

## NC State University - Radiation Transport Group

### THOR User's Manual

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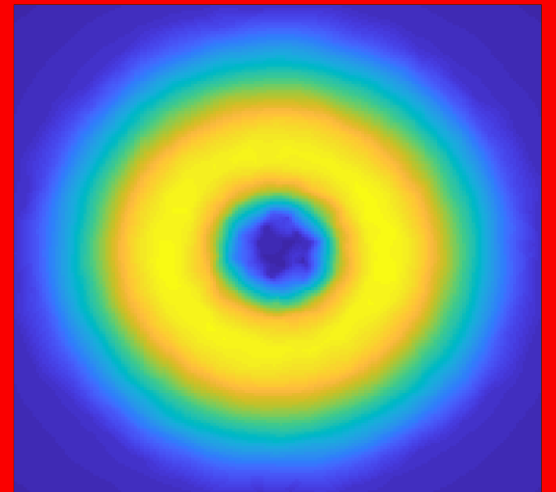
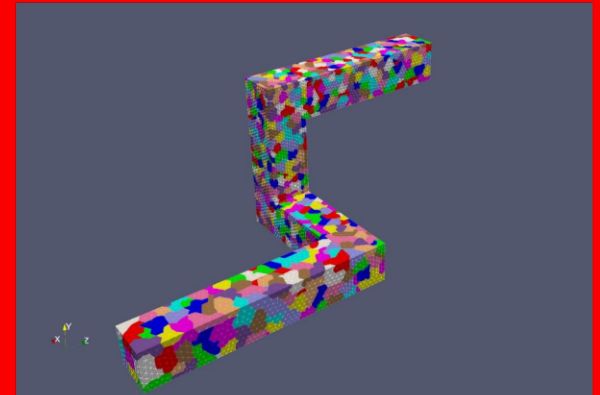
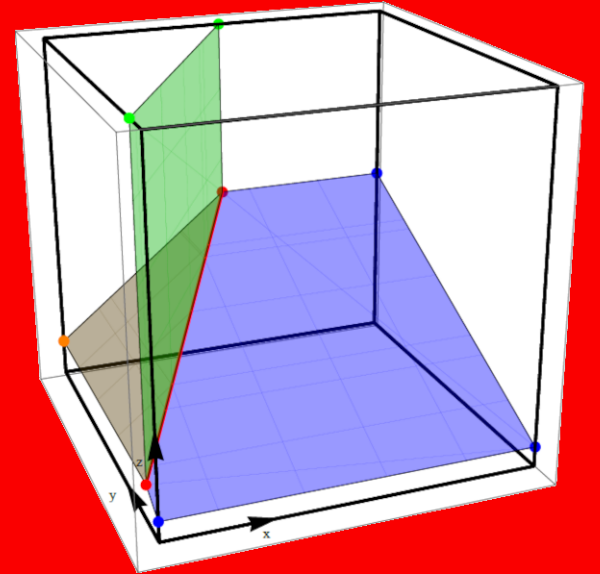
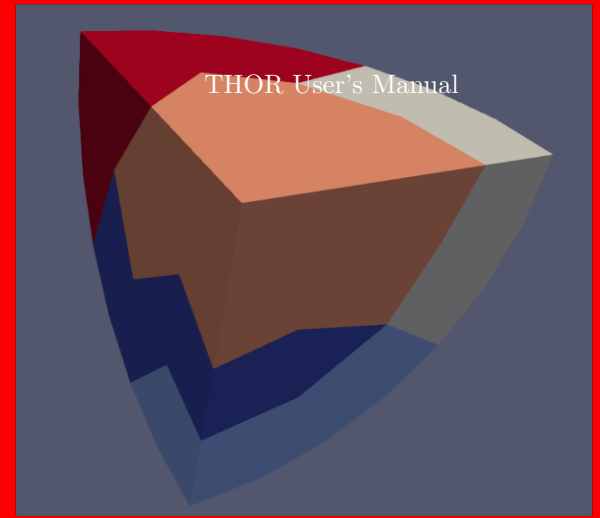
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03/16/2022



## Revision Log

Revision	Date	Affected Pages	Revision Description
0	03/16/2022	All	Initial Release

# Acronyms

# Contents

<b>1</b>	<b>System Requirements</b>	<b>2</b>
<b>2</b>	<b>Getting Started</b>	<b>3</b>
2.1	Accessing THOR on github . . . . .	3
2.2	Cloning THOR repository from github . . . . .	3
2.3	Updating the devel branch . . . . .	5
2.4	Obtaining lapack dependencies . . . . .	5
2.5	Compiling THOR . . . . .	6
2.6	Running THOR for the first time . . . . .	7
2.7	Pre/post Processors . . . . .	8
2.7.1	Setting up THOR_MESH_Generator . . . . .	8
<b>3</b>	<b>Godiva Tutorial</b>	<b>11</b>
3.0.1	Godiva Mesh . . . . .	12
3.0.2	Cross section data . . . . .	13
3.0.3	THOR input file and executing THOR . . . . .	13
<b>4</b>	<b>Input Format</b>	<b>16</b>
4.0.1	THOR Transport Solver . . . . .	16
4.0.2	Compilation and Invocation . . . . .	16
4.0.3	Standard Input . . . . .	17
4.0.4	THOR Mesh Format . . . . .	20
4.0.5	THOR Cross Section Format . . . . .	20
4.0.6	List of all Inputs and Outputs of THOR transport solver . . . . .	21

4.0.7	THOR Mesh Generator . . . . .	21
4.0.7.1	Compilation and Invocation . . . . .	21
4.0.7.2	Format of the THOR's mesh generator standard input . . . . .	22
4.0.7.3	Formatting instructions for regions, source, and boundary id reassignment files	23
4.0.8	Testing . . . . .	23
<b>References</b>		<b>24</b>

The purpose of the User Manual is to provide the novice user with the necessary instructions to install, compile, and execute the Tetrahedral-grid High Order Radiation (THOR) transport code.

# 1. System Requirements

- UNIX-like operating system.
- mpi
- GNU make
- (Conditional on setup method) Version control software git

The THOR team has good experiences with setting up your environment following the [MOOSE](#) setup instructions.

## 2. Getting Started

### 2.1 Accessing THOR on github

THOR is hosted at North Carolina State University's github repository. Open a browser and navigate to:

```
github.ncsu.edu
```

Log in with your unity ID and password. For accessing THOR you have to be a member of the THOR project. Please contact the code owner [Yousry Azmy](#) to be added to the project's membership. Once you have obtained access to THOR, click on the THOR link on THOR's github page, then click fork, and then your username; below we refer to this username as `<git_usr>`. This creates your own personal THOR repository that is separate from the main repository. You have write access to this repository, while you most likely do not have write access to the main repository.

### 2.2 Cloning THOR repository from github

This subsection describes how to clone, i.e. copy a fresh version, of THOR from the github repository to your local computer. The first step is to set up ssh keys. This can be accomplished by following the directions provided [here](#).

Now, the process of cloning is described. First open a terminal window on your local computer where you wish to clone THOR. In the following description terminal commands are indicated by `>>`; by `/home/<usr>` the home directory is indicated but it is understood that `<usr>` must be replaced by the actual user name on the local computer; this tutorial also assumes that `git` is installed on the local computer and is accessible to `<usr>`. In addition, we recognize that username on the local computer, `<usr>`, can be different from the same user's github username, `<git_usr>`, hence they are distinguished in the following instructions by different notation. It is a good idea to create a folder for all github projects, e.g. by

```
>> cd /home/<usr> ; mkdir projects
```

Navigate to the `projects` directory.

```
>> cd projects
```

Clone THOR by typing:



```
>> git clone git@github.ncsu.edu:NCSU-Rad-Transport/THOR.git
```

Alternatively, a user who will not communicate frequently with THOR's github repository can avoid establishing ssh keys and clone THOR directly by issuing the following line-command on the local computer:

```
git clone https://github.ncsu.edu/NCSU-Rad-Transport/THOR.git
```

Now, navigate into the THOR directory and check if things are properly set up.

```
>> cd THOR
```

First check the current branch:

```
>> git branch
```

It should return

```
>> * devel
```

indicating that `devel` is the current branch. The `devel` branch (short for development) contains the most up to date version of THOR. The current branch can be changed by:

```
>> git checkout <branch>
```

where `<branch>` is the branch name to switch to. Next, the remotes are set up. The remotes are addresses to remote repositories and serve as shorthand when information is pulled from or pushed to one of the remotes. The convention is to call the remote of the master repository *upstream*, and to call the remote of the personal repository *origin*. To check the remotes, type:

```
>> git remote -v
```

This should show the following:

```
>> origin git@github.ncsu.edu:NCSU-Rad-Transport/THOR.git (fetch)
>> origin git@github.ncsu.edu:NCSU-Rad-Transport/THOR.git (push)
```

To set up the remote according to the THOR convention, type:

```
>> git remote rm origin
>> git remote add upstream git@github.ncsu.edu:NCSU-Rad-Transport/THOR.git
>> git remote add origin git@github.ncsu.edu:<git_usr>/THOR.git
```

Checking the remote again should show:

```
>> upstream git@github.ncsu.edu:NCSU-Rad-Transport/THOR.git (fetch)
>> upstream git@github.ncsu.edu:NCSU-Rad-Transport/THOR.git (push)
>> origin git@github.ncsu.edu:<git_usr>/THOR.git (fetch)
>> origin git@github.ncsu.edu:<git_usr>/THOR.git (push)
```

## 2.3 Updating the devel branch

As THOR is developed, the local devel branch or any other user-created branches will become outdated. This section demonstrates how to obtain the most up-to-date version of THOR. It is a good idea to check the current status of the repository, by navigating to the THOR directory and typing:

```
>> git status
```

This command reveals if there are any modified or uncommitted files. Updating the current branch is prohibited if there are any modified files. Let us first assume that there is a modified file called `/path/to/modified_file.txt`. There are two options:

- You can stash the file by:

```
>> git stash
```

This removes the modification and stores it in the stash. To un-stash the modifications, do:

```
>> git stash pop
```

- You can commit the file by:

```
>> git add /path/to/modified_file.txt
>> git commit -m "A message for this commit"
```

After either committing or stashing, the branch can be updated by:

```
>> git pull --rebase upstream devel
```

Updating can lead to merge conflicts when local changes conflict with changes in the upstream version of the devel branch. Conflict resolution is beyond the scope of this primer. Please consult git literature or google for guides on conflict resolution.

## 2.4 Obtaining lapack dependencies

THOR depends on certain lapack routines. These are provided with THOR as a submodule. The lapack submodule can be initialized by:

```
>> git submodule update --init
```

The lapack submodule is not expected to change at all. However, if it does, the THOR repository keeps track of the associated version of the lapack repository, so after updating as described in Sect. 2.3, the user may run:

```
>> git submodule update
```

to obtain the latest lapack submodule. If as expected lapack hasn't changed an empty line will be displayed.

## 2.5 Compiling THOR

This section describes how to compile THOR and its dependencies. If THOR is not set up from github, then this is your entry point for the *Getting started* tutorial. Simply unzip the THOR directory where you want it to reside; this tutorial assumes that THOR is unzipped in the `/home/<usr>/projects` directory.

The first step is to compile the lapack dependency. To this end, navigate to:

```
>> cd /home/<usr>/projects/THOR/contrib/scripts
```

Edit the file `make.inc` to specify the MPI Fortran compiler available on the local machine. Also, if necessary, enter command line that modify the environment to enable the compilation process to find the path to required executables; these typically have the form `>> load module pathname`, where `pathname` is a directory on the local computer where these necessary executables reside. Execute the `build_lapack.sh` script by (first command may not be necessary, it only ensures that `build_lapack.sh` is executable):

```
>> chmod +x build_lapack.sh
>> ./build_lapack.sh <n>
```

where `<n>` is the number of processors. For example, on Idaho National Laboratory's Sawtooth HPC the compiler is set in `make.inc` via the statement `FORTTRAN = mpif90`, and the environment is modified with the command line

```
>> module load mvapich2/2.3.3-gcc-8.4.0".
```

A successful lapack build will conclude the scrolled output on the screen with a table of the form:

```
--> LAPACK TESTING SUMMARY <--
Processing LAPACK Testing output found in the TESTING directory
```

SUMMARY	nb test run	numerical error	other error
REAL	1291905	0 (0.000%)	0 (0.000%)
DOUBLE PRECISION	1292717	0 (0.000%)	0 (0.000%)
COMPLEX	749868	0 (0.000%)	0 (0.000%)
COMPLEX16	749588	1 (0.000%)	1 (0.000%)
--> ALL PRECISIONS	4084078	1 (0.000%)	1 (0.000%)

Now, THOR can be compiled. Navigate to the THOR source folder:

```
>> cd /home/<usr>/projects/THOR/THOR/src
```

and, as before, edit the file `Makefile` to utilize the available MPI Fortran compiler and if necessary modify the environment to enable `make` to locate the compiler. Then type:

```
>> make
```

Successful compilation of THOR will conclude with the line:

```
mv ./thor-1.0.exe ../
```

The THOR executable (named in the above line) can be found here:

```
>> ls /home/<usr>/projects/THOR/THOR/
```

that should produce:

```
doc  examples  hello_world  scripts  src  tests  thor-1.0.exe  unit
```

## 2.6 Running THOR for the first time

Navigate to the `hello_thor` directory:

```
>> cd /home/<usr>/projects/THOR/THOR/hello_world
```

Check the content of this folder:

```
>> ls
```

It should show the following files:

```
>> ls -l
total 696
-rwxrwxr-x 1 azmyyy azmyyy 603 Jun 26 19:56 hello_world.in
-rw-rw-r-- 1 azmyyy azmyyy 42700 Jun 26 19:56 hello_world.o
-rwxrwxr-x 1 azmyyy azmyyy 150040 Jun 26 19:56 hello_world.thrm
-rwxrwxr-x 1 azmyyy azmyyy 23 Jun 26 19:56 hello_world.xs
```

These files have the following significance:

- `hello_world.in` is a sample input file to THOR. This file is used to execute THOR.
- `hello_world.thrm` is the corresponding mesh file that is referenced within `hello_world.in`. At this point, it is only important that it is present and has the proper THOR mesh format. Creation of THOR mesh files is covered later in this manual.
- `hello_world.xs` is the corresponding cross section file, also referenced within `hello_world.in`, and again at this point, it is only important that it is present.
- `hello_world.o` is the corresponding output file created by redirecting THOR's standard output. This file can be used to compare THOR's printed output with what it should be upon correct termination of this run.

THOR is invoked at a minimum with the executable name and the standard input file that is specified after the `-i` modifier.

```
>> ../thor-1.0.exe -i hello_world.in
```

For parallel execution type:

```
>> mpiexec -n <n> ../thor-1.0.exe -i hello_world.in
```

where <n> is the number of processors. Several files should have been created:

- hello\_world.flux
- hello\_world.fluxeven
- hello\_world.fluxodd
- hello\_world.in\_out.csv
- intermediate\_output\_even.dat
- intermediate\_output\_odd.dat

The significance of these files will be discussed later. THOR's standard output should start with a banner and conclude with:

```
-----
      Region averaged reaction rates
-----

-- Region --    0 Volume=    1.500000E+01

      Group          Flux          Fission      Absorption      Fiss Src
      1    9.515584E-01  1.284604E+00  8.564026E-01  1.284604E+00
Total    9.515584E-01  1.284604E+00  8.564026E-01  1.284604E+00

-----
      Execution of THOR completed successfully
-----
```

## 2.7 Pre/post Processors

### 2.7.1 Setting up THOR\_MESH\_Generator

THOR mesh generator converts *exodus* and *gms* mesh formats to THOR's native mesh format. It also permits uniform refinement of meshes provided in *exodus* files. Conversion from *exodus* format and uniform refinement uses the *libmesh* [5] *meshtool*. Therefore, *libmesh* has to be set up first. To this end, navigate to the `scripts` directory:

```
>> cd /home/<usr>/projects/THOR/contrib/scripts
```

and execute `build_libmesh.sh`:

```
>> chmod +x build_libmesh.sh
>> ./build_libmesh.sh <n>
```

where the first command makes `build_libmesh.sh` executable (if it is not already) and `<n>` is the number of processors. It must be provided even if it is simply 1. Executing this script may take a long time to complete installing *libmesh*, however, it will show progress on the screen. If the git-clone command in the `build_libmesh.sh` script does not work, replace it with the command:

```
git clone https://github.com/libMesh/libmesh.git
```

Finally, the `THOR_LIBMESH_DIRECTORY` environment variable has to be set. This environment variable must point to the directory that the `meshtool-opt` executable is located. For the standard installation, one should execute:

```
>> export THOR_LIBMESH_DIRECTORY=/home/<usr>/projects/THOR/contrib/libmesh/build
```

The next step is to make the `THOR_MESH_GENERATOR` application. Navigate to its source folder:

```
>> cd /home/<usr>/projects/THOR/pre-processors/THOR_Mesh_Generator/src
```

and type:

```
>> make
```

The executable

```
/home/<usr>/projects/THOR/pre-processors/THOR_Mesh_Generator/Thor_Mesh_Generator.exe
```

should have been created.

```
*****
```

The tests as described below did not execute as described.

Instead, I did the following and still execution of the tests did not work properly:

```
1. In ~/PROJECTS/THOR/pre-processors/THOR_Mesh_Generator:
ln -s Thor_Mesh_Generator\MP.exe Thor_Mesh_Generator.exe
```

```
2. In ~/PROJECTS/THOR/pre-processors/THOR\_Mesh_Generator/scripts:
chmod u+x test\_all.sh
```

```
3. ./test\_all.sh
```

This ran but did not give the output below and reported execution errors. It is not clear if these reported errors are part of the testing since some of the cases are labeled Bad, or the error indicates erroneous installation of *libmesh*.

```
*****
```

To ensure that the `THOR_MESH_GENERATOR` application compiled correctly, execute the regression tests. Change directory to:

```
>> cd /home/<usr>/projects/THOR/pre-processors/THOR_Mesh_Generator/scripts
```

and execute:

```
>> python run_thor_tests.py
```

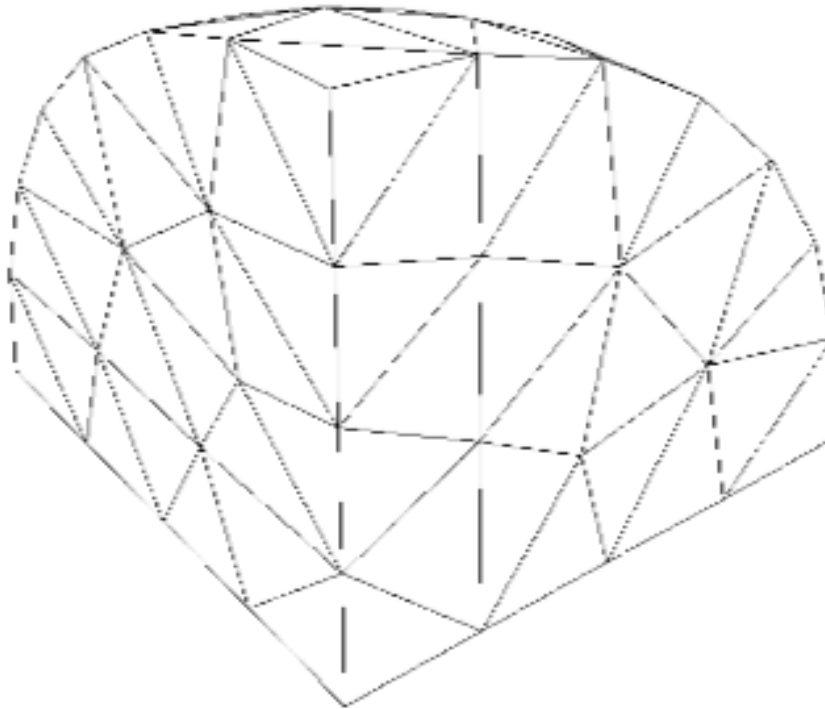
You should see screen output similar to this:

```
-----
Test  1 tests/bad_gmesh_non_tet_element:bad_gmesh_non_tet_element success
Test  2 tests/bad_gmesh_no_elements_block:bad_gmesh_no_elements_block success
Test  3 tests/homogeneous_domain:homogeneous success
Test  4 tests/homogeneous_domain:homogeneous_from_exodus success
Test  5 tests/homogeneous_domain:homogeneous_r1 success
Test  6 tests/homogeneous_domain:homogeneous_from_exodus_r1 success
Test  7 tests/bad_gmesh_no_nodes_block:bad_gmesh_no_nodes_block success
Test  8 tests/bad_gmesh_no_format_block:bad_gmesh_no_format_block success
Test  9 tests/bad_gmesh_non_tri_face:bad_gmesh_non_tri_face success
Test 10 tests/convert_old_to_new_THOR:convert_old_to_new_THOR success
Test 11 tests/unv_sphere_in_shell_in_box:unv_sphere_in_shell_in_box success
Test 12 tests/Basic_Cube_Mesh_test:basic_cube_mesh success
Test 13 tests/split_hex_and_prism:split_hex success
Test 14 tests/split_hex_and_prism:split_prism success
Test 15 tests/bad_unv_sphere_in_shell_in_box_no_2411:bad_unv_sphere_in_shell_in_box_no_2411 success
-----
Successes:  15          Failures:  0
```

All or at least the vast majority of tests should pass, so **Failures** should be close to zero.

### 3. Godiva Tutorial

Godiva is an un-shielded, pulsed, nuclear burst reactor. It is essentially a homogeneous sphere of highly enriched uranium with a diameter of 30 cm, that was operated by inserting a piston of fissile material [2]. In this tutorial the critical benchmark configuration described in Ref. [1] is considered. The geometry that is modeled by THOR is a homogeneous sphere of radius 8.71 cm discretized by tetrahedra similar to Fig. 3.1 (with the exception that Fig. 3.1 shows one-eighth of the domain). The energy domain is discretized with six energy groups, and cross sections are provided by [1].



**Figure 3.1:** Coarse mesh for Godiva problem, picture courtesy [3]

This tutorial first explains how a tetrahedral mesh is created for the Godiva problem, then the cross sections data input is discussed, and finally the standard input to THOR is covered. The input files discussed below for the Godiva tutorial are located in:



```
>> /home/<usr>/projects/THOR/THOR/examples/Godiva_tutorial
```

### 3.0.1 Godiva Mesh

The workflow described here is suitable if the user has access to the Cubit mesh generator. A Cubit journal file is provided in directory:

```
>> /home/<usr>/projects/THOR/THOR/examples/Godiva_tutorial
```

It creates the exodus file **godiva.e**. To verify whether Cubit is available to the user on the target computer execute the command line:

```
>> which cubit
```

Note that even if Cubit is installed on the target computer it might not be available to the user unless its path is defined in the user's search paths. To execute the journal file, the last line must be modified by replacing **<path>** with:

```
/home/<usr>/projects/THOR/THOR/examples/Godiva_tutorial ..  
/create_mesh/from_cubit
```

then execute Cubit with the command line:

```
>> cubit <godiva_mesh_CUBIT.jou
```

For users that do not have access to Cubit, the **godiva.e** file is provided in the **from\_cubit** directory. The exodus file **godiva.e** is converted to THOR's native mesh format by executing the THOR mesh generator using the standard input file **convert\_godiva.in** that is included in the **from\_cubit** directory with the command line:

```
>> /home/<usr>/projects/THOR/pre-processors/THOR_Mesh_Generator/Thor_Mesh_Generator_MP.exe -i conver
```

In this case this input file contains the following lines:

```
./godiva_1_c.e  
./godiva_1_c.thrm
```

specifying the input exodus file and the THOR mesh formatted output file. The mesh conversion infers the file format of the input by the filename extension; currently **.e** and **.gmsh** for exodus and gmsh formats, respectively, are supported. The conversion is performed by typing:

```
/home/<usr>/projects/THOR/pre-processors/THOR_Mesh_Generator/ ...  
Thor_Mesh_Generator.exe -i convert_godiva.in
```

After successful completion of the conversion, the following printout should appear:

```
=====
```

```
TTTTTTT HH      HH 00000 RRRRRR
TTT      HH      HH 0000000 RRRRRRR
TTT      HH      HH 00    00 RR  RR
TTT      HHHHHHHHH 00    00 RRRRRR
TTT      HHHHHHHHH 00    00 RRRR
TTT      HH      HH 00    00 RR  RR
TTT      HH      HH 0000000 RR  RR
TTT      HH      HH 00000  RR  RR
```

Tetrahedral High Order Radiation Transport Code

By North Carolina State University

Version 1.0 - Update 2020

```
=====
```

```
=====
```

```
Executing -->THOR_MESH_GENERATOR<-- application
```

```
=====
```

```
----- Input -----
```

```
Infile mesh file name: ./godiva_1_c.msh
Output mesh file name: ./godiva_1_c.thrm
Mesh refinement level: 0
```

```
No region ID edits are provided
No source ID edits are provided
No boundary ID edits are provided
```

```
=====
Application -->THOR_MESH_GENERATOR<-- terminated successfully
=====
```

The file `godiva_1_c.thrm` should result from this execution for use by THOR. This concludes the mesh generation step for this tutorial.

### 3.0.2 Cross section data

The THOR cross section file for the Godiva benchmark is provided by:

```
/home/<usr>/projects/THOR/THOR/examples/Godiva_tutorial/cross_sections/godiva.xs
```

THOR uses a custom cross section format that is explained in detail in Sec. 4.0.1.

### 3.0.3 THOR input file and executing THOR

The THOR input file is

```
/home/<usr>/projects/THOR/THOR/examples/Godiva_tutorial/...
THOR/godiva.i
```

THOR uses a keyword-based input that is listed in Sect. 4.0.1. The Godiva tutorial input file is now explained in detail; note that this listing includes comments separated by # at the end of each line that are not present in the original input file.

```
Godiva 6 group example          # title line

start problem                   # beginning of problem block
  execution = yes               # problem is solved
  lambda = 0                   # constant approximation (SC)
  type = keig                  # solve an eigenvalue problem
  keigsolver = pi ; piacc = none # Power iterations w/o
                                # acceleration, note separation
                                # by ; to have same-line syntax
  sweep = precomp              # sweep path precomputed
  page_refl = save             # reflected bnd fluxes stored
  kconv = 1E-8                 # stopping tolerance on k-eff
  innerconv = 1E-12            # stopping tolerance inner it
  outerconv = 1E-7             # stopping tolerance outer it
  maxinner = 4 ; maxouter = 5000 # max # inner/outer it
end problem                    # terminate the problem block

start inout                     # start input/output file block
  mesh_file = ../create_mesh/from_CUBIT/godiva_1_c.thrm # mesh file
  xs_file = ../cross_sections/godiva.xs # XS file
  flux_file = godiva.flux       # plain flux output file
  vtk = flux mat               # print flux and mat ids to vtk
  print_xs = yes               # XS are echoed
end inout                      # terminate this block

start cross_sections
  ngroups = 6                  # the number of energy groups
  upscattering = no           # upscattering is not present
                                # scat. XS is lower triangular
end cross_sections

start quadrature
  qdtype = levelsym           # level-symmetric quadrature
  qdorder = 4                 # order 4
end quadrature

start regionmap
  1                           # the region map maps block
                                # ids to materials; in this
                                # case the only block is
end regionmap                 # assigned material 1 that is
                                # present in XS file

end file                       # the input file is terminated
                                # by "end file"
```

The Godiva tutorial is solved with THOR via the command line:

```
>> /home/<usr>/projects/THOR/THOR/thor-1.0.exe -i godiva.i
```

Completion of execution of the Godiva tutorial is indicated by the printout:

```
-----
Execution of THOR completed successfully
-----
```

THOR provides the following output that is discussed in this tutorial:

- The final estimate of the multiplication factor is printed under “Execution Summary”, “Final eigenvalue”. In this case the value is 0.9611. This is not close to critical because the mesh that is created does not conserve the volume of the critical sphere. When creating a tetrahedral mesh, Cubit places the nodes on the sphere’s surface, but that necessarily leads to the volume of the discretized geometry to be smaller than the original volume. In this case, the computational volume is 2742.4 cm<sup>3</sup>, while the actual volume is 2767.8 cm<sup>3</sup>.
- A summary of group-wise, region-averaged reaction rates is provided for each region identifier separately under “Region averaged reaction rates”. The volume of each region, and group-wise fluxes, fission, absorption, and fission source rates are listed.
- Two vtk formatted files, `flux.vtk` contains spatial flux maps, and `mat.vtk` contains the material map. These files can be opened with the paraview post-processing tool that is available [here](#).

The reaction rate summary is given by:

```
-----
Region averaged reaction rates
-----

-- Region --   1 Volume=   2.742401E+03

Group      Flux      Fission    Absorption    Fiss Src
  1  8.424948E-01  1.399140E-01  4.946487E-02  1.399140E-01
  2  1.573081E+00  2.333728E-01  9.509448E-02  2.333728E-01
  3  9.814645E-01  1.379655E-01  5.968679E-02  1.379655E-01
  4  1.631588E+00  2.215631E-01  1.006868E-01  2.215631E-01
  5  1.198341E+00  1.915944E-01  9.089592E-02  1.915944E-01
  6  1.792724E-01  4.652155E-02  2.413688E-02  4.652155E-02
Total    6.406241E+00  9.709313E-01  4.199657E-01  9.709313E-01
```

The results can be improved greatly by enforcing volume conservation on the Cubit mesh. This is accomplished manually in this tutorial by changing the radius in the Cubit journal file to .

## 4. Input Format

### 4.0.1 THOR Transport Solver

add a description of input format for

- standard input
- XS
- mesh
- source

This section describes the input format of the THOR transport solver.

### 4.0.2 Compilation and Invocation

Navigate to:

```
>> cd /home/<usr>/projects/THOR/THOR/src
```

and make the application:

```
>> make
```

The executable `THOR_TRANSPORT_SOLVER.exe` should have been created here:

```
>> /home/<usr>/projects/THOR/THOR/THOR_TRANSPORT_SOLVER.exe
```

The THOR transport solver is invoked by

```
>> mpiexec -n <n> THOR_TRANSPORT_SOLVER.exe -i standard_input -t
```

where `-i` precedes the name of the standard input file `standard_input` and `-t` is an optional parameter for additional timing. **Remark:** `-i standard_input` must currently follow the executable name `THOR_TRANSPORT_SOLVER.exe`. This will be fixed in future versions.

### 4.0.3 Standard Input

THOR input is organized in blocks. The blocks are

- The first line must contain a user selected name for the problem.
- **problem**: general parameters to define the problem to be solved
- **inout**: Names of inputs and outputs files and toggles for specific output.
- **cross\_sections**: parameters pertaining to the cross section data.
- **quadrature**: parameters pertaining to the angular quadrature.
- **postprocess**: parameters pertaining to postprocessing outputs. **cartesian\_map** sets up an overlaid Cartesian mesh that fluxes and reactions rates are averaged over. The Cartesian mesh is defined by the minimum and maximum coordinates for each direction (x, y, z) and number of subdivisions between. **point\_value\_locations** allows extraction of flux values at user provided points.
- **regionmap**: mapping from region id to cross section id. Region ids are an integer assigned to each tetrahedral element that are used to group elements into regions or blocks (see Sect. 4.0.4). Cross section ids are indices that identify sets of cross sections provided in the cross section input file (see Sect. 4.0.5).

Blocks are delineated with **start** and **end** keywords like this:

```
start <block_name>
  key1 = value1 ; key2 = value2
  key3 = value3
end <block_name>
```

Each blocks contains several keyword-value pairs. Multiple assignments can be placed on the same line if they are separated by `;`. The keyword-value pairs can be provided in an arbitrary order. All keywords are listed in Table ??.

The **regionmap** block does not contain keyword-value pairs. Instead, it maps region ids to cross section ids. We denote by **min\_reg** and **max\_reg** the smallest and largest region ids in the mesh file. The number of entries in the **regionmap** field must then be **num\_entries** = **max\_reg** - **min\_reg** + 1. The assignment is best illustrated for an example. Let us assume that **min\_reg** = -1 and **max\_reg** = 2 and we want to assign

```
-1 -> 12
0 -> 1
1 -> 1
2 -> 3
```

Then the **regionmap** block is given by:

```
start regionmap
  12 1 1 3
end regionmap
```

Unused region ids can be accommodated by padding the entries in the **regionmap** field.

**Table 1:** Keywords of THOR Transport Solver Application

Keyword	Type	Options	Explanation
<b>Block:</b> problem			
type	string	keig/fsrc	Problem type. Either eigenvalue (keig) or fixed source (fsrc)
keigsolver	string	pi/jfnk	Solve type for keig. Either power iteration (pi) or Jacobian-Free Newton
lambda	integer	-	Expansion order, negative number indicates reduced set
inflow	string	yes/no	If fixed inflow boundary conditions are provided for fsrc problems
piacc	string	errmode/none	Type of power iteration acceleration: none or error mode extrapolation
sweep	string	precomp	Must be set to precomp at this point. (Redundant keyword)
page_sweep	string	yes/no	If the sweep path is saved or is paged to scratch file when not needed
page_refl	string	page/save/inner	If significant angular fluxes are paged to/from scratch file (page), stored
page_iflw	string	bygroup/all	If inflow information is loaded to memory completely (all) or for each gr
kconv	Real	-	Stopping criterion for eigenvalue
innerconv	Real	-	Stopping criterion for group flux during inner iteration
outerconv	Real	-	Stopping criterion for group flux during outer/power iteration
maxinner	integer	-	Maximum number of inner iterations
maxouter	integer	-	Maximum number of outer/power iterations
jfnk_krsze	integer	-	Maximum size of Krylov subspace during jfnk
jfnk_maxkr	integer	-	Maximum number of Krylov iterations
jfnk_method	string	outer/flat/flat_wds	Type of jfnk formulation, see []
initial_guess	string	yes/no	If an initial guess file should be read
save_restart	string	yes/no	If a restart file should be written
ipiter	integer	-	Number of initial power iterations for jfnk
print_conv	string	yes/no	If convergence monitor is written to file <b>thor.convergence</b>
density_factor	string	no/byvolume/fromfile	Density factor options: use no density factors (no), use density factors a
execution	string	yes/no	If yes problem is executed, if no then input is only read and checked.

**Table 2:** Keywords of THOR Transport Solver Application

Keyword	Type	Options	Explanation
<b>Block:</b> inout			
mesh_file	string	-	Name of the mesh file
inflow_file	string	-	Name of the boundary inflow information file
source_file	string	-	Name of the volumetric source file
flux_file	string	-	Name of the THOR formatted output flux file
xs_file	string	-	Name of the cross section file
density_factor_file	string	-	Name of the density factor file
quad_file	string	-	Name of the angular quadrature file
vtk	string (multiple)	flux/mat/reg/src	Which information is written to vtk format
vtk_flux_file	string	-	Name of the vtk flux file
vtk_mat_file	string	-	Name of the vtk material file
vtk_reg_file	string	-	Name of the vtk region file
vtk_src_file	string	-	Name of the vtk volumetric source file
restart_file	string	-	Name of the restart file written when saving
inguess_file	string	-	Name of the initial guess file if initial guess is used
cartesian_map_file			Name of the file that the cartesian map is written to
print_xs	string	yes/no	If cross sections are echoed to standard output
<b>Block:</b> cross_sections			
ngroups	integer	-	Number of energy groups
pnorder	integer	-	Spherical harmonics order used for scattering
pnread	integer	-	Spherical harmonics expansion provided
upscattering	string	yes/no	Read upscattering data from cross section file
multiplying	string	yes/no	If the cross section file contains fission information
scatt_mult_included	string	yes/no	If the scattering data includes the $2l + 1$ Legendre moments
<b>Block:</b> quadrature			
qdtype	string	levelsym/legcheb/fromfile	Quadrature type: level-symmetric, Legendre, or from file
qdorder	integer	-	Order of the angular quadrature
<b>Block:</b> postprocess			
cartesian_map	Real/integer (9 entries)	-	xmin, xmax, nx, ymin, ymax, ny, zmin, zmax, nz
point_value_locations	Real (3 N)	-	N is the number of points, (x,y,z) coordinates



#### 4.0.4 THOR Mesh Format

Line 1: number of vertices

Line 2: number of elements

Line 3: unused enter 1

Line 4: unused enter 1

Block 1: vertex coordinates, number of lines = number of vertices; each line is as follows: vertex\_id (integer) x-coordinate (Real) y-coordinate (Real) z-coordinate (Real)

Block 2: region and source id assignments, number of lines = number of elements; each line is as follows: element\_id region\_id source\_id (all integers). For setting up Monte Carlo on the tet mesh, this block can be ignored.

Block 3: element descriptions, the vertex\_ids that form each element. Number of lines = number of elements; each line is as follows: element\_id vertex\_id1 vertex\_id2 vertex\_id3 vertex\_id4 (all integers).

Next line: number of boundary face edits

Block 4: boundary face descriptions. All exterior faces associated with their boundary condition id, number of lines = number of boundary face edits; each line is as follows: element\_id local\_tetrahedron\_face\_id boundary\_condition\_id.

**Explanation:** local\_tetrahedron\_face\_id: natural local id of tetrahedrons face which is the id of the vertex opposite to this face. Note: indexed 0-3. boundary\_condition\_id: value = 0: vacuum BC value = 1: reflective BC value = 2: fixed inflow

Next line: number of adjacency list entries

Block 5: adjacency list, number of lines = number of adjacency list entries; each line is as follows: element\_id face\_id neighbor\_id neighbor\_face\_id. Explanation: The element\_id is the current element. The neighbor across the face indexed by face\_id has the element id neighbor\_id and the its own local index for the said common face is neighbor\_face\_id.

#### 4.0.5 THOR Cross Section Format

Line 1: number of materials

Block 1: each entry in this block contains cross sections for a single material. Each entry contains L

Entry line 1: material\_id

Entry line 2: fission\_spectrum\_1 fission\_spectrum\_2 fission\_spectrum\_G

Entry line 3: energy\_group\_boundary\_1 energy\_group\_boundary\_3 energy\_group\_boundary\_G

Entry line 4: fission\_xs\_1 fission\_xs\_2 fission\_xs\_3 fission\_xs\_G

Entry line 5: nu\_bar\_1 nu\_bar\_2 nu\_bar\_G

Entry line 6: total\_xs\_1 total\_xs\_2 total\_xs\_G

```

Entry line 7: sig_scatter_{0, 1->1} sig_scatter_{0, 2->1} sig_scatter_{0, G->1}
Entry line 8: sig_scatter_{0, 1->2} sig_scatter_{0, 2->2} sig_scatter_{0, G->2}

:
Entry line G + 6: sig_scatter_{0, 1->G} sig_scatter_{0, 2->G} sig_scatter_{0, G->G}
Entry line G + 7: sig_scatter_{1, 1->1} sig_scatter_{1, 2->1} sig_scatter_{1, G->1}
Entry line G + 8: sig_scatter_{1, 1->2} sig_scatter_{1, 2->2} sig_scatter_{1, G->2}

:
Entry line 2 * G + 6: sig_scatter_{1, 1->G} sig_scatter_{1, 2->G} sig_scatter_{1, G->G}
:

```

- $G$  = total number of groups.
- $L$  = scattering expansion.
- `fission_spectrum_g`: fraction of neutrons born in fission that appear in energy group  $g$ .
- `energy_group_boundary_g`: currently unused, can be filled with 0s. Upper bound of energy group  $g$ .
- `fission_xs_g`: fission cross section (NOTE: not  $\nu_{\text{bar}} * \text{fission\_xs}$ ) in group  $g$ .
- `nu_bar_g`: average number of neutrons released by fission caused by a neutron in energy group  $g$ .
- `total_xs_g`: total cross section in energy group  $g$ .
- `sig_scatter_l, g->g`:  $l$ -th Legendre polynomial moment of the scattering cross section from group  $g$  to  $g$ . The  $(2 * l + 1)$  factor may be included in the value of the cross section or not, THOR can handle both cases. It needs to be specified separately every time.

## 4.0.6 List of all Inputs and Outputs of THOR transport solver

### 4.0.7 THOR Mesh Generator

This section discusses the input of THOR's mesh generator tool. The mesh generator tool provides the following capabilities:

- Convert exodus [6] formatted meshes to THOR format.
- Convert gmsh [4] formatted meshes to THOR format.
- Convert universal file format meshes to THOR format.
- Split the elements of hexahedra and wedge (triangular prisms) meshes into tetrahedra before converting to THOR mesh format.
- Conversion of legacy THOR mesh format to new format.

#### 4.0.7.1 Compilation and Invocation

Navigate to:

```
>> cd /home/<usr>/projects/THOR/pre-processors/THOR_Mesh_Generator/src
```

and make the application:

```
>> make
```

The executable `Thor_Mesh_Generator.exe` should have been created here:

```
>> /home/<usr>/projects/THOR/pre-processors/THOR_Mesh_Generator/Thor_Mesh_Generator.exe
```

The application is invoked with:

```
>> /path/to/Thor_Mesh_Generator.exe -i standard_input
```

where `standard_input` is the standard input file.

**Remark:** Conversion of exodus files to gmsh files relies on using libMesh's `mesh-tool` [5]. You must compile libMesh and set the environment variable `THOR_LIBMESH_DIRECTORY`.

#### 4.0.7.2 Format of the THOR's mesh generator standard input

The standard input file of the THOR mesh generator contains the following lines:

```
input_mesh_file
output_mesh_file
region_id_file
source_id_file
boundary_id_file
```

where `input_mesh_file` and `output_mesh_file` are required parameters, while the remaining parameters are optional. If a parameter is omitted, the line should just be left blank (that means that e.g. `source_id_file` will always be provided on line 4 regardless of whether `region_id_file` was provided).

The purpose of these files is as follows:

- `input_mesh_file`: name of the input mesh file. Note that the file ending matters: exodus is `.e` and gmsh is `.msh`, because the mesh format is inferred from it.
- `output_mesh_file`: name of the THOR mesh format files. Use file ending `.thrm`.
- `region_id_file`: name of file that contains instructions to reassign region (sometimes called block) ids. Formatting instructions for this file is provided in Sect. 4.0.7.3.
- `source_id_file`: name of file that contains instructions to reassign source ids. Formatting instructions for this file is provided in Sect. 4.0.7.3.
- `boundary_id_file`: name of file that contains instructions to reassign boundary ids. Formatting instructions for this file is provided in Sect. 4.0.7.3.

#### 4.0.7.3 Formatting instructions for regions, source, and boundary id reassignment files

IDs are integers that are assigned to each tetrahedral element to group it into a region, a source region or assigned to boundary faces to group it into a set of faces for boundary assignment. The file format for id reassignment files is:

```
n
old_id_<1> new_id_<1>
:
:
old_id_<n> new_id_<n>
```

The meaning is as follows:

- **n**: the number of instructions in the file (i.e. the number of lines following this line).
- **old\_id\_<j>**: the j-th old id that will be replaced by **new\_id\_<j>**.
- **new\_id\_<j>**: the j-th new id that will replace **old\_id\_<j>**

#### Remarks:

- Note, each instruction needs to be on a different line.
- The boundary id characterizes the boundary condition. It must be 0, 1, or 2, where 0 is a vacuum, 1 is a reflective, and 2 is a fixed inflow boundary.

### 4.0.8 Testing

THOR provides a convenient test harness that can be executed by the user by navigating to the **scripts** directory that exists as subdirectory for all applications; for example, for the THOR transport solver the test file is located in:

```
>> /home/<usr>/projects/THOR/THOR/scripts
```

The tests are executed via the **run\_thor\_tests.py** python script. The execution of the python script requires *python3*. Tests are executed by:

```
>> cd /home/<usr>/projects/THOR/THOR/scripts
>> python run_thor_tests.py
```

The output of the test script should look like this:

```
-----
Test  1 name success
:
Test  N name success
-----
Successes:  N           Failures:  0
```

# Bibliography

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