

# **Tutorial 5: Explanation of CPP Options, ocean.in, and r4dvar.in**

# **Tutorial 5: Explanation of CPP Options, ocean.in, and psas.in**

# R4D-Var Tutorial Wiki Page

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## Reprenter-based Data Assimilation (R4D-Var)

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## Introduction

[edit]

During this exercise you will apply the dual form strong/weak constraint, 4-Dimensional Variational (**4D-Var**) data assimilation based on the indirect representer algorithm to ROMS configured for the U.S. west coast and the California Current System (CCS). This configuration, referred to as **WC13**, has 30 km horizontal resolution, and 30 levels in the vertical. While 30 km resolution is inadequate for capturing much of the energetic mesoscale circulation associated with the CCS, **WC13** captures the broad scale features of the circulation quite well, and serves as a very useful and efficient illustrative example of R4D-Var.

## Model Set-up

The **WC13** model domain is shown in Fig. 1 and has open boundaries along the northern, western, and southern edges of the model domain.

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# WC13 C-preprocessing Options

## (Basic Configuration)

### Momentum Equations Options:

#define UV_ADV	Including advection terms
#define UV_COR	Including Coriolis term
#define DJ_GRADPS	splines density Jacobian PGF
#define UV_QDRAG	quadratic bottom friction
#define UV_VIS2	harmonic horizontal mixing
#define MIX_S_UV	mixing along s-levels

### Tracers Equations Options:

#define TS_U3HADVECTION	3 <sup>rd</sup> -order Upstream H. advection
#define TS_C4VADVECTION	4 <sup>th</sup> -order Centered V. advection
#define TS_DIF2	harmonic horizontal mixing
#define MIX_GEO_TS	mixing along geo-potentials
#define SALINITY	including salinity
#define NONLIN_EOS	nonlinear equation of state
#define ANA_BTFLUX	analytical bottom Temp flux
#define ANA_BSFLUX	analytical bottom Salt flux

### Vertical Turbulent Mixing Parameterization:

#define GLS_MIXING	Generic Length Scale Mixing
#ifdef GLS_MIXING	(K-omega)
#define N2S2_HORAVG	smoothing of buoyancy/shear
#define KANTHA_CLAYSON	stability function
#endif	

### Atmospheric Boundary Layer Parameterization:

#define BULK_FLUXES	Air/sea COARE bulk fluxes
#define DIURNAL_SRFLUX	imposing local diurnal cycle
#define SOLAR_SOURCE	solar radiation source term
#define LONGWAVE_OUT	compute outgoing long wave rad
#define EMINUSP	compute E-P

### Model Configuration Options:

#define SOLVE3D	solve 3D primitive equations
#define CURVGRID	curvilinear grid
#define MASKING	land/sea masking
#define SPHERICAL	spherical grid
#define PROFILE	time profiling
#define SPLINES	parabolic splines reconstruction

### Lateral Boundary Conditions:

#define EASTERN_WALL	closed eastern wall condition
#define WEST_FSCHAPMAN	free-surface, Chapman
#define WEST_M2FLATHER	2D momentum, Flather
#define WEST_M3CLAMPED	3D momentum, clamped
#define WEST_TCLAMPED	tracers, clamped condition
#define NORTH_FSCHAPMAN	free-surface, Chapman
#define NORTH_M2FLATHER	2D momentum, Flather
#define NORTH_M3CLAMPED	3D momentum, clamped
#define NORTH_TCLAMPED	tracers, clamped
#define SOUTH_FSCHAPMAN	free-surface, Chapman
#define SOUTH_M2FLATHER	2D momentum, Flather
#define SOUTH_M3CLAMPED	3D momentum, clamped
#define SOUTH_TCLAMPED	tracer, clamped

#define SPONGE	enhanced viscosity/diffusion areas
----------------	------------------------------------

# WC13 C-preprocessing Options

## (R4D-Var Configuration)

### Algorithm:

#define W4DVAR	dual form strong/weak constraint indirect representer algorithm
#define POSTERIOR_EOFS	estimate posterior error analysis error
#define POSTERIOR_ERROR_I	estimate initial conditions posterior analysis error

### Control Vector:

#define ADJUST_BOUNDARY	open boundary conditions increments
#define ADJUST_STFLUX	surface tracer flux increments
#define ADJUST_WSTRESS	surface wind stress increments

### Error Covariance Modeling:

#define VCONVOLUTION	Vertical correlation modeling
#define IMPLICIT_VCON	Implicit vertical diffusion operator
#define BALANCE_OPERATOR	Multivariate balance constraint
#ifdef BALANCE_OPERATOR	
#define ZETA_ELLIPTIC	SSH elliptic equation method
#endif	

### Prior:

#define FORWARD_READ	read basic state linearization in TLM and ADM files
#define FORWARD_WRITE	writing basic state by the NLM
#define FORWARD_MIXING	processing basic state vertical mixing coefficients
#define NL_BULK_FLUXES	surface kinematic fluxes from nonlinear model
#define RPM_RELAXATION	include relaxation term using previous Picard iteration solution

### I/O:

#define OUT_DOUBLE	double precision data in output NLM, TLM, RPM, and ADM
--------------------	--

# WC13 C-preprocessing Options

## (4D-PSAS Configuration)

### Algorithm:

```
#define W4DPSAS          dual form strong/weak constraint 4D-PSAS
#define POSTERIOR_EUFS   estimate posterior error analysis error
#define POSTERIOR_ERROR_I estimate initial conditions posterior analysis error
```

### Control Vector:

```
#define ADJUST_BOUNDARY  open boundary conditions increments
#define ADJUST_STFLUX   surface tracer flux increments
#define ADJUST_WSTRESS  surface wind stress increments
```

### Error Covariance Modeling:

```
#define VCONVOLUTION    Vertical correlation modeling
#define IMPLICIT_VCON   Implicit vertical diffusion operator
#define BALANCE_OPERATOR Multivariate balance constraint
#ifdef BALANCE_OPERATOR
#define ZETA_ELLIPTIC   SSH elliptic equation method
#endif
```

### Prior:

```
#define FORWARD_READ    read basic state linearization in TLM and ADM files
#define FORWARD_WRITE   writing basic state by the NLM
#define FORWARD_MIXING  processing basic state vertical mixing coefficients
#define NL_BULK_FLUXES  surface kinematic fluxes from nonlinear model
#define RPM_RELAXATION  include relaxation term using previous Picard iteration solution
```

### I/O :

```
#define OUT_DOUBLE      double precision data in output NLM, TLM, RPM, and ADM
```

# Include File: wc13.h

```
/*
** svn $Id: wc13.h 476 2010-06-26 20:25:30Z arango $
*****
** Copyright (c) 2002-2010 The ROMS/TOMS Group
** Licensed under a MIT/X style license
** See License_ROMS.txt
*****
**
** Options for the California Current System, 1/3 degree resolution.
**
** Application flag:   WC13
** Input script:      ocean_wc13.in
**                    s4dvar.in
**
** Available Drivers options: choose only one and activate it in the
**                             build.sh script (MY_CPP_FLAGS definition)
**
** AD_SENSITIVITY      Adjoint Sensitivity Driver
** AFT_EIGENMODES      Adjoint Finite Time Eigenmodes
** ARRAY_MODES         Stabilized representer matrix array modes
** CLIPPING            Stabilized representer matrix clipped analysis
** CORRELATION          Background-error Correlation Check
** GRADIENT_CHECK      TLM/ADM Gradient Check
** FORCING_SV          Forcing Singular Vectors
** FT_EIGENMODES       Finite Time Eigenmodes
** IS4DVAR              Incremental, strong constraint 4DVAR
** NLM_DRIVER           Nonlinear Basic State trajectory
** OPT_PERTURBATION     Optimal perturbations
** PICARD_TEST          Picard Iterations Test
** R_SYMMETRY           Representer Matrix Symmetry Test
** SANITY_CHECK         Sanity Check
** SO_SEMI              Stochastic Optimals: Semi-norm
```

# ROMS Standard Input Parameters

```
NtileI == 2          ! I-direction partition
NtileJ == 2          ! J-direction partition
. . .
NTIMES == 192        ! Number of time-steps (4 days)
  DT == 1800.0d0     ! Number of time-steps (48 steps per day)
. . .
Nouter = 1           ! Number of 4D-Var outer loops
Ninner = 50          ! Number of 4D-Var inner loops
. . .
LDEFOUT == T         ! Switch to create new history files
  NHIS == 48         ! Steps between writing of NLM data (daily)
  NDEFHIS == 0       ! Steps between creation of new NLM files
. . .
LcycleTLM == F       ! Switch to recycle records in TLM file
  NTLM == 48         ! Steps between writing of TLM data (daily)
  NDEFTLM == 0       ! Steps between creation of new TLM files
LcycleADJ == T       ! Switch to recycle records in ADM file
  NADJ == 192        ! Steps between writing of ADM data (strong constraint)
  NDEFADJ == 0       ! Steps between creation of new ADM files
  NSFF == 48         ! Steps between adjustment of surface fluxes (daily)
  NOBC == 48         ! Steps between adjustment of open boundary (daily)
. . .
APARNAM = r4dvar.in ! R4D-Var standard input parameters
```



# ROMS Standard Input Parameters

```
NtileI == 2          ! I-direction partition
NtileJ == 2          ! J-direction partition
. . .
NTIMES == 192        ! Number of time-steps (4 days)
  DT == 1800.0d0     ! Number of time-steps (48 steps per day)
. . .
Nouter = 1           ! Number of 4D-Var outer loops
Ninner = 50          ! Number of 4D-Var inner loops
. . .
LDEFOUT == T         ! Switch to create new history files
  NHIS == 48         ! Steps between writing of NLM data (daily)
  NDEFHIS == 0       ! Steps between creation of new NLM files
. . .
LcycleTLM == F       ! Switch to recycle records in TLM file
  NTLM == 48         ! Steps between writing of TLM data (daily)
  NDEFTLM == 0       ! Steps between creation of new TLM files
LcycleADJ == T       ! Switch to recycle records in ADM file
  NADJ == 192        ! Steps between writing of ADM data (strong constraint)
  NDEFADJ == 0       ! Steps between creation of new ADM files
  NSFF == 48         ! Steps between adjustment of surface fluxes (daily)
  NOBC == 48         ! Steps between adjustment of open boundary (daily)
. . .
APARNAM = psas.in    ! 4D-PSAS standard input parameters
```

# Standard Input File: ocean\_wc13.in

```
RSTNAME == wc13_rst.nc
HISNAME == wc13_his.nc
TLMNAME == wc13_tlm.nc
TLFNAME == wc13_tlf.nc
ADJNAME == wc13_adj.nc
AVGNAME == wc13_avg.nc
DIANAME == wc13_dia.nc
STANAME == wc13_sta.nc
FLTNAME == wc13flt.nc

! Input ASCII parameter filenames.

    APARNAM = r4dvar.in
    SPOSNAM = stations.in
    FPOSNAM = floats.in
    BPARNAM = bioFasham.in
    SPARNAM = sediment.in
    USRNAME = MyFile.dat

!
! GLOSSARY:
! =====
!
!-----
! Application title (string with a maximum of eighty characters) and
! C-preprocessing flag.
!-----
!
! TITLE          Application title.
!
! MyAppCPP       Application C-preprocessing option.
!
```

# 4D-Var Parameters: Normalization

```
Nmethod == 0 ! normalization method
Nrandom == 5000 ! randomization iterations
. . .
LdefNRM == F F F F ! Create a new normalization files
LwrtNRM == F F F F ! Compute and write normalization
. . .
CnormI(isFsur) = T ! 2D variable at RHO-points
CnormI(isUbar) = T ! 2D variable at U-points
CnormI(isVbar) = T ! 2D variable at V-points
CnormI(isUvel) = T ! 3D variable at U-points
CnormI(isVvel) = T ! 3D variable at V-points
CnormI(isTvar) = T T ! NT tracers
. . .
CnormB(isFsur) = T ! 2D variable at RHO-points
CnormB(isUbar) = T ! 2D variable at U-points
CnormB(isVbar) = T ! 2D variable at V-points
CnormB(isUvel) = T ! 3D variable at U-points
CnormB(isVvel) = T ! 3D variable at V-points
CnormB(isTvar) = T T ! NT tracers
. . .
CnormF(isUstr) = T ! surface U-momentum stress
CnormF(isVstr) = T ! surface V-momentum stress
CnormF(isTsur) = T T ! NT surface tracers flux
. . .
NRMnameM == wc13_nrm_m.nc ! model error (weak constraint)
NRMnameI == wc13_nrm_i.nc ! initial conditions
NRMnameB == wc13_nrm_b.nc ! open boundary conditions
NRMnameF == wc13_nrm_f.nc ! surface forcing (wind stress and net heat flux)
```

# 4D-Var Parameters: Decorrelation Scales

## Horizontal and vertical stability and accuracy factors (< 1):

!		IC	Model	OBC	Sur For	
	Hgamma =	0.5	0.5	0.5	0.5	! horizontal operator
	Vgamma =	0.0005	0.0005	0.0005	0.0005	! vertical operator

## Initial conditions correlations (m):

HdecayI(isFsur)	==	50.0d+3				! free-surface
HdecayI(isUbar)	==	50.0d+3				! 2D U-momentum
HdecayI(isVbar)	==	50.0d+3				! 2D V-momentum
HdecayI(isUvel)	==	50.0d+3				! 3D U-momentum
HdecayI(isVvel)	==	50.0d+3				! 3D V-momentum
HdecayI(isTvar)	==	50.0d+3	50.0d+3			! 1:NT tracers
VdecayI(isUvel)	==	30.0d0				! 3D U-momentum
VdecayI(isVvel)	==	30.0d0				! 3D V-momentum
VdecayI(isTvar)	==	30.0d0	30.0d0			! 1:NT tracers

## Surface forcing correlations (m):

HdecayF(isUstr)	==	100.0d+3				! surface U-momentum stress
HdecayF(isVstr)	==	100.0d+3				! surface V-momentum stress
HdecayF(isTsur)	==	100.0d+3	100.0d+3			! 1:NT surface tracer flux

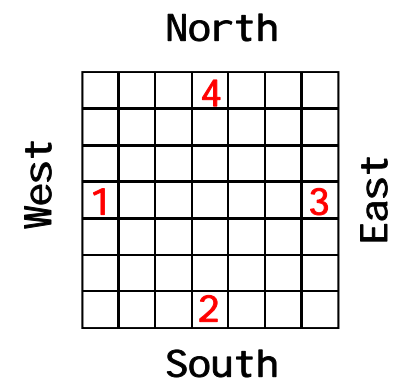
# 4D-Var Parameters: Decorrelation Scales

## Open boundary conditions correlations (m):

!		1: west	2: south	3: east	4: north	
HdecayB(i sFsur)	==	100.0d+3	100.0d+3	100.0d+3	100.0d+3	! free-surface
HdecayB(i sUbar)	==	100.0d+3	100.0d+3	100.0d+3	100.0d+3	! 2D U-momentum
HdecayB(i sVbar)	==	100.0d+3	100.0d+3	100.0d+3	100.0d+3	! 2D V-momentum
HdecayB(i sUvel)	==	100.0d+3	100.0d+3	100.0d+3	100.0d+3	! 3D U-momentum
HdecayB(i sVvel)	==	100.0d+3	100.0d+3	100.0d+3	100.0d+3	! 3D V-momentum
HdecayB(i sTvar)	==	4*100.0d+3	4*100.0d+3			! 1:NT tracers
VdecayB(i sUvel)	==	30.0d0	30.0d0	30.0d0	30.0d0	! 3D U-momentum
VdecayB(i sVvel)	==	30.0d0	30.0d0	30.0d0	30.0d0	! 3D V-momentum
VdecayB(i sTvar)	==	4*30.d0	4*30.d0			! 1:NT tracers

## Boundary edges to adjust (logical switches):

!		1	2	3	4	
Lobc(i sFsur)	==	T	T	F	T	! free-surface
Lobc(i sUbar)	==	T	T	F	T	! 2D U-momentum
Lobc(i sVbar)	==	T	T	F	T	! 2D V-momentum
Lobc(i sUvel)	==	T	T	F	T	! 3D U-momentum
Lobc(i sVvel)	==	T	T	F	T	! 3D V-momentum
Lobc(i sTvar)	==	T	T	F	T	\
		T	T	F	T	



# 4D-Var Parameters: Balance Operator

## SSH, elliptic solver:

Nbico == 200 ! bi conjugate gradient iteration

## SSH, integration of hydrostatic equation:

LNM\_depth == 1000.0d0 ! level of no motion (m, positive)

LNM\_flag = 1 ! Integration flag

[0] integrate from bottom to surface

[1] integrate from LNM to surface or  
from local depth, if shallower

## Balanced salinity empirical T-S relationship:

dTdz\_min == 0.001d0 ! minimum dT/dz (Celsius/m)

ml\_depth == 100.0d0 ! mixed-layer depth (m; positive)

## State Variables switches:

balance(i sSal t) = T

! salinity

balance(i sFsur) = T

! free-surface

balance(i sVbar) = F

! 2D momentum (ubar, vbar)

balance(i sVvel) = T

! 3D momentum (u, v)

# Other 4D-Var Parameters

## Lanczos algorithm parameters:

GradErr = 1.0d-4

! Upper bound on the relative error of the gradient

HevecErr = 1.0d-1

! Maximum error bound on Hessian eigenvectors

LhessianEV = T

! Compute approximated hessian eigen pairs

## Preconditioning:

Lprecond = F

! Limited-Memory Preconditioner: Spectral

Lritz = T

! Limited-Memory Preconditioner: Ritz

NritzEV = 0

! If preconditioning, number of eigenvectors if NritzEV = 0, use HevecErr

## Weak constraint:

LhotStart = T

! Hot start in subsequent outer loops

Npostl = 50

! Posterior error analysis Lanczos iterations

Nvct = 50

! Stabilized representer matrix eigenvector to process

tl\_M2diff == 0.0d0

! RPM relaxation (m2/s), 2D momentum

tl\_M3diff == 0.0d0

! RPM relaxation (m2/s), 3D momentum

tl\_Tdiff == 0.0d0 0.0d0

! RPM relaxation (m2/s), tracers

# Other 4D-Var Parameters

## Lanczos algorithm parameters:

GradErr = 1.0d-4

! Upper bound on the relative error of the gradient

HevecErr = 1.0d-1

! Maximum error bound on Hessian eigenvectors

LhessianEV = T

! Compute approximated hessian eigen pairs

## Preconditioning:

Lprecond = F

! Limited-Memory Preconditioner: Spectral

Lritz = T

! Limited-Memory Preconditioner: Ritz

NritzEV = 0

! If preconditioning, number of eigenvectors if NritzEV = 0, use HevecErr

## Weak constraint:

LhotStart = T

! Hot start in subsequent outer loops

NpostI = 50

! Posterior error analysis Lanczos iterations

Nvct = 50

! Stabilized representer matrix eigenvector to process

~~tl\_M2diff -- 0.0d0~~

~~! RPM relaxation (m2/s), 2D momentum~~

~~tl\_M3diff -- 0.0d0~~

~~! RPM relaxation (m2/s), 3D momentum~~

~~tl\_Tdiff -- 0.0d0 0.0d0~~

~~! RPM relaxation (m2/s), tracers~~



# R4D-Var Parameters File: r4dvar.in

```
! If weak constraint 4DVar, set number of iterations in the Lanczos
! algorithm used to estimate the posterior analysis error covariance
! matrix.

      NpostI = 50

! If weak constraint 4DVar, set diffusive relaxation coefficients (m2/s)
! used to relax representer tangent linear solution to previous Picard
! iteration linearized trajectory.

      tl_M2diff == 0.0d0          ! 2D momentum
      tl_M3diff == 0.0d0          ! 3D momentum

      tl_Tdiff == 0.0d0 0.0d0     ! NT tracers

! Switches (T/F) to create and write error covariance normalization
! factors for model, initial conditions, boundary conditions, and
! surface forcing. If TRUE, these factors are computed and written
! to NRMname(1:4) NetCDF files. If FALSE, they are read from NRMname(1:4)
! NetCDF file. The computation of these factors is very expensive and
! need to be computed only once for a particular application provided
! that grid land/sea masking, and decorrelation scales remains
! the same. Notice that four values are needed (1=initial conditions,
! 2=model, 3=boundary conditions, 4=surface forcing) per each nested
! grid, [1:4,1:Ngrids].

      LdefNRM == F F F F          ! Create a new normalization files
      LwrtNRM == F F F F          ! Compute and write normalization

! Switches to compute the correlation normalization coefficients for
! model error covariance.
```

# R4D-Var Job Script: job\_r4dvar.sh

1. Set path definition to one directory up in the tree.

```
set Dir = `dirname ${PWD}`
```

2. Set string manipulations perl script.

```
set SUBSTITUTE = ${ROMS_ROOT}/ROMS/Bin/substitute
```

3. Copy nonlinear model initial conditions file.

```
cp -p ${Dir}/Data/wc13_ini.nc wc13_ini.nc
```

4. Copy reparameter model initial conditions file, same as nonlinear model

```
cp -p ${Dir}/Data/wc13_ini.nc wc13_irp.nc
```

5. Copy reparameter model initial conditions file.

```
cp -p ${Dir}/Data/wc13_ini.nc wc13_irp.nc
```

6. Set model, initial conditions, boundary conditions and surface forcing error covariance standard deviations files.

```
set STDnameM = ../Data/wc13_std_m.nc  
set STDnameI = ../Data/wc13_std_i.nc  
set STDnameB = ../Data/wc13_std_b.nc  
set STDnameF = ../Data/wc13_std_f.nc
```

7. Set initial conditions, boundary conditions and surface forcing error covariance normalization factors files.

```
set NRMnameM = ../Data/wc13_nrm_m.nc  
set NRMnameI = ../Data/wc13_nrm_i.nc  
set NRMnameB = ../Data/wc13_nrm_b.nc  
set NRMnameF = ../Data/wc13_nrm_f.nc
```

8. Set observations file.

```
set OBSname = wc13_obs.nc
```

7. Get a clean copy of the observation file. This is really important since this file is modified.

```
cp -p ${Dir}/Data/${OBSname} .
```

8. Modify 4D-Var template input script and specify above files.

```
set R4DVAR = r4dvar.in  
if (-e $R4DVAR) then  
  /bin/rm $R4DVAR  
endif  
cp s4dvar.in $R4DVAR
```

```
$SUBSTITUTE $R4DVAR ocean_std_m.nc $STDnameM  
$SUBSTITUTE $R4DVAR ocean_std_i.nc $STDnameI  
$SUBSTITUTE $R4DVAR ocean_std_b.nc $STDnameB  
$SUBSTITUTE $R4DVAR ocean_std_f.nc $STDnameF  
$SUBSTITUTE $R4DVAR ocean_nrm_m.nc $NRMnameM  
$SUBSTITUTE $R4DVAR ocean_nrm_i.nc $NRMnameI  
$SUBSTITUTE $R4DVAR ocean_nrm_b.nc $NRMnameB  
$SUBSTITUTE $R4DVAR ocean_nrm_f.nc $NRMnameF  
$SUBSTITUTE $R4DVAR ocean_obs.nc $OBSname  
$SUBSTITUTE $R4DVAR ocean_hss.nc wc13_hss.nc  
$SUBSTITUTE $R4DVAR ocean_icz.nc wc13_icz.nc  
$SUBSTITUTE $R4DVAR ocean_mod.nc wc13_mod.nc  
$SUBSTITUTE $R4DVAR ocean_err.nc wc13_err.nc
```

# 4D-PSAS Job Script: `job_psas.sh`

1. Set path definition to one directory up in the tree.

```
set Dir = `dirname ${PWD}`
```

2. Set string manipulations perl script.

```
set SUBSTITUTE = ${ROMS_ROOT}/ROMS/Bin/substitute
```

3. Copy nonlinear model initial conditions file.

```
cp -p ${Dir}/Data/wc13_ini.nc wc13_ini.nc
```

4. ~~Copy representer model initial conditions file, same as nonlinear model~~

```
cp -p ${Dir}/Data/wc13_ini.nc wc13_irp.nc
```

5. Copy representer model initial conditions file.

```
cp -p ${Dir}/Data/wc13_ini.nc wc13_irp.nc
```

6. Set model, initial conditions, boundary conditions and surface forcing error covariance standard deviations files.

```
set STDnameM = ../Data/wc13_std_m.nc  
set STDnameI = ../Data/wc13_std_i.nc  
set STDnameB = ../Data/wc13_std_b.nc  
set STDnameF = ../Data/wc13_std_f.nc
```

7. Set initial conditions, boundary conditions and surface forcing error covariance normalization factors files.

```
set NRMnameM = ../Data/wc13_nrm_m.nc  
set NRMnameI = ../Data/wc13_nrm_i.nc  
set NRMnameB = ../Data/wc13_nrm_b.nc  
set NRMnameF = ../Data/wc13_nrm_f.nc
```

8. Set observations file.

```
set OBSname = wc13_obs.nc
```

7. Get a clean copy of the observation file. This is really important since this file is modified.

```
cp -p ${Dir}/Data/${OBSname} .
```

8. Modify 4D-Var template input script and specify above files.

```
set PSAS = psas.in  
if (-e $PSAS) then  
  /bin/rm $PSAS  
endif  
cp s4dvar.in $PSAS
```

```
$SUBSTITUTE $PSAS ocean_std_m.nc $STDnameM  
$SUBSTITUTE $PSAS ocean_std_i.nc $STDnameI  
$SUBSTITUTE $PSAS ocean_std_b.nc $STDnameB  
$SUBSTITUTE $PSAS ocean_std_f.nc $STDnameF  
$SUBSTITUTE $PSAS ocean_nrm_m.nc $NRMnameM  
$SUBSTITUTE $PSAS ocean_nrm_i.nc $NRMnameI  
$SUBSTITUTE $PSAS ocean_nrm_b.nc $NRMnameB  
$SUBSTITUTE $PSAS ocean_nrm_f.nc $NRMnameF  
$SUBSTITUTE $PSAS ocean_obs.nc $OBSname  
$SUBSTITUTE $PSAS ocean_hss.nc wc13_hss.nc  
$SUBSTITUTE $PSAS ocean_icz.nc wc13_icz.nc  
$SUBSTITUTE $PSAS ocean_mod.nc wc13_mod.nc  
$SUBSTITUTE $PSAS ocean_err.nc wc13_err.nc
```

# R4D-Var Job Script File: `job_r4dvar.sh`

```
#!/bin/csh -f
#
# svn $Id: job_r4dvar.sh 474 2010-06-25 20:19:44Z arango $
#####
# Copyright (c) 2002-2010 The ROMS/TOMS Group
# Licensed under a MIT/X style license
# See License_ROMS.txt
#####
#
# Strong/Weak constraint R4D-Var job script:
#
# This script NEEDS to be run before any run:
#
# (1) It copies a new clean nonlinear model initial conditions
# file. The nonlinear model is initialized from the
# background or reference state.
# (2) It copies representer model initial condition, same as
# nonlinear model.
# (3) Specify model, initial conditions, boundary conditions, and
# surface forcing error covariance input standard deviations
# files.
# (4) Specify model, initial conditions, boundary conditions, and
# surface forcing error covariance input/output normalization
# factors files.
# (5) Copy a clean copy of the observations NetCDF file.
# (6) Create 4D-Var input script "r4dvar.in" from template and
# specify the error covariance standard deviation, error
# covariance normalization factors, and observation files to
# be used.
#
#####
```

# Compile: build.sh

1. Set a local environmental variable to define the path to the directories where all this project's files are kept.

```
setenv MY_ROOT_DIR /home/arango/ocean/toms/repository
setenv MY_PROJECT_DIR ${PWD}
```

2. Location of your ROMS source code.

```
setenv MY_ROMS_SRC ${MY_ROOT_DIR}/branches/arango
```

3. Build script invoked CPP options.

```
setenv MY_CPP_FLAGS "-DW4DVAR"
setenv MY_CPP_FLAGS "${MY_CPP_FLAGS} -DPOSTERIOR_EOFS"
setenv MY_CPP_FLAGS "${MY_CPP_FLAGS} -DPOSTERIOR_ERROR_I"
```

4. Compiler selection environment variables.

```
setenv USE_MPI on
setenv USE_MPI_F90 on
setenv FORT pgi
```

5. Use custom library paths.

```
#setenv USE_MY_LIBS on
```

## Libraries for PGI

```
if ($?USE_MY_LIBS) then
  switch ($FORT)
  case "pgi"
    setenv ARPACK_LIBDIR /opt/pgi/soft/serial/ARPACK
    if ($?USE_MPI) then
      setenv PARPACK_LIBDIR /opt/pgi/soft/mpi/parpack
    endif

    if ($?USE_NETCDF4) then
      if ($?USE_MPI) then
        setenv NETCDF_INCDIR /opt/pgi/soft/mpi/netcdf4/include
        setenv NETCDF_LIBDIR /opt/pgi/soft/mpi/netcdf4/lib
        setenv HDF5_LIBDIR /opt/pgi/soft/mpi/hdf5/lib
      else
        setenv NETCDF_INCDIR /opt/pgi/soft/serial/netcdf4/include
        setenv NETCDF_LIBDIR /opt/pgi/soft/serial/netcdf4/lib
        setenv HDF5_LIBDIR /opt/pgi/soft/serial/hdf5/lib
      endif
    else
      setenv NETCDF_INCDIR /opt/pgi/soft/serial/netcdf3/include
      setenv NETCDF_LIBDIR /opt/pgi/soft/serial/netcdf3/lib
    endif
  breaksw
endif
```

# Compile: build.sh

1. Set a local environmental variable to define the path to the directories where all this project's files are kept.

```
setenv MY_ROOT_DIR /home/arango/ocean/toms/repository
setenv MY_PROJECT_DIR ${PWD}
```

2. Location of your ROMS source code.

```
setenv MY_ROMS_SRC ${MY_ROOT_DIR}/branches/arango
```

3. Build script invoked CPP options.

```
setenv MY_CPP_FLAGS "-DW4DPSAS"
setenv MY_CPP_FLAGS "${MY_CPP_FLAGS} -DPOSTERIOR_EOFs"
setenv MY_CPP_FLAGS "${MY_CPP_FLAGS} -DPOSTERIOR_ERROR_I"
```

4. Compiler selection environment variables.

```
setenv USE_MPI on
setenv USE_MPI_F90 on
setenv FORT pgi
```

5. Use custom library paths.

```
#setenv USE_MY_LIBS on
```

## Libraries for PGI

```
if ($?USE_MY_LIBS) then
  switch ($FORT)
  case "pgi"
    setenv ARPACK_LIBDIR /opt/pgi/soft/serial/ARPACK
    if ($?USE_MPI) then
      setenv PARPACK_LIBDIR /opt/pgi/soft/mpi ch/PARPACK
    endif

    if ($?USE_NETCDF4) then
      if ($?USE_MPI) then
        setenv NETCDF_INC_DIR /opt/pgi/soft/mpi ch/netcdf4/include
        setenv NETCDF_LIBDIR /opt/pgi/soft/mpi ch/netcdf4/lib
        setenv HDF5_LIBDIR /opt/pgi/soft/mpi ch/hdf5/lib
      else
        setenv NETCDF_INC_DIR /opt/pgi/soft/serial/netcdf4/include
        setenv NETCDF_LIBDIR /opt/pgi/soft/serial/netcdf4/lib
        setenv HDF5_LIBDIR /opt/pgi/soft/serial/hdf5/lib
      endif
    else
      setenv NETCDF_INC_DIR /opt/pgi/soft/serial/netcdf3/include
      setenv NETCDF_LIBDIR /opt/pgi/soft/serial/netcdf3/lib
    endif
  breaksw
endif
```

# Build Script: build.sh

```
#
# Sometimes it is desirable to activate one or more CPP options to run
# different variants of the same application without modifying its header
# file. If this is the case, specify each options here using the -D syntax.
# Notice also that you need to use shell's quoting syntax to enclose the
# definition. Both single or double quotes works. For example, to write
# time-averaged fields set:
#
#   setenv MY_CPP_FLAGS "-DAVERAGES"

setenv MY_CPP_FLAGS "-DW4DVAR"
setenv MY_CPP_FLAGS "${MY_CPP_FLAGS} -DPOSTERIOR_EOFs"
setenv MY_CPP_FLAGS "${MY_CPP_FLAGS} -DPOSTERIOR_ERROR_I"

# Other user defined environmental variables. See the ROMS makefile for
# details on other options the user might want to set here. Be sure to
# leave the switched meant to be off set to an empty string or commented
# out. Any string value (including off) will evaluate to TRUE in
# conditional if-stamentents.

setenv USE_MPI           on
setenv USE_MPIF90        on
setenv FORT               pgi

#setenv USE_OpenMP       on

#setenv USE_DEBUG         on
setenv USE_LARGE          on
setenv USE_NETCDF4        on

# Activate Data Access Protocol (like OPeNDAP) support for input
# NetCDF files. This is only possible for NetCDF library version
```