

The SLAM+(ERSEM+POLCOMS) model coupling

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1 Introduction

This document contains a short description and user guide of the SLAM (Sandeel Larval Advection model) model, coupled offline with the POLCOMS+ERSEM model provided by Plymouth Marine Laboratory (PML) in the MEECE project. The SLAM model underlying this implementation is published mainly in Can. J. Fish. Aquat. Sci. 65: 1498-1511 (2008) by Christensen *et al.*

The SLAM in the present implementation is set up within the IBMlib (Individual based model library) framework maintained by DTU Aqua. The present code package, however, contains only a self contained subset of the IBMlib framework. The basic idea of IBMlib is sketched and is very well in line with the goal of MEECE to ease the traditionally cumbersome process of coupling biological models to arbitrary biogeochemical models. The present code package is a preliminary demonstration coupling of the SLAM and POLCOMS+ERSEM models; the production runs in MEECE WP 3+4 will be performed with an upgraded and extended version of the present code package, and interested MEECE participants are encouraged to obtain latest version before embarking on simulations on their own with the SPAM model setup.

OFFLINE RUN

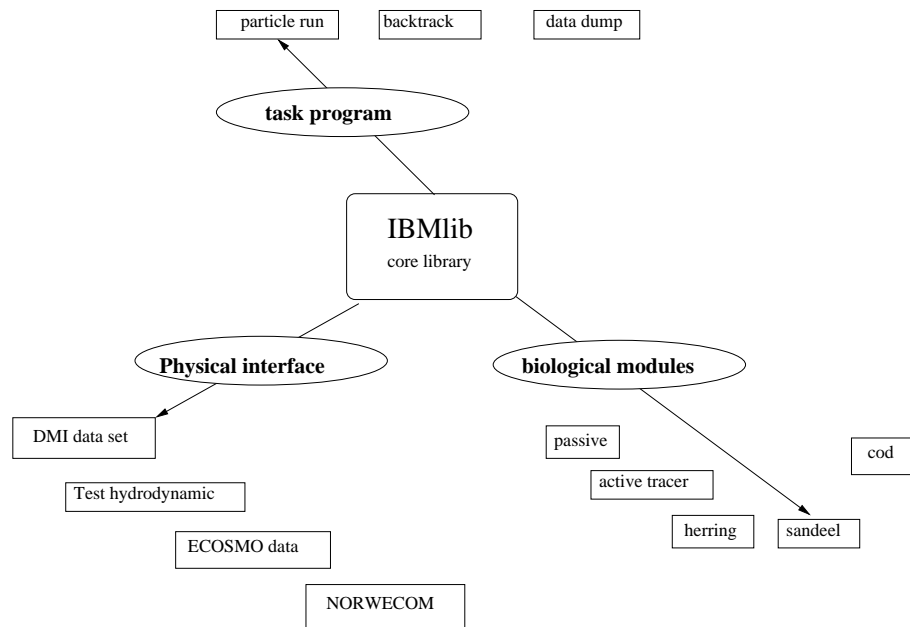


Figure 1: Concept diagram of the IBMlib framework.

2 Functionality

SLAM+(POLCOMS+ERSEM) setup allows you to perform flexible individual-based modelling of early life stages within the space-time window covered by a supplied POLCOMS+ERSEM data set. Sandeel larvae and/or eggs can be released according to any space-time pattern specified in the input file. The model describes growth of egg/larvae in relation to local physical conditions obtained from the supplied POLCOMS+ERSEM data set. The model can easily be extended by alternative biological processes, e.g. alternative active larval behavior.

3 License and intellectual property rights

At some point it is planned to make the IBMlib public available, but this awaits the clarification of some of the license issues of integrated external packages as well as code hosting and distribution issues; until then, the code is available to MEECE participants within the MEECE project under provisions specified in the MEECE consortium agreement and the intellectual property rights belong to DTU Aqua.

4 Installation

4.1 Portability

The code is written in strict Fortran 90 and does not invoke operating system specific features nor does the code rely on non-standard compiler specific services. Therefore the code (possibly with some superficial preparations) should compile and execute on any platform; however, the code has only been applied on Linux platforms until now.

4.2 Requirements

In order to build the SLAM executable, the following resources must be available:

- A Fortran 90+ compiler and C++ compiler. The compiler and link flags in the Makefile applies to Intels *ifort* compiler, however the transcription of of these flags should be straightforward by consulting the documentation of alternative compilers; most flags are standardized.
- NetCDF. The code assumes the Fortran 90 implementation (the fortran-only variant) is installed (at compilation time by use association of of the netcdf module - usually netcdf.mod, somewhere in the include path - and at link time by linking to libnetcdf.a and libnetcdf.a). The NetCDF is freely available at www.unidata.ucar.edu/software/netcdf. It is recommended to use NetCDF 4+ (which requires HDF5 and zlib installed), but NetCDF 3.6+ should also work.
- GNU make. The current makefile is written for GNU make, however, with minor adaptation, it should also work with other implementations of make.
- tar and gzip. The code is packed with tar and compressed with gzip.

4.3 Code installation under Linux

The code package comes as *SLAM4MEECE.tgz*. Put it in a suitable location and type the following commands

```
tar xvfz SLAM4MEECE.tgz
make tracker
```

This builds the SLAM executable **tracker** which you - on UNIX like systems - may place somewhere in the executable paths for convenience.

5 Running the code

On UNIX like systems, the SLAM model is invoked from the command line simply as

```
tracker <inputfile> [ > <outputfile>]
```

(assuming tracker is somewhere in the executable paths). Here “<inputfile>” means the filename of an ASCII text file containing parameters for the simulation. The structure of inputfile” is outlined below. Optionally, an output file “<outputfile>” can be specified to capture the simulation output; if an output file is not specified, it is written to standard output.

6 Input files

6.1 Simulation parameters

The SLAM model code comes along with an input parser that reads input with minimalistic markup in the form

TAG = VALUE|SET OF VALUES

from the ASCII text input file specified on the command line above. *TAG* is a name of something (like starting time or a parameter) and *VALUE* is the value SLAM should use for *TAG*. *SETOFVALUES* means that this can also be a simple list (separated by spaces) of values. A small pieces of an input file could look like this

```
! --- Simulation parameters for ...
  y = 3445      ! an optional trailing comment
  x = 45.
  zvector = 1 56 99
  aname   = whatever
```

A complete example of an input file is given below. Line order and white spaces does not matter. The following rules also apply for the markup:

- Comments: everything after (and including) “!” is skipped
- Empty lines: just ignored
- Malformed lines: just ignored (e.g. if you forgot “=”)
- Everything after “=” is considered part of the value for tag
- Required tag (or value) is missing. This will generate a runtime error, SLAM will stop, unless a default applied to the tag.

The input parser is given in the file *input_parser.f* in the code package
 An example of a minimal input file (i.e. the mandatory input tags) is
 given below (provided in file "input/test_input.txt" in the code package
SLAM4MEECE.tgz)

```
!-----
!           Main simulation control file
!-----

start_time  = 2005 03 01 0    ! year  month  day  second_of_day
end_time    = 2005 03 18 0    ! year  month  day  second_of_day

hydroDBpath = hydroDB        ! file path to POLCOMS+ERSEM data set
grid_desc   = grid_desc.txt  ! sub grid descriptor

advec_intg_method = euler    ! advection scheme: euler/rk27rk4
particle_time_step = 1800    ! in seconds for time integration of motion

! ----- biology spatial control -----
! r(1:4) start:  year month day sec_of_day
! r(5:8) end:    year month day sec_of_day
! r(9:11)        lon_min lat_min z_min  (z=0 -> surface)
! r(12:14)       lon_max lat_max z_max  (z=1 -> bottom)
! r(15)          max_number_of_tracers
! r(16:)         other input item to particle state

emitbox = 2005 03 02 0      2005 03 02 3600  2 54 1   3 55 1   100 e
emitbox = 2005 03 02 3600  2005 03 02 7200  4 54 1   5 56 1   100 1 9.66

! ----- biology growth -----
! parameters corresponds to Can. J. Fish. Aquat. Sci. 65: 1498-1511 (2008)

egg_hatch_begin = 78.484  0.10984 ! days since fertilization (A,k0 > 0)
egg_hatch_mid   = 108.75   0.12488 ! days since fertilization (B,k1 > 0)
egg_hatch_end   = 176.14   0.11358 ! days since fertilization (C,k2 > 0)

larvae_temp_coeff    = -1.725  0.136142  0.00    ! mm/day/Celcius^n
larvae_temp_func     = 1       ! 0: func=polynomial  1: func=exp(polynomial)
```

```

larvae_hatch_len      = 7.7253      ! mm
larvae_length_expo    = 0.315544 ! beta (small larvae growth exponent)
larvae_metamorph_len  = 40.0        ! mm

```

- start_time. The beginning of the simulation, specified as (year month day second_of_day)
- end_time. The end time of the simulation, specified as (year month day second_of_day)
- hydroDBpath. File path to POLCOMS+ERSEM data set
- grid_desc. The file name of the sub grid descriptor (see below).
- advec_intg_method. The integration algorithm for time forward integration of advected sandeel larvae (options Euler forward, Runge-Kutta 2 or 4)
- particle_time_step. Nominal time step for the integration algorithm in seconds.
- emitbox. Gives a window in space and time, where sandeel larvae/eggs are released. This may be specified as many emitbox entries as desired, in this way they may act in parallel and quite complex release patterns can be set up. The first 4 integers are (year month day second_of_day) where the release begins (of that box); the next 4 integers are (year month day second_of_day) where the release stops (of that box). The next six numbers specify a spatial box (in latitude and longitude) where sandeel larvae/eggs are released. Dry (land-locked) sectors of the spatial box are omitted when releasing larvae/eggs. The first three numbers are the spatial lower SW corner of the release box given as (longitude,latitude,vertical position), the next three numbers are the spatial upper NE corner of the release box given as (longitude,latitude,vertical position). The vertical position can be specified as absolute depth (counted negative below the water surface) or as relative depth $z: 0 < z < 1$. $z = 0$ corresponds to the sea surface, $z = 1$ corresponds to the sea bed. Biological particles are released uniformly in time and space within the space-time window specified, so that the total number of particles released adds up to the integer given as number 15.

All other parameters following number 15 in emit box are passed to the biological module. There can be an "e", which means sandeel eggs are released by this emitbox; the can be an "l" which means sandeel larvae are released by this emitbox - in the latter case, a number giving the initial length (in mm) of the released sandeel larvae should be provided.

- stochastic egg growth parameters (egg_hatch_begin, egg_hatch_mid, egg_hatch_end), see Can. J. Fish. Aquat. Sci. 65: 1498-1511 (2008) for details.
- larval growth parameters (larvae_temp_coeff, larvae_temp_func, larvae_hatch_len, larvae_leng, larvae_metamorph_len) see Can. J. Fish. Aquat. Sci. 65: 1498-1511 (2008) for details.

6.2 Sub grid descriptors

The SLAM model allows to operate on a sub grid of the POLCOMS+ERSEM spatial domain in MEECE. The sub grid descriptors allows to cut out a subgrid, when the relevant biological habitat is smaller than the domain of the hydrodynamic model, to speed up calculations and run the simulation on a laptop. The sub grid descriptor is specified in same format as the simulation parameters above.

```
! -----
! ----- full grid definition -----
! -----

lambda_start_fullgrid = -19.833333333
lambda_end_fullgrid   = 13.000000000
dlambda_fullgrid      = 0.166666667

phi_start_fullgrid    = 40.111111111
phi_end_fullgrid      = 64.888888889
dphi_fullgrid         = 0.111111111

nz                    = 40

! -----
! ----- subgrid definition -----
! -----
```



```

lambda_start_subgrid = -3.0
lambda_end_subgrid   = 10.0
phi_start_subgrid    = 52.0
phi_end_subgrid      = 59.0

```

- (lambda_start_fullgrid, lambda_end_fullgrid) is the longitude range of the POLCOMS+ERSEM spatial domain in degrees East.
- lambda_fullgrid is the longitude grid spacing of the POLCOMS+ERSEM spatial domain.
- (phi_start_fullgrid, phi_end_fullgrid) is the latitude range of the POLCOMS+ERSEM spatial domain in degrees North.
- dphi_fullgrid is the latitude grid spacing of the POLCOMS+ERSEM spatial domain.
- (lambda_start_subgrid, lambda_end_subgrid) is the longitude range of the SLAM spatial sub domain in degrees East.
- (phi_start_subgrid, phi_end_subgrid) is the latitude range of the SLAM spatial sub domain in degrees North.

The sub grid will be coherent to the full POLCOMS+ERSEM grid (i.e. overlapping grid nodes) (provided in file "input/grid_desc.txt" in the code package *SLAM4MEECE.tgz*)

7 Output

The simulation writes logging information to standard output, and at the end of the simulation, the task example writes the state of the sandeel egg/larvae ensemble to standard output. It is normally highly specialized which output is desired, and therefore, to save disk space, it is most efficient to insert write statements in "main_program.f" selecting exactly the information needed. The IBMlib features several writing subroutines.

8 Programming

The IBMlib code is written in strict Fortran 90 and the coding style of IBMlib adheres to modern object-oriented programming principles to the extent they are supported by Fortran. IBMlib API will be provided later in MEECE. The most important interfaces in IBMlib are

- The physical interface: provided by `physical_fields.f`. Here the biological modules can access the local physical/biological environment.
- The particle state interface: provided by `physical_state.f`. This interface lets IBMlib update the biological ensembles. All biological mechanisms are behind the particle state interface
- The task interface: this is interface provided by IBMlib to the main program.