

The CWR Estuary and Lake Computer Model ELCOM

User Guide

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ELCOM User-Guide

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Estuary and Lake Computer Model

ELCOM: User Manual

These notes are designed to aid an uninitiated user in performing an ELCOM simulation. This document gives a basic description of ELCOM's operation and instructions on how to create the input files required for a standard simulation.

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1. Introduction

1.1 The ELCOM model

ELCOM (Estuary and Lake Computer Model) is a numerical modeling tool that applies hydrodynamic and thermodynamic models to simulate the temporal behavior of stratified water bodies with environmental forcing. The hydrodynamic simulation method solves the unsteady, viscous Navier-Stokes equations for incompressible flow using the hydrostatic assumption for pressure. Modeled and simulated processes include baroclinic and barotropic responses, rotational effects, tidal forcing, wind stresses, surface thermal forcing, inflows, outflows, and transport of salt, heat and passive scalars. Through coupling with the CAEDYM (Computational Aquatic Ecosystem DYNamics Model) water quality module, ELCOM can be used to simulate three-dimensional transport and interactions of flow physics, biology and chemistry. The hydrodynamic algorithms in ELCOM are based on the Euler-Lagrange method for advection of momentum with a conjugate-gradient solution for the free-surface height (Casulli and Cheng, 1992). Passive and active scalars (i.e. tracers, salinity and temperature) are advected using a conservative ULTIMATE QUICKEST discretization (Leonard 1991).

1.2 Implementation language

ELCOM is implemented in Fortran 90 (with F95 extensions) so that three-dimensional space can be mapped into a single vector for fast operation using array-processing techniques. Only the computational cells that contain water are represented in the single vector so that memory usage is minimized. This allows Fortran 90 compiler parallelization and vectorization without platform-specific modification of the code. A future extension of ELCOM will include dynamic pressure effects to account for nonlinear dynamics of internal waves that may be lost due to the hydrostatic approximation.

1.3 Outline of a simulation run

ELCOM is designed to be used in conjunction with a graphical user interface (GUI) called CWR-Modeller, which supports post-processing and analysis of the data. However, ELCOM is not executed by CWR-Modeller, but is initiated with simple command-line execution statements.

Setting up and running an ELCOM simulation requires the following steps:

1. compile the PRE executable (see [section 2.2](#)),
2. prepare `bathymetry.dat` file that provides topography and grid information (see [section 2.3](#)),
3. prepare `bc.dat` file that designates sets of cells for applying boundary conditions (see [section 2.4](#)),
4. run the pre-processor to produce `sparsedata.unf` and `usedata.unf` files [section 2.5](#),
5. compile the ELCOM executable (see [section 3.3](#)),
6. configure the simulation `run_elcom` file (see [section 3.6](#))
7. specify output in a datablock file (`*.db`) for designating data output operations (see [section 3.7.2](#)),
8. prepare temporal boundary condition files (`*.dat`) for data applied to the boundary condition cells (see [section 3.7.3](#)),
9. prepare any user-defined subroutines for initial conditions (see [section 3.8](#))
10. run the executable, directing the screen output to a file(see [section 3.3](#))
11. post-process data

ELCOM uses and creates a large number of files. The following directory structure is suggested to help streamline ELCOM's use:

/ELCOM/source/PRE/	contains pre-processor source code
/ELCOM/source/ELCOM/	contains ELCOM source code
/ELCOM/BIN/	contains ELCOM, PRE, db2nc executables, and must be in the UNIX search path
/ELCOM/bathymetry/	contains bathymetry.dat and bc.dat files for various user simulations
/ELCOM/runs/RUN_A/	contains run_elcom file for a simulation, as well as ELCOM status output files
/ELCOM/runs/RUN_A/infiles/	contains user input files for a simulation (ie. sparsedata.unf, usedata.unf, temporal boundary condition files, and datablock.db)
/ELCOM/runs/RUN_A/unffiles/	contains ELCOM * .unf output files to be converted to * .nc files using db2nc
/ELCOM/runs/RUN_A/txtfile/	contains output text files from ELCOM

The /ELCOM/BIN/ directory must be in the UNIX search as defined in your user .cshrc file. For example, the following line should be included in your .cshrc file:

```
setenv PATH "${PATH}:${HOME}/ELCOM/bin"
```

1.4 Conventions and definitions

In this document, a number of conventions will be used. For clarity, the computational domain will generally be referred to as a "lake". However, the simulation method is designed for applicability to estuaries and coastal ocean environments

x,y,z space

Full Cartesian space which covers all of the bathymetry.

i,j,k space

Discrete integer space of computational cells which covers all of the bathymetry.

land

A computational cell that cannot contain water.

open

A cell used for an open boundary condition.

interior

A cell that may contain water.

x_rows

Number of grid cells in the x direction.

y_columns

Number of grid cells in the y direction.

z_layers

Number of grid cells in the (vertical) z direction.

land_value

Number used in the bathymetry file to represent the land cells.

open_value

Number used in the bathymetry file to represent the open boundary cells.

dx

Uniform grid spacing (delta) in the x direction (in metres).

dy

Uniform grid spacing (delta) in the y direction (in metres).

dz

Grid spacing (delta) in the z direction, may be uniform or non-uniform (in metres).

[Continue on to Pre-processor](#)

2. Pre-Processor

2.1 Introduction

The ELCOM pre-processor is used to convert the user's three-dimensional data of lake bathymetry, boundary forcing sections and grid structure into one-dimensional vectors used by ELCOM and the three-dimensional data format required for CWR-Modeller. Configuration for running the pre-processor comes from a `run_pre` file. The `run_pre` file is a text file that contains the names and location of user-supplied input files and pre-processor output files. The pre-processor requires two user-supplied text files: `bathymetry.dat` and `bc.dat`. The former describes the bathymetry of the lake and grid configuration. The latter defines sets of boundary cells. The pre-processor output consists of two Fortran unformatted (binary) files: `sparsedata.unf` and `usedata.unf`. The former contains information on three-dimensional space and is used by CWR-Modeller and ELCOM. The latter is used by ELCOM, and contains the one-dimensional spatial relationships required for the simulation. These files may be renamed, but the user must keep track of which files are "sparsedata" files and which files are "usedata" files.

2.2 Compiling the pre-processor

Before compiling the pre-processor source code directory should contain the following files:

- `pre_main.f90`
- `pre_globals.f90`
- `pre_process.f90`
- `pre_bathymetry.f90`
- `pre_common.f90`
- `makefile`
- `run_pre`

The `*.f90` files should not need modification by the user. However, the `makefile` file will need modification by the user to indicate the path to the bin directory. The pre-processor is compiled using a make utility and the file `makefile`. Typical UNIX syntax for compiling is simply:

```
make
```

This produces an executable file (`PRE`) which is run to generate the pre-processor output binary files. The `PRE` executable file is then installed using the UNIX syntax:

```
make install
```

The pre-processor source code directory can then be cleaned up using the following UNIX command:

```
make clean
```

2.3 Setting up the bathymetry file (bathymetry.dat)

An up-to-date copy of `bathymetry.dat` can be downloaded from [Appendix A](#).

2.3.1 Introduction

The bathymetry file is designed to allow the user to input the geometry of the simulation domain [at the resolution that the simulation will be conducted!](#). The format of the bathymetry file is rigid with regards to the amount of data provided and the order in which it is listed. Each item in the bathymetry file must be present even if it is not applicable to the present simulation. Comment lines may be added into the file at any point before the actual bathymetry values. Each line of the `bathymetry.dat` file must have at least two "items" separated by a space or a tab. An "item" may be an integer, a real number or a character string. Comment

lines must also have at least two items, the first of which is an exclamation point (the Fortran 90 comment symbol). The order in which the data is presented is fixed so that CWR-Modeller can process the [bathymetry.dat file](#)². A common crash of the ELCOM preprocessor produces a Fortran error message regarding insufficient input data. This is typically because (1) one or more non-optional items were omitted from the input file or (2) a data line or comment line has less than two items. The file `bathymetry.dat` is organized as a series of text lines of keywords, data and comments that follow one of the formats shown below.

keyword	data
data	keyword
! -	comment
keyword	keyword
data(j=1,y_columns)	

The keywords are ELCOM-defined words that indicate the type of user-provided data presented: (1) before the keyword, (2) after the keyword, or (3) on succeeding lines of the file. Exclamation points are used to indicate lines with comments. For proper functioning with CWR-Modeller, the `bathymetry.dat` file must be organized with the arguments presented in a fixed order (however any number of comment lines may be inserted). Additional comments may be provided after the keywords to help the user remember the keyword purpose (the code only reads the first two items on a line). Detailed information on the keywords and the format of the file are presented in the following sections.

2.3.2 Physical and computational space coordinates

In the literature, bathymetry is generally presented as the vertical distance from some horizontal baseline (often a mean water height) and may be either positive downward or positive upward. The convention used in ELCOM is an (x,y,z) space that follows the right-hand rule and has z positive in the upward direction ([Figure 1](#)). The user is free to choose any suitable $z=0$ baseline for the bathymetry as long as positive z upward is maintained as the convention for measurement. The pre-processor will adjust the user-defined $z=0$ baseline to an internal code baseline that is suitable for the ELCOM simulation. For the most part, the user does not need to know or consider this internal baseline adjustment. However, some debugging/monitoring output is presented in terms of the code coordinate system rather than the user coordinate system, so care must be taken in interpreting debugging output that is intended for use by code developers. The position of physical space origin ($x=0, y=0$) is set implicitly by the user through the layout of the bathymetry data in the file `bathymetry.dat`. This will be covered in more detail below. Physical space x and y values may not be negative (i.e., the x,y origin must be in a corner of the domain).

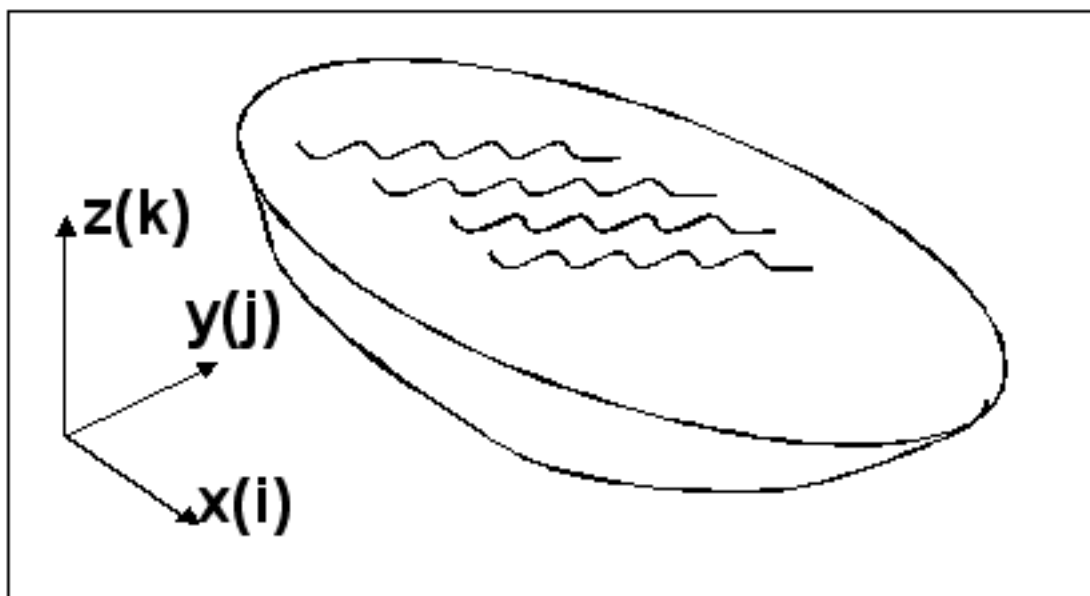


Figure 1. Schematic of coordinate system

ELCOM performs its simulation on computational cells of a three-dimensional Cartesian mesh. The user needs to understand the layout of the mesh so that boundary condition cells can be properly identified. In the present version of the code, the mesh has uniform spacing in each of the horizontal dimensions (x and

y) and allows non-uniform spacing in the vertical (z) direction. [Figure 2](#) shows a plan view (x-y plane) of a lake with a discrete (uniform) Cartesian mesh overlay. The computational domain is divided into "cells", whose centers are represented in [Figure 2](#) by solid dots. Dashed lines represent the faces between the cells. Each cell center has an (i,j,k) coordinate that designates its location in discrete computational space. The i,j,k values are the integer indices of a three-dimensional array. The right-hand rule is used, with (i,j,k) being non-zero, positive integer values; $k=1$ is associated with the lower-most cell layer in the domain. The (i,j) pairs of a horizontal plane represent a matrix, so that by conventional mathematical notation the (1,1) position is always in the upper left corner. The 'i' index is the row index and increases down the page. The 'j' index is the column index and increases across the page. The (i,j) indices for corner cells are shown in [Figure 2](#). For convenience and consistency, the origin (0,0) of the x,y (physical space) coordinates is taken to be at the (1,1) location in the computational space i,j indices.

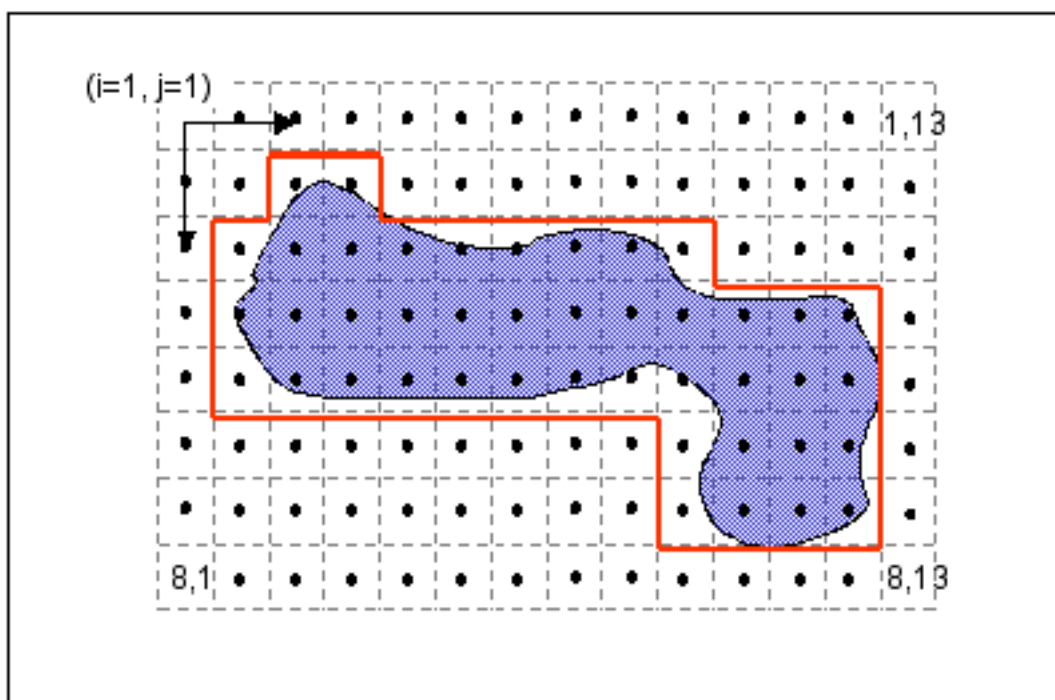


Figure 2. Plan view of Cartesian mesh

Using the upper left corner as the origin may initially be confusing to users familiar with mirror-image storage (used by many graphics programs) where the *lower* left corner is stored in position $i=1, j=1$. The problem with mirror-image storage is that a simple printout of bathymetry data always appears as an upside-down mirror image of the real world when viewed with a text processor or spreadsheet. The present convention gives the user a "What You See Is What You Get" view of the bathymetry simply by opening the file with a word processor or importing into a spreadsheet. This convention has been found to be intuitive for users who are not graphics experts. However, caution needs to be taken when converting raw bathymetry data from graphics programs that use a mirror-image standard. Likewise the direction of the positive u (x direction) velocity is confusing as it is positive down the page.

[Figure 3](#) shows the *physical* space outline of the lake as the border of the shaded area. The discrete outline of a lake is shown with a heavy solid line. In developing a `bathymetry.dat` file for an ELCOM simulation, the user is required to provide bathymetry values (i.e. vertical distance measured from the user's $z=0$ baseline) for each of the cells within the discrete boundary of the lake. The region outside the lake boundary is defined as the "land" region. In [Figure 3](#), the cell center dots have been replaced with numbers that represent the distance from the user's $z=0$ baseline to the bottom of the lake. In this case, all of the "land" regions have been given a value of 99. Note that the boundary of the bathymetry data set shown in [Figure 3](#) has land values around the entire boundary. In general, it is required that either a land cell or open boundary cell is required around the outside of the bathymetry data. For the purposes of the `bathymetry.dat` file, the user must provide bathymetry values for *all* cells in a rectangular array, including those cells that only contain land. ELCOM will only reserve computational memory for (1) the cells which may contain water, and (2) the first layer of land cells surrounding the water cells.

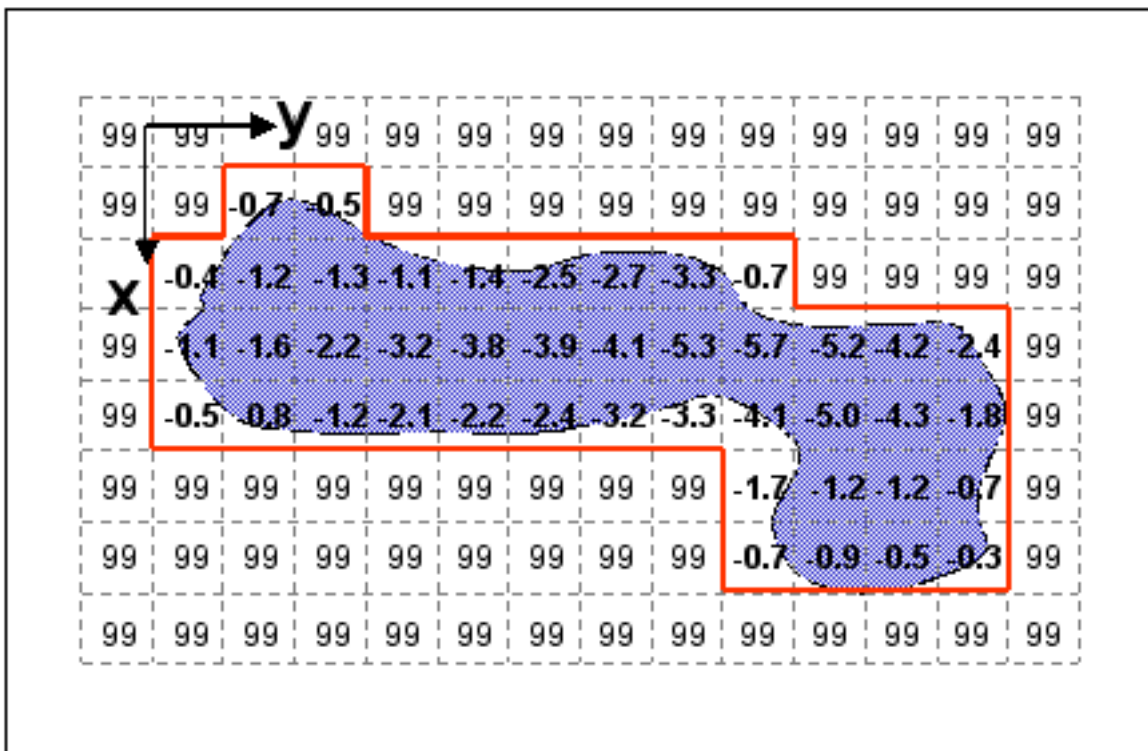


Figure 3. Plan view of bathymetry data

Figure 4 shows an elevation view (in the x - z plane) of a lake with a Cartesian overlay. A continuous, *physical-space* lake bottom is shown as a dashed line, with the resulting discrete lake bottom shown in a solid line. ELCOM computes the bathymetry on the face of each cell (the dashed lines) from user-supplied bathymetry data at the cell centers (on the solid dots). In this figure, the user $z=0$ baseline is chosen as the top of the domain, so that all the bathymetry values should be negative. Note that the $k=1$ index of (i,j,k) space corresponds with the lowermost layer, regardless of the user $z=0$ baseline position.

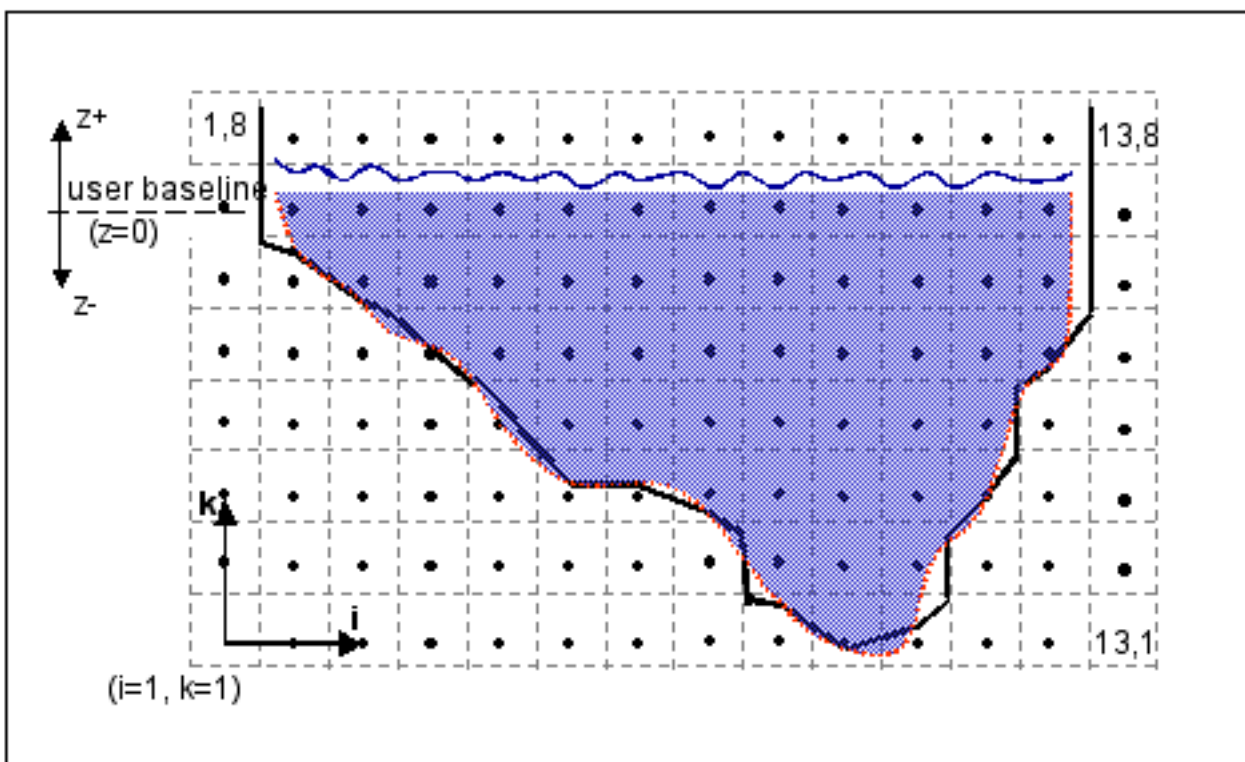


Figure 4. Elevation view of bathymetry data

2.3.3 Creating a bathymetry.dat file

An up-to-date copy of `bathymetry.dat` can be downloaded from [Appendix A](#). This file can be used as a template for creating your own `bathymetry.dat` file. The judicious use of extra comment lines and comments after the keywords, leads to a file that is much more readable. The `bathymetry.dat` file can be renamed, but the corresponding name in the `run_pre` must also be changed. The pre-processor will run with some items out of order, but the GUI will be unable to process the bathymetry file.

The remainder of this section will detail the use of each keyword in the `bathymetry.dat` file.

File header

keywords = FILE, VERSION

The first two non-comment lines of the `bathymetry.dat` file must begin with the keywords `FILE` and `VERSION`. The data following `FILE` must be the name of the file. The data following `VERSION` must be `1.1.a` (the present version of the `bathymetry.dat` file). Thus, the first two non-comment lines *must* appear as shown:

```
FILE    bathymetry.dat
VERSION 1.1.a
```

Description lines

keywords = TITLE, ANALYST, ORGANIZATION, COMMENT

Several keywords are provided to allow the user to put identifying information at the top of the bathymetry file. This information is printed to the standard output at the start of the run to provide identification of the debugging output data. The format for the header lines is `'string' keyword`, where the string must be a non-null character array enclosed in single quotations marks. If the string is left blank or not enclosed in quotes, the pre-processor will crash with a Fortran read error. The user must have only one `COMMENT` line. A typical set of header keywords is shown below.

```
'Lake Nowhere'      TITLE
'I.M. Underpaid'    ANALYST
'Very Large Company' ORGANIZATION
'test run number 1' COMMENT
```

Overwriting files

keyword = overwrite

This command is now superceded by the `run_pre` file.

```
yes overwrite
no  overwrite
```

Size of input bathymetry data

keywords = `x_rows`, `y_columns`, `z_layers`, `n_max`

In order to correctly read the bathymetry data from file `bathymetry.dat`, the user must tell the pre-processor the size of grid to expect. The value for `x_rows` is the number of rows in a two dimensional plan view of the bathymetry (e.g. figure 4). The `y_columns` is the number of columns in the two-dimensional plan view. The number of layers in the vertical direction is identified with the keyword `z_layers`. The keyword `n_max` is used to set an estimate of the number of grid cells in the domain that are required for simulation (i.e. interior points plus the land points in one layer surrounding the interior). For most simulations, `n_max` can be set to 0, and the code will start with an estimate of `x_rows x y_columns x z_layers`. However, in some cases, this may create a temporary data array larger than can be processed

using the available memory and the pre-processor will crash. The crash may produce a machine error message stating "out of memory," or a more enigmatic "segmentation fault". If insufficient memory is available, the user can estimate the actual number of grid points required for the simulation and rerun the preprocessor.

```
102 x_rows
84  y_columns
22  z_layers
0   n_max
```

Land cells

keyword = land_value

The land region and is designated by a unique bathymetry value (defined using the keyword "land_value") that is greater than any value found within the lake. The user is free to choose the land_value as any convenient number that meets this criterion. Thus, if the user chooses a $z=0$ baseline at the very top of the domain, the land_value may be set to zero (and all the bathymetry values should be negative numbers). If the user chooses the bottom of the domain as the $z=0$ baseline, then the land value may be any positive number that is greater than the largest number appearing within the lake boundary (it is usually convenient to use 9999). In this latter case, all the bathymetry values will be positive numbers. Note that any cell whose bathymetry is not equal to the land value is considered a cell within the domain — thus typographical errors can produce spurious interior cells of the domain. The user is not required to have a $z=0$ baseline within the domain: it may be convenient to use an external reference such as sea level. This may result in the lowest point in the domain having a positive non-zero bathymetry value.

```
99 land_value
```

Open boundary cells

keyword = open_value

Open boundary cells are boundaries with tidal forcing from an adjacent body of water that is not in the simulation domain. The user-defined bathymetry value that indicates an open boundary cell is defined with the keyword "open_value."

```
88 open_value
```

North vector

keywords = north_x, north_y

The user may create a bathymetry data file with any orientation (i.e. there is no requirement that north be at the top of the page). Since wind forcing on the free surface requires a direction, the user must designate the north direction. The north vector is determined as shown in [Figure 5](#). Note that the north vector is given in terms of a vector in physical (x,y,z) space, not the discrete (i,j,k) space. Because the origin is in the upper left corner, a standard "north up" data set has a north vector of $(-1, 0)$.

```
-1.0 north_x
0.5  north_y
```

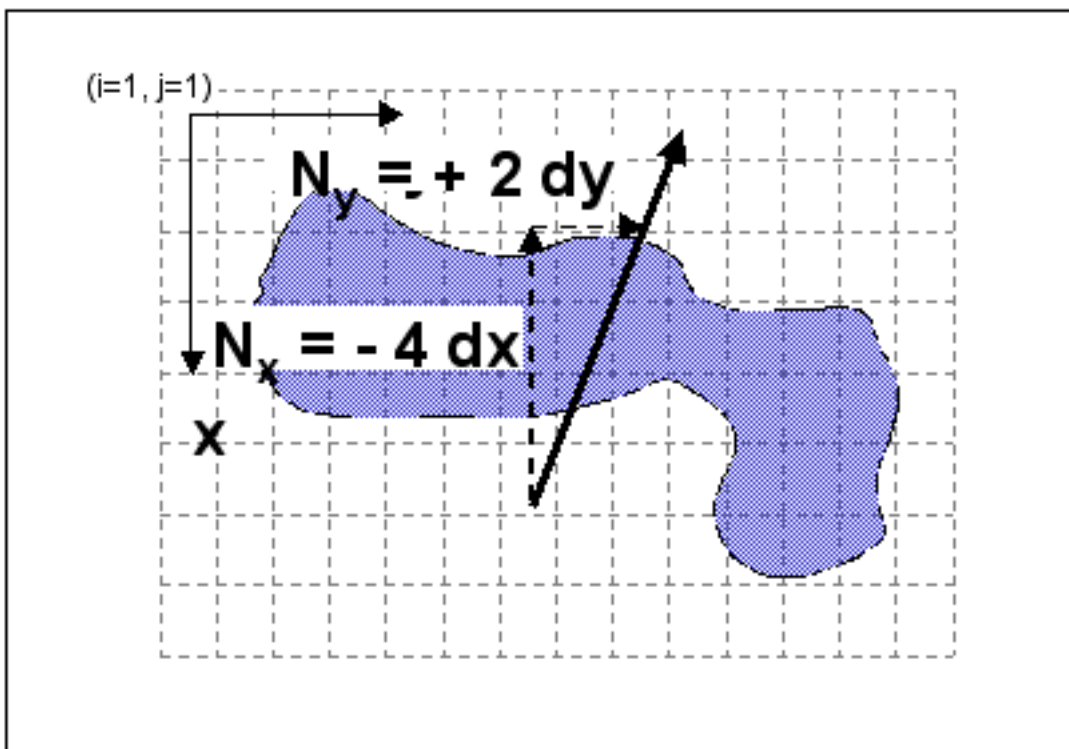


Figure 5. Calculation of the north vector

Geographic position

keywords = latitude, longitude, altitude

The user provides the geographic position of the simulation domain using the keywords `latitude`, `longitude` and `altitude`. Longitude and altitude are provided for future use in expanding CWR-Modeller to provide more complete graphics. Presently they have no effect on the simulation or output and may be safely set to 0. The latitude is used to provide appropriate Coriolis forcing in the ELCOM simulation. North latitudes are entered as positive numbers while south latitudes are entered as negative numbers.

```
-32.0 latitude
0.0 longitude
0.0 altitude
```

Horizontal grid spacing

keywords = `x_grid`, `y_grid`

The present version of ELCOM uses uniform (non-stretched) grid spacing in the x and y direction. Separate grid spacings are implemented in each horizontal dimension so that $dx \neq dy$ is valid. However, the user should be careful in implementing a solution grid with different dx and dy grid spacings since this can effect the solution accuracy. The x and y grid spacings are designated with the key words "`x_grid`" and "`y_grid`".

```
100.0 x_grid
100.0 y_grid
```

Layer thickness

keyword = `dz`

In the vertical direction, ELCOM allows the user to design a structure of horizontal layers whose thickness varies with depth (i.e. $dz = f(z)$). Note that the layer thickness may not vary with x or y coordinates. In the `bathymetry.dat` file, the successive layer thicknesses (in descending order from the top of the domain)

are designated by the keyword \bar{dz} . While the user is free to set the dz values in an arbitrary manner, the accuracy and stability of the simulation method may be effected by the degree of non-uniformity of the layers. In general, the simulation method will perform best (i.e. with greatest accuracy) using a grid with uniform dz . For non-uniform dz , the degradation of accuracy is a function of the rate at which the dz changes. Thus, a grid that has a dz that varies slowly will perform very well, while a grid with abrupt changes in grid size will be less accurate. For example, [Figure 6](#) shows two possible configurations which provide vertical resolution of 0.25 metres in the upper portion of a lake and 2.0 metre resolution in the lower portion. Both use the same number of grid cells, but the dz distribution on the left of [Figure 6](#) will show better performance.

```
0.5 dz
0.5 dz
0.5 dz
0.51 dz
0.54 dz
0.58 dz
0.65 dz
0.75 dz
0.90 dz
1.13 dz
1.41 dz
1.65 dz
1.88 dz
2.0 dz
```

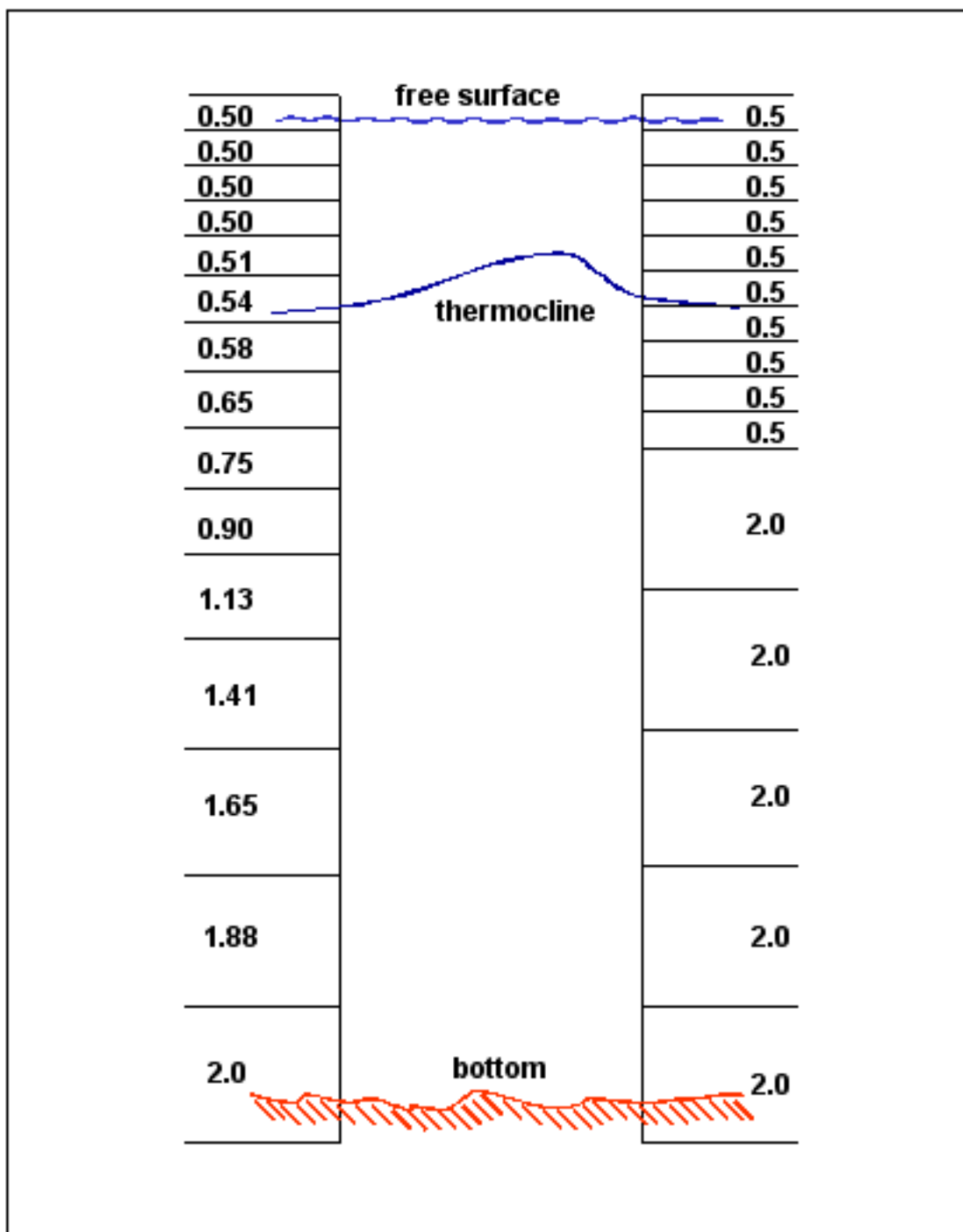


Figure 6. Vertical grid spacings

Bathymetry values

keywords = BATHYMETRY DATA

The bathymetry values must be presented in the file `bathymetry.dat` following a line with the two keywords `BATHYMETRY DATA`. Each line of bathymetry values must contain exactly `y_columns` of data and there must be exactly `x_rows` of data lines following `BATHYMETRY DATA`. Each value on the line may be separated from the next by one or more spaces or tabs. Thus, for a domain with `x_rows` = 8 and `y_columns` = 13, and `land_value` = 99 (i.e. similar to figure (3)), the bathymetry values would be entered into the `bathymetry.dat` file as shown:

```

BATHYMETRY DATA
99 99 99 99 99 99 99 99 99 99 99 99 99
99 99 -0.7 -0.5 99 99 99 99 99 99 99 99 99
99 -0.4 -1.2 -1.3 -1.1 -1.4 -2.5 -2.7 -3.3 -0.7 99 99 99
99 -1.1 -1.6 -2.2 -3.2 -3.8 -3.9 -4.1 -5.3 -5.7 -5.2 -2.4 99
99 -0.5 -0.8 -1.2 -2.1 -2.2 -2.4 -3.2 -3.3 -4.1 -5.0 -1.8 99
99 99 99 99 99 99 99 99 99 -1.7 -1.2 -0.7 99
99 99 99 99 99 99 99 99 99 -0.7 -0.9 -0.3 99
99 99 99 99 99 99 99 99 99 99 99 99 99

```

2.4 Setting up the boundary cell set file (bc.dat)

An up-to-date copy of `bc.dat` can be downloaded from [Appendix A](#).

2.4.1 Introduction

The boundary cell set file designates logical grouping of grid cells over which boundary conditions are enforced in ELCOM. One set of points may have a river inflow, another set may have a dam offtake, another set may represent groundwater inflows, etc. Each set of boundary points is defined by a unique *reference number*, (used also by ELCOM temporal boundary condition files), *type* (one of several keywords), *title* (user reference only), and a series of cell *i,j,k* coordinates. This file gives the user complete flexibility in the assignment of diverse and complex boundary conditions around a three-dimensional topography.

2.4.2 Boundary cell set categories

The keywords for the boundary cell sets fall into six categories, defined as follows:

`land`

Cell boundary which is considered to be land, on which a Dirichlet "no-slip" condition is imposed on all components of the velocity, and a Neumann "zero-gradient" condition is imposed on transported scalars.

`slip`

Cell boundary which is considered to be land, but on which the velocity components tangential to the boundary have a Neumann "free-slip" condition applied, while the normal component of the velocity is a Dirichlet "no-flux" condition (this may be thought of more precisely as normal Dirichlet, tangential Neumann, or NDTN). A Neumann "zero-gradient" condition is imposed on transported scalars.

`flow`

Cell boundary through which an inflow or outflow will be imposed as a Dirichlet "fixed velocity" condition. This also enforces a Dirichlet condition on transported scalars.

`section`

Cell boundary which is grouped into a set for application of an environmental forcing or scalar forcing boundary condition, but does not change the underlying form of the velocity boundary condition (e.g. the free surface may be divided into sections over which different wind conditions may be enforced — the type velocity boundary condition is not changed, but its magnitude is adjusted on different

`open`

Cell boundary which is "open" to inflows and outflows from another body of water where the velocity is not fixed *a priori* (i.e. not a Dirichlet boundary) but is a function of the baroclinic and barotropic forcing near the boundary (presently not implemented).

`interior`

An interior cell on which a source or sink of mass or tracer is enforced, e.g. a sewage outfall or drinking water offtake (presently not implemented).

A boundary cell set consists of a collection of cell *faces* that are the boundaries between *interior* cells and

land cells. The *land* boundary condition is the default for all faces of cells identified by the *land_value* in the *bathymetry.dat* file and for all boundary cell faces below the bathymetry data. As such, the user will rarely need to use a *land* boundary condition in the *bc.dat* file. For a lake without any inflows, outflows or special boundary conditions, the *bc.dat* file need not be present.

2.4.3 Boundary cell set face direction

The cell faces of the boundaries are identified by the (i,j,k) coordinates of the interior cell and the direction from the cell center to the boundary. The conventions used are as follows:

<i>xp</i>	the y-z cell face in the positive x direction from the cell center
<i>yp</i>	the x-z cell face in the positive y direction from the cell center
<i>xm</i>	the y-z cell face in the negative x direction from the cell center
<i>ym</i>	the x-z cell face in the negative y direction from the cell center
<i>top</i>	the x-y cell face above the
<i>all_sides</i>	all possible side boundary faces (i.e. all y-z and x-z faces)
<i>all</i>	all possible side boundary faces plus the bottom boundary

2.4.4 Boundary cell set keywords

The keywords used to identify boundary cell sets are a concatenation of the categories and the face direction. The valid keywords are listed below:

Boundary Cell Set Keywords			
<i>land_xp</i>	<i>flow_xp</i>	<i>slip_xp</i>	<i>section_xp</i>
<i>land_xm</i>	<i>flow_xm</i>	<i>slip_xm</i>	<i>section_xm</i>
<i>land_yp</i>	<i>flow_yp</i>	<i>slip_yp</i>	<i>section_yp</i>
<i>land_ym</i>	<i>flow_ym</i>	<i>slip_ym</i>	<i>section_ym</i>
<i>land_bottom</i>	<i>flow_bottom</i>	<i>slip_bottom</i>	<i>section_bottom</i>
<i>land_all_sides</i>	<i>flow_all_sides</i>	<i>slip_all_sides</i>	<i>section_all_sides</i>
<i>land_all</i>	<i>flow_all</i>	<i>slip_all</i>	<i>section_all</i>
			<i>section_top</i>
interior (not implemented)			
open (not implemented)			

2.4.5 Organization of the boundary cell set file (bc.dat)

At the top of the boundary cell set file (*bc.dat*) the user may put any number of comment lines. Each comment line must be started with an exclamation point (!) and have at least 3 items on the line (i.e. two words or numbers separated from the exclamation point and each other by a spaces or tabs). After the initial comments, the boundary cell set file consists of a series of definitions of boundary sets. Each set consists of a *set identification line* followed by one or more lines of *cell definitions*. Each line must have three items on it or the pre-processor will crash with a Fortran file read error. Blank lines are not allowed in the boundary cell set file.

Each set identification line is of the form: `integer keyword title`. The first item, `integer`, is a *reference number* for the boundary cell set that is used in preparing temporal boundary condition data files for the ELCOM simulation run (see section 3.5). Each boundary cell set must have a unique `integer` reference number. The `keyword` is one of the valid keywords listed in table 14. The `title` is a user-designated name for the boundary condition set. The `title` may have more than one word, although only the first word will be used by the code. After the set identification line, there will be one or more lines of cell definitions. A cell definition consists of three sets of number in the general form `istart:iend`
`jstart:jend kstart:kend`. The `i`, `j`, and `k` are integer values of the cell positions in space for a rectangular box of cells on which the relevant boundary cell keyword will be applied. For example, a simple boundary cell set definition might be

```
107 flow_xp river_inflow
    3:10    4:5        10:14
```

The preprocessor would designate as a single set, all positive x boundaries in the range $3 \leq i \leq 10$ and $4 \leq j \leq 5$ and $10 \leq k \leq 14$. A boundary cell set may have multiple lines of cell definitions:

```
13 flow_bottom groundwater_inflow
    5:10    1:5        10:14
    5:10    7:11       10:14
    3:8     12:17      1:5
```

Note that each number pair in a cell definition must be in increasing order.

There are two simpler cases for implementing boundary cell sets (1) where `i`, `j` or `k` is a single value rather than a range, (2) where all possible values in `i`, `j` or `k` are to be included. In the former case, only a single number needs to be provided (rather than a pair), in the latter case, only the colon (:) needs to be listed. For example, the cell set for an inflow boundary along a flat side where `j = 2` might be written as:

```
2791 flow_ym river_inflow
    3:15    2          :
```

This would group all the possible *y-minus* cell faces for cells at any `k` level, where `j = 2` and `i` is bounded by 3 and 15. This generalizes such that a boundary cell set may be written as

```
214 slip_all freeslip_domain
    :       :         :
```

which enforces a free-slip boundary condition on all boundary faces (sides and bottom) of the domain.

There is no error cross-checking within the pre-processor for cell faces which have been defined as belonging to more than one boundary cell set. In some cases, multiple definitions may be desirable. For example, one might define a river inflow which covers the entire depth of the river for purposes of the velocity boundary condition; and then define a separate section at the bottom of the river that is used to place a tracer. The `bc.dat` file might then contain

```
2791 flow_ym    river_inflow
    3:15    2          :
73  section_ym tracer_inflow
    3:15    2          1:5
```

In other cases, multiple set definitions for a particular face can produce unintended results. For example, if we define an inflow boundary and a global definition of free-slip boundaries on all the cells:

```
2791 flow_ym    river_inflow
    3:15    2          :
214 slip_all    freeslip_domain
    :       :         :
```

We then find that the cells (3:15, 2, :) get defined both with imposed velocity (Dirichlet) boundary conditions and free-slip (normal-Dirichlet, tangential-Neumann) boundary conditions. Obviously this is inconsistent and the flow simulation cannot enforce both boundary conditions. Checking for which combinations of multiple definitions are consistent and which are not is an error capturing routine that is not yet implemented. Thus, the above boundary conditions will actually be enforced, but one cannot tell *a priori* which boundary condition it will be without detailed examination of source code. Note that the order the items are defined may affect the operation of an inconsistent definition. Thus if the above inconsistent definition were reversed and written as:

```
214 slip_all freeslip_domain
    :      :
2791 flow_ym river_inflow
    3:15      2 :
```

The results in ELCOM may be different than the previous implementation — *they would be wrong in a different manner*. As a general rule, the user should not provide multiple definitions of velocity boundary conditions on a single cell face. The `land`, `flow` and `slip` keywords all effect the type of velocity boundary condition; it is incumbent on the user to be sure that the cell faces defined with these types do not have multiple definitions.

The following table illustrates a `bc.dat` file with a few subtle uses of the boundary cell set definitions. A flow boundary has been declared on the x-minus faces of cells with $i=2$. Thus, to get free-slip boundary conditions on all the other face, we use the keyword `slip_all` for cells with $i>2$. Then to get free slip on the y-minus and y-plus faces at $i=2$ requires two more definitions (These are not inconsistent with the `flow_xm` definition since they are on different faces of the same cell). This file also shows the use of the `section_top` to segregate the free surface into different sets onto which different wind data can be enforced.

```
! - - - -
! - bc.dat file
! - comment lines need three items

34 flow_xm      outflow_number_1
   2           :
75 section_top  wind_region_a
   :           1:10
118 section_top wind_region_b
   :           11:21
25 slip_all_sides free_slip_section
   3:34
26 slip_ym      free_slip_near_inflow_1
   2:
27 slip_yp      free_slip_near_inflow_2
   2:
```

Table: Example of boundary cell set file (bc.dat)

2.5 Running the pre-processor with `run_pre`

An up-to-date copy of `run_pre` can be downloaded from [Appendix A](#).

The `run_pre` file is a text file that contains the names and location of user-supplied input files and pre-processor output files. The pre-processor requires two user-supplied text files: `bathymetry.dat` and `bc.dat`. The former describes the bathymetry of the lake and grid configuration. The latter defines sets of boundary cells. The pre-processor output consists of two Fortran unformatted (binary) files: `sparsedata.unf` and `usedata.unf`. The former contains information on three-dimensional space and is used by CWR-Modeller and ELCOM. The latter is used by ELCOM, and contains the one-dimensional spatial relationships required for the simulation. These files may be renamed, but the user must keep track of which files are "sparsedata" files and which files are "usedata" files. Once the `run_pre` file is in the

directory from which you intend to run the pre-processor and the paths and names of the two input and two output files are specified in `run_pre`, then the pre-processor can be run by invoking the name of the executable, namely:

`PRE`

Notes

1. The pre-processor cannot interpolate from a coarse bathymetry to a finer grid or filter from a fine grid to a coarse grid. The user must perform this task before setting up the bathymetry file.
2. The pre-processor will correctly operate on files with any line order of the data:keyword commands, however the GUI can only process `bathymetry.dat` file that fit the data order prescribed in this section.

[Continue on to ELCOM](#)

```

FILE          bathymetry.dat
VERSION       1.1.a
! -----!
! headers: must be contained in quotes ' '.
! -
! 'a very small lake'      TITLE
! 'I.M. Underpaid'        ANALYST
! 'Very Large Corporation' ORGANIZATION
! 'example file'          COMMENT
! -----!
yes          overwrite files
! -----!
! number of grid cells
! -
8           x_rows
13          y_columns
14          z_layers
0           n_max
! -----!
! land and open boundary bathymetry values
! -
99          land_value
88          open_value
! -----!
! geographic position
! -
-4.0        north_x (positive x towards bottom of page)
2.0         north_y (positive y towards right on line)
-32.0       latitude
0.0         longitude
0.0         altitude
! -----!
! grid spacing
! -
100.0       x_grid_size
100.0       y_grid_size
! -----!
! - dz vector starting from top going down
!
0.5         dz top
0.5         dz
0.5         dz
0.54        dz
0.58        dz
0.65        dz
0.75        dz
0.90        dz
1.13        dz
1.41        dz
1.65        dz
1.88        dz
2.0         dz bottom
! -----!
! x in rows (increasing down), y in columns (increasing across)
! BATHYMETRY DATA
99  99  99  99  99  99  99  99  99  99  99  99  99
99  99  -0.7  0.5  99  99  99  99  99  99  99  99  99
99  -0.4  -1.2  -1.3  -1.1  -1.4  -2.5  -2.7  -3.3  -0.7  99  99  99
99  -1.1  -1.6  -2.2  -3.2  -3.8  -3.9  -4.1  -5.3  -5.7  -5.2  -2.4  99
99  -0.5  -0.8  -1.2  -2.1  -2.2  -2.4  -3.2  -3.3  -4.1  -5.0  -1.8  99
99  99  99  99  99  99  99  99  99  -1.7  -1.2  -0.7  99
99  99  99  99  99  99  99  99  99  -0.7  -0.9  -0.3  99
99  99  99  99  99  99  99  99  99  99  99  99  99

```

```
! - - - - -  
! boundary condition file for  
! A Very Large Lake  
! File prepared by I.M. Underpaid - 24.05.00  
! - - - - -
```

```
! ----- !
!
FILE          run_pre
VERSION      1.1.a
!
! ----- !
! input files
!
'bathymetry.dat'      BATHYMETRY_FILE
'bc.dat'              BOUNDARY_CONDITION_FILE
!
! ----- !
! output files
!
bathout.txt           DATAOUT_FILE
usedata.unf           ELCOM_USEDATA_FILE
sparsedata.unf        ELCOM_SPARSEDATA_FILE
!
! ----- !
! permission to overwrite existing files
!
yes      overwrite
!
! ----- !
```

3. ELCOM Operation

3.1 Introduction

Running an ELCOM simulation requires the user to provide several types of data: (1) initial conditions, (2) temporal boundary conditions, (3) configuration parameters, and (4) output control. ELCOM obtains this data from the user through user-prepared input files, pre-processor prepared input files, CWR-Modeller-prepared input files, and user-prepared subroutines. A separate user guide is provided for the use of CWR-Modeller, so this aspect of configuring ELCOM will not be covered in detail within this document. The following section will describe the types of data that CWR-Modeller is used to set, the files the user is required to prepare without CWR-Modeller, and the files that will be produced as ELCOM output.

3.2 Files for ELCOM

ELCOM is a command-line executable that communicates with the user through files. To simplify the task of setting up a simulation run, CWR-Modeller, pre-processor, and `makefile` have been developed. The pre-processor creates the files `sparsedata.unf` and `usedata.unf` which provide all the geometrical information for a simulation run. The `makefile` compiles the source code into an executable (ELCOM). CWR-Modeller creates the `datablock.db` file that controls simulation output. The following table provides a summary of file types, whether their presence is required or optional, the purpose of the file, and responsibility for preparation of the file. The file names in the `courier` bold font without brackets are fixed named files (i.e. the file must be present with exactly this name). File names such as `[data_name].unf` use square brackets to indicate that various names determined by ELCOM or CWR-Modeller may be in the bracketed section (these are specified in the following sections of this document). File names such as `<datablock>.db` use the `<>` brackets to indicate file names (or sections of names) which the user can set, either through CWR-Modeller or by changing the file name at the command line.

Files for ELCOM				
File Name	Presence	Usage	Preparation	Number of Files
<code>run_elcom</code>	<i>required</i>	configuration control	user	1
<datablock>.db	optional	output control	CWR-Modeller	1
<environment>.dat	optional	boundary condition	user	10
<code>elcom_user.f90</code>	<i>required</i>	initial conditions	user	1
<elcom_user_*.f90	optional	forced update	user	no limit
<sparsedata>.unf	<i>required</i>	geometry data	pre-processor	1
<usedata>.unf	<i>required</i>	geometry data	pre-processor	1
<group_name>.unf	output	datablock output	ELCOM	no limit
<code>[data_name].unf</code>	output	3D data type output	ELCOM	no limit
<code>[time_]<save>.unf</code>	output	simulation space output	ELCOM	no limit
<restart_out>.unf	output	restart file	ELCOM	1
ELCOM	<i>required</i>	executable	make	1
<code>makefile</code>	optional	compiling and linking	none	1
<code>[*].f90</code>	optional	source code	none	see makefile

3.3 Compiling and running ELCOM

If the ELCOM executable (ELCOM) is not available, one must be created by compiling the source code. The source code consists of all the files listed below, plus any user-prepared files.

ELCOM Fortran90 Modules

elcom_bc.f90	elcom_caedym1.f90
elcom_caedym2.f90	elcom_cgm_setup.f90
elcom_curvilinear.f90	elcom_datablock.f90
elcom_errors.f90	elcom_flowbound.f90
elcom_free_surface.f90	elcom_globals.f90
elcom_heat.f90	elcom_inputs.f90
elcom_mixing.f90	elcom_momentum.f90
elcom_no_caedym1.f90	elcom_no_caedym2.f90
elcom_openbound.f90	elcom_outputs.f90
elcom_quick.f90	elcom_scalar_solve.f90
elcom_scalar_utilities.f90	elcom_setup.f90
elcom_solver.f90	elcom_testcase.f90
elcom_time_step.f90	elcom_update.f90
elcom_user.f90	elcom_user_examples.f90
elcom_utilities.f90	elcom_velocities.f90
elcom_wind.f90	main/elcom_main.o
fd_no_caedym.f90	

ELCOM is compiled by typing the command

```
make
```

at the command line in the same directory as the source code and the `makefile`. This executes the `makefile` and produces the executable `ELCOM`. The `ELCOM` executable file is then installed using the UNIX syntax:

```
make install
```

The ELCOM source code directory can then be cleaned up using the following UNIX command:

```
make clean
```

ELCOM is executed at the command line by typing `ELCOM` (capital letters are required) within the directory *where the `run_elcom` and all the appropriate input data files* are located. The monitoring/debugging output is sent to the standard output of the computer (the display terminal on most UNIX workstations). Unless the simulation run is very short, it is recommended that the user redirect the output to a file with a UNIX execution redirection such as:

```
ELCOM > run_output
```

3.4 Simulation modules

ELCOM is organized into simulation modules that can be turned off/on depending on the type of simulation that is desired. A module is turned off/on by setting the value of control keywords. This is done within the `run_elcom` file. These controls provide a simple method for modifying the simulation. For example: a boundary condition file can be set up to include both meteorological data and inflow data. As a simple test, the user decides to turn off the inflow. One approach would be to return to the pre-processor, eliminate the inflow boundary cell set, re-run the pre-processor, then modify the temporal boundary condition file to remove the inflow data. As a simpler approach, the user need only set `iflow = 0`, and the inflow data and boundary cell sets will be ignored.

This section will give a brief overview of the modules within `ELCOM` and their corresponding keywords which are set within the `run_elcom` file.

3.4.1 Stratification

Density stratification in ELCOM can be through temperature, salinity, or a combination of the two by using the controls `itemperature` and `isalinity` in combination with `idensity = 1`. If the user desires, temperature and salinity can be treated as passive tracers (i.e. eliminating the density term from the momentum equation) by setting `idensity = 0`. If temperature and/or salinity are turned on, they must be initialized throughout the simulation domain through the `default_***` controls or the `user_init_***` controls and their associated subroutines.

3.4.2 Tracers

Up to 10 passive tracers may be transported in a simulation. They may be initialized in the domain using the `user_init_tracer` control and subroutine, or they may flow into the domain through an inflow boundary cell set using the keywords `TRACER_**` in a temporal boundary condition file.

3.4.3 Surface thermodynamics

The surface thermodynamics module controls the heat transfer across the free surface. To use this module, the user is required to provide the following data in temporal boundary condition files: solar radiation, wind speed, air temperature, relative humidity.

3.4.4 Wind

The wind effects ELCOM simulations in three ways: (1) evaporative cooling in the surface thermodynamics module, (2) application of a wind stress to the free surface, or (3) application of a wind induced momentum source to the wind mixed layer. The control `iwind` only sets the functioning of the wind input. Therefore, with `iwind = 0`, any wind value read in temporal boundary condition files will not provide a wind stress boundary condition, but will be used in the surface thermodynamics module (for `iheat_input > 0`). With `iwind = 1` or `2`, any wind value read in temporal boundary condition files will provide a wind stress boundary condition or a momentum source to the wind mixed layer. As well, the wind value will be used in the surface thermodynamics module (for `iheat_input > 0`).

If the keywords `WIND_SPEED` and `WIND_DIR` in the temporal boundary condition files are associated with the boundary cell set 0 (the entire free surface), then a wind field of uniform speed and direction is applied across the entire free surface. Alternatively, the user may declare multiple boundary cell sets on the free surface (using the keyword `section_top` in the file `bc.dat` for the preprocessor). The different sections are given reference numbers in `bc.dat`, which are then used in the temporal boundary condition files to identify the `WIND_SPEED` and `WIND_DIR` for each section.

3.4.5 Inflow/outflow

Inflow and outflow boundaries must be initially declared in the preprocessor `bc.dat` file. The user must then ensure that temporal boundary conditions are provided for the flow rate `n` in the temporal boundary condition files. For inflow conditions, all transported scalars must also be given boundary values temporal boundary condition files. If a transported scalar is not given an inflow value, the inflow volume will have a scalar concentration of 0.

3.4.6 User initialization

The user is provided with the ability to set up complex initial conditions through the adding code to subroutines in the [elcom_user.f90](#) file. The subroutines in this file are only called if the appropriate `user_init_***` control in the `run_elcom` file is non-zero. The stub routines supplied with ELCOM are designed to be executed with the `user_init_***` controls = 1. However, this can be modified by the user so that different initializations are performed for different values of `user_init_***` (see section 4 for more detail).

3.4.7 User update

The user is provided with the ability to design new subroutines for modifying the velocity or any transported scalars in the simulation. These are invoked through the subroutine `user_update` in the file `elcom_user.f90`. Details of writing code for user-developed subroutines is provided in section 5. This code is turned on or off with the control `user_update` in file `run_elcom`

3.5 Configuration controls

3.5.1 Uniform initial conditions

For simple initial conditions that (1) distribute a scalar temperature or salinity over the entire simulation volume or (2) provide a flat free surface across the entire domain, the user can set the desired values through the `run_elcom` file using keywords of the form `default_****`. Any default value entered will be overwritten by the user initialization subroutines (when invoked).

Time step

The time step (keyword = `del_t`) is the number of seconds that the simulation advances in time for each simulation step.

Number of steps

The total number of time steps to be computed in the present simulation run is set using the keyword `iter_max` in the `run_elcom` file.

Start date

The start date is entered in CWR Julian Day format which specifies the year and day using a real number with 7 digits to the left of the decimal place as `yyyddd.ddd`. The first 4 digits are the year and the subsequent 3 digits are the Julian Day. Hours, minutes, seconds are represented as fractions of a day (ie. 1am on January 1st, 2000 is represented as `2000000.041667`).

3.5.2 CGM tolerance

The `CGM_TOL` is the residual at which iteration of the conjugate gradient method (used for the free-surface solution) is stopped.

3.5.3 Turbulence

ELCOM solves the unsteady Reynolds-averaged Navier-Stokes equations (RANS) with four types of turbulence closure. The base closure scheme, which is always applied (`iclosure = 0`) requires the user to set fixed values for horizontal (`h_vis`) and vertical (`v_vis`) eddy viscosity. The base closure scheme is best applied only in unstratified flows. For stratified flow, the base closure scheme can be supplemented by unstable mixing (`iclosure = 1`) algorithms. The unstable mixing algorithm mixes momentum and transported scalars in vertical cells where an unstable gradient of potential density exists (i.e. $Ri < 0.0$). For stratified flows driven by surface wind, the wind-mixed-layer closure scheme with energy transport (`iclosure = 6`) is recommended. This approach computes the turbulent kinetic energy due to wind stirring, convective overturns and wind shear, then uses this to compute vertical mixing. Note that `iwind = 2` must be used with `iclosure = 6`.

3.5.4 Output frequency

Outputs of data that are not controlled by the datablock file are set through the `iter_out_***` and `start_output_***` keywords. The former controls the time step interval for providing output, the latter controls the first step on which output will be provided.

3.6 Configuration using file `run_elcom`

A `run_elcom` template can be obtained from [Appendix A](#).

ELCOM looks for the file `run_elcom` in the local directory (where the executable reside) to set the configuration of the simulation run. In future versions, the `run_elcom` file will be produced through CWR-Modeller. The `run_elcom` configuration file controls execution of optional modules and models within ELCOM, allowing the user to set the parameters that govern the physical and numerical simulation. The format of the file is flexible in that lines may occur in any order, and not all items must be present. Comment lines (preceded by `!`) may occur anywhere in the file. Blank lines are not allowed in the file (the code stops reading at a blank line). Each line (including comment lines) must have at least two items separated by a blank space. The comment marker (`!`) counts as an item, so a valid line is `! !` or `! -`, while `!!` (no space between) is not valid. The general format for the file is a series of lines, with each line listing a user setting, then a keyword that defines the parameter. A typical `run_elcom` file can be viewed [here](#). The individual keywords and settings are explained briefly below:

File headers:

```
FILE
    file name for this file (must be run_elcom)
VERSION
    current version of run_elcom file (1.1.a)
TITLE
    user title for this run (must be inside single quotes ' ')
ANALYST
    user name (must be inside single quotes ' ')
ORGANIZATION
    user company name (must be inside single quotes ' ')
COMMENT
    user comment — only 1 line (must be inside single quotes ' ')
```

CASE_KEYWORD

allows user to add code to be executed for a particular simlatoon case. Must be a string without blank spaces enclosed in single quotes.

Simulation module controls

```
iwind controls the wind module
    0 = no wind
    1 = wind model 1: wind stress boundary conditions; used only with iclosure = 0 or 1
    2 = wind model 2: direct wind momentum input; used only with iclosure = 6 -
    RECOMMENDED

iheat_input controls the surface thermodynamics
    0 = no surface thermodynamics
    1 = surface thermodynamics model 1

iflow controls inflow/outflow (Dirichlet boundaries) overrides boundary condition files
    0 = no inflow/outflow
    1 = inflow /outflow model 1

itemperature controls whether temperature is a transported scalar
    0 = no temperature simulation
    1 = temperature as transported scalar
```

`isalinity` controls whether salinity is a transported scalar

- 0 = no salinity in simulation
- 1 = salinity as a transported scalar

`idensity` controls whether density (buoyancy term) is included in simulation

- 0 = no buoyancy term
- 1 = density computed from UNESCO equation of state and used for buoyancy term

`ntracer` sets the number of tracers which can be transported in a simulation

- 0 = no tracers
- 1 to 10 = number of tracers initialized in `user_init_tracer` or supplied through inflows or `user_update`

`ICAEDYM` controls whether the CAEDYM water quality module is used

- 0 = CAEDYM off
- 1 = CAEDYM on

`icoriolis` controls whether or not the latitude (from `bathymetry.dat`) is used to compute a Coriolis term

- 0 = Coriolis off
- 1 = Coriolis on

`idatablock` controls which version of `datablock.db` is used

- 1 = version 1.0
- 2 = version 1.1

`icurvilinear` controls function of curvilinear approximation module

- 0 = curvilinear off
- 1 = output only
- 2 = use curvilinear and output

Initialization and update options

`irestart` use of restart file for initial conditions

- 0 = do not user restart file
- 1 = use restart file specified by `restart_in_file`

`user_init_u_vel` controls user written subroutine for initializing x velocity

- 0 = ignore subroutine `user_u_init`
- 1 = execute subroutine `user_u_init` with `USER_START_U = 1`

`user_init_v_vel` controls user written subroutine for initializing y velocity

- 0 = ignore subroutine `user_v_init`
- 1 = execute subroutine `user_v_init` with `USER_START_V = 1`

`user_init_w_vel` controls user written subroutine for initializing z velocity

- 0 = ignore subroutine `user_w_init`
- 1 = execute subroutine `user_w_init` with `USER_START_W = 1`

`user_init_temperature` controls user written subroutine for initializing temperature

- 0 = ignore subroutine `user_temperature_init`
- 1 = execute subroutine `user_temperature_init` with `USER_START_TEMPERATURE = 1`

`user_init_salinity` controls user written subroutine for initializing salinity

- 0 = ignore subroutine `user_salinity_init`
- 1 = execute subroutine `user_salinity_init` with `USER_START_SALINITY = 1`

`user_init_tracer` controls user written subroutine for initializing tracer

0 = ignore subroutine `user_tracer_init`
 1 = execute subroutine `user_trace_init` with `USER_START_TRACER = 1`

`user_init_viscosity` controls user written subroutine for initializing viscosity

0 = ignore subroutine `user_viscosity_init`
 1 = execute subroutine `user_viscosity_init` with `USER_START_VIS = 1`

`user_init_height` controls user written subroutine for initializing free surface height

0 = ignore subroutine `user_height_init`
 1 = execute subroutine `user_height_init` with `USER_START_HEIGHT = 1`

`user_init_extinction` controls user written subroutine for initializing the light extinction (or attenuation) coefficient

0 = ignore subroutine `user_extinction_init`
 1 = execute subroutine `user_extinction_init` with `USER_START_EXTINCTION = 1`

`user_update` controls user written subroutine for temporal updating of scalars

0 = ignore subroutine `user_update`
 1 = execute subroutine `user_update`

Turbulence modeling

`iclosure` controls the type of closure scheme used

0 = constant eddy viscosity
 1 = closure model #1 (constant eddy viscosity plus mixing of statically unstable stratification)
 2 = disabled
 3 = disabled
 4 = disabled
 5 = disabled
 6 = wind-mixed layer model including mixing energy transport; must be used with `iwind = 2` - RECOMMENDED

`vis_u` x-dir eddy viscosity (metres²/second)

`vis_v` y-dir eddy viscosity (metres²/second)

`vis_w` z-dir eddy viscosity (metres²/second)

`vis_u_tracer` x-dir eddy viscosity for tracers (metres²/second)

`vis_v_tracer` y-dir eddy viscosity for tracers (metres²/second)

`vis_w_tracer` z-dir eddy viscosity for tracers (metres²/second)

`vis_u_salt` x-dir eddy viscosity for salt (metres²/second)

`vis_v_salt` y-dir eddy viscosity for salt (metres²/second)

`vis_w_salt` w-dir eddy viscosity for salt (metres²/second)

`vis_u_temperature` x-dir eddy viscosity for heat (metres²/second)

`vis_v_temperature` y-dir eddy viscosity for heat (metres²/second)

`vis_w_temperature` z-dir eddy viscosity for heat (metres²/second)

`prandtl` turbulent Prandtl number (diffusion of heat κ_T)

`salt_schmidt` turbulent Schmidt number (diffusion of salinity κ_T)

`tracer_schmidt` turbulent Schmidt number (diffusion of tracer κ_T)

Model Settings and Controls

`drag_btm_cd` drag coefficient on bottom cells

`model_grav_damp_x` damping of gravity for x baroclinic term

`model_grav_damp_y` damping of gravity for y baroclinic term

Scalar maximum and minimum for debugging

`min_temperature`
`max_temperature`
`min_salinity`
`max_salinity`
`min_tracer1, min_tracer2, ... min_tracer10`
`max_tracer1, max_tracer2, ... max_tracer10`

Debugging controls

`debug_check` controls level of error checking in code

0 = no error checking
 1 = checking for errors, stop on fatal errors
 2 = highest level of checking for errors, stop on fatal errors
 3 = highest level of checking for errors, stop on warning

`debug_print` controls level of debugging printout

0 = none
 1 = minimal
 2 = moderate
 10 = extensive

`debug_point` sparse data index point for debug output

`debug_baroclinic_x` error trapping

0 = off
 1 = on

`debug_baroclinic_y` error trapping

0 = off
 1 = on

Output frequency

`iter_out_monitor` time step interval for output of monitoring data
`iter_out_save` time step interval for output of saved full simulation data
`iter_out_restart` time step interval for output of restart data

Output start time

`start_output_monitor` time step to start providing monitoring data
`start_output_save` time step to start providing simulation save files

Time controls

`start_date_CWR` start day in CWR Julian Day format yyyyddd (real)
`del_t` size of time step (in seconds)
`iter_max` maximum number of time steps in the simulation

Iterative (conjugate gradient method) solution controls

`CGM_TOL` Tolerance (residual) for stopping conjugate-gradient solution of free-surface equation.
`CGM_MIN` Minimum number of iterations that must be computed for the conjugate gradient method to solve the free-surface (over-rides the `CGM_TOL`)
`CGM_MAX` Maximum number of iterations that must be computed for the conjugate gradient method to solve the free-surface

Default (uniformly distributed) values

`default_height` Default water height (in metres). May be over-ridden by user-prepared subroutine if `user_init_height = 1`. Note that default height must be lower than the maximum height of the domain or an error will occur.

`default_salinity` Default salinity (in practical salinity units [psu], which is essentially equivalent to parts per thousand). May be over-ridden by user-prepared subroutine if `user_init_salinity = 1`.

`default_temperature` Default water temperature (in Celsius degrees). May be over-ridden by user-prepared subroutine if `user_init_temperature = 1`.

`default_wind_speed` Default wind speed (in ms^{-1}). May be over-ridden by user-prepared input file (see following section)

`default_wind_direction` Default wind direction (in degrees measured counter-clockwise from North in direction wind is coming *from*). May be over-ridden by user-prepared input file (see following section)

`default_extinction` Default light extinction coefficient (in Celsius degrees). May be over-ridden by user-prepared subroutine if `user_init_extinction = 1`.

`default_bc` Default boundary conditions.

- 1 = no-slip all
- 2 = free-slip all
- 3 = drag all
- 4 = neumann all
- 5 = free-slip sides, no-slip bottom
- 6 = free-slip sides, drag bottom
- 7 = free-slip sides, neumann bottom
- 8 = values defined in `mixing_elcom.f90`

Input file names (require exact file name, including extension)

`3D_data_file` Name of the file containing the 3D data produced by the pre-processor (`sparsedata.unf`)

`preprocessor_file` Name of the file containing the 1D data (maps) produced by the pre-processor (`usedata.unf`).

`boundary_condition_file` Name(s) of the files containing boundary condition information

`CAEDYM_file` Name of the file containing CAEDYM configuration information

`datablock_file` Name of the file containing datablock configuration (produced by GUI)

File Directories (require quote marks around; note that incomplete quotes can cause code to hang)

`infile_dir` input files directory (ie. `datablock.db`, `sparsedata.unf`, `usedata.unf`, boundary condition files)

`outfile_txt_dir` output *.unf files directory. Unformatted binary output files specified by `datablock.db` file.

`outfile_txt_dir` ELCOM diagnostic ascii format files.

Output files names (do not include extension)

`restart_save_file` kernel of name used for simulation save file (time stamp and `.unf` are appended by ELCOM)

`restart_out_file` kernel of name used for simulation restart output file (`.unf` is appended by ELCOM)

3.7 Input files

3.7.1 sparsedata.unf and usedata.unf

The files `sparsedata.unf` and `usedata.unf` are required files for ELCOM. These are the Fortran unformatted (binary) files produced by the pre-processor. The user may rename these files, but must keep track of which file is the `sparsedata` file (keyword = `3D_data_file`) and which is the `usedata` file (keyword = `preprocessor_file`). ELCOM will look for the files specified in the `run_elcom` file. If the incorrect files are used, the simulation will crash.

3.7.2 Datablock.db

A `datablock.db` template can be obtained from [Appendix A](#)

The file `datablock.db` is a text file produced by CWR-Modeller, and contains the configuration data used to set up the users output. The name for this file is specified by the user from CWR-Modeller. Alternatively, the `datablock.db` file may be setup manually using a text editor.

A `datablock.db` file is structured as a series of blocks of instructions. Each block requires an exact number of lines, with specific instructions on each line. Each line contains a single instruction followed by a comment. The comment must begin with a "!" character. In manually editing `datablock.db` files, it is important to keep in mind that while blocks of instructions may be added or removed, lines within a block can only be modified, but not removed. The following will discuss the structure of each block of instructions.

3.7.2.1 Datablock Description

The first 4 lines of a `datablock.db` are the datablock description. In order, they are the block descriptor, the current datablock version number, the number of output groups, and the number of output sets. Output groups and sets are described below.

3.7.2.2 Group Description

Following the datablock description are the group description blocks. A group represents a group of data to be output. A data group is defined as an output filename, datatypes to be output (ie. `TEMPERATURE`, `SALINITY`, etc...), the location within the simulation domain of the points at which data is to be output, and the time at which data is to be output. If several groups are desired, the group descriptions must follow each other without breaks between them.

A group description block contains 11 lines, which are as follows:

1. `BEGIN GROUP_DESCRIPTION`
begin block statement
2. group reference number
an integer labelling the group, this can have any value
3. user label
a name containing no blank spaces labelling the group (used by CWR modeller only)
4. output filename
should include `.unf` suffix
5. set reference number
an integer corresponding to a description of a set of points at which the data are to be output (sets will be described after groups)
6. number of datatypes in group
an integer indicating the desired number of datatypes to be output (ie. `TEMPERATURE`, `SALINITY`, etc...)
7. should read `.TRUE.`, do not modify

8. output interval
an integer indicating the desired number of timesteps between outputs
9. output start time
an integer indicating which time step to begin outputting data
10. output end time
an integer indicating which time step to stop outputting data, 0 indicates entire run
11. should read 0, do not modify

3.7.2.3 Set Description

A set is a set of points at which data are to be output. This can include a vertical column of points, a contiguous series of column of points (curtain), a horizontal sheet of points, or any group of points in the three dimensional domain. The valid set types are

```

PROFILE
CURTAIN_2D
SHEET_2D
ALL_3D
GENERAL_3D

```

Profiles and curtains require only the (x,y) locations of the base of each column to describe their location. Sheets can be referenced to a layer, a height above user datum, a height above bottom, or a height below the free-surface. ALL_3D simply outputs all points in the domain, and GENERAL_3D at points whos coordinates are individually specified.

A set description block contains 16 lines, which are as follows:

1. BEGIN SET_DESCRIPTION
begin set statement
2. set reference number
an integer labelling the set, this can have any value
3. user label
a name containing no blank spaces labelling the set (used by CWR modeller only)
4. type of set
on of the 5 valid set type names described above
5. number of cells in set
for type PROFILE this is the desired number of columns, for type CURTAIN2D this is the number of columns making up the desired curtain, for type GENERAL_3D this is the desired number of points, for types SHEET_2D and ALL_3D this value is 0
6. sheet type
for type SHEET_2D this indicates the SHEET datum, for all other set types the keyword is NULL. Valid sheet type keywords are:
 - LAYER horizontal layer referenced as layer number up from bottom as described in bathymetry.dat file
 - HEIGHT horizontal layer referenced as vertical distance in metres up from datum as defined in bathymetry.dat file
 - BOTTOM two dimensional surface defined as a vertical distance in metres up from the bottom
 - SURFACE two dimensional surface defined as a vertical distance down from the free-surface (note: the location of this sheet will evolve in time as the free-surface evolves)
 - FLAT a two dimensional sheet made up of calculated properties in each water column, the calculation type is specified below

7. sheet calc

calculation type for SHEET_2D, for all other set types the keyword is NULL.
Valid sheet calcs keywords are:

NEAREST Assigns each sheet value to the grid value vertically nearest the specified sheet height. Required for sheet types LAYER, HEIGHT, BOTTOM, and SURFACE. Not used with sheet type FLAT
MAX Assigns each sheet value to the maximum value in a column. Used only with sheet type FLAT
MIN Assigns each sheet value to the minimum value in a column. Used only with sheet type FLAT
AVG Assigns each sheet value to the average value in a column. Used only with sheet type FLAT
TOTAL Assigns each sheet value to the integrated value in a column. Used only with sheet type FLAT

8. sheet value

Layer number for sheet type LAYER. Vertical distance in metres for sheet types HEIGHT, BOTTOM, and SURFACE. 0 for sheet type FLAT.

9-16.

Not implemented, do not modify.

3.7.2.4 Group data types

This block begins with the desired number of data types, followed a listing of the desired data types.

1. BEGIN GROUP_DATA_TYPES

Begin data type description

2. group reference number

An integer referencing the group number defined in the group description

3- data types

One data type descriptor per line. The possible data type descriptors are:

Output Datatype Keywords		
U_VELOCITY,	V_VELOCITY,	W_VELOCITY,
TEMPERATURE,	SALINITY,	DENSITY,
EXTINCTION_COEF,	FLUC_PRESSURE,	TOTAL_PRESSURE,
TRACER_1,	TRACER_2,	TRACER_3,
TRACER_4,	TRACER_5,	TRACER_6,
TRACER_7,	TRACER_8,	TRACER_9,
TRACER_10,		
HEIGHT,	RI_NUM,	
SHEAR,	MIX_ENERGY,	
AUX_1,	AUX_2,	AUX_3,
AUX_4,	AUX_5,	
U_VIS,	V_VIS,	W_VIS,
TRACER_XVIS,	TRACER_YVIS,	TRACER_ZVIS,
SALT_XVIS,	SALT_YVIS,	SALT_ZVIS,
TEMP_XVIS,	TEMP_YVIS,	TEMP_ZVI,
S,	CELLS,	NORTH,
CURVATURE,	EPSILON,	LAMBDA,
GRAPH_X,	GRAPH_Y,	
WIND_U,	WIND_V,	
RADIAL,	TANGENT,	
GRAPH_R,	GRAPH_S	

3.7.2.5 Group Time List

This block is required to initialize a time stamp for each group. It requires only two lines as follows:

```
1. BEGIN GROUP_TIME_LIST
    Begin group time list
2. group reference number
    An integer referencing the group number defined in the group description
```

3.7.2.6 Set Cell Data

This block describes the spatial location of each datablock set. The locations are described as row, column, and/or layer number.

```
1. BEGIN SET_CELL_DATA
    Begin set cell data
2. set reference number
    An integer referencing the set number defined in the set description
3- cell locations
    This input varies for different set types as follows:
    PROFILE Each profile requires a row of input. Each row requires 4
    integers: 1) x-row, 2) y-column, 3) 0, 4) 0
    CURTAIN_2D Each profile making up the curtain requires a row of
    input. Each row requires 4 integers: x-row y-column 0 0
    SHEET_2D No Input
    ALL_3D No input
    GENERAL_3D Each point requires a row of input. Each row requires
    3 integers: x-row y-column z-layer
```

3.7.3 Temporal boundary condition (1D) files

A temporal boundary condition file template can be obtained from [Appendix A](#)

The one-dimensional temporal boundary condition files (keyword = `boundary_condition_file`) are set up by the user to provide environmental forcing information that changes as a function of time. These files require a format that consists of four header lines followed by columns of data. The data supplied are values for particular instants of time that are applied at particular points in space (i.e. over a boundary cell set). Each column is data of a particular type (defined by a keyword in the header). The first column must be time in CWR format (yyyddd.dd). Each row is data at a particular time. The file may have as many comment lines (preceded by !) at the top of the file as desired. Comment lines within the header data must have as many items as there are columns (i.e. if there are 3 columns a valid comment line is ! - -). No comment lines are allowed within the columns of data. The columns must be completely dense with data (i.e. there may not be "missing" data such that a row has less data than the number of columns).

The first header line contains one number: the number of data sets (columns) in the boundary condition file *not including the time column*.

The second line in the header also contains the number 0, indicating that the CWR time format will be used in the data below, and the first column in the data file should be a `TIME` column. Note that the user does *not* have to supply boundary condition data at the same time interval that the simulation uses as the computational time step (`del_t`): the simulation interpolates (linearly) between the temporal boundary condition data to find the appropriate data value at an instant in time.

The third line in the header must have an integer for each column. This is the reference number for the applicable set of boundary cell set (defined in the file `bc.dat` during the setup of the pre-processor) that each column of data will be applied to. For example, the example `bc.dat` file shown [here \(example bc.dat at bottom of section 2D\)](#) provides three sets of boundary points with reference numbers 34, 75, 118, 25, 26, 27. These numbers are used in the third line of the header to identify the points to which a

column of data applies. Note that a given reference number may be used for multiple data columns in multiple files, so that (for example) inflow volumes and temperatures for a particular set of boundary points may be supplied in one file, while tracer inflow for the same set of points may be supplied in a different file. Any temporal data boundary condition applied to the *entire* free surface (which is not required to be defined as a boundary cell set in `bc.dat`) is given a reference number of 0 (zero).

The fourth line in the header contains the keyword that describes the data in the column. The keywords are limited to a set specifically defined and understood by ELCOM and are listed in the table below.

The remaining lines in the file contain the data for the specified time (or time intervals).

Temporal Boundary Condition Data-types		
HEIGHT	m	water height for open boundary condition
INFLOW	m ³ s ⁻¹	inflow volume
OUTFLOW	m ³ s ⁻¹	inflow volume
WIND_SPEED	ms ⁻¹	wind speed
WIND_DIR	degrees	wind direction (<i>degrees measured clockwise from north that the wind comes from</i>)
SOLAR_RAD	Wm ²	solar radiation that reaches water surface
ATM_PRESS	Pascals	atmospheric pressure
CLOUDS		fractional cloud cover (<i>non-dimensional number between 0 and 1 inclusive</i>)
REL_HUM		relative humidity (<i>non-dimensional number between 0 and 1 inclusive</i>)
SALINITY	PSU	salinity (<i>practical salinity units which can be treated as parts per thousand within the numerical resolution of the model</i>)
WTR_DENSITY	kgm ³	density of inflowing water
AIR_TEMP	degC	air temperature
WTR_TEMP	degC	water temperature for inflowing water
U_VEL	ms ⁻¹	velocity in x-direction NOT_IMPLEMENTED
V_VEL	ms ⁻¹	velocity in y-direction NOT_IMPLEMENTED
W_VEL	ms ⁻¹	velocity in z-direction NOT_IMPLEMENTED
ANGLE	degrees	inflow/outflow angle (<i>degrees measured clockwise from north the water is flowing to</i>)
TRACER_1	concentration	tracer input
TRACER_2	concentration	tracer input
TRACER_3	concentration	tracer input
TRACER_4	concentration	tracer input
TRACER_5	concentration	tracer input
TRACER_6	concentration	tracer input

TRACER_7	concentration	tracer input
TRACER_8	concentration	tracer input
TRACER_9	concentration	tracer input
TRACER_10	concentration	tracer input

3.8 Non-Uniform Initial Conditions -- `elcom_user.f90`

Non-uniform initial conditions for certain variables can be set up within the `elcom_user.f90` file. As described in the [Compiling ELCOM](#) section, `elcom_user.f90` is a module of the ELCOM source code. Within `elcom_user.f90` are a number of subroutines which the user can modify to implement initial conditions. The `elcom_user.f90` file will contain a number of examples which the user can emulate to create their own initial conditions. A list of the initialization subroutines contained within `elcom_user.f90` is given below.

User Initialization Subroutines in <code>ELCOM_user.f90</code>		
<code>user_u_init,</code>	<code>user_v_init,</code>	<code>user_w_init,</code>
<code>user_temperature_init,</code>	<code>user_salinity_init,</code>	
<code>user_tracer_init,</code>	<code>user_extinction_init,</code>	<code>user_height_init,</code>
<code>user_uvis_init,</code>	<code>user_vvis_init,</code>	<code>user_wvis_init,</code>
<code>user_tracer_xvis_init,</code>	<code>user_tracer_yvis_init,</code>	<code>user_tracer_zvis_init,</code>
<code>user_salt_xvis_init,</code>	<code>user_salt_yvis_init,</code>	<code>user_salt_zvis_init,</code>
<code>user_temp_xvis_init,</code>	<code>user_temp_yvis_init,</code>	<code>user_temp_zvis_init</code>

For a particular subroutine to be used, the appropriate `user_init_***` keyword must be set to 1 in the `run_elcom` file. Within each initialization subroutine, various possible initial conditions are separated from each other by a series of `IF (CASE_KEYWORD == 'string1') THEN` or `ELSEIF (CASE_KEYWORD == 'string2') THEN` statements. A particular initial condition is implemented by setting the `CASE_WORD` equal to the appropriate string variable in the `IF` logical argument. The `CASE_WORD` is assigned in the `run_elcom` file.

3.9 Output files

3.9.1 Datablock output files

An example of a current `run_db2nc` file can be found in [Appendix A](#).

The datablock output files are named and configured through the [datablock.db file](#) which is configured either manually or through CWR modeller. ELCOM produces Fortran unformatted (binary) files as specified by the datablock output configuration. In general, one datablock output file contains one or more data types (e.g. velocity, temperature) applied to as set of points (chosen through CWR-Modeller) at a specified time interval. These files can be reprocessed into NETCDF files through CWR-Modeller, or manually using the command-line program `db2nc`

The program `db2nc` is used for converting datablock unformatted by binary files created by ELCOM to NETCDF files which can then be processed and viewed in a program such as Matlab. The `db2nc` executable should be placed either in the working directory or as recommended in the `/bin` directory which is in the operating system's search path. `db2nc` is run through a configuration file called `run_db2nc`.

The `run_db2nc` file contains 8 lines as follows:

1. Comment line, up to 64 characters
2. `.unf` file name and path
3. `sparsedata.unf` file name and path

4. `.nc` netcdf output file name and path
5. Start time step, -1 for beginning
6. End time step, -1 for end of file
7. Overwriting of `.nc` file? `replace` for overwriting, `noreplace` for no overwriting
8. Output diagnostics to monitor? `echo` to output to monitor, `noecho` to not output to monitor

Once the `run_db2nc` file is correctly configured, `db2nc` can be executed at the command line by typing:

```
db2nc run_db2nc
```

Notes:

- If you get the path for the `sparsedata.unf` file wrong in your `run_db2nc` file, `db2nc` will still run through and create a slightly smaller `.nc` file, complete with correct scalar data, however, all spatial information (x,y,z, etc) will be assigned a value of zero.
- The `run_db2nc` can have any file name, however the correct file name must be invoked at run time.
- The `run_db2nc` can only be configured for one file at a time. For batch processing of several `.unf` files, it is recommended that a `run_db2nc` file be created for each `.unf` to be converted and `db2nc` be invoked for each `run_db2nc` using a batch file.

3.9.2 Save and Restart output files

Save and Restart output files are Fortran unformatted (binary) files that contain run configuration information used to allow the user to restart ELCOM runs from a certain point. As described in the [run_elcom](#) section, the kernel name and output frequency of save and restart files are defined in the `run_elcom` file.

The save files are created starting at the time step specified by the value assigned to the `run_elcom` keyword `start_output_save`, and subsequently at intervals defined by `iter_out_save`. The file name of the save files begin with the user-specified kernel, given by `restart_save_file`, appended by a time stamp, appended by `.unf`. *Note that the time stamp does not currently include the simulation year and there exists the possibility that save files will be overwritten after a year of simulation time.*

The restart files are created at intervals defined by `iter_out_restart`, starting from the first time step. Subsequent restart files overwrite existing restart files. The file name of the restart files begin with the user-specified kernel, given by `restart_out_file`, appended by `.unf`. All start and restart files are written in the `outfiles` directory specified in the `run_elcom` file.

To restart an ELCOM simulation using a save or restart file, the user must place a copy of the desired save or restart file in the `infiles` directory as specified in the `run_elcom` file. In the `run_elcom` file the name of the desired save or restart file must be assigned to the `restart_infile` keyword, and keyword `i_restart` must be assigned the value 1.

[Continue on to templates](#)

```

! ----- !
!
FILE          run_elcom
VERSION       1.1.a
!
! ----- !
!
'A Very Large Lake'          TITLE
'I.M. Underpaid'            ANALYST
'Centre for Water Research'  ORGANIZATION
'ELCOM guide template'      COMMENT
! ----- !
! simulation modules (0=OFF,1=ON)
!
1      iwind          ! wind forcing (1=stress, 2=momentum, 3=special)
0      iheat_input    ! surface thermodynamics
0      iflow          ! inflow/outflow
1      itemperature   ! temperature
1      isalinity      ! salinity
1      idensity       ! density (baroclinic forcing)
0      ntracer        ! number of tracers
0      iCAEDYM        ! water quality module
0      icoriolis      ! turn off/on coriolis forcing
0      icurvilinear   ! curvilinear terms 1=output only 2=use & output
2      idatablock     ! datablock version control
! ----- !
! initialization options
!
0      irestart       ! start from restart file (yes=1/no=0)
0      user_init_u_vel ! 1 = use stub routine in elcom_user.f90
0      user_init_v_vel
0      user_init_w_vel
0      user_init_temperature
0      user_init_salinity
0      user_init_tracer
0      user_init_viscosity
0      user_init_height
0      user_init_extinction
0      user_update
'SURFACE_SLOPE'    CASE_KEYWORD
! ----- !
! turbulence modeling
! (0=no mixing, 1=static stability mixing, 6=mixed layer)
!
0      iclosure       ! closure model
!
! - bottom drag
!
1.0e-4 drag_btm_cd   ! drag coefficient on bottom cells
!
! - eddy viscosities
!
0.0    vis_u          ! x-dir eddy viscosity
0.0    vis_v          ! y-dir eddy viscosity
3.0e-3 vis_w          ! z-dir vertical viscosity
0.0    vis_x_tracer   ! x-dir eddy vis for tracer
0.0    vis_y_tracer   ! y-dir eddy vis for tracer
0.0    vis_z_tracer   ! z-dir eddy vis for tracer
0.0    vis_x_salt     ! x-dir eddy vis for salt
0.0    vis_y_salt     ! y-dir eddy vis for salt
0.0    vis_z_salt     ! z-dir eddy vis for salt
0.0    vis_x_temp     ! x-dir eddy vis for temperature
0.0    vis_y_temp     ! y-dir eddy vis for temperature
0.0    vis_z_temp     ! z-dir eddy vis for temperature
1.0    prandtl        ! turbulent Prandtl number

```



```
1.0    salt_schmidt    ! turbulent salinity Schmidt number
1.0    tracer_schmidt  ! tracer Schmidt number
!
! ----- !
! model settings and controls !
!
1.0    model_grav_damp_x ! damping of gravity for x baroclinic term
1.0    model_grav_damp_y ! damping of gravity for y baroclinic term
!
! ----- !
! scalar max and min for debugging !
!
0      min_temperature
40     max_temperature
0      min_salinity
35     max_salinity
0      min_tracer_1
1.1    max_tracer_1
0      min_tracer_2
1.1    max_tracer_2
0      min_tracer_3
0      max_tracer_3
0      min_tracer_4
0      max_tracer_4
0      min_tracer_5
0      max_tracer_5
0      min_tracer_6
0      max_tracer_6
0      min_tracer_7
0      max_tracer_7
0      min_tracer_8
0      max_tracer_8
0      min_tracer_9
0      max_tracer_9
0      min_tracer_10
0      max_tracer_10
!
! ----- !
! special setup (see elcom_setup.f90) !
!
0      itest_case      ! test case number
0      iflow_special   ! special flow setup
!
! ----- !
! debug output controls !
!
1      debug_check      ! level of error checking
2      debug_print      ! level of debug printing
0      debug_point      ! sparse data index point for debug output
1      debug_baroclinic_x ! error trapping on = 1, off = 0
1      debug_baroclinic_y ! error trapping on = 1, off = 0
!
! ----- !
! frequency of outputs !
!
10     iter_out_monitor  ! monitoring output to screen
2000   iter_out_save     ! saving separate restart files
500    iter_out_restart  ! overwriting restart output
!
! ----- !
! start of outputs !
!
1      start_output_monitor ! monitor output to screen
1      start_output_save    ! saved restart files
!
! ----- !
```

```
! time controls
!
2000001.0      start_date_cwr  ! CWR Julian day
950           del_t           ! Time step
100           iter_max        ! Number of time steps
!
! ----- !
! conjugate gradient iteration controls
!
1.e-16 CGM_TOL      ! CGM tolerance
30      CGM_MIN     ! minimum iterations of CGM
1000    CGM_MAX     ! maximum iterations of CGM
!
! ----- !
! default values (superseded by other inputs)
!
0.5      default_height
0.0      default_salinity
12.5     default_temperature
0.0      default_wind_speed
0.0      default_wind_direction
0.3      default_extinction
0        default_bc
!
! ----- !
! input file names (include exact file extension)
!
sparsedata.unf      3D_data_file
usedata.unf         preprocessor_file
met.dat             boundary_condition_file
datablock.db        datablock_file
!
! ----- !
! file directories (require quote marks around)
! (note that incomplete quotes can cause code to hang)
!
'infiles'           infile_dir
'unfiles'           outfile_unf_dir
'txtfiles'          outfile_txt_dir
!
! ----- !
! output file names (do not include file extension)
!
save                restart_save_file
restart_final       restart_out_file
```

```
DATABLOCK DESCRIPTION FILE
1.1.a ! datablock version
13 ! number of output groups
13 ! number of output sets
BEGIN GROUP_DESCRIPTION
100 ! GUI group reference number
tchain ! user label
tchain.unf ! output filename
100 ! GUI set reference number
13 ! number of datatypes in group
.TRUE. ! is_interval? logical: true if time step interval is used for output
5 ! output interval
0 ! output start time
0 ! output end time, 0 indicates entire run
0 ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
1 ! group reference number
curtain_group1 ! user label
curtain1.unf ! output filename
2 ! set reference number
4 ! number of datatypes in group
.TRUE. ! is_interval? logical: true if time step interval is used for output
5 ! output interval
0 ! output start time
0 ! output end time, 0 indicates entire run
0 ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
3 ! group reference number
curtain_group2 ! user label
curtain2.unf ! output filename
4 ! set reference number
9 ! number of datatypes in group
.TRUE. ! is_interval? logical: true if time step interval is used for output
5 ! output interval
0 ! output start time
0 ! output end time, 0 indicates entire run
0 ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
7 ! group reference number
all_group ! user label
all3d.unf ! output filename
6 ! set reference number
9 ! number of datatypes in group
.TRUE. ! is_interval? logical: true if time step interval is used for output
5 ! output interval
0 ! output start time
0 ! output end time, 0 indicates entire run
0 ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
9 ! group reference number
general_group ! user label
general3d.unf ! output filename
8 ! set reference number
9 ! number of datatypes in group
.TRUE. ! is_interval? logical: true if time step interval is used for output
5 ! output interval
0 ! output start time
0 ! output end time, 0 indicates entire run
0 ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
11 ! group reference number
sheet1_group ! user label
sheet1.unf ! output filename
10 ! set reference number
10 ! number of datatypes in group
.TRUE. ! is_interval? logical: true if time step interval is used for output
```

```
5      ! output interval
0      ! output start time
0      ! output end time, 0 indicates entire run
0      ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
13     ! group reference number
sheet2_group      ! user label
sheet2.unf       ! output filename
12              ! set reference number
10             ! number of datatypes in group
.TRUE.         ! is_interval? logical: true if time step interval is used for output
5             ! output interval
0             ! output start time
0             ! output end time, 0 indicates entire run
0             ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
15     ! group reference number
sheet3_group      ! user label
sheet3.unf       ! output filename
14              ! set reference number
10             ! number of datatypes in group
.TRUE.         ! is_interval? logical: true if time step interval is used for output
5             ! output interval
0             ! output start time
0             ! output end time, 0 indicates entire run
0             ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
17     ! group reference number
sheet4_group      ! user label
sheet4.unf       ! output filename
16              ! set reference number
10             ! number of datatypes in group
.TRUE.         ! is_interval? logical: true if time step interval is used for output
5             ! output interval
0             ! output start time
0             ! output end time, 0 indicates entire run
0             ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
19     ! group reference number
sheet5_group      ! user label
sheet5.unf       ! output filename
18              ! set reference number
10             ! number of datatypes in group
.TRUE.         ! is_interval? logical: true if time step interval is used for output
5             ! output interval
0             ! output start time
0             ! output end time, 0 indicates entire run
0             ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
21     ! group reference number
sheet6_group      ! user label
sheet6.unf       ! output filename
20              ! set reference number
10             ! number of datatypes in group
.TRUE.         ! is_interval? logical: true if time step interval is used for output
5             ! output interval
0             ! output start time
0             ! output end time, 0 indicates entire run
0             ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
23     ! group reference number
sheet7_group      ! user label
sheet7.unf       ! output filename
22              ! set reference number
10             ! number of datatypes in group
.TRUE.         ! is_interval? logical: true if time step interval is used for output
```

```
5      ! output interval
0      ! output start time
0      ! output end time, 0 indicates entire run
0      ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN GROUP_DESCRIPTION
25     ! group reference number
sheet8_group      ! user label
sheet8.unf       ! output filename
24     ! set reference number
10     ! number of datatypes in group
.TRUE. ! is_interval? logical: true if time step interval is used for output
5      ! output interval
0      ! output start time
0      ! output end time, 0 indicates entire run
0      ! size of time list, non-zero only if is_interval = .FALSE.
BEGIN SET_DESCRIPTION
100    ! GUI set reference number
tchain  ! user label
PROFILE_1D      ! type of set
3       ! number of cells in set
NULL     ! sheet type
NULL     ! sheet calc
0.0     ! sheet value
NULL     ! dynamic name #1
NULL     ! dynamic calc #1
NULL     ! dynamic direction #1
0.0     ! dynamic value #1
NULL     ! dynamic name #2
NULL     ! dynamic calc #2
NULL     ! dynamic direction #2
0.0     ! dynamic value #2
BEGIN SET_DESCRIPTION
2       ! set reference number
curtain_set    ! user label
CURTAIN_2D    ! type of set
3          ! number of cells in set
NULL       ! sheet type
NULL       ! sheet calc
0.0       ! sheet value
NULL      ! dynamic name #1
NULL      ! dynamic calc #1
NULL      ! dynamic direction #1
0.0      ! dynamic value #1
NULL      ! dynamic name #2
NULL      ! dynamic calc #2
NULL      ! dynamic direction #2
0.0      ! dynamic value #2
BEGIN SET_DESCRIPTION
4       ! set reference number
curtain_set2 ! user label
CURTAIN_2D  ! type of set
11         ! number of cells in set
NULL       ! sheet type
NULL       ! sheet calc
0.0       ! sheet value
NULL      ! dynamic name #1
NULL      ! dynamic calc #1
NULL      ! dynamic direction #1
0.0      ! dynamic value #1
NULL      ! dynamic name #2
NULL      ! dynamic calc #2
NULL      ! dynamic direction #2
0.0      ! dynamic value #2
BEGIN SET_DESCRIPTION
6       ! set reference number
all_set  ! user label
```

```
ALL_3D ! type of set
0 ! number of cells in set
NULL ! sheet type
NULL ! sheet calc
0.0 ! sheet value
NULL ! dynamic name #1
NULL ! dynamic calc #1
NULL ! dynamic direction #1
0.0 ! dynamic value #1
NULL ! dynamic name #2
NULL ! dynamic calc #2
NULL ! dynamic direction #2
0.0 ! dynamic value #2
BEGIN SET_DESCRIPTION
8 ! set reference number
general_set ! user label
GENERAL_3D ! type of set
4 ! number of cells in set
NULL ! sheet type
NULL ! sheet calc
0.0 ! sheet value
NULL ! dynamic name #1
NULL ! dynamic calc #1
NULL ! dynamic direction #1
0.0 ! dynamic value #1
NULL ! dynamic name #2
NULL ! dynamic calc #2
NULL ! dynamic direction #2
0.0 ! dynamic value #2
BEGIN SET_DESCRIPTION
10 ! set reference number
sheet1_set ! user label
SHEET_2D ! type of set
0 ! number of cells in set
LAYER ! sheet type
NEAREST ! sheet calc
3 ! sheet value
NULL ! dynamic name #1
NULL ! dynamic calc #1
NULL ! dynamic direction #1
0.0 ! dynamic value #1
NULL ! dynamic name #2
NULL ! dynamic calc #2
NULL ! dynamic direction #2
0.0 ! dynamic value #2
BEGIN SET_DESCRIPTION
12 ! set reference number
sheet2_set ! user label
SHEET_2D ! type of set
0 ! number of cells in set
HEIGHT ! sheet type
NEAREST ! sheet calc
4.5 ! sheet value
NULL ! dynamic name #1
NULL ! dynamic calc #1
NULL ! dynamic direction #1
0.0 ! dynamic value #1
NULL ! dynamic name #2
NULL ! dynamic calc #2
NULL ! dynamic direction #2
0.0 ! dynamic value #2
BEGIN SET_DESCRIPTION
14 ! set reference number
sheet3_set ! user label
SHEET_2D ! type of set
0 ! number of cells in set
```

```
BOTTOM          ! sheet type
NEAREST        ! sheet calc
0.5            ! sheet value
NULL           ! dynamic name #1
NULL           ! dynamic calc #1
NULL           ! dynamic direction #1
0.0           ! dynamic value #1
NULL           ! dynamic name #2
NULL           ! dynamic calc #2
NULL           ! dynamic direction #2
0.0           ! dynamic value #2
BEGIN SET_DESCRIPTION
16             ! set reference number
sheet4_set     ! user label
SHEET_2D       ! type of set
0              ! number of cells in set
SURFACE        ! sheet type
NEAREST        ! sheet calc
0.5            ! sheet value
NULL           ! dynamic name #1
NULL           ! dynamic calc #1
NULL           ! dynamic direction #1
0.0           ! dynamic value #1
NULL           ! dynamic name #2
NULL           ! dynamic calc #2
NULL           ! dynamic direction #2
0.0           ! dynamic value #2
BEGIN SET_DESCRIPTION
18             ! set reference number
sheet5_set     ! user label
SHEET_2D       ! type of set
0              ! number of cells in set
FLAT           ! sheet type
MAX            ! sheet calc
0.0           ! sheet value
NULL           ! dynamic name #1
NULL           ! dynamic calc #1
NULL           ! dynamic direction #1
0.0           ! dynamic value #1
NULL           ! dynamic name #2
NULL           ! dynamic calc #2
NULL           ! dynamic direction #2
0.0           ! dynamic value #2
BEGIN SET_DESCRIPTION
20             ! set reference number
sheet6_set     ! user label
SHEET_2D       ! type of set
0              ! number of cells in set
FLAT           ! sheet type
MIN            ! sheet calc
0.0           ! sheet value
NULL           ! dynamic name #1
NULL           ! dynamic calc #1
NULL           ! dynamic direction #1
0.0           ! dynamic value #1
NULL           ! dynamic name #2
NULL           ! dynamic calc #2
NULL           ! dynamic direction #2
0.0           ! dynamic value #2
BEGIN SET_DESCRIPTION
22             ! set reference number
sheet7_set     ! user label
SHEET_2D       ! type of set
0              ! number of cells in set
FLAT           ! sheet type
AVG            ! sheet calc
```

```
0.0          ! sheet value
NULL        ! dynamic name #1
NULL        ! dynamic calc #1
NULL        ! dynamic direction #1
0.0         ! dynamic value #1
NULL        ! dynamic name #2
NULL        ! dynamic calc #2
NULL        ! dynamic direction #2
0.0         ! dynamic value #2
BEGIN SET_DESCRIPTION
24          ! set reference number
sheet8_set  ! user label
SHEET_2D    ! type of set
0           ! number of cells in set
FLAT        ! sheet type
TOTAL       ! sheet calc
0.0         ! sheet value
NULL        ! dynamic name #1
NULL        ! dynamic calc #1
NULL        ! dynamic direction #1
0.0         ! dynamic value #1
NULL        ! dynamic name #2
NULL        ! dynamic calc #2
NULL        ! dynamic direction #2
0.0         ! dynamic value #2
BEGIN GROUP_DATA_TYPES
100         ! GUI group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
TEMPERATURE
DENSITY
AUX_1
AUX_2
AUX_3
AUX_4
AUX_5
MIX_ENERGY
SHEAR
RI_NUM
BEGIN GROUP_DATA_TYPES
1           ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
U_VIS
BEGIN GROUP_DATA_TYPES
3           ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
U_VIS
BEGIN GROUP_DATA_TYPES
7           ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
```



```
TRACER_2
U_VIS
BEGIN GROUP_DATA_TYPES
9      ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
U_VIS
BEGIN GROUP_DATA_TYPES
11     ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
HEIGHT
U_VIS
BEGIN GROUP_DATA_TYPES
13     ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
HEIGHT
U_VIS
BEGIN GROUP_DATA_TYPES
15     ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
HEIGHT
U_VIS
BEGIN GROUP_DATA_TYPES
17     ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
HEIGHT
U_VIS
BEGIN GROUP_DATA_TYPES
19     ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
```

```
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
HEIGHT
U_VIS
BEGIN GROUP_DATA_TYPES
21      ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
HEIGHT
U_VIS
BEGIN GROUP_DATA_TYPES
23      ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
HEIGHT
U_VIS
BEGIN GROUP_DATA_TYPES
25      ! group reference number
U_VELOCITY
V_VELOCITY
W_VELOCITY
SALINITY
TEMPERATURE
DENSITY
TRACER_1
TRACER_2
HEIGHT
U_VIS
BEGIN GROUP_TIME_LIST
100     ! GUI group reference number
BEGIN GROUP_TIME_LIST
1       ! group reference number
BEGIN GROUP_TIME_LIST
3       ! group reference number
BEGIN GROUP_TIME_LIST
7       ! group reference number
BEGIN GROUP_TIME_LIST
9       ! group reference number
BEGIN GROUP_TIME_LIST
11      ! group reference number
BEGIN GROUP_TIME_LIST
13      ! group reference number
BEGIN GROUP_TIME_LIST
15      ! group reference number
BEGIN GROUP_TIME_LIST
17      ! group reference number
BEGIN GROUP_TIME_LIST
19      ! group reference number
BEGIN GROUP_TIME_LIST
21      ! group reference number
BEGIN GROUP_TIME_LIST
```

```
23      ! group reference number
BEGIN GROUP_TIME_LIST
25      ! group reference number
BEGIN SET_CELL_DATA
100     ! GUI set reference number
4 3 0 0
4 6 0 0
4 11 0 0
BEGIN SET_CELL_DATA
2       ! set reference number
3 7 0 0
4 7 0 0
5 7 0 0
BEGIN SET_CELL_DATA
4       ! set reference number
4 2 0 0
4 3 0 0
4 4 0 0
4 5 0 0
4 6 0 0
4 7 0 0
4 8 0 0
4 9 0 0
4 10 0 0
4 11 0 0
4 12 0 0
BEGIN SET_CELL_DATA
6       ! set reference number
BEGIN SET_CELL_DATA
8       ! set reference number
4 6 2
4 6 3
4 10 1
4 10 3
BEGIN SET_CELL_DATA
10      ! set reference number
BEGIN SET_CELL_DATA
12      ! set reference number
BEGIN SET_CELL_DATA
14      ! set reference number
BEGIN SET_CELL_DATA
16      ! set reference number
BEGIN SET_CELL_DATA
18      ! set reference number
BEGIN SET_CELL_DATA
20      ! set reference number
BEGIN SET_CELL_DATA
22      ! set reference number
BEGIN SET_CELL_DATA
24      ! set reference number
```

```
! ----- - -
! A Very Large Lake
! Top hat 5m/s wind
! File prepared by I.M. Underpaid - 24.05.00
! ----- - -
6 data sets
0 seconds between data
      0          0          0          0          0          0          0
! ----- - -
  TIME      WIND_SPEED  WIND_DIR  SOLAR_RAD  AIR_TEMP  REL_HUM  CLOUDS
2000001.00      0         0.0       100.00     20.0     0.94     0.20
2000001.15      0         0.0       100.00     20.0     0.94     0.20
2000001.50      5         0.0       100.00     20.0     0.94     0.20
2000002.00      5         0.0       100.00     20.0     0.94     0.20
2000003.10      5         0.0       100.00     20.0     0.94     0.20
2000003.55      0         0.0       100.00     20.0     0.94     0.20
2000004.00      0         0.0       100.00     20.0     0.94     0.20
2000005.00      0         0.0       100.00     20.0     0.94     0.20
2000006.00      0         0.0       100.00     20.0     0.94     0.20
2000007.00      0         0.0       100.00     20.0     0.94     0.20
2000008.00      0         0.0       100.00     20.0     0.94     0.20
2000009.00      0         0.0       100.00     20.0     0.94     0.20
```

```
! Curtain translation
../curtain1.unf
../infile/sparsedata.unf
../ncfiles/curtain1.nc
-1
-1
replace
echo
```

A. Pre-processor Templates

- [bathymetry.dat](#)
- [bc.dat](#)
- [run_pre](#)

B. ELCOM Templates

- [run_elcom](#)
- [datablock.db](#)
- [temporal boundary conditions](#)
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