

Wikiprint Book

Title: Command Line Interface (CLI) for WRF4G

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Command Line Interface (CLI) for WRF4G

wrf4g_framework

Usage

wrf4g_framework {start|stop|restart|reload|status}

Synopsis

wrf4g_framework manages WRF4G framework components: GridWay and MySQL (in case it is needed). It loads the framework configuration from `$WRF4G_LOCATION/etc/framework4g.conf`.

wrf4g_prepare

Usage

wrf4g_prepare [--dry-run] [--reconfigure] [--verbose] [--help]

Synopsis

Given a file (experiment.wrf4g) describing the experiment, prepare it by creating the realization and chunks needed to perform the experiment.

Options

<code>--dry-run</code>	Perform a trial run with no changes made.
<code>--reconfigure</code>	Reconfigure experiment. With this option we can change the start and end date of the experiments and add new physics. Values are taken from a modified <code>experiment.wrf4g</code> .
<code>--verbose</code>	Verbose mode. Explain what is being done
<code>--help</code>	Shows this help

wrf4g_submit

Usage

wrf4g_submit [--dry-run] [--{exp Experiment|rea Realization|frea File|chunk Chunk|nchunk Number_of_chunks|nrea Number_of_realizations}] [--rerun] [--run-just-one][--priority P] [--verbose] [--force] [--tdep afterok | afternotok | afterany] [--help]

Synopsis

Submit an experiment or a realization. If the user doesn't specify an experiment or a realization with the options, the name of the experiment will be retrieved from `experiment.wrf4g` in case it exists in the current folder.

Options

<code>--version</code>	Show program's version number and exit
<code>-h, --help</code>	Show this help message and exit
<code>-n, --dry-run</code>	Perform a trial run with no changes made
<code>-e name, --exp=name</code>	Name of the experiment to submit
<code>-r name, --rea=name</code>	Name of the realization to submit
<code>-F FILE, --frea=FILE</code>	File containing the name of the realization to submit.
<code>-a, --rerun</code>	Force to run although the realization or experiment has finished
<code>-o, --run-just-one</code>	Run just the first chunk of the first realization. Only for testing purposes.
<code>-C N, --nchunk=N</code>	Run the next N chunks not finished of each realization
<code>-R N, --nrea=N</code>	Run the next N realizations not finished of the experiment

```

-p P, --priority=P    P is the priority of the experiment or realization that is
                      going to be launched (P is an integer between 0 and 20)
-v, --verbose        Verbose mode. It explains what is being done
-f, --force          Don't ask the user if he wants to submit an experiment
                      already submitted
-d TYPE_DEP, --tdep=TYPE_DEP
                      Specify dependencies between Chunks. afterok: The
                      chunk may be scheduled for execution only after jobs
                      jobid have finished without any error. afternotok: The
                      chunk may be scheduled for execution only after jobs
                      jobid have finished with errors. afterany: The chunk
                      may be scheduled for execution after jobs jobid have
                      finished, with or without errors.

```

wrf4g_status

Usage

```
wrf4g_status [--{exp Experiment|rea Realization}] [--long] [--ncharacters N] [--help]
```

Synopsis

Print the experiment or realization status.

Options

```

--version            Show program's version number and exit
-h, --help          Show this help message and exit
-e name, --exp=name Name of the experiment.
-r name, --rea=name Name of the realization
-l, --long          Show a detailed status.
-n NUMBER_OF_CHARACTERS, --ncharacters=NUMBER_OF_CHARACTERS
                    Print n characters of the Experiment's name
                    (default value is 20 characters)

```

Output field description

Summarized output: Shows realizations' status of each experiment (Default).

```

[user@mycomputer~]$ wrf4g_status
Experiment P    W    R    D    F
test      0    0    0    1    0
uc_phys   0    4    1    0    0
uc_single 1    0    0    0    0

```

- P: Prepared
- W: Waiting
- R: Running
- F: Failed

Long output: Shows a detailed realization status of every experiment

```

[user@mycomputer~]$ wrf4g_status --long
Realization      GW  Stat Chunks Comp.Res  WN      Run.Sta      ext  %
test             2   D   3/3   mycomputer sipcl8   Finished    0 100.00
uc_phys__phys1   3   R   1/3   mycomputer sipcl8   WRF         - 0.00
uc_phys__phys2   6   W   1/3   -           -         Submitted   - 0.00

```

uc_phys__phys3	9	W	1/3	-	-	Submitted	- 0.00
uc_phys__phys4	12	W	1/3	-	-	Submitted	- 0.00
uc_phys__phys5	15	W	1/3	-	-	Submitted	- 0.00
uc_single	-	P	0/3	-	-	Prepared	- 0.00

- Realization: Realization name.
- Status: It can take the following values: P(Prepared), S(Submitted), R(Running), F(Failed) and D(Done).
- Chunks [Chunk currently running/Total Chunks]: A realization is split into chunks. Each chunk is sent as a job.
- Computer resource: Computing Resource where the job is running. (It has to be one of the resources listed by wrf4g_resources)
- WN: Computing node where the job is running.
- Run.Sta: Job status in the WN (Downloading data, running ungrib, real, wrf, ...)
- ext: Exit Code. If exit code is different from 0, there has been an error. Error codes are explained in `$WRF4G_LOCATION/lib/bash/wrf4g_exit_codes.sh`
- % : percentage of simulation finished.

wrf4g_resources

Usage

```
wrf4g_resources [-h] [-c delay] [-nfx] [-m job_id] [host_id]
```

Synopsis

Print information about all the resources configured in `framework4g.conf` (default)

Options

```
-h          print this help
-c delay    refresh host information every delay seconds
-n          do not print the header
-f          full format
-x          xml format
-m job_id   print hosts matching the requirements of a given job
host_id     only monitor this host_id, printing also queue information
```

Output field description

```
[user@mycomputer~]$ wrf4g_resources
HID PRIO OS          ARCH  NODES(U/F/T) LRMS          HOSTNAME
0   1   GNU/Linux2.6.32 x86_6  0/1/1 FORK      mycomputer
1   1   GNU/Linux2.6.18 x86_6  88/0/88 PBS        asna_t1
2   1   GNU/Linux2.6.18 x86_6  168/0/168 PBS        asna_b2
```

FIELD INFORMATION

```
HID          host unique identification assigned by the GridWay system
PRIO         priority assigned to the host
OS           operating system
ARCH         architecture
NODES(U/F/T) number of slots: U = used by GridWay, F = free, T = total
LRMS         local resource management system, the jobmanager name
HOSTNAME     FQDN of this resource
```

QUEUE FIELD INFORMATION

```
QUEUENAME    name of this queue
SL(F/T)      slots: F = Free, T = Total
WALLT        queue wall time
CPUT         queue cpu time
COUNT       queue count number
```

MAXR	max. running jobs
MAXQ	max. queued jobs
STATUS	queue status
DISPATCH	queue dispatch type
PRIORITY	queue priority

wrf4g_priority

Usage

```
wrf4g_priority [--dry-run] [--(exp Experiment |rea Realization |frea File)] [--priority P] [--verbose] [--help]
```

Synopsis

Change the priority of an experiment/realization. The priority must be in range [0,20] and default value is 0. When a chunk gets a priority of 20, it becomes urgent. This chunk is dispatched as soon as possible, passing all the scheduling policies.

Options

--version	Show program's version number and exit
-h, --help	Show this help message and exit
-n, --dry-run	Perform a trial run with no changes made
-e name, --exp=name	Name of the experiment
-r name, --rea=name	Name of the realization
-F FILE, --frea=FILE	File containing the name of the realization to change the priority
-v, --verbose	Verbose mode. Explain what is being done
-p PRIORITY, --priority=PRIORITY	The priority must be in range [0,20].

wrf4g_kill

Usage

```
wrf4g_kill [--dry-run] [--(exp Experiment |rea Realization|frea File)] [--verbose] [--help]
```

Synopsis

wrf4g_kill command kill the jobs that belong to an experiment or realization. Additionally, the experiment and realization chunks, which have not done, will go back to "Prepared" status

Options

--version	Show program's version number and exit
-h, --help	Show this help message and exit
-n, --dry-run	Perform a trial run with no changes made
-e name, --exp=name	Name of the experiment
-r name, --rea=name	Name of the realization
-F FILE, --frea=FILE	File containing the name of the realization
-v, --verbose	Verbose mode. Explain what is being done.