Building BLSS

HADES

ARES



BORG

Structure

- ARES3 is actually the instrastructure package. Maybe will be renamed later
- On top of it there is an "extra/" subdirectory that contains extra modules (like foreground handling, HADES, BORG)
- the build system is using CMake (http://cmake.org)
- You will need CMake \geq 3.6
- New script available since October 25th 2016: build.sh

Building

- Go your ares build directory and run "./build.sh -h"
- You will get the help:

```
lavaux@reims:~/PROJECTS/ares$ ./build.sh -h
Ensure the current directory is ARES
This is the build helper. The arguments are the following:
 -cmake CMAKE BINARY
                       instead of searching for cmake in the path,
 use the indicated binary
                        build without openmp support (default with)
 -without-openmp
                        build with MPI support (default without)
 -with-mpi
 -c compiler COMPILER
                        specify the C compiler to use (default to cc)
 -cxx compiler COMPILER specify the CXX compiler to use (default to c++)
 -build dir DIRECTORY
                        specify the build directory (default to "build/" )
 -debua
                        build for full debugging
 -no-debug-log
                        remove all the debug output to increase speed on parallel
                        filesystem.
                        add timing instructions and report in the log files
 -perf
 -extra flags FLAGS
                        extra flags to pass to cmake
After the configuration, you can further tweak the configuration using comake (if
available on your system).
```

- Type "./build.sh" for default options, no MPI
- When done, go to the build directory and type "make"

Building... takes quite some time



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 - hades3: implements HADES and BORG algorithms. No powerspectrum inference or foreground cleaning here.
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 - borg_forward: possibility to replay borg mcmc to get more defailed informations.

Let's try ares !

• We run "./src/ares3"



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Error message: command line arguments

ARES command line

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 - first argument is generally:
 - INIT: to initialize a run from scratch
 - RESUME: to resume a previous interrupted run, requires restart files in that case
 - second argument is the configuration file
- First argument can also be:
 - SPECIAL_RESUME: this requires a monolithic restart files. This is used for MPI runs that requires a change in the number of nodes (e.g. you changed the supercomputer you are running the chain on). The chain is reseeded from the state of the random number generator.
 - RESUME_RESEED: to reseed a chain. Note! the seed in the configuration file is not used there.

HADES command line

• hades3 requires the same two arguments!

BORG_FORWARD

- borg_forward takes an mcmc element of a BORG chain, a configuration file identical to the one used to produce the chain (except for some options) and recompute the simulation
- Can produce:
 - density field, velocity field
 - particle set (positions and/or velocities)
- It is possible to change the following options:
 - number of time steps