

SeisSol_ParamStudies Manual

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https://github.com/christadler/SeisSol_ParamStudies

This is a bunch of generic scripts and templates that allows you to use SeisSol for parameter studies. It was developed for DT-Geo (DTC-E4, Block 3) to generate a catalogue for AltoTiberina

How to Use the SeisSol_ParamStudies Workflow

1. Set-up one test case (Examples/)
 - a. Use the right directory structure
 - b. Adapt and test the slurm job file
2. Decide on the parameters that should change (Inputs/parameter_study_list.csv)
 - a. Create a csv file for all runs
3. Create .jinja Templates (Inputs/Templates/)
 - a. for all input files that should use parameters from the .csv file
 - b. all input files that don't change should be copied to Inputs/Files/
4. Adapt Scripts/template_parameters.py
 - a. Adapt **tp_param_list**, **tp_list**, **tp_mapping** & add **user_name** + user info below
5. Finally adapt Inputs/Templates/SeisSol_slurm_JobFarm.slurm.jinja (with help)

... now run Scripts/main.py to generate one folder for each run in Outputs/
copy the Outputs/ directory to SuperMucNG and submit the slurm_xxx.scripts

1. Set-up Test Case (and copy it to the Examples/ dir)

On SuperMucNG set up your file structure:
(In your *workdir*, Output should go to *scratch*)

- inputs/
 - <xyz>.yaml
 - parameters.par
- mesh/
- seissol_bin/
 - e.g. SeisSol_Release_dskx_4_elastic
- JobSubmission.slurm

inputs/parameters.par example:

```
MeshFile =  
'../../mesh/mesh topo1km 7 faults smoothed v3 small1  
z'
```

```
OutputFile =  
'/hpf/fs/scratch/0A/di35ban/AltoTiberina/outputs/000  
1/0001'
```

Slurm script example:

```
cd <path in work dir>
```

```
mkdir -p <path in scratch for output>
```

```
srun -n $SLURM_NTASKS  
../seissol bin/<binary file> parameters.pa
```

```
# These files are used for post-processing
```

```
myfile=<path in scratch for output>/<name>-surface  
e.xdmf
```

```
echo "Myfile: $myfile"
```

```
cd <path in scratch for output>
```

```
<post processing>
```

2. Build Inputs/parameter_study_list.csv

	A	B	C	D	E	F	G	H	I
1	id	nucleation	d_c	y	mu_s	R	xha_x	xha_y	xha_z
2	0		1 0.4	0.6	0.35	0.65	295214.0	4817590.0	-8238.0
3	1		1 0.4	0.6	0.35	0.7	295214.0	4817590.0	-8238.0
4	2		1 0.4	0.6	0.35	0.75	295214.0	4817590.0	-8238.0
5	3		1 0.4	0.6	0.35	0.8	295214.0	4817590.0	-8238.0
6	4		1 0.4	0.6	0.35	0.85	295214.0	4817590.0	-8238.0
7	5		1 0.4	0.6	0.4	0.7	295214.0	4817590.0	-8238.0
8	6		1 0.4	0.6	0.4	0.75	295214.0	4817590.0	-8238.0
9	7		1 0.4	0.6	0.4	0.8	295214.0	4817590.0	-8238.0
10	8		1 0.4	0.6	0.4	0.85	295214.0	4817590.0	-8238.0
11	9		1 0.4	0.6	0.45	0.75	295214.0	4817590.0	-8238.0
12	10		1 0.4	0.6	0.45	0.8	295214.0	4817590.0	-8238.0
13	11		1 0.4	0.6	0.45	0.85	295214.0	4817590.0	-8238.0
14	12		1 0.4	0.6	0.45	0.9	295214.0	4817590.0	-8238.0
15	13		1 0.4	0.6	0.5	0.7	295214.0	4817590.0	-8238.0
16	14		1 0.4	0.6	0.5	0.8	295214.0	4817590.0	-8238.0
17	15		1 0.4	0.6	0.5	0.85	295214.0	4817590.0	-8238.0
18	16		1 0.4	0.6	0.5	0.9	295214.0	4817590.0	-8238.0
19	17		1 0.4	0.6	0.55	0.7	295214.0	4817590.0	-8238.0
20	18		1 0.4	0.6	0.55	0.8	295214.0	4817590.0	-8238.0
21	19		1 0.4	0.6	0.55	0.85	295214.0	4817590.0	-8238.0

Example from AltoTiberina:

Each line represents one run.

Column A holds the “id”, starting with zero. It will be used as identifier/”prefix”.

The following columns are named according to the parameters of the templates.

3a. Create .jinja Templates in Inputs/Templates

3b. Copy all other Input Files to Inputs/Files/

To create the .jinja Template, just replace the variable with {{<var_name>}}

Examples:

```
components: !FunctionMap
  map:
    sig_zz: |
      return 2670.0*(1.0-{{y}})*9.8*min(-200.0,z);
```

sig_zz.yaml.jinja

```
!LuaMap
returns: [forced_rupture_time]
function: |
  function f (x)
    xha = { {{xha_x}}, {{xha_y}}, {{xha_z}} }
```

forced_rupture_time.yaml.jinja

```
function f (x)
  R = {{R}}
  return {
    S = (1.0/R-1.0)
  }
end

- !GroupFilter
# This is for the Gubbio fault
groups: [66,67,68,69,70,71]
components: !LuaMap
  returns: [S]
  function: |
    function f (x)
      R = {{R}}
      return {
        S = (1.0/R-1.0)
      }
```

initial_stress.yaml.jinja

4. Adapt Scripts/template_parameters.py

```
tp_param_list = ["nucleation", "d_c", "y", "mu_s", "R"]  
#  
# list of all jinja template files (except slurm tmpl for stage 2)  
# to be filled by user  
# stored in dir "../Input/Templates/*SeisSol_slurm_JobFarm.slurm.jinja"  
# "AltoTiberina_forced_rupture_time.yaml" (xha) is currently not used  
tp_list = ["AltoTiberina_fault.yaml",  
           "AltoTiberina_initial_stress.yaml",  
           "AltoTiberina_sig_zz.yaml",  
           "AltoTiberina_forced_rupture_time.yaml"]  
  
# mapping of parameters to jinja templates  
# to be filled by user  
# "AltoTiberina_forced_rupture_time.yaml" : ["xha"],  
tp_mapping = { "AltoTiberina_fault.yaml" : ["mu_s", "d_c"],  
               "AltoTiberina_initial_stress.yaml": ["R"],  
               "AltoTiberina_sig_zz.yaml": ["y"],  
               "AltoTiberina_forced_rupture_time.yaml": ["xha_x", "xha_y", "xha_z"]  
             }
```

template_parameters.py

tp_param_list

all parameters/columns
from the .csv-file

tp_list

all jinja templates
that should be used

tp_mapping

mapping of parameters to
jinja templates

user_name

(and choose TEST or PRODUCTION)

tp_slurm_account

tp_slurm_email

tp_output_file_dir

5. Finally adapt `Inputs/Templates/SeisSol_slurm_JobFarm.slurm.jinja`

should not be necessary for testing the AltoTiberina catalogue

... now run `Scripts/main.py` to generate
one folder for each run in `Outputs/`

copy the `Outputs/` directory to SuperMucNG, add the mesh and `.asagi-files` (as described in
slide “Set-Up Test Case” and submit the `slurm_XXX.scripts`