

# **SHAREMED - WP3.1**

Developing a relocatable high-resolution short-term forecasting system for the marine environment

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# Document Control Sheet

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## Introduction

This report describes the high-resolution, relocatable modelling system for shortterm forecasts developed within Sharemed and provides the basic instructions to implement this numerical tool in different areas of the Mediterranean Sea. The relocatable system is based on the coupled hydrodynamic-biogeochemical model MITgcm-BFM, and it is initialized and forced by the Copernicus Marine Environment Monitoring Service (CMEMS-CMS) and other available operational products (e.g., satellite data, in-situ measurements). The package has been designed for expert users, with a suite of scripts and procedures for downloading initial, boundary and forcing data and for model set up, once the user has selected a specific geographical domain in the Mediterranean. The whole code is publicly available on specific GitHub repositories, while further details on model setup, implementation and provided in а dedicated online upgrades are wiki page (https://github.com/inogs/MITgcm\_BFM\_chain/wiki) in which the description of the modeling system can be updated according to its evolution and further developments.

The forecasting system was designed to be easily relocatable: it was firstly implemented in operational mode in the northern Adriatic study site, then it was further tested and applied in a cluster of transnational coastal areas in the Mediterranean Sea (i.e., Sicily Channel -WP4.5- and North Western Med and subbasins -WP5.1-). All the local implementations were customized by adapting the python and bash scripts developed for the pre- and post-processing phases of each operational system, which shapes up to be an effective regional downscaling of the Copernicus Marine Service (CMS) Analysis and Forecast products.

#### General description of the relocatable forecasting system

The relocatable forecasting system was developed by OGS and it is built upon the coupled hydrodynamic-biogeochemical model based on the MITgcm and BFM open-source codes (Figure 1), both of which keep up to the state-of-the-art in the scope of ocean modelling. It is nested into the CMEMS Analysis and Forecasting

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system, and it includes the whole operational chain for running the simulations on a daily basis.

The general documentation for the coupled (hydrodynamic-biogeochemical) model MITgcm-BFM can be found at the following links:

- hydrodynamic model: <u>http://mitgcm.org</u>
- biogeochemical model: <u>https://cmcc-foundation.github.io/</u> <u>www.bfm-community.eu/</u>
- coupler (MITgcm-BFM): <u>https://gmd.copernicus.org/articles/10/1423/2017/</u>

The "relocability" means that an experienced user should be able to implement and run the forecasting system on his/her own HPC platform without too many efforts. Indeed, the operational chain includes pre-processing, launch and post-processing steps, i.e.:

- selection, "cut" and interpolation of the initial (ICs) and open boundary conditions (OBCs) from CMEMS;
- integration of river contributions into the open boundary condition (OBCs: sections/"slices" derived from CMEMS);
- definition of surface and bottom biogeochemical fluxes;
- creation of model namelists containing all the runtime parameters;
- various scripts in order to manipulate meteorological forcings originating from different sources/services.

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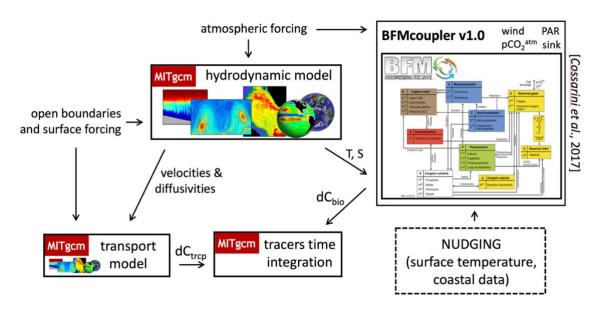


Figure 1 - Description of the MITgcm-BFM coupling used in SHAREMED.

The forecasting system was initially developed by OGS on the northern Adriatic Sea (at 1/128° resolution) and it is running operationally since several years, providing 72hour forecasts (<u>http://medeaf.ogs.it/adriatic</u>). The implementation of the relocatable system relies on code sharing through GitHub repositories and, during the project, different customized configurations were developed at each study site (Figure 2).

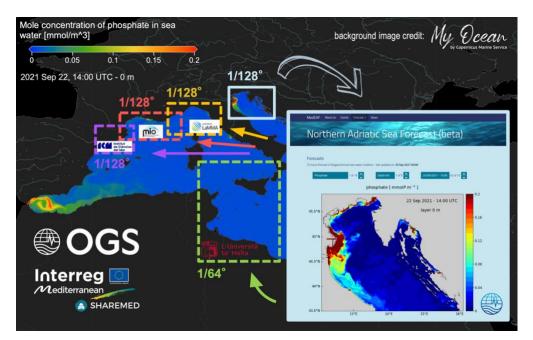


Figure 2 - The different configurations of the relocatable forecasting system in SHAREMED.

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# Instructions to compile and build the coupled MITgcm-BFM executable

In order to build the MITgcm-BFM executable, the user must follow the steps below.

Define your ModelBuild directory and clone from the GitHub repository the tool for building the model:

DIR=ModelBuild

git clone git@github.com:inogs/MITgcmBFM-build.git \$DIR

Download the source code of BFM, of the **coupler** and of MITgcm:

cd \$DIR git checkout main ./downloader MITgcm bfm.sh

Edit builder\_MITgcm\_bfm.sh, to set debug mode or not, compiler, modules, etc...

For example:

MIT\_COMPILER=intel (gfortran)
export MODULEFILE=\$PWD/compilers/machine\_modules/g100.intel
(fluxus.gfortran, yourmachine.yourcompiler)

Put your HPC-platform-specific *INC\_FILE* for BFM in:

```
$BFMDIR/compilers/$INC_FILE (e.g., x86_64.LINUX.gfortran.inc)
```

Compile and build BFM:

./builder\_MITgcm\_bfm.sh -o bfm

Decide the number of cores and the domain decomposition of your model run, configuring your SIZE.h (e.g., choosing among the various options of the northern Adriatic benchmark experiment). E.g., for 95 cores:

cp presets/NORTH\_ADRIATIC/SIZE.h\_095p presets/NORTH\_ADRIATIC/SIZE.h

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Configure/customize the setup:

./configure\_MITgcm\_bfm.sh

If needed, update the MYCODE directory with your own modified routines (different OBCs, meteorological forcing files, etc...).

Put your HPC-platform-specific *INC\_FILE* for MITgcm in:

\$PWD/compilers/\$INC\_FILE (e.g., x86\_64.LINUX.gfortran.inc).

Compile and build MITgcm: use the same builder used for compiling BFM (i.e., builder MITgcm bfm.sh)

./builder\_MITgcm\_bfm.sh -o MITgcm

The executable (mitgcmuv) is in \$PWD/MITGCM\_BUILD.

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# Instructions to install the operational chain

In this section we describe the main dependencies in the code pipeline and we provide a primer for the production workflow, from the download of the code to the installation of the chain.

Clone the operational chain from the GitHub repository.

```
git clone git@github.com:inogs/MITgcm_BFM_chain.git
```

The installation of the chain relies on the script called mit\_setup\_chain.ksh. In order to run it, the user must set properly the mit\_profile\_machine.src\_inc, where *machine* is the name of the host, as set up in the MIT\_HOSTNAME variable.

With this configuration script, the user must set some variables and run some commands, depending on the *machine* used (i.e., the available computing facility). The structure of this file is explained below, so that the user/installer can configure it without difficulties.

#### Variables

- MIT\_NETCDF\_PATH: is the path of folder containing the NetCDF library on the machine (one can locate easily, for instance, loading the proper module, and spotting the folder in the LD\_LIBRARY\_PATH\_env variable);
- MIT\_DATE\_PATH: the path of date command (almost certainly will be /usr/bin/date);
- MIT\_NCKS\_PATH: the path of ncks on the machine;
- MIT\_NCFTPGET\_PATH: the path of ncftpget on the machine;
- MITGCMUV\_PATH: the path of the executable produced during the compilation of the model (see section above). This will be located in the MITGCM\_BUILD subfolder of the model folder;
- MOTUCLIENT\_UNPACK\_DIR: a path where the tool motuclient will be unpacked (usually set to \$MIT\_BINDIR/motuclient);

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 PYTHON3\_PATH: the path of a valid python3 interpreter. It is used just to create virtual environment, the interpreter included in this environment will be used.

#### Functions

- mit\_mkdir <dir>: creates the directory <dir> with all its parents and checks for the success;
- symlink\_if\_not\_present <path> <link>: symbolic links the <path> to
   <link> if <link> is not already present. If it is so, the function does nothing. If
   <path> it not present, it returns an error state;
- clone\_or\_update <git\_link> <path> <branch>: if <path> is not present, it clones the <git\_link> repo to <path> checking out branch <branch> (notice: if <path> is present it does nothing, the 'update' part has still to be implemented);
- download\_and\_unpack <url> <path> <filename> <extraction\_dir>: downloads file at <url> placing it to <path>/<filename>. Then it creates the directory <extraction\_dir> and extracts the contents to it. If some step fail, the function returns an error status;
- <pyhton3 path> <requirements>: set python environment uses <python3\_path> to create an environment according to the variable \$MIT VENV 1 (set in mit profile.src inc). Then it activates the environment, and uses pip to install all the requirements specified in <requirements> (the python3 path is set for convenience in the variable PYTHON3 PATH. The requirements are listed in the file \$MIT HOME/requirements.txt).

Note that all the functions above are idempotent up to an upgrade coming from online repositories. This means that a script containing these function can be safely executed n times without spoiling the setup.

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#### Commands

Although the exact command depends on the specific machine (i.e., modules to load, particular env variable to set) the structure of the file is the following:

# env variable section mit batch prex "export LD LIBRARY PATH=\${MIT LIBDIR}:\${MIT NETCDF PATH}:\${LD LIBRARY PATH}" mit batch prex "mit mkdir \$MIT BINDIR" # symlink section mit batch prex "symlink if not present \$MIT DATE PATH \$MIT BINDIR/date" mit\_batch\_prex "symlink if not present \$MIT NCKS PATH \$MIT BINDIR/ncks" mit batch prex "symlink if not present \$MIT NCFTPGET PATH \$MIT BINDIR/ncftpget" mit batch prex "symlink if not present \$MITGCMUV PATH \$MIT BINDIR/mitgcmuv" # git section mit batch prex "clone or update \$MIT BITSEA REPO \$MIT BITSEA \$MIT BITSEA BRANCH" mit batch prex "clone or update \$MIT POSTPROC REPO \$MIT POSTPROCDIR \$MIT POSTPROC BRANCH" mit\_batch\_prex "clone\_or\_update \$MIT\_BC\_IC\_REPO \$MIT BC IC FROM OGSTM DIR \$MIT BC IC BRANCH" # motuclient section mit batch prex "download and unpack \$MIT MOTUCLIENT URL \$MIT BINDIR motuclient.tar.gz \$MOTUCLIENT UNPACK DIR" mit batch prex "symlink if not present \$MOTUCLIENT UNPACK DIR/src/python \$MIT BINDIR/motuclient-python" # python3 section mit batch prex "set python environment \$PYTHON3 PATH \$MIT HOME/requirements.txt" # module section # ad libitum 

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After setting mit\_profile\_\_machine.src\_inc, create a script (e.g., chain\_env.sh) with the following content (export and alias).

NOTE: given the example below, you could put it in /yourpath/.

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```
# MIT_HOME is your MITgcm_BFM_chain directory, the root of the git
repository
export MIT_HOME=/yourpath/MITgcm_BFM_chain
# your machine name
export MIT_HOSTNAME=fluxus
# your work-launch (e.g., "scratch" area) directory in a parallel
filesystem
export MIT_WORKDIR=/yourpath/WORK
export MIT_VERSION_NUMBER=1
export MIT_STAGE=devel
alias mitcd="cd $MIT_HOME ; pwd"
```

Then do:

source chain\_env.sh
cd \$MIT\_HOME/bin/src

For a first check that everything works, compile and install all the scripts:

make make install

```
Add a line referring to your mit_profile__<yourmachinename>.inc in the SCRIPTS list of the Makefile ($MIT_HOME/bin/src/Makefile) and perform another (updated) installation:
```

```
make
make install
```

Once the job scheduler options are set (consult your IT manager for these details), you can launch your model run (with bulletin date: 19/10/2021, as an example).

```
cd $MIT_HOME/bin
./mit_start.ksh --pass --job-multiple --force --try-resume --without-
kill --rundate 20211019
```

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mit\_start.ksh generates the log directory, where the single jobs and/or the single executables can be submitted or launched one by one for each of the three "stages" of a forecast run: preprocessing, model run, postprocessing.

To conclude, the compilation phase will produce an executable (MITgcm-BFM) called mitgcmuv. For running the model:

#### • option 1 - full operational chain

mit\_start.ksh should run the whole model chain (if there are no setup/runtime errors). The job produced by mit\_start.ksh (which needs the machine-specific scheduler options) will look for the MITgcm-BFM executable located in \$MIT\_HOME/HOST/\$MIT\_HOSTNAME/bin/ (together with the other executables: date, ncks...).

The executable can be renamed "mitgcmuv\_<number\_of\_cores>", where the number of cores is specified in mit mpi intelmpi.src inc.

#### • option 2 - test simulation

copy the mitgcmuv executable in your run directory and run the simulation using a customized job for your local scheduler.

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# **Model output**

Model output is dumped in the native MITgcm format, in png images or in NetCDF products, organized in the same way as the CMS modelling products for the Mediterranean Sea. Figure 3 shows an example of model output published on a dedicated web page (https://medeaf.oqs.it/adriatic).

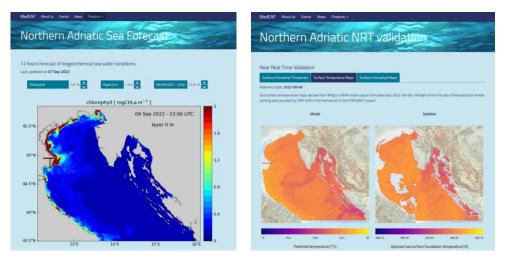


Figure 3 - Examples of model output published on dedicated web pages: chlorophyll forecast fields (left) and SST comparisons with satellite data (right).

NetCDF products can be also stored on specific web services, like the THREDDS server shown in Figure 4.

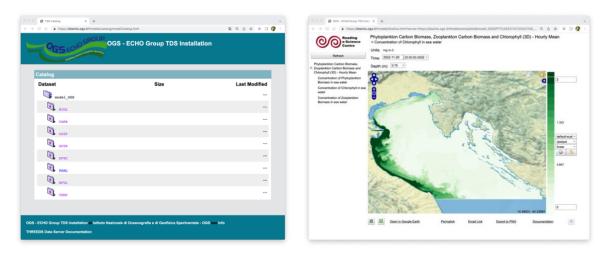


Figure 4 - Example of model output stored on a dedicated THREDDS server and visible through the "preview" tools available.

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# Conclusions

In the framework of SHAREMED, we developed a short-term forecasting system for ocean hydrodynamics and biogeochemistry in a cluster of transnational coastal areas in the Mediterranean Sea.

Model development has been carried out in the northern Adriatic Sea (AS). The system was subsequently tested in the Sicily Channel (SC), and finally transferred to the northwestern Mediterranean Sea (NWM), where the domain has been further subdivided into three transnational sub-basins: the Ligurian Sea, the Gulf of Lion and the Catalan Sea.

At present, the system is running in operational or pre-operational mode on a daily basis in all the study sites, providing 2- to 3-day forecasts of the main physical and biogeochemical variables (e.g., currents, thermohaline properties, nutrients, chlorophyll, dissolved oxygen). The results are published on dedicated SHAREMED portals and site-specific webpages, while the source code of the modelling system and of the operational chain described in the present report is available on public GitHub repositories.

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