layers need to be listed line by line Format: {minZ, maxZ}

The first layer listed is layer 1, the second is layer 2, etc Stochastic
families can be assigned to theses layers (see stochastic shape family
section) */

h: 0.050 /* Minimum fracture length scale(meters) Any fracture with a
feature, such as and intersection, of less than h will be rejected. */

//=====//
/* Fracture Network Parameters:
*/
```

```

tripleIntersections: 1 /* Options:      0: Off 1: On    */

printRejectReasons: 0 /* Useful in debugging, This option will print all
fracture rejection reasons as they occur.  0: disable 1: print all rejection
reasons to screen */

visualizationMode: 0 /* Options: 0 or 1 Used during meshing: 0: creates a
fine mesh, according to h parameter; 1: produce only first round of
triangulations. In this case no modeling of flow and transport is possible.
*/

seed: 92731535 //Seed for random generator.

domainSizeIncrease: {0,0,0} //temporary size increase for inserting fracture
centers outside domain //increases the entire width by this ammount. So,
{1,1,1} will increase //the domain by adding .5 to the +x, and subbtracting
.5 to the -x, etc

keepOnlyLargestCluster: 0 /* 0 = Keep any clusters which connects the
specified boundary faces in boundaryFaces option below

    1 = Keep only the largest cluster which connects the specified
    boundary faces in boundaryFaces option below */

ignoreBoundaryFaces: 1 /* 0 = use boundaryFaces option below

    1 = ignore boundaryFaces option and keep all clusters and will still
    remove fractures with no intersections */

boundaryFaces: {1,1,0,0,0,0} /* DFN will only keep clusters with
connections to domain boundaries which are set to 1:

    boundaryFaces[0] = +X domain boundary boundaryFaces[1] = -X domain
    boundary boundaryFaces[2] = +Y domain boundary boundaryFaces[3] = -Y
    domain boundary boundaryFaces[4] = +Z domain boundary boundaryFaces[5] =
    -Z domain boundary

    Be sure to set ignoreBoundaryFaces to 0 when using this feature.

rejectsPerFracture: 10 /*If fracture is rejected, it will be re-translated
to a new position this number of times.

    This helps hit distribution targets for stochastic
    families (Set to 1 to ignore this feature)

//=====
//                               Shape and Probability Parameters
//=====

//user rectangles and user Ellipses defined in their cooresponding files

```

```

famProb: {.5,.5} /* Probability of occurrence of each family of randomly
distributed rectangles and ellipses. User-ellipses and user-rectangles
insertion will be attempted with 100% likelihood, but with possibility they
may be rejected. The famProb elements should add up to 1.0 (for %100). The
probabilities are listed in order of families starting with all stochastic
ellipses, and then all stochastic rectangles.

    For example: If then there are two ellipse families, each with
    probabiliy .3, and two rectangle families, each with probabiliy .2,
    famProb will be: famProb: {.3,.3,.2,.2} Notice: famProb elements add to 1
    */

/*=====*/
//=====
//          Elliptical Fracture Options
//      NOTE: Number of elements must match number of ellipse families //
(first number in nShape input parameter)
//=====
/*=====*/

//Number of ellipse families nFamEll: 0 //Having this option = 0 will ignore
all rectangle family variables

eLayer: {0,0} /* Defines which domain the family belings to. Layer 0 is the
entire domain. Layers numbered > 0 coorespond to layers defined above 1
corresponst to the first layer listed, 2 is the next layer listed, etc */

//edist is a mandatory parameter if using statistically generated ellipses
edistr: {2,3} /* Ellipse statistical distribution options: 1 - lognormal
distribution 2 - truncated power law distribution 3 - exponential
distribution 4 - constant */

ebetaDistribution: {1,1} /* Beta is the rotation around the polygon's
normal vector, with the polygon centered on x-y plane at the orgin

    0 - uniform distribution [0, 2PI]    1 - constant angle
    (specefied below by "ebeta")        */

e_p32Targets: {.1,.1} /* Elliptical families target fracture intensity per
family. When using stopCondition = 1 (defined at the top of the input
file), families will be inserted untill the families desired fracture
intensity has been reached. Once all families desired fracture intensity
has been met, fracture generation will be complete. */

//=====
// Parameters used by all stochastic ellipse families // Mandatory
Parameters if using statistically generated ellipses

easpect: {1,1} /* Aspect ratio. Used for lognormal and truncated power law
distribution. */

enumPoints: {12, 12} /*Number of vertices used in creating each elliptical
fracture family. Number of elements must match number of ellipse families

```

```
(first number in nShape) */

eAngleOption: 0      /* All angles for ellipses: 0 - degrees 1 - radians
(Must use numerical value for PI) */

etheta: {-45, 45,} /*Ellipse fracture orientation. The angle the normal
vector makes with the z-axis */

ephi: {0,0} /* Ellipse fracture orientation. The angle the projection of
the normal onto the x-y plane makes with the x-axis */

ebeta: {0, 0} /* rotation around the normal vector */

ekappa: {8,8} /*Parameter for the fisher distribnShaprutions. The bigger,
the more similar (less diverging) are the elliptical familiy's normal
vectors */

//=====
// Options Specific For Ellipse Lognormal Distribution (edistr=1): //
Mandatory Parameters if using ellispes with lognormal distribution

//      NOTE: Number of elements must match number of //
ellipse families (first number in nShape)

eLogMean: {2} //Mean value For Lognormal Distribution.

eLogMax: {100} eLogMin: {1}

esd: {.5} // Standard deviation for lognormal distributions of ellipses

//=====
//      Options Specific For Ellipse Exponential Distribution (edistr=3): //
Mandatory Parameters if using ellispes with exponential distribution

eExpMean: {2} //Mean value for Exponential Distribution      eExpMax: {3}
//Mean value for Exponential Distribution      eExpMin: {1} //Mean value
for Exponential Distribution

//=====
//      Options Specific For Constant Size of ellipses (edistr=4):

econst: {10, 10, 10} // Constant radius, defined per family

//=====
// Options Specific For Ellipse Truncated Power-Law Distribution (edistr=2)
// Mandatory Parameters if using ellispes with truncated power-law dist.

// NOTE: Number of elements must match number //      of ellipse families
(first number in nShape)

emin: {1} // Minimum radius for each ellipse family. // For power law
distributions.

emax: {6} // Maximum radius for each ellipse family. // For power law
distributions.
```



```

ealpha: {2.4} // Alpha. Used in truncated power-law // distribution
calculation

/*=====*/
/*=====*/ /*
Rectangular Fractures Options      /* /* NOTE: Number of elements must
match number of rectangle families */ /*      (second number in nShape
parameter)      */
/*=====*/
/*=====*/

//Number of rectangle families nFamRect: 0 //Having this option = 0 will
ignore all rectangle family variables

rLayer: {0,0} /* Defines which domain the family belongs to. Layer 0 is the
entire domain. Layers numbered > 0 correspond to layers defined above 1
corresponds to the first layer listed, 2 is the next layer listed, etc */

/*rdist is a mandatory parameter if using statistically generated rectangles
*/ rdistr: {2,3} /* Rectangle statistical distribution options: 1 -
lognormal distribution 2 - truncated power law distribution 3 - exponential
distribution 4 - constant */

rbetaDistribution: {1,1} /* Beta is the rotation/twist about the z axis
with the polygon centered on x-y plane at the origin before rotation into 3d
space

                0 - uniform distribution [0, 2PI]      1 - constant angle
                (specified below by "rbeta")

                */

r_p32Targets: {.1,.1} /* Rectangle families target fracture intensity per
family. When using stopCondition = 1 (defined at the top of the input
file), families will be inserted until the families desired fracture
intensity has been reached. Once all families desired fracture intensity has
been met, fracture generation will be complete. */

//=====
// Parameters used by all stochastic rectangle families // Mandatory
Parameters if using statistically generated rectangles

raspect: {1,1} /* Aspect ratio */

rAngleOption: 0 /* All angles for rectangles: 0 - degrees 1 - radians
(must be numerical value, cannot use "Pi") */

rtheta: {-45,45} /*Rectangle fracture orientation. The angle the normal
vector makes with the z-axis */

rphi: {0,45} /* Rectangle fracture orientation. The angle the projection of
the normal onto the x-y plane makes with the x-axis */

```

```

rbeta: {0,0}    /* rotation around the normal vector */

rkappa: {8,8}  /*Parameter for the fisher distributions. The bigger, the
more similar (less diverging) are the rectangle familiy's normal vectors */

//=====
// Options Specific For Rectangle Lognormal Distribution (rdistr=1): //
Mandatory Parameters if using rectangles with lognormal distribution

rLogMean: {1.6} /*For Lognormal Distribution. Mean radius (1/2 rectangle
length) in lognormal distribution for rectangles. */

rLogMax: {100} rLogMin: {1}

rsd: {.4}      /* Standard deviation for lognormal distributions of
rectangles */

//=====
// Options Specific For Rectangle Truncated Power-Law Distribution
(rdistr=2): // Mandatory Parameters if using rectangles with power-law
distribution

rmin: {1,1}     /* Minimum radius for each rectangle family. For power
law distributions. */

rmax: {6,5}     /* Maximum radius for each rectangle family. For power law
distributions. */

ralpha: {2.4,2.5} // Alpha. Used in truncated power-law // distribution
calculation

/*=====*/
/* Options Specific For Rectangle Exponential Distribution (edistr=3):
*/ /* Mandatory Parameters if using rectangles with exponential
distribution */

rExpMean: {2} //Mean value for Exponential Distribution rExpMax: {100}
rExpMin: {1}

/*=====*/
/* Options Specific For Constant Size of rectangles (edistr=4):
*/

rconst: {4,4} // Constant radius, defined per rectangular family

/*=====*/
/*=====*/
/* User-Specified Ellipses
*/ /* Mandatory Parameters if using user-ellipses
*/ /* NOTE: Number of elements must match number of user-ellipse families
*/ /*(third number in nShape parameter)
*/
/*=====*/
/* NOTE: Only one user-ellipse is placed into the domain per defined
user-ellipse, with possibility of being rejected */

```

```

userEllipsesOnOff: 0    //0 - User Ellipses off //1 - User Ellipses on

UserEll_Input_File_Path: ./TestCases/test/uEllInput.dat

/*=====*/
/*=====*/
/* User-Specified Ellipses
*/ /* Mandatory Parameters if using user-ellipses
*/ /* NOTE: Number of elements must match number of user-ellipse families.
*/ /* NOTE: Only one user-ellipse is placed into the domain per defined
*/ /* user-ellipse, with possibility of being rejected
*/
/*=====*/
/*=====*/

userEllByCoord: 0 /* 0 - User ellipses defined by coordinates off 1 - User
ellipses defined by coordinates on */

EllByCoord_Input_File_Path:
/home/jharrod/GitProjects/DFNGen/DFNC++Version/inputFiles/
userPolygons/ellCoords.dat


/*=====*/
/* User-Specified Rectangles
*/ /* Mandatory Parameters if using user-rectangles
*/ /* NOTE: Number of elements must match number of user-ellipse families
*/ /* (fourth number in nShape parameter)
*/
/*=====*/
/* NOTE: Only one user-rectangle is placed into the domain per defined
user-rectangle, with possibility of being rejected */


userRectanglesOnOff: 1    //0 - User Rectangles off //1 - User Rectangles on

UserRect_Input_File_Path: /home/nknapp/dfnWorks-Version2.0/
tests/define_4_user_rects.dat

/*=====*/
/* If you would like to input user specified rectangles according to their
coordinates, you can use the parameter userDefCoordRec. In that case, all
of the user specified rectangles will have to be according to coordinates.
*/

userRecByCoord: 0 // 0 - user defined rectangles not used // 1 - user
defined rectangles used and defined by input file:

RectByCoord_Input_File_Path: ./inputFiles/userPolygons/rectCoords.dat


/*WARNING: userDefCoordRec can cause LaGriT errors because the polygon
vertices are not put in clockwise or counter-clockwise order. If errors
(Usually seg fault during meshing if using LaGriT), try to reorder the points
till u get it right. Also, coordinates must be co-planar */

/*=====*/

```

```

// Aperture [m] /* Mandatory parameter, and can be specified in several
ways: - 1)meanAperture and stdAperture for using LogNormal distribution. -
2)apertureFromTransmissivity, first transmissivity is defined, and then,
using a cubic law, the aperture is calculated; - 3)constantAperture, all
fractures, regardless of their size, will have the same aperture value; -
4)lengthCorrelatedAperture, aperture is defined as a function of fracture
size*/

//NOTE: Only one aperture type may be used at a time

aperture: 3 //choise of aperture option described above

//(**** 1)meanAperture and stdAperture for using LogNormal
distribution.*****) meanAperture: -3 /*Mean value for aperture using
normal distribution */ stdAperture: 0.8 //Standard deviation

/*(***** 2)apertureFromTransmissivity, first transmissivity is defined, and
then, using a cubic law, the aperture is calculated;*****/
apertureFromTransmissivity: {1.6e-9, 0.8} /* Transmissivity is calculated as
transmissivity = F*R^k, where F is a first element in
aperturefromTransmissivity, k is a second element and R is a mean radius of
a polygon. Aperture is calculated according to cubic law as
b=(transmissivity*12)^1/3 */

/*(***** 3)constantAperture, all fractures, regardless of their size, will
have the same aperture value; *****/

constantAperture: 1e-5 //Sets constant aperture for all fractures

/*(***** 4)lengthCorrelatedAperture, aperture is defined as a function of
fracture size *****/

lengthCorrelatedAperture: {5e-5, 0.5} /*Length Correlated Aperture Option:
Aperture is calculated by: b=F*R^k, where F is a first element in
lengthCorrelatedAperture, k is a second element and R is a mean radius of a
polygon.*/

//=====
//Permeability /* Options: 0: Permeability of each fracture is a function of
fracture aperture, given by k=(b^2)/12, where b is an aperture and k is
permeability 1: Constant permeabilty for all fractures */

permOption: 1 //See above for options

constantPermeability: 1e-12 //Constant permeability for all fractures

//=====

outputAcceptedRadiiPerFamily:1 /* output radii files for each family
containing the final radii chosen */

disableFram:0 /* 0 if FRAM (feature rejection algorithm for meshing) is
disabled, 1 otherwise */

outputFinalRadiiPerFamily:1 /* output radii files for each family containing
the final radii chosen */

```

```
insertUserRectanglesFirst:1 /* 1 if user defined rectangles should be
inserted first, 0 otherwise */

forceLargeFractures:0 /* Force large fractures to be included in the network
*/

radiiListIncrease: 0.1 /* Increase the length of the initially generated
radii list (before rejections) by this percentage */

removeFracturesLessThan: 0 /*Used to change the lower cutoff of fracture
size*/
```



**DFNFLOW**

*dfnFlow* involves using flow solver such as PFLOTTRAN or FEHM. PFLOTTRAN is recommended if large number of fractures ( $> O(1000)$ ) are involved in a network. Using the function calls that are part of *dfnworks\_python*, one can create the mesh files needed to run PFLOTTRAN. This will involve creating unstructured mesh file *\*uge* as well as the boundary *\*ex* files. Please see the PFLOTTRAN user manual at <http://www.pfлотran.org> under unstructured *explicit* format usage for further details. An example input file for PFLOTTRAN is provided in the repository. Please use this as a starting point to build your input deck.

Below is a sample input file. Refer to the PFLOTTRAN user manual at <http://www.pfлотran.org> for input parameter descriptions.

```
# Jan 13, 2014
# Nataliia Makedonska, Satish Karra, LANL
#=====

SIMULATION
  SIMULATION_TYPE SUBSURFACE
  PROCESS_MODELS
    SUBSURFACE_FLOW flow
    MODE RICHARDS
  /
/
END
SUBSURFACE

DFN

#===== discretization =====
GRID
  TYPE unstructured_explicit full_mesh_vol_area.uge
  GRAVITY 0.d0 0.d0 0.d0
END

#===== fluid properties =====
FLUID_PROPERTY
  DIFFUSION_COEFFICIENT 1.d-9
END

DATASET Permeability
  FILENAME dfn_properties.h5
END

#===== material properties =====
MATERIAL_PROPERTY soil1
  ID 1
```

```

POROSITY 0.25d0
TORTUOSITY 0.5d0
CHARACTERISTIC_CURVES default
PERMEABILITY
    DATASET Permeability
/
END

#===== characteristic curves =====
CHARACTERISTIC_CURVES default
SATURATION_FUNCTION VAN_GENUCHTEN
    M 0.5d0
    ALPHA 1.d-4
    LIQUID_RESIDUAL_SATURATION 0.1d0
    MAX_CAPILLARY_PRESSURE 1.d8
/
PERMEABILITY_FUNCTION MUALEM_VG_LIQ
    M 0.5d0
    LIQUID_RESIDUAL_SATURATION 0.1d0
/
END

#===== output options =====
OUTPUT
    TIMES s 0.01 0.05 0.1 0.2 0.5 1
# FORMAT TECPLOT BLOCK
PRINT_PRIMAL_GRID
FORMAT VTK
MASS_FLOWRATE
MASS_BALANCE
VARIABLES
    LIQUID_PRESSURE
    PERMEABILITY
/
END

#===== times =====
TIME
    INITIAL_TIMESTEP_SIZE 1.d-8 s
    FINAL_TIME 1.d0 d==
    MAXIMUM_TIMESTEP_SIZE 10.d0 d
    STEADY_STATE
END

# REFERENCE_PRESSURE 1500000.

#===== regions =====
REGION All
COORDINATES
    -1.d20 -1.d20 -1.d20
    1.d20 1.d20 1.d20
/
END

REGION inflow
    FILE pboundary_left_w.ex
END

```



```

REGION outflow
  FILE pboundary_right_e.ex
END

#===== flow conditions =====
FLOW_CONDITION initial
  TYPE
    PRESSURE dirichlet
  /
  PRESSURE 1.01325d6
END

FLOW_CONDITION outflow
  TYPE
    PRESSURE dirichlet
  /
  PRESSURE 1.d6
END

FLOW_CONDITION inflow
  TYPE
    PRESSURE dirichlet
  /
  PRESSURE 2.d6
END

#===== condition couplers =====
# initial condition
INITIAL_CONDITION
  FLOW_CONDITION initial
  REGION All
END

BOUNDARY_CONDITION INFLOW
  FLOW_CONDITION inflow
  REGION inflow
END

BOUNDARY_CONDITION OUTFLOW
  FLOW_CONDITION outflow
  REGION outflow
END

#===== stratigraphy couplers =====
STRATA
  REGION All
  MATERIAL soil1
END

END_SUBSURFACE

```



## DFNTRANS

dfnTrans is a method for resolving solute transport using control volume flow solutions obtained from dfnFlow on the unstructured mesh generated using dfnGen. We adopt a Lagrangian approach and represent a non-reactive conservative solute as a collection of indivisible passive tracer particles. Particle tracking methods (a) provide a wealth of information about the local flow field, (b) do not suffer from numerical dispersion, which is inherent in the discretizations of advection–dispersion equations, and (c) allow for the computation of each particle trajectory to be performed in an intrinsically parallel fashion if particles are not allowed to interact with one another or the fracture network. However, particle tracking on a DFN poses unique challenges that arise from (a) the quality of the flow solution, (b) the unstructured mesh representation of the DFN, and (c) the physical phenomena of interest. The flow solutions obtained from dfnFlow are locally mass conserving, so the particle tracking method does not suffer from the problems inherent in using Galerkin finite element codes.

dfnTrans starts from reconstruction of local velocity field: Darcy fluxes obtained using dfnFlow are used to reconstruct the local velocity field, which is used for particle tracking on the DFN. Then, Lagrangian transport simulation is used to determine pathlines through the network and simulate transport. It is important to note that dfnTrans itself only solves for advective transport, but effects of longitudinal dispersion and matrix diffusion, sorption, and other retention processes are easily incorporated by post-processing particle trajectories. The detailed description of dfnTrans algorithm and implemented methodology is in Makedonska, N., Painter, S. L., Bui, Q. M., Gable, C. W., & Karra, S. (2015). Particle tracking approach for transport in three-dimensional discrete fracture networks. *Computational Geosciences*, 19(5), 1123–1137.

All source files of C code of dfnTrans are in ParticleTracking/ directory of dfnWorks 2.0. It compiles under linux/mac machines using makefile. In order to run transport, first, all the parameters and paths should be set up in the PTDFN Control file, PTDFN\_control.dat. Then, the following command should be run: ./dfnTrans <PTDFN\_control.dat

PTDFN\_control.dat file sets all necessary parameters to run particle tracking in dfnWorks. Below is one PTDFN\_control.dat example that includes a short explanation of each parameter setting:

```

/*****/ /*
CONTROL FILE FOR PARTICLE TRACKING IN DISCRETE FRACTURE NETWORK */
/*****/

// 1. Define paths of files that were generated by DFNGEN and contain all
the information about computational grid.

/***** INPUT FILES: grid *****/
/**** input files with grid of DFN, mainly it's output of dfnGen *****/
param: params.txt //generated by dfnGen, contains the number of fractures in
the DFN

poly: poly_info.dat //generated by dfnGen, contains the normal vectors of
each polygon; used for rotation each fracture

inp: full_mesh.inp // AVS file of full mesh of DFN: positions of vertices

```

```

and connectivity list

stor: tri_fracture.stor // produced by LaGriT: the matrix of all connected
nearest neighbors faces of Voronoi polygons and their geometrical
coefficients

boundary: allboundaries.zone //list of nodes that located on domain
boundaries /* boundary conditions: reading the nodes that belong to in-flow
and out-flow boundaries. Should be consistent with those applied to obtain
steady state pressure solution (PFLOTTRAN) */ /*1 - top; 2 - bottom; 3 -
left_w; 4 - front_s; 5 - right_e; 6 - back_n */ in-flow-boundary: 3
out-flow-boundary: 5

/* The allboundaries.dat file can be modified according to task. For
example, inflow boundaries are left and top, and out flow is a bottom face
of the domain. In this case, left and top list of boundary nodes should be
combined together in allboundaries.dat and only one number given (for
example 1). dfnTrans code can't read more than one defined number for
in-flow or out-flow boundaries. Also, number of boundary zone is not
limited by 6. */

// 2. Input of Flow Solver results

/***** INPUT FILES: PFLOTTRAN flow solution *****/
PFLOTTRAN: yes //yes - the PFLOTTRAN flow solver is used

PFLOTTRAN_vel: darcyvel.dat // the Darcy flux for each Control Volume face of
the full mesh of DFN

PFLOTTRAN_cell: cellinfo.dat

/*connectivity list and area of each Control Volume face of the full mesh of
DFN. In case of using one of recent versions of PFLOTTRAN, where
PFLOTTRAN_vel file consists areas of connection, PFLOTTRAN_uge file will not
be used.*/

PFLOTTRAN_uge: lagrit_pflotran_dat/lagrit_pflotran_corrected.uge

// The current version of DFNWorks uses PFLOTTRAN /***** INPUT
FILES: FEHM flow solution *****/ /*currently we are using
PFLOTTRAN , but the code would work with FEHM, too */ FEHM: no

FEHM_fin: tri_frac.fin // fluxes and pressure outputs of FEHM

// 3. Settings of OUTPUT files with particle tracking results.

/***** OUTPUT FILES *****/
/* initial grid info structure output, useful for debugging */ out_grid: yes
// yes means that the output ASCII files, named "nodes" and "fractures",
will be generated. If you don't want this output - type "no" instead of
"yes".

/* flow field: 3D Darcy velocities: output ASCII file "Velocity3D" has an
each nodes position and its Darcy velocity, reconstructed from fluxes */
out_3dflow: yes

/* out initial positions of particles into separate ASCII file "initpos" */
out_init: yes

```

```

/* out particle trajectories tortuosity file, torts.dat */ out_tort: no

/***** output options for particles trajectories *****/
/* output frequency is set according to trajectories curvature. We check the
curvature of particles trajectory each segment, from intersection to
intersection. If it's like a straight line, then the output is less
frequent (in case of "out_curv:yes", if "no", the output file will contain
every time step of the simulation) */ out_curv: no

/* output into avs file (GMV visualization, Paraview visualization) */
out_avs: no

/* output into trajectories ascii files (veloc+posit+cell+fract+time) */
out_traj: no

/***** output directories *****/
out_dir: ptresults1 /* path and name of directory where all the particle
tracking results will be written, including those defined above*/

out_path: trajectories /*name of directory where all particle trajectories
will be saved, in out_dir path */

/* name of resultant file (in out_dir path), which contains total travel
time and final positions of particles */ out_time: partime

// 4. Options for Particles Initial Positions in DFN

/***** PARTICLES INITIAL POSITIONS *****/
/*****init_nf: if yes - the same number of
particles (init_partn) will be placed on every boundary fracture edge on
in-flow boundary, equidistant from each other *****/ init_nf: yes init_partn:
5

/*****init_eqd: if yes - particles will be placed on the same distance from
each other on all over in-flow boundary edges *****/ //
The difference between options init_eqd: and init_nf: is the following. In
case of init_eqd the total length of fracture edges on in-flow boundaries
will be calculated. Then, according to init_npart given number of particles,
the particles will be distributed equidistant over all fracture edges on
in-flow boundaries. In init_nf option, the init_partn number of particles
will be equidist in each edge of fracture on in-inflow boundaries. In this
case distance between two neighboring particles in one fracture will not be
the same as distance between two particles in other fracture. init_eqd: no
//maximum number of particles that user expects on one boundary edge
init_npart: 1

/* all particles start from the same region at in-flow boundary, in a
range {in_xmin, in_xmax,in_ymin, in_ymax, in_zmin, in_zmax} *****/
// In this option, the region on in-flow boundary should be defined
according to x, y, and z coordination of the domain. Then particles will be
placed equidistant in those part of fracture edges that cross the defined
region. If there are no fracture edge found there, the program will be
terminated, and user should redefine the region. init_oneregion: no
in_partn: 10 in_xmin: -20.0 in_xmax: 20.0 in_ymin: -20.0 in_ymax: 20.0
in_zmin: 499.0 in_zmax: 501.0

/**** all particles are placed randomly over all fracture surface (not only

```

```

on boundary edges!) *****/ // In this option
the particles will be placed on the center of randomly chosen Control Volume
cell over all cells in DFN mesh (not only on in-flow boundary).
init_random: no // total number of particles in_randpart: 110000

// 5. Flow and Fracture Parameters

/***** FLOW AND FRACTURE PARAMETERS *****/
porosity: 0.25 // porosity

density: 997.73 //fluid density

satur: 1.0

thickness: 1.0 //DFN aperture (used in case of no aperture file provided)

aperture: yes //DFN aperture

aperture_type: frac //aperture is giving per cell (type "cell") // or per
fracture (type "frac") // In the current version of DFNWorks the only an
aperture per fracture option is given.

aperture_file: aperture.dat // The ASCII file aperture.dat is produced by
dfnGen

/***** TIME *****/
timesteps: 2000000

//units of time (years, days, hours, minutes) time_units: seconds

/**** flux weighted particles*/ // in all the options of initial positions
of particles, particles can be weighted either by input flux )in case of
placing particles in in-flow boundary) of by current cell aperture (in case
of randomly defined initial positions). flux_weight: yes

/* random generator seed */ seed: 337799

// 6. Control Plane Output /***** Control Plane/Cylinder
Output *****/ // Here is another option for output. The
control Planes can be defined on any position along the flow direction. For
example, if fluid flow goes from top to bottom along Z direction
(flowdir=2), the imaginary control planes will be parallel to x-y plane and
placed each 1 m (delta_Control: 1). Each particle will have it's data output
(location, current velocity) every time it crosses control plane.

/** virtual Control planes will be build in the direction of flow. Once
particle crosses the control plane, it's position, velocity, time will
output to an ascii file. ****/ ControlPlane: no

/* the path and directory name with all particles output files */
control_out: outcontroldir

/* Delta Control Plane - the distance between control planes */
delta_Control: 1

/* ControlPlane: direction of flow: x-0; y-1; z-2 */ flowdir: 1

```

```
/*****  
END
```





## PYDFNWORKS: THE DFNWORKS PYTHON PACKAGE

The pydfnworks package allows the user to easily run dfnWorks from the command line and call dfnWorks within other python scripts. Because pydfnworks is a package, users can call individual methods from the package easily.

The pydfnworks must be setup by the user using the following command in the directory dfnWorks-Version2.0/pydfnworks/ :

python setup.py install (if the user has admin privileges), OR:

python setup.py install --user (if the user does not have admin privileges):

The documentation below includes all the methods and classes of the pydfnworks package.

### 5.1 DFNWORKS

```
class pydfnworks.DFNWORKS (jobname='', local_jobname='', dfnGen_file='', output_file='', local_dfnGen_file='', ncpu='', dfnFlow_file='', local_dfnFlow_file='', dfnTrans_file='', inp_file='full_mesh.inp', uge_file='', vtk_file='', mesh_type='dfn', perm_file='', aper_file='', perm_cell_file='', aper_cell_file='', dfnTrans_version='', num_frac='')
```

Class for DFN Generation and meshing

#### Attributes:

- `_jobname`: name of job, also the folder where output files are stored
- `_ncpu`: number of CPUs used in the job
- `_dfnGen_file`: the name of the dfnGen input file
- `_dfnFlow_file`: the name of the dfnFlow input file
- `_local` prefix: indicates that the name contains only the most local directory
- `_vtk_file`: the name of the VTK file
- `_inp_file`: the name of the INP file
- `_uge_file`: the name of the UGE file
- `_mesh_type`: the type of mesh
- `_perm_file`: the name of the file containing permeabilities
- `_aper_file`: the name of the file containing apertures
- `_perm_cell_file`: the name of the file containing cell permeabilities
- `_aper_cell_file`: the name of the file containing cell apertures

- `_dfnTrans_version`: the version of dfnTrans to use
- `_freeze`: indicates whether the class attributes can be modified
- `_large_network`: indicates whether C++ or Python is used for file processing at the bottleneck

of inp to vtk conversion

**argparse** = <module 'argparse' from '/n/swdev/packages/Ubuntu-14.04-x86\_64/anaconda-python/2.4.1/lib/python2.7/ar

**check\_input** (*input\_file*='', *output\_file*='')

#### Input Format Requirements:

- Each parameter must be defined on its own line (separate by newline)
- A parameter (key) MUST be separated from its value by a colon ':' (ie. -> key: value)
- Values may also be placed on lines after the 'key'
- Comment Format: On a line containing // or / \*, nothing after \* / or // will be processed but text before a comment will be processed

#### Kwargs:

- `input_file` (name): name of dfnGen input file
- `output_file` (name): stripped down input file for DFNGen

**commandline\_options** ()

Read command lines for use in dfnWorks.

#### Options:

- `-name` : Jobname (Mandatory)
- `-ncpu` : Number of CPUS (Optional, default=4)
- `-input` : input file with paths to run files (Mandatory if the next three options are not specified)
- `-gen` : Generator Input File (Mandatory, can be included within this file)
- `-flow` : PFLORAN Input File (Mandatory, can be included within this file)
- `-trans`: Transport Input File (Mandatory, can be included within this file)
- `-cell`: True/False Set True for use with cell based aperture and permeability (Optional, default=False)
- `-large_network`: True/False Set True to use CPP for file processing bottleneck (Optional, default=False)

**copy\_dfnTrans\_files** ()

create link to DFNTRANS and copy input file into local directory

**create\_network** ()

Execute dfnGen and print whether the generation of the fracture network failed or succeeded. The params.txt file must be there for success.

**dfnFlow** ()

Run the dfnFlow portion of the workflow.

**dfnGen** ()

#### Run the dfnGen workflow:

1. `make_working_directory`: Create a directory with name of job
2. `check_input`: Check input parameters and create a clean version of the input file

- 3. `create_network`: Create network. DFNGEN v2.0 is called and creates the network
- 4. `output_report`: Generate a PDF summary of the DFN generation
- 5. `mesh_network`: calls module `dfnGen_meshing` and runs LaGriT to mesh the DFN

**dfnTrans** ()

Copy input files for `dfnTrans` into working directory and run `DFNTrans`

**dump\_time** (*local\_jobname*, *section\_name*, *time*)

keeps log of cpu run time, current formulation is not robust

**get\_num\_frac** ()

Get the number of fractures from the `params.txt` file.

**class input\_helper** (*params*, *minFracSize*)

Functions to help parse the input file and check input parameters.

**Attributes:**

- `params` (list): list of parameters specified in the input file.
- `minFracSize` (float): the minimum fracture size.

**checkFamCount** ()

Makes sure at least one polygon family has been defined in `nFamRect` or `nFamEll` OR that there is a user input file for polygons.

**checkMean** (*minParam*, *maxParam*, *meanParam*, *warningFile*='')

Warns the user if the minimum value of a parameter is greater than the family's mean value, or if the maximum value of the parameter is less than the family's mean value.

**checkMinFracSize** (*valList*)

Corrects the minimum fracture size if necessary, by looking at the values in `valList`.

**checkMinMax** (*minParam*, *maxParam*, *shape*)

Checks that the minimum parameter for a family is not greater or equal to the maximum parameter.

**curlyToList** (*curlyList*)

'{1,2,3}' → [1,2,3]

**error** (*errString*)

print an error

**Args:** `errString` (str): a string describing the error

**extractParameters** (*line*, *inputIterator*)

Returns line without comments or white space.

**findKey** (*line*, *unfoundKeys*, *warningFile*)

Input: line containing a parameter (key) preceding a ":"

**Returns:**

- `key` – if it has not been defined yet and is valid
- `None` – if key does not exist
- `exits` – if the key has already been defined to prevent duplicate confusion

**findVal** (*line*, *key*, *inputIterator*, *unfoundKeys*, *warningFile*)

Extract the value for key from line.

**getGroups** (*line*, *valList*, *key*)

extract values between { and }

**hasCurlys** (*line*, *key*)

Checks to see that every { has a matching }.

**isNegative** (*num*)  
“returns True if num is negative, false otherwise

**listToCurly** (*strList*)  
[1,2,3] -> ‘{1,2,3}’ for writing output

**processLine** (*line, unfoundKeys, inputIterator, warningFile*)  
Find the key in a line, and the value for that key.

**scale** (*probList, warningFile*)  
scales list of probabilities (famProb) that doesn’t add up to 1 ie [.2, .2, .4] -> [0.25, 0.25, 0.5]

**valHelper** (*line, valList, key*)  
pulls values from curly brackets

**valueOf** (*key, writing=False*)  
Use to get key’s value in params. writing always false

**verifyFlag** (*value, key=’, inList=False*)  
Verify that value is either a 0 or a 1.

**verifyFloat** (*value, key=’, inList=False, noNeg=False*)  
Verify that value is a positive float.

**verifyInt** (*value, key=’, inList=False, noNeg=False*)  
Verify that value is a positive integer.

**verifyList** (*valList, key, verificationFn, desiredLength, noZeros=False, noNegs=False*)  
verifies input list that come in format {0, 1, 2, 3}

**Input:**

- valList - list of values (flags, floats, or ints) corresponding to a parameter
- key - the name of the parameter whose list is being verified
- verificationFn - (either verifyflag, verifyfloat or verifyint) checks each list element
- desiredLength - how many elements are supposed to be in the list
- noZeros - (optional) True for lists than cannot contain 0’s, false if 0’s are ok
- noNegs - (optional) True for lists than cannot contain negative numbers, false otherwise

**Output:**

- returns negative value of list length to indicate incorrect length and provide meaningful error message
- prints error and exits if a value of the wrong type is found in the list
- returns None if successful

**warning** (*warnString, warningFile=’*’*)*  
print a warning to a file (currently does not work)

**zeroInStdDevs** (*valList*)  
returns True is there is a zero in valList of standard deviations

**DFNWORKS.lagrit2pflotran** (*inp\_file=’, mesh\_type=’, hex2tet=False*)  
Takes output from LaGriT and processes it for use in PFLOTRAN.

**Kwargs:**

- inp\_file (str): name of the inp (AVS) file produced by LaGriT
- mesh\_type (str): the type of mesh
- hex2tet (boolean): True if hex mesh elements should be converted to tet elements, False otherwise.

**DFNWORKS.legal** ()  
Print the legal LANL statement for dfnWorks.

**DFNWORKS.make\_working\_directory()**

make working directories for fracture generation

**DFNWORKS.mesh\_network** (*production\_mode=True, refine\_factor=1, slope=2*)

Mesh Fracture Network using ncpus and lagrit meshing file is separate file: dfnGen\_meshing.py

**DFNWORKS.os** = <module 'os' from '/n/swdev/packages/Ubuntu-14.04-x86\_64/anaconda-python/2.4.1/lib/python2.7/os.py'

**DFNWORKS.output\_report** (*radiiFile='radii.dat', famFile='families.dat', trans-  
File='translations.dat', rejectFile='rejections.dat', output\_name=''*)

Create PDF report of generator

#### Notes:

- Set the number of histogram buckets (bins) by changing numBuckets variable in his graphing functions
- Also change number of x-values used to plot lines by changing numXpoints variable in appropriate funcs
- Set show = True to show plots immediately and still make pdf
- NOTE future developers of this code should add functionality for radiiList of size 0.

**DFNWORKS.parse\_pflotran\_vtk** (*grid\_vtk\_file=''*)

Using C++ VTK library, convert inp file to VTK file, then change name of CELL\_DATA to POINT\_DATA.

**DFNWORKS.parse\_pflotran\_vtk\_python** (*grid\_vtk\_file=''*)

Replace CELL\_DATA with POINT\_DATA in the VTK output.

**DFNWORKS.pflotran** ()

Run pflotran Copy PFLOTTRAN run file into working directory and run with ncpus

**DFNWORKS.pflotran\_cleanup** ()

Concatenate PFLOTTRAN output files and then delete them

**DFNWORKS.print\_run\_time** (*local\_jobname*)

Read in run times from file and and print to screen with percentages

**DFNWORKS.re** = <module 're' from '/n/swdev/packages/Ubuntu-14.04-x86\_64/anaconda-python/2.4.1/lib/python2.7/re.py'

**DFNWORKS.run\_dfnTrans** ()

run dfnTrans simulation

**DFNWORKS.subprocess** = <module 'subprocess' from '/n/swdev/packages/Ubuntu-14.04-x86\_64/anaconda-python/2.4.1/lib/python2.7/subprocess.py'

**DFNWORKS.sys** = <module 'sys' (built-in)>

**DFNWORKS.write\_perms\_and\_correct\_volumes\_areas** (*inp\_file='', uge\_file='',  
perm\_file='', aper\_file=''*)

Write permeability values to perm\_file, write aperture values to aper\_file, and correct volume areas in uge\_file

**DFNWORKS.zone2ex** (*uge\_file='', zone\_file='', face=''*)

Convert zone files from LaGriT into ex format for LaGriT inputs: uge\_file: name of uge file zone\_file: name of zone file face: face of the plane corresponding to the zone file

zone\_file='all' processes all directions, top, bottom, left, right, front, back

## 5.2 dfnGen

### 5.2.1 Processing generator input

```
pydfnworks.gen_input.check_input(self, input_file='', output_file='')
```

#### Input Format Requirements:

- Each parameter must be defined on its own line (separate by newline)
- A parameter (key) MUST be separated from its value by a colon ':' (ie. -> key: value)
- Values may also be placed on lines after the 'key'
- Comment Format: On a line containing // or / \*, nothing after \* / or // will be processed but text before a comment will be processed

#### Kwargs:

- input\_file (name): name of dfnGen input file
- output\_file (name): stripped down input file for DFNGen

```
class pydfnworks.distributions.distr(params, numEdistrib, numRdistrib, minFracSize)
```

Verifies the fracture distribution input parameters for dfnGen.

#### Attributes:

- params (list): parameters for dfnGen
- numEdistrib (int): number of ellipse family distributions
- numRdistrib (int): number of rectangle family distributions
- minFracSize (double): minimum fracture size

#### betaDistribution (prefix)

Verifies both the "ebetaDistribution" and "rBetaDistribution". If either contain any flags indicating constant angle (1) then the corresponding "ebeta" and/or "rbeta" parameters are also verified.

**Args:** prefix (str): Indicates shapes that the beta distribution describes. 'e' if they are ellipses, 'r' if they are rectangles.

#### constantDist (prefix)

Verifies parameters for constant distribution of fractures

#### distr (prefix)

Verifies "edistr" and "rdistr" making sure one distribution is defined per family and each distribution is either 1 (log-normal), 2 (Truncated Power Law), 3 (Exponential), or 4 (constant). Stores how many of each distrib are in use in numEdistrib or numRdistrib lists.

#### exponentialDist (prefix)

Verifies parameters for exponential distribution of fractures.

#### lognormalDist (prefix)

Verifies all logNormal Parameters for ellipses and Rectangles.

#### tplDist (prefix)

Verifies parameters for truncated power law distribution of fractures.

## 5.2.2 Running the generator

`pydfnworks.generator.create_network(self)`

Execute dfnGen and print whether the generation of the fracture network failed or succeeded. The params.txt file must be there for success.

`pydfnworks.generator.dfnGen(self)`

**Run the dfnGen workflow:**

- 1. `make_working_directory`: Create a directory with name of job
- 2. `check_input`: Check input parameters and create a clean version of the input file
- 3. `create_network`: Create network. DFNGEN v2.0 is called and creates the network
- 4. `output_report`: Generate a PDF summary of the DFN generation
- 5. `mesh_network`: calls module dfnGen\_meshing and runs LaGriT to mesh the DFN

`pydfnworks.generator.make_working_directory(self)`

make working directories for fracture generation

## 5.2.3 Graphing generator output

`pydfnworks.gen_output.output_report(self, radiiFile='radii.dat', famFile='families.dat', transFile='translations.dat', rejectFile='rejections.dat', output_name='')`

Create PDF report of generator

**Notes:**

- Set the number of histogram buckets (bins) by changing numBuckets variable in his graphing functions
- Also change number of x-values used to plot lines by changing numXpoints variable in appropriate funcs
- Set show = True to show plots immediately and still make pdf
- NOTE future developers of this code should add functionality for radiiList of size 0.

## 5.3 dfnFlow

`pydfnworks.flow.dfnFlow(self)`

Run the dfnFlow portion of the workflow.

`pydfnworks.flow.inp2gmw(self, inp_file='')`

Convert inp file to gmw file, for general mesh viewer .

**Kwargs:** `inp_file` (str): name of inp file

`pydfnworks.flow.lagrit2pflotran(self, inp_file='', mesh_type='', hex2tet=False)`

Takes output from LaGriT and processes it for use in PFLOTRAN.

**Kwargs:**

- `inp_file` (str): name of the inp (AVS) file produced by LaGriT
- `mesh_type` (str): the type of mesh
- `hex2tet` (boolean): True if hex mesh elements should be converted to tet elements, False otherwise.

```
pydfnworks.flow.parse_pflotran_vtk(self, grid_vtk_file='')
```

Using C++ VTK library, convert inp file to VTK file, then change name of CELL\_DATA to POINT\_DATA.

```
pydfnworks.flow.parse_pflotran_vtk_python(self, grid_vtk_file='')
```

Replace CELL\_DATA with POINT\_DATA in the VTK output.

```
pydfnworks.flow.pflotran(self)
```

Run pflotran Copy PFLOTTRAN run file into working directory and run with ncpus

```
pydfnworks.flow.pflotran_cleanup(self)
```

Concatenate PFLOTTRAN output files and then delete them

```
pydfnworks.flow.write_perms_and_correct_volumes_areas(self, inp_file='', uge_file='',  
                                                         perm_file='', aper_file='')
```

Write permeability values to perm\_file, write aperture values to aper\_file, and correct volume areas in uge\_file

```
pydfnworks.flow.zone2ex(self, uge_file='', zone_file='', face='')
```

Convert zone files from LaGriT into ex format for LaGriT inputs: uge\_file: name of uge file zone\_file: name of zone file face: face of the plane corresponding to the zone file

zone\_file='all' processes all directions, top, bottom, left, right, front, back

## 5.4 dfnTrans

```
pydfnworks.transport.copy_dfnTrans_files(self)
```

create link to DFNTRANS and copy input file into local directory

```
pydfnworks.transport.dfnTrans(self)
```

Copy input files for dfnTrans into working directory and run DFNTrans

```
pydfnworks.transport.run_dfnTrans(self)
```

run dfnTrans simulation

## 5.5 LaGriT (meshing)

### 5.5.1 Mesh DFN

```
pydfnworks.meshdfn.mesh_network(self, production_mode=True, refine_factor=1, slope=2)
```

Mesh Fracture Network using ncpus and lagrit meshing file is separate file: dfnGen\_meshing.py

### 5.5.2 LaGrit scripts

```
pydfnworks.lagrit_scripts.create_lagrit_scripts(visual_mode, ncpu, refine_factor=1,  
                                                  production_mode=True)
```

Creates LaGriT script to be run for each polygon

#### Inputs:

- visual\_mode (True / False)
- ncpu: number of cpus
- refine\_fractor: used rectangles
- production\_mode (True/False)

```
pydfnworks.lagrit_scripts.create_merge_poly_files(ncpu, num_poly, visual_mode)
```



**Section 4** [Create merge\_poly file] Creates a lagrit script that reads in each mesh, appends it to the main mesh, and then deletes that mesh object Then duplicate points are removed from the main mesh using EPS\_FILTER The points are compressed, and then written in the files full\_mesh.gmv, full\_mesh.inp, and an FEHM dump is preformed.

```
pydfnworks.lagrit_scripts.create_parameter_mgli_file(num_poly, h, slope=2, re-
                                                    fine_dist=0.5)
```

create parameter\_mgli\_files Outputs parameteri.mgli files used in running LaGriT Scripts

**Inputs:**

- num\_poly: Number of polygons
- h: meshing length scale
- slope: Slope of coarsening function, default = 2, set to 0 for uniform mesh resolution
- refine\_dist: distance used in coarsing function, default = 0.5,

```
pydfnworks.lagrit_scripts.create_user_functions()
```

Create user\_function.lgi files for meshing

```
pydfnworks.lagrit_scripts.define_zones(h, domain)
```

Creates and runs LaGriT script to define domain size

### 5.5.3 Run meshing in parallel

```
pydfnworks.run_meshing.merge_the_meshes(num_poly, ncpu, n_jobs, visual_mode)
```

Merges all the meshes together, deletes duplicate points, dumps the .gmV and fehM files

```
pydfnworks.run_meshing.mesh_fracture(fracture_id, visual_mode, num_poly)
```

Child function for parallelized meshing of fractures

```
pydfnworks.run_meshing.mesh_fractures_header(num_poly, ncpu, visual_mode)
```

Header function for Parallel meshing of fractures

Creates a queue of fracture numbers ranging form 1, num\_poly

Each fractures is meshed using mesh\_fracture called within the worker function.

If any fracture fails to mesh properly, then a folder is created with that fracture information and the fracture number is written into failure.txt.

**Returns:**

- True: If failure.txt is empty meaning all fractures meshed correctly
- False: If failure.txt is not empty, then at least one fracture failed.

```
pydfnworks.run_meshing.worker(work_queue, done_queue, visual_mode, num_poly)
```

Worker function for parallelized meshing

## 5.6 Helper methods

### 5.6.1 Mesh helper methods

```
pydfnworks.mesh_dfn_helper.check_dudded_points(dudded)
```

Parses Lagrit log\_merge\_all.txt and checks if number of dudded points is the expected number

**Returns:**

- True if the number of dudded points is correct
- False if the number of dudded points is incorrect

`pydfnworks.mesh_dfn_helper.cleanup_dir()`

Removes files from meshing

`pydfnworks.mesh_dfn_helper.output_meshing_report (visual_mode)`

Prints information about the final mesh to file

`pydfnworks.mesh_dfn_helper.parse_params_file()`

Reads params.txt file and parse information

**Returns:**

- `num_poly`: (int) Number of Polygons
- `h`: (float) meshing length scale `h`
- `dudded_points`: (int) Expected number of dudded points in Filter (LaGriT)
- `visual_mode`: True/False
- `domain`: dict: x,y,z domain sizes

## 5.6.2 Print legal statement

`pydfnworks.legal.legal (self)`

Print the legal LANL statement for dfnWorks.

## 5.6.3 Other helper methods

`pydfnworks.helper.commandline_options()`

Read command lines for use in dfnWorks.

**Options:**

- `-name` : Jobname (Mandatory)
- `-ncpu` : Number of CPUS (Optional, default=4)
- `-input` : input file with paths to run files (Mandatory if the next three options are not specified)
- `-gen` : Generator Input File (Mandatory, can be included within this file)
- `-flow` : PFLOLAN Input File (Mandatory, can be included within this file)
- `-trans`: Transport Input File (Mandatory, can be included within this file)
- `-cell`: True/False Set True for use with cell based aperture and permeability (Optional, default=False)
- `-large_network`: True/False Set True to use CPP for file processing bottleneck (Optional, default=False)

`pydfnworks.helper.dump_time (local_jobname, section_name, time)`

keeps log of cpu run time, current formulation is not robust

`pydfnworks.helper.get_num_frac()`

Get the number of fractures from the params.txt file.

**class** `pydfnworks.helper.input_helper (params, minFracSize)`

Functions to help parse the input file and check input parameters.

**Attributes:**

- `params` (list): list of parameters specified in the input file.
- `minFracSize` (float): the minimum fracture size.

**checkFamCount** ( )

Makes sure at least one polygon family has been defined in `nFamRect` or `nFamEll` OR that there is a user input file for polygons.

**checkMean** (*minParam*, *maxParam*, *meanParam*, *warningFile*='')

Warns the user if the minimum value of a parameter is greater than the family's mean value, or if the maximum value of the parameter is less than the family's mean value.

**checkMinFracSize** (*valList*)

Corrects the minimum fracture size if necessary, by looking at the values in *valList*.

**checkMinMax** (*minParam*, *maxParam*, *shape*)

Checks that the minimum parameter for a family is not greater or equal to the maximum parameter.

**curlyToList** (*curlyList*)

'{1,2,3}' → [1,2,3]

**error** (*errString*)

print an error

**Args:** *errString* (str): a string describing the error

**extractParameters** (*line*, *inputIterator*)

Returns line without comments or white space.

**findKey** (*line*, *unfoundKeys*, *warningFile*)

Input: line containing a parameter (key) preceding a ":"

**Returns:**

- `key` – if it has not been defined yet and is valid
- `None` – if key does not exist
- `exits` – if the key has already been defined to prevent duplicate confusion

**findVal** (*line*, *key*, *inputIterator*, *unfoundKeys*, *warningFile*)

Extract the value for *key* from *line*.

**getGroups** (*line*, *valList*, *key*)

extract values between { and }

**hasCurlys** (*line*, *key*)

Checks to see that every { has a matching }.

**isNegative** (*num*)

"returns True if *num* is negative, false otherwise

**listToCurly** (*strList*)

[1,2,3] → '{1,2,3}' for writing output

**processLine** (*line*, *unfoundKeys*, *inputIterator*, *warningFile*)

Find the key in a line, and the value for that key.

**scale** (*probList*, *warningFile*)

scales list of probabilities (*famProb*) that doesn't add up to 1 ie [.2, .2, .4] → [0.25, 0.25, 0.5]

**valHelper** (*line*, *valList*, *key*)

pulls values from curly brackets

**valueOf** (*key*, *writing=False*)

Use to get key's value in params. writing always false

**verifyFlag** (*value*, *key=''*, *inList=False*)

Verify that value is either a 0 or a 1.

**verifyFloat** (*value*, *key=''*, *inList=False*, *noNeg=False*)

Verify that value is a positive float.

**verifyInt** (*value*, *key=''*, *inList=False*, *noNeg=False*)

Verify that value is a positive integer.

**verifyList** (*valList*, *key*, *verificationFn*, *desiredLength*, *noZeros=False*, *noNegs=False*)

verifies input list that come in format {0, 1, 2, 3}

**Input:**

- *valList* - list of values (flags, floats, or ints) corresponding to a parameter
- *key* - the name of the parameter whose list is being verified
- *verificationFn* - (either *verifyflag*, *verifyfloat* or *verifyint*) checks each list element
- *desiredLength* - how many elements are supposed to be in the list
- *noZeros* - (optional) True for lists than cannot contain 0's, false if 0's are ok
- *noNegs* - (optional) True for lists than cannot contain negative numbers, false otherwise

**Output:**

- returns negative value of list length to indicate incorrect length and provide meaningful error message
- prints error and exits if a value of the wrong type is found in the list
- returns None if successful

**warning** (*warnString*, *warningFile=''*)

print a warning to a file (currently does not work)

**zeroInStdDevs** (*valList*)

returns True is there is a zero in *valList* of standard deviations

`pydfnworks.helper.print_run_time` (*local\_jobname*)

Read in run times from file and and print to screen with percentages

## **SCRIPTS**

The pydfnworks package has four Python scripts that assist with compiling, testing, and running the software. These scripts are all in the folder dfnWorks-Version2.0/pydfnworks/bin/ .

### **6.1 fix\_paths.py: fix the paths in the tests folder**

Syntax: `python fix_paths.py`

### **6.2 compile.py: compile dfnWorks components**

The compile.py script is called by run.py, but can also be called on its own. This script compiles the C and C++ components of dfnWorks. Without arguments, the script performs the compiling. With the argument 'clean,' the script cleans up C and C++ object files, before compiling.

### **6.3 test.py: test dfnWorks**

The test.py script can be used to easily test functionality of dfnWorks. It runs a test suite (input files contained in the tests folder) that test the exponential, power law, and lognormal distributions as well as user-defined ellipses and user-defined rectangle fracture inputs. Syntax: `python test.py [TEST_NAME_PART]`, where [TEST\_NAME\_PART] is all or part of the name of the tests the user wants to run.

### **6.4 run.py: run dfnWorks**

The run.py script allows the user to execute a single run of dfnWorks. The syntax for this script is:

```
python run.py -name [JOBNAME] -input [INPUT_FILE_NAME] -ncpu [NUMBER OF CPUS TO USE] -input [INPUT_FILE_NAME] -large_network
```

Please refer to the tutorial section (the next section for a detailed description of how to run dfnWorks.



## TUTORIAL

This document contains a short, five example, tutorial for dfnWorks. The five test cases provided are for:

- `4_user_defined_rectangles`: The user defines the precise location and size of four rectangular fractures.
- `4_user_defined_ellipses`: The user defines the precise location and size of four elliptical fractures, where the ellipses are approximated as polygons.
- `exponential_dist`: The user specifies the parameters for two families of fractures with an exponential distribution of fracture size.
- `lognormal_dist`: The user specifies the parameters for two families of fractures with a lognormal distribution of fracture size.
- `truncated_power_law_dist`: The user specifies the parameters for two families of fractures with a truncated power-law distribution of fracture size.

All required input files for these examples are contained in the folder `dfnWorks-Version2.0/tests/`. The focus of this document is to provide visual confirmation that new users of dfnWorks have the code set up correctly, can carry out the following runs and reproduce the following images. All images are rendered using Paraview, which can be obtained for free at [http : //www.paraview.org/](http://www.paraview.org/). The first two examples are simpler than the last three so it is recommended that the user proceed in the order presented here.

### 7.1 Turn on X forwarding if on server

Ensure that X forwarding is turned on if you are running dfnWorks from an ssh connection. This requires that the ssh login have the `-X` option:

```
$ ssh -X SERVER_NAME
```

### 7.2 Go to the dfnWorks-Version2.0 repository

```
$ cd ~/dfnWorks-Version2.0/
```

### 7.3 Fix paths in test directory

Fix the pathnames for all files in the folder `/tests/`. This can be done automatically by running the script `fix_paths.py`:

```
$ cd /pydfnworks/bin/  
$ python fix_paths.py
```

## 7.4 Set the PETSC, PFLOTRAN, Python, and LaGriT paths correctly

Before executing dfnWorks, the following paths must be set:

- DFNWORKS\_PATH: the dfnWorks-Version2.0 repository folder
- PETSC\_DIR and PETSC\_ARCH: PETSC environmental variables
- PFLOTRAN\_DIR: The PFLOTRAN directory
- python\_dfn: The location of the Python distribution to use
- lagrit\_dfn: The location of the LaGriT executable

```
$ vi /pydfnworks/pydfnworks/paths.py
```

For example:

```
os.environ['DFNWORKS_PATH'] = '/home/nknapp/dfnWorks-Version2.0/'
```

## 7.5 Setup the Python package pydfnworks

Go up a directory:

```
$ cd ..
```

If the user has admin privileges:

```
$ python setup.py install
```

If the user DOES NOT have admin privileges:

```
$ python setup.py install --user
```

## 7.6 Executing dfnWorks

To run one of the test cases enter the following command:

- `python run.py -name [JOBNAME] -input [INPUT_FILE] -ncpu [NUMBER_OF_CPUS] -large_network`

The arguments are:

- [JOBNAME]: The name of the run, which is also the absolute path name of the folder which will contain the run's output.
- [INPUT\_FILE]: An input file with three lines that have input files for dfnGen, dfnFlow, and dfnTrans, respectively. Any of the files with ending .txt in the directory tests can be used as examples of input files.
- [NUMBER\_OF\_CPUS]: The number of CPUs that the user would like to use for the parallel computation of the meshing and flow solutions.
- large\_network (optional): Only use this flag if the user should use CPP for file processing.

For example, to run the 4\_user\_defined rectangles test below, on 4 CPUs, the command line input would be:

```
$ cd bin
$ python run.py -name ~/4_user_defined_rectangles
-input ../tests/4_user_defined_rectangles.txt
-ncpu 4
```



This command will run the `4_user_defined_rectangles` test and create a new folder `4_user_defined_rectangles` where all output files will be located. Descriptions of each output file are in the documentation. In the following sections, we provide descriptions of the output you should expect for each of the five examples.

## 7.7 Summary

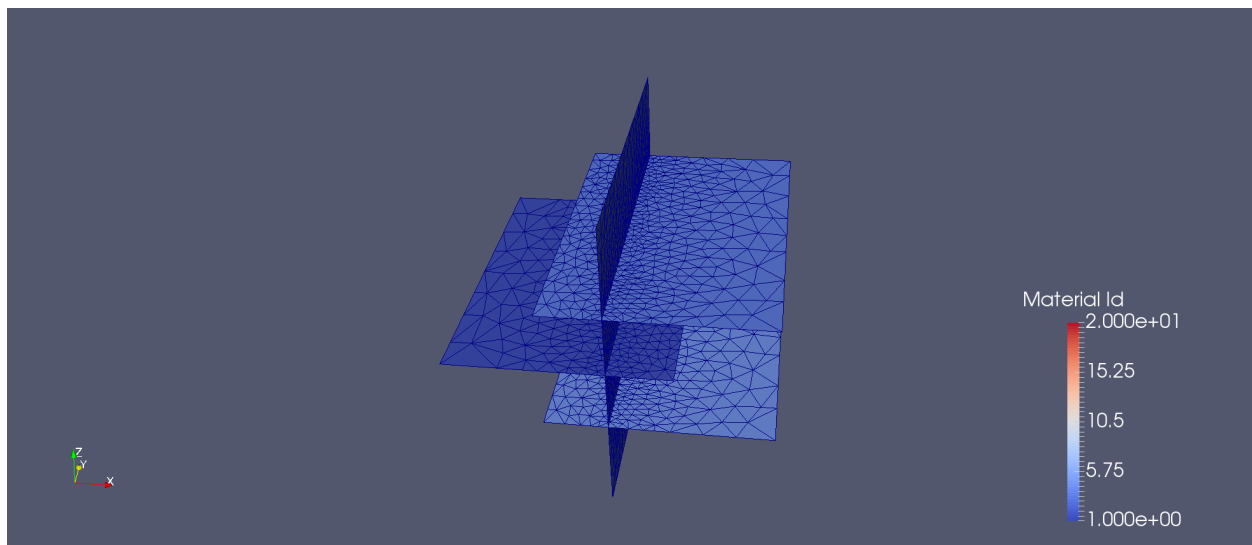
Here is a quick summary of the steps described above, with commands for the `4_user_defined_rectangles` test case. **You MUST edit the paths in `paths.py`, and change `~` to the location of the repository, for these commands to work!**

```
$ ssh -X SERVER_NAME
$ cd ~/dfnWorks-Version2.0/pydfnworks/bin/
$ python fix_paths.py
$ vi /pydfnworks/pydfnworks/paths.py
$ cd ..
$ python setup.py install --user
$ cd bin
$ python run.py -name ~/4_user_defined_rectangles
-input ../tests/4_user_defined_rectangles.txt
-ncpu 4
```

## 7.8 4\_user\_defined\_rectangles

This test case consists of four user defined rectangular fractures within a cubic domain with sides of length one meter. The input file specifying the ellipses is in `dfnWorks-Version2.0/tests`, and is named `define_4_user_rectangles.dat`. To run the test on 4 cpus, enter the following command line input:

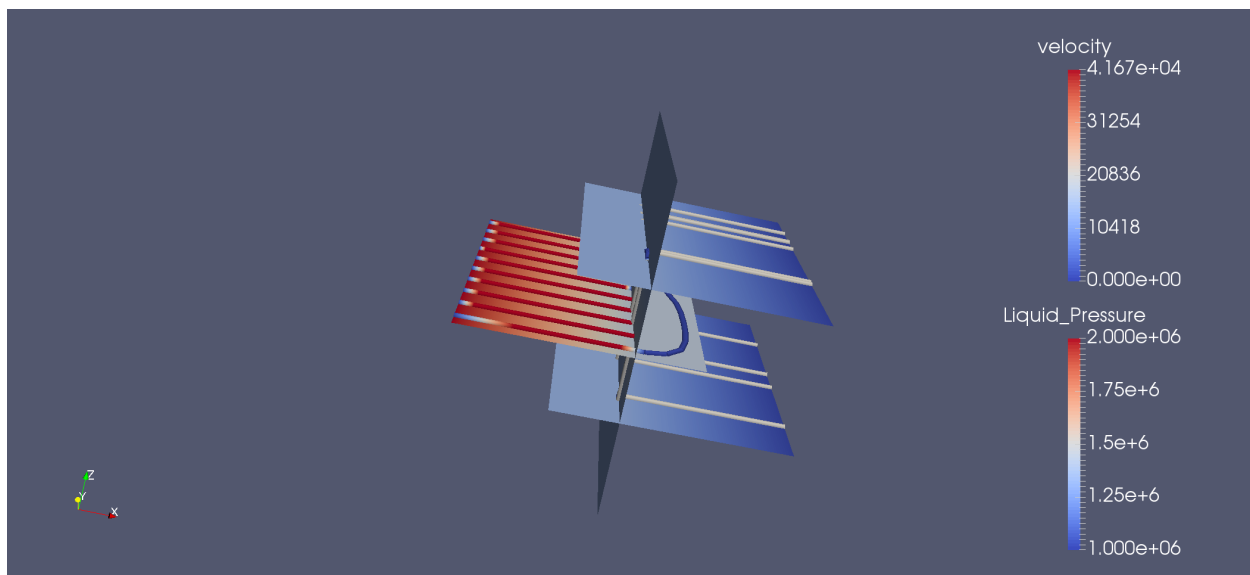
The network of four fractures, each colored by material ID. The computational mesh is overlaid on the fractures. This image is created by loading the file `full_mesh.inp`, located in the folder `4_user_defined_rectangles/LaGriT/`, into Paraview.



The network of four fractures, colored by pressure solution. High pressure (red) Dirichlet boundary conditions are applied on the edge of the single fracture along the boundary  $x = -0.5$ , and low pressure (blue) boundary conditions are applied on the edges of the two fractures at the boundary  $x = 0.5$ . This image is created by loading the file `4_user_defined_rectangles/PFLOTRAN/parsed_vtk/dfn_explicit-001.vtk` into Paraview.



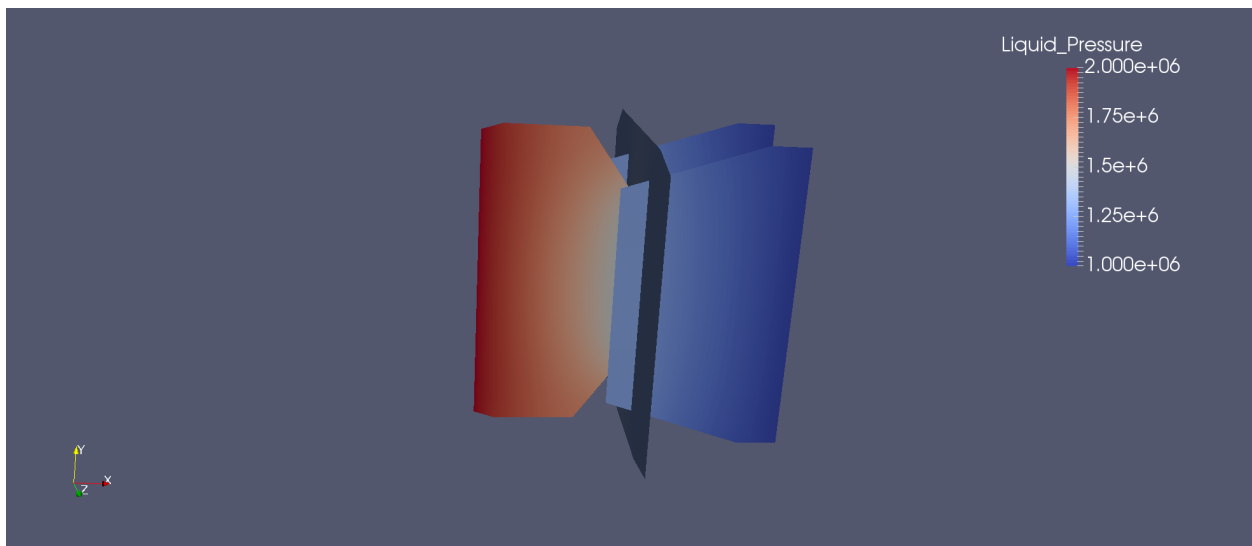
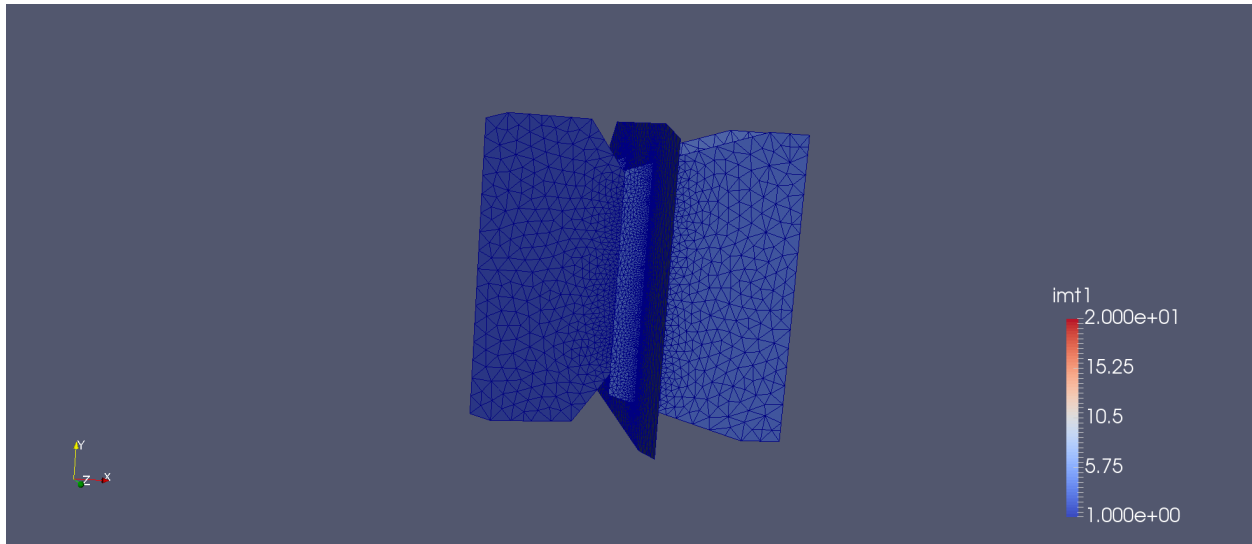
Particle trajectories on the network of four fractures. Particles are inserted uniformly along the inlet fracture on the left side of the image. Particles exit the domain through the two horizontal fractures on the right side of the image. Due to the stochastic nature of the particle tracking algorithm, your pathlines might not be exactly the same as in this image. Trajectories are colored by the current velocity magnitude of the particle's velocity. Trajectories can be visualized by loading the files `part_*.inp`, in the folder `4_user_rectangles/dfnTrans/trajectories/`. We have used the extract surface and tube filters in paraview for visual clarity.

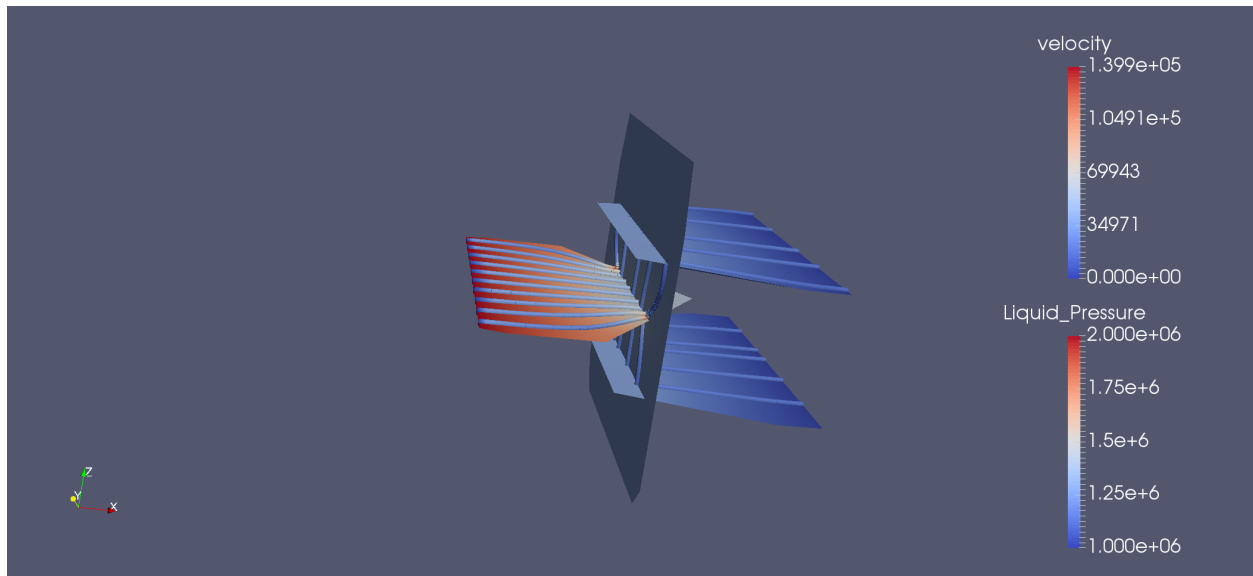


In the other tests, only a brief description and pictures are provided.

## 7.9 4\_user\_defined\_ellipses

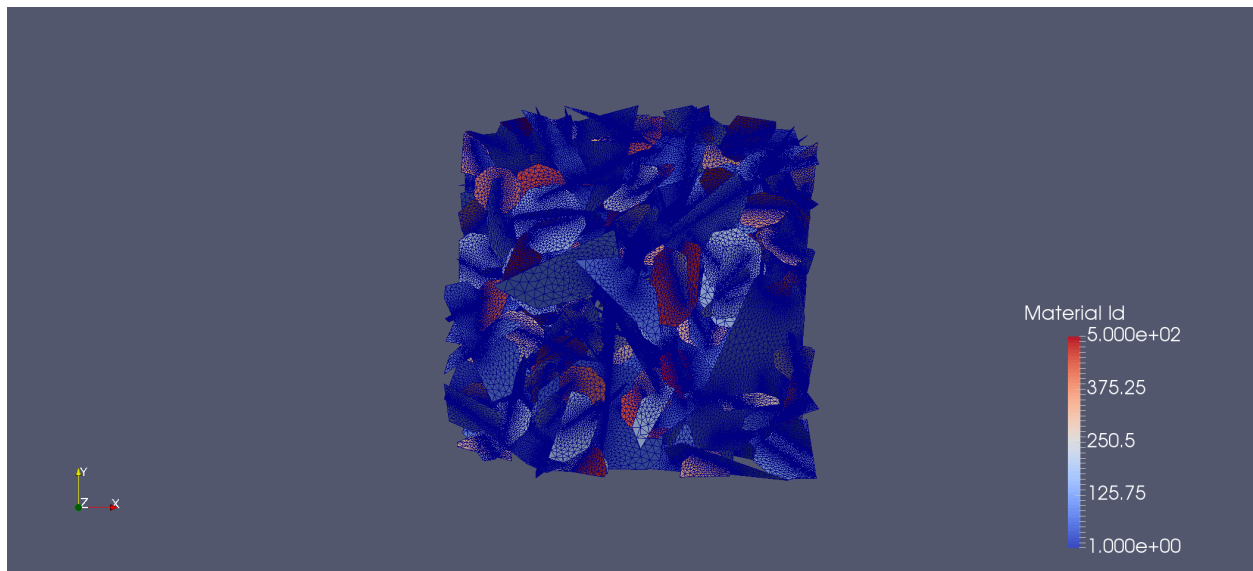
This test case consists of four user defined elliptical fractures within a cubic domain with sides of length one meter. In this case the ellipses are approximated using 5 vertices. The input file specifying the ellipses is in `dfnWorks-Version2.0/tests`, and is named `define_4_user_ellipses.dat`.





## 7.10 truncated\_power\_law\_dist

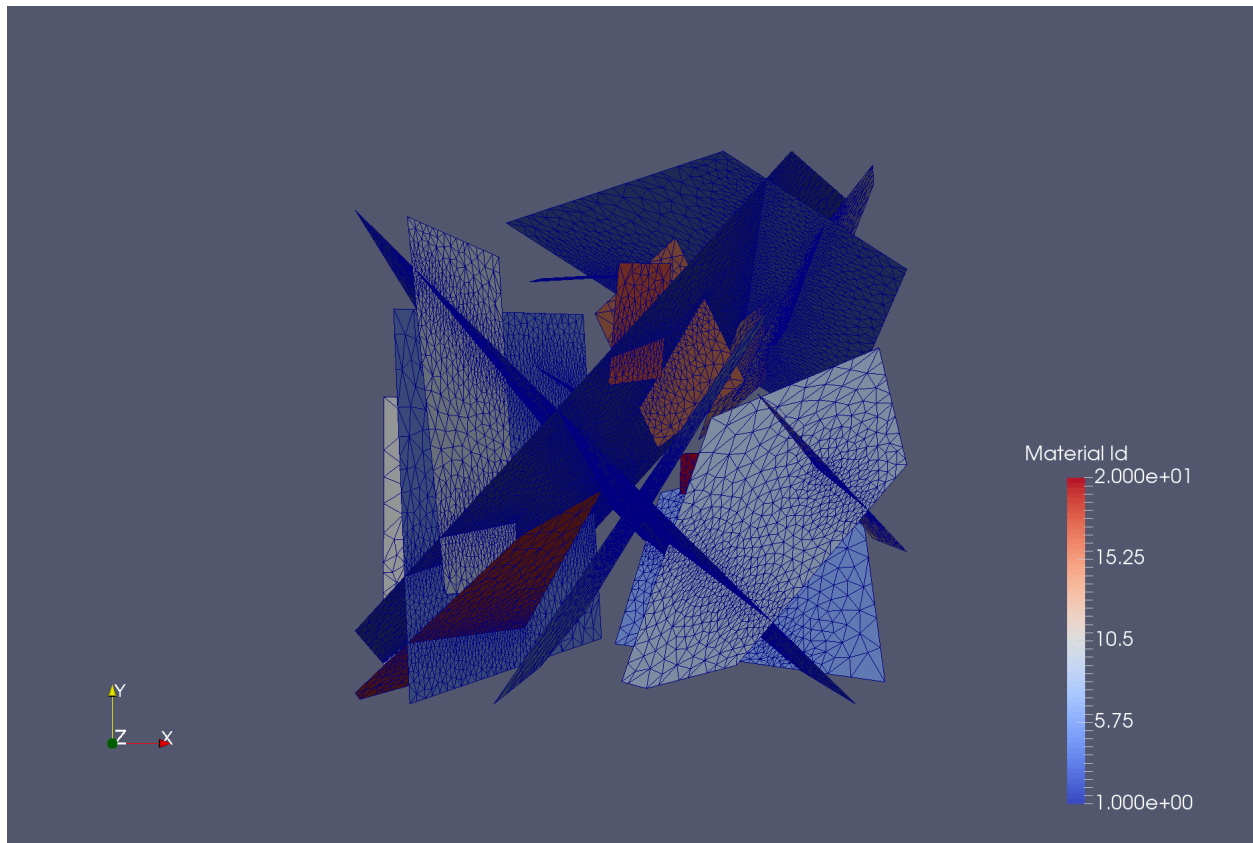
This test case consists of two families whose sizes have a truncated power law distribution with a minimum size of 0.5m and a maximum size of 50m. The domain size is cubic with an edge length of 4m. The other input parameters can be found in tests/gen\_truncated\_power\_law\_dist.dat.

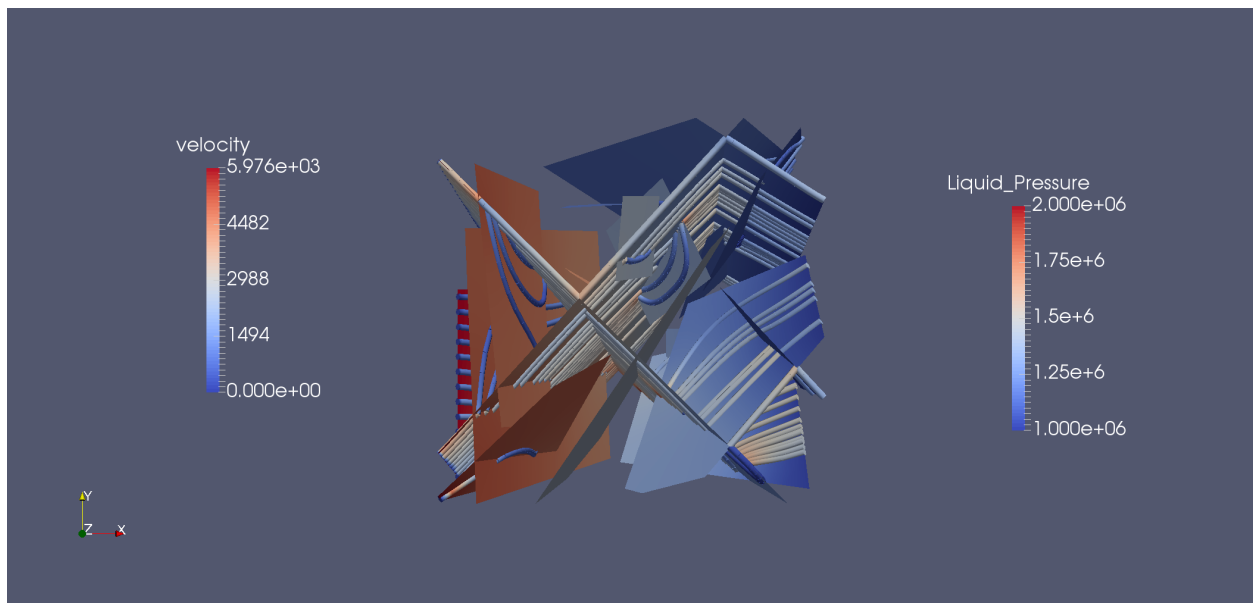
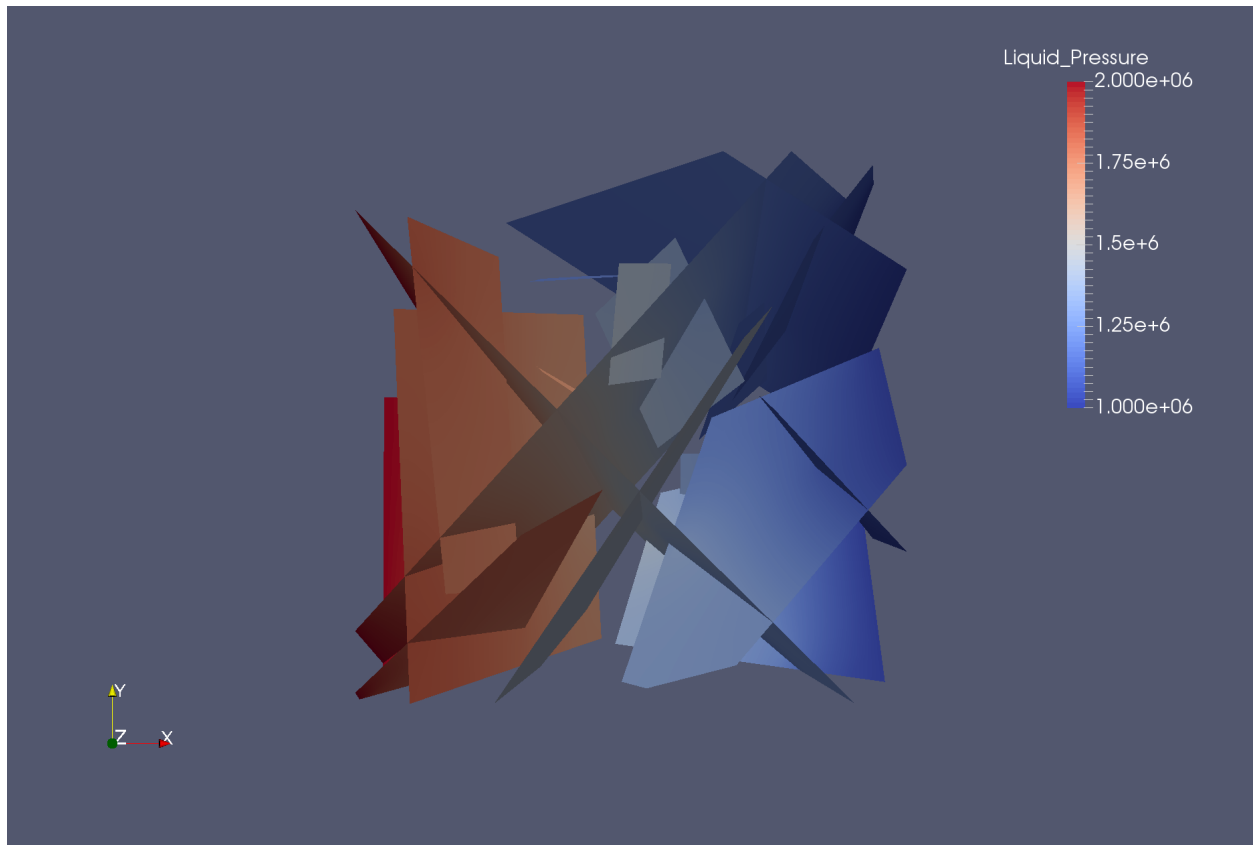




## 7.11 exponential\_dist

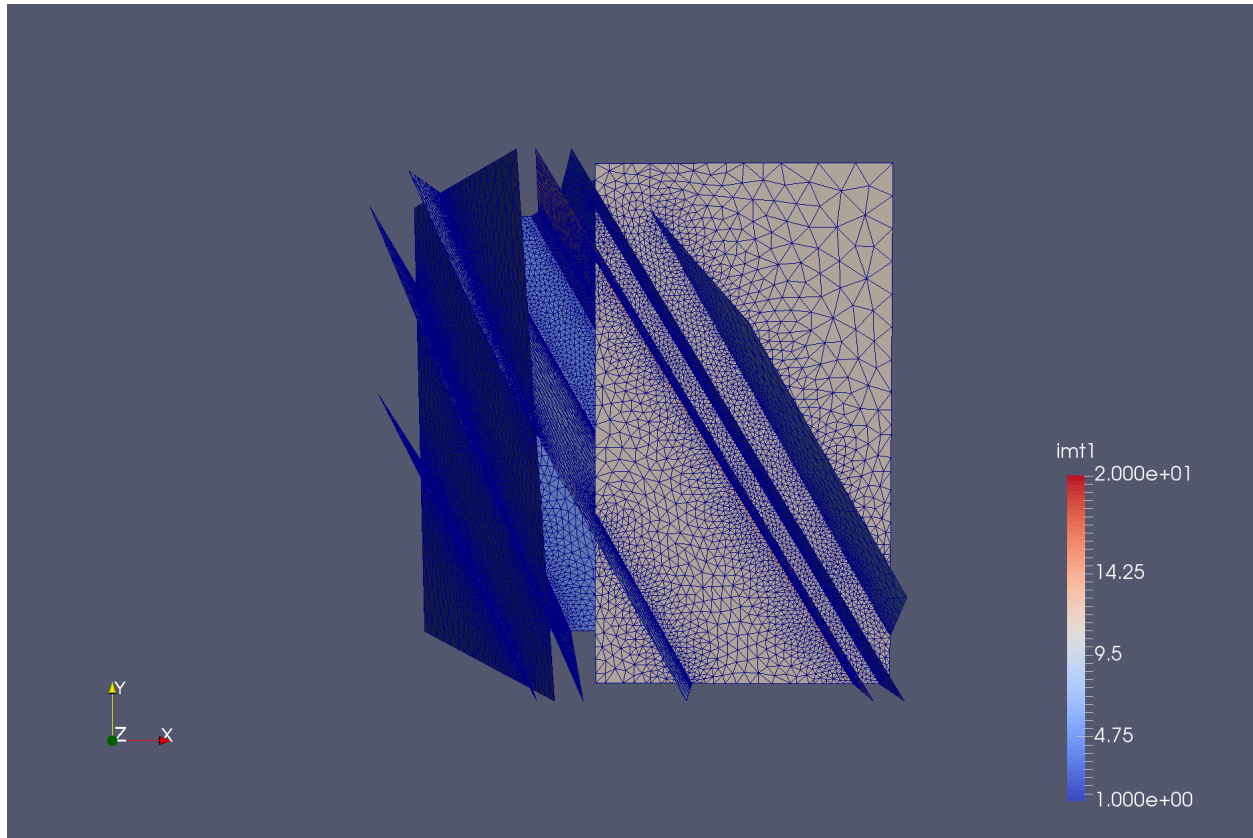
This test case consists of a family of fractures whose size is exponentially distributed with a minimum size of 1m and a maximum size of 50m. The domain is cubic with an edge length of 10m. All input parameters for the generator can be found in tests/gen\_exponential\_dist.dat.



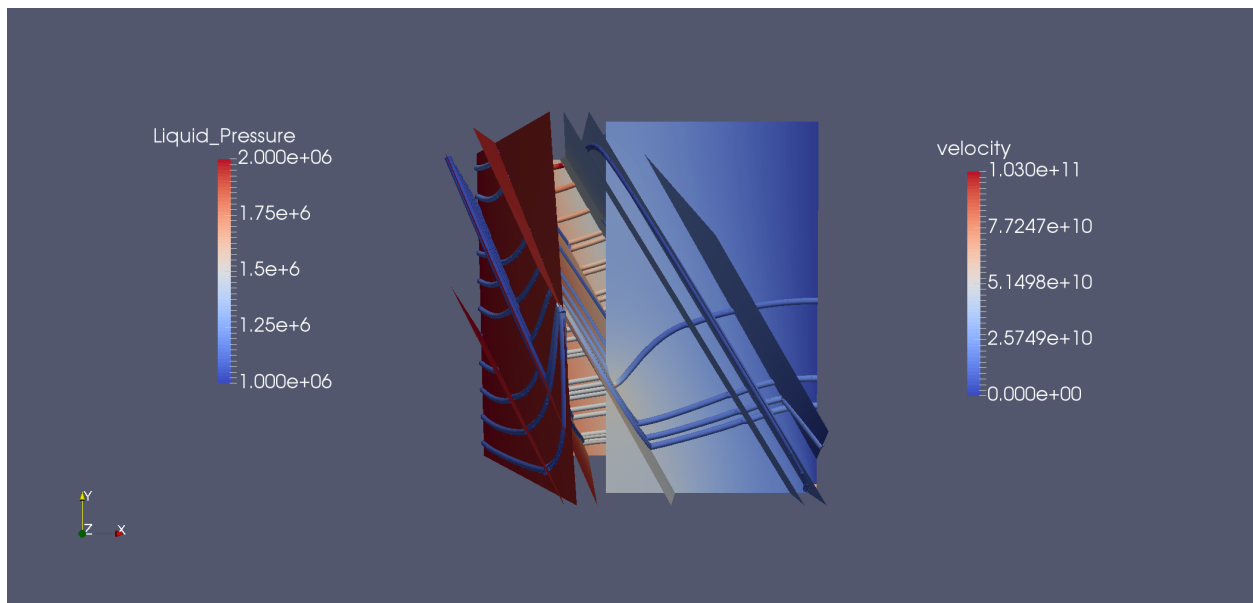
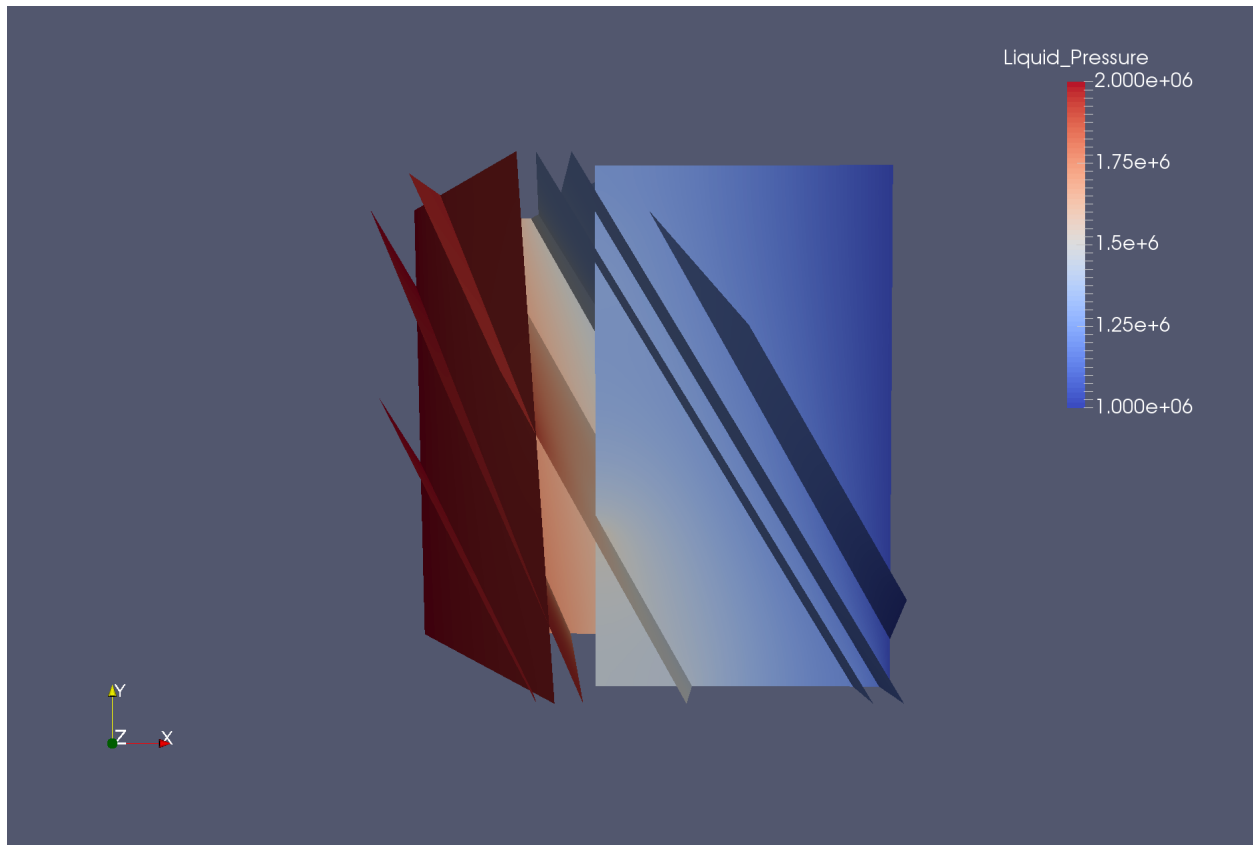


## 7.12 lognormal\_dist

This test case consists of two fracture families whose sizes have a lognormal distribution with a minimum size of 0.5m and a maximum size of 50m. The domain size is cubic with an edge length of 10m. All input parameters for the generator can be found in tests/gen\_lognormal\_dist.dat.









## **EXAMPLE APPLICATIONS**

### **8.1 Carbon dioxide sequestration**

DFNWORKS provides the framework necessary to perform multiphase simulations (such as flow and reactive transport) at the reservoir scale. A particular application, highlighted here, is sequestering CO<sub>2</sub> from anthropogenic sources and disposing it in geological formations such as deep saline aquifers and abandoned oil fields. Geological CO<sub>2</sub> sequestration is one of the principal methods under consideration to reduce carbon footprint in the atmosphere due to fossil fuels (Bachu, 2002; Pacala and Socolow, 2004). For safe and sustainable long-term storage of CO<sub>2</sub> and to prevent leaks through existing faults and fractured rock (along with the ones created during the injection process), understanding the complex physical and chemical interactions between CO<sub>2</sub>, water (or brine) and fractured rock, is vital. DFNWORKS capability to study multiphase flow in a DFN can be used to study potential CO<sub>2</sub> migration through cap-rock, a potential risk associated with proposed subsurface storage of CO<sub>2</sub> in saline aquifers or depleted reservoirs. Moreover, using the reactive transport capabilities of PFLOTTRAN coupled with cell-based transmissivity of the DFN allows one to study dynamically changing permeability fields with mineral precipitation and dissolution due to CO<sub>2</sub>–water interaction with rock.

### **8.2 Shale energy extraction**

Hydraulic fracturing (fracking) has provided access to hydrocarbon trapped in low-permeability media, such as tight shales. The process involves injecting water at high pressures to reactivate existing fractures and also create new fractures to increase permeability of the shale allowing hydrocarbons to be extracted. However, the fundamental physics of why fracking works and its long term ramifications are not well understood. Karra et al. (2015) used DFNWORKS to generate a typical production site and simulate production. Using this physics based model, they found good agreement with production field data and determined what physical mechanisms control the decline in the production curve.

### **8.3 Nuclear waste repository**

The Swedish Nuclear Fuel and Waste Management Company (SKB) has undertaken a detailed investigation of the fractured granite at the Forsmark, Sweden site as a potential host formation for a subsurface repository for spent nuclear fuel (SKB, 2011; Hartley and Joyce, 2013). The Forsmark area is about 120 km north of Stockholm in northern Uppland, and the repository is proposed to be constructed in crystalline bedrock at a depth of approximately 500 m. Based on the SKB site investigation, a statistical fracture model with multiple fracture sets was developed; detailed parameters of the Forsmark site model are in SKB (2011). We adopt a subset of the model that consist of three sets of background (non-deterministic) circular fractures whose orientations follow a Fisher distribution, fracture radii are sampled from a truncated power-law distribution, the transmissivity of the fractures is estimated using a power-law model based on the fracture radius, and the fracture aperture is related to the fracture size using the cubic law (Adler et al., 2012). Under such a formulation, the fracture apertures are uniform on each fracture, but vary among fractures.

The network is generated in a cubic domain with sides of length one-kilometer. Dirichlet boundary conditions are imposed on the top (1 MPa) and bottom (2 MPa) of the domain to create a pressure gradient aligned with the vertical axis, and noflow boundary conditions are enforced along lateral boundaries.

Sources:

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- Hartley, L., Joyce, S., 2013. Approaches and algorithms for groundwater flow modeling in support of site investigations and safety assessment of the Forsmark site, Sweden. *J. Hydrol.* 500, 200–216.
- Karra, S., Makedonska, N., Viswanathan, H., Painter, S., Hyman, J., 2015. Effect of advective flow in fractures and matrix diffusion on natural gas production. *Water Resour. Res.*, under review.
- Pacala, S., Socolow, R., 2004. Stabilization wedges: solving the climate problem for the next 50 years with current technologies. *Science* 305, 968–972.
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