

1 Passo fazer o download do arquivo

[http://ftp.cptec.inpe.br/pesquisa/bam/bam\\_serial-2.0.tar.gz](http://ftp.cptec.inpe.br/pesquisa/bam/bam_serial-2.0.tar.gz)

2 Passo fazer o descompactação do arquivo

```
tar -zxvf bam\_serial-2.0.tar.gz
```

3 Passo Compilar o GetICnAGCM

```
cd bam\_serial-2.0/GetICnAGCM/Sources
```

```
make
```

```
make linux_gnu_serial
```

4 Passo Compilar o pre-processamento

```
cd bam\_serial-2.0/pre/sources
```

```
make
```

```
make linux_gnu_serial
```

5 Passo Compilar o modelo

```
cd bam\_serial-2.0/model/source
```

```
make
```

```
make linux_gnu_serial
```

6 Passo Compilar o pos-processamento

```
cd bam\_serial-2.0/pos/source
```

```
make
```

```
make linux_gnu_serial
```

7 Passo Rodar o GetICnAGCM para gerar a análise GAN...

```
cd bam\_serial-2.0/GetICnAGCM/GrADSScripts
```

```
grads -lbc "run GetERAWebICn_daily.gs 12z01dec2011 6 1"
```

```
cd bam\_serial-2.0/GetICnAGCM/RunScripts
```

```
./RunICnAGCM.XT6 62 28 2011120112 EIT Q T
```

8 Passo Rodar o pre-processamento

```
cd bam\_serial-2.0/run
```

```
./runPre
```

```
./runPre 62 28 2011120112 NMC 1 T
```

9 Passo Rodar o modelo

```
cd bam\_serial-2.0/run
```

```
./runModel
```

```
./runModel 1 1 1 bam 62 28 2011120112 2011120212 2011120212 2011120212 NMC sstmt 2
```

10 Passo Rodar o pos-processamento

```
cd bam_serial-2.0/run
```

```
./runPos
```

```
./runPos 1 1 1 posbam 126 28 2011120112 2011120212 2011120212 NMC COLD 1
```

Experimentos:

The bam model presents the following structure:

1. GetICnAGCM: (program that creates the initial condition of the atmosphere using as input data the reanalysis of the interim ERA)
2. bam  
pre (programs that pre-process the data for the resolution of the bam model grid.)  
model (program that performs the forecast for a given period.)

Pos (programs that pro-process the data of the bam model grid for the resolution of the output grid )

Before you run the GetlCnAGCM program you must edit the GetlCnAGCM / Configure.ksh file  
Example

```
:  
Vim GetlCnAGCM/Configure.ksh  
....  
....  
....  
export DirHome=/gws/nopw/j04/klngaman/paulo.kubota/GetlCnAGCM  
export DirData=/gws/nopw/j04/klngaman/paulo.kubota/GetlCnAGCM;mkdir -p ${DirData}  
export DirDataBam=/gws/nopw/j04/klngaman/paulo.kubota/bam  
export DirGrADS=/opt/grads/2.0.a9/bin  
.....  
....
```

The environment variables DirHome, DirData, DirDataBam, DirGrADS need to be modified for each user:

To run GetlCnAGCM you use the scripts:

```
./RunlCnAGCM.XT6 126 28 2011120112 EIT Q T
```

Submits the Geration of CPTEC AGCM Spectral Initial Condition Based on Input Grided Data of Topography, Mean Sea Level Pressure, Temperature, Zonal and Meridional Winds and Specific Humidity, for ERA 40 and ERA Interim Reanalysis and for NCEP Final Analysis.

```
#####  
# Arguments: #  
# #
```

```

#      1st : Trc: Triangular Truncation      #
#      2nd :   Lv: Number of Vertical Layers  #
#      3rd : DatelCn: Date of Initial Condition #
#      DatelCn: yyyyymmddhh                 #
#              yyyy = Year with 4 Digits     #
#              mm  = Month with 2 Digits     #
#              dd  = Day with 2 Digits       #
#              hh  = Hour with 2 Digits      #
#      4th : Case:                           #
#      E40 - ERA 40 at CPTEC                 #
#      E4W - ERA 40 at ECMWF                 #
#      EIT - ERA Interim at ECMWF           #
#      FNL - NCEP Final Analysis at NCAR    #
#      GFS - NCEP Analysis                   #
#      5th : GD: Grid:                       #
#              Quadratic (Q) or Linear (L)   #
#      6th : T: Level:                       #
#              Sigma (T) or Hibrid (F)      #
#
#
#*****#

```

Before running the programs of the BAM model it is necessary to modify the EnvironmentalVariablesMCGA file that is the bam/run directory.

In this step edit the file EnvironmentalVariablesMCGA

Vim EnvironmentalVariablesMCGA

```

.....
.....

```

```

echo ${DK}; export DK=/gws/nopw/j04/klngaman/paulo.kubota/bam
if [ -z "${DK}" ] ;then
export DK=`cd ../pwd`      # Default, if no filename specified.
fi
echo ${DK2}; export DK2=/gws/nopw/j04/klngaman/paulo.kubota/bam
.....
.....

```

The DK and DK2 environment variables define the paths where the input and output data are to run the model. The DK is where is the input data and the DK2 is where the model will write the files.

Para executar o pré-processamento do modelo bam é necessário executar o scripts que está no diretório bam/run/

```
./runPre 62 28 1979120112 NMC 0 T F 213 42
```

```

#*****#
#
# script to run CPTEC OPERATION #
# runPre TRC LV LABELI PREFIX total SmoothTopo GDASOnly RESIN KMIN #
# example : #
# runPre 94 28 2004032600 NMC 1 T F 62 42 #
# #
# #
# runPre 62 28 1979120112 NMC 1 T F 213 42 #
# #
# TRC => 4 digits spectral resolution #
# LV => 3 digits vertical resolution #
# LABELI => initial data YYYYMMDDHH #
# PREFIX => NMC #
# total => 1 tudo 0 parcial #
# SmoothTopo => Flag to Performe Topography Smoothing #
# GDASOnly => Flag to Only Produce Input CPTEC Analysis File (T or F) #
# RESIN => Spectral Horizontal Resolution of Input Data #
# 4 digits spectral resolution #
# KMIN => Number of Layers of Input Data #
# 3 digits vertical resolution #
# #
#*****#

```

```

#*****#
#
#
# script to run CPTEC Global Model on PC Clusters under MPI Scali #
# and Sun Grid Engine without OpenMP #
#
#

```

```

# assumptions: assume present at the same directory: #
#   ParModel_MPI (Global Model Executable file) #
#   MODELIN (Global Model input Namelist file) #
#
#
# usage: runModel cpu_mpi cpu_node task_omp name TRC LV LABELI LABELR LABELW LABELF NMC NMSST initlz #
#
#
#
#   runModel 72 4 6 bam 62 28 2004032600 2004042600 2004042600 2004042600 NMC sstmttd 2 #
#
#
# where: #
# cpu_mpi : integer, the desired number of mpi processes #
# cpu_node : integer, the desired number of mpi processes per shared memory node #
# task_omp : integer, number task omp by processos mpi #
# name : character, the job name (for SGE) #
# TRC : Spectral resolution (Number of wave) #
# LV : Vertical resolution(Number of layer in atmosphere) #
# LABELI : initial data YYYYMMDDHH initial condition #
#
# LABELR : Read restart data YYYYMMDDHH #
# LABELW : Write restart data YYYYMMDDHH data da inicializacao do restart do modelo #
# LABELF : Final data YYYYMMDDHH #
# NMC : prefix of files NMC #
#
# NMSST : type of sst (sstwkl, sstaoi, sstmttd, sstdyl) #
# initlz =2, ! initlz = 2 diabatic initialization and normal mode initialization #
# ! = 1 diabatic initialization and without normal mode initialization #
# ! = 0 without diabatic initialization and without normal mode initialization #
# ! [TOTAL RESTART ](adiabatic with no normal mode initialization) #
# ! < 0 same as >0 with sib variables read in instead of initialized #
# ! =-1 diabatic initialization and without normal mode initialization #
# ! with sib variables read in instead of initialized #
# ! =-2 diabatic initialization and normal mode initialization #
# ! with sib variables read in instead of initialized #
#
# hold: any, present or not;
#
# if absent, script finishes after queueing job; #
# if present, script holds till job completion
#

```

```
#####
```

```
#####
```

```
#
```

```
#
```

```
# script to run CPTEC Post-processing on PC Clusters under MPI Scali
```

```
#
```

```
# and Sun Grid Engine without OpenMP
```

```
#
```

```
#
```

```
#
```

```
# assumptions: assume present but NOT at the same directory:
```

```
#
```

```
#     $FEXE/PostGrib (Post-processing Executable file)
```

```
#
```

```
#     $FSCR/POSTIN-GRIB (Post-processing input Namelist file)
```

```
#
```

```
#
```

```
#
```

```
# usage:runPos cpu_mpi cpu_node task_omp name TRC LV LABELI LABELW LABELF NMC START FIRST #
```

```
#     runPos 1 1 1 posbam 62 28 2004032600 2004042600 2004042600 NMC COLD 1 #
```

```
# where:
```

```
#
```

```
# cpu_mpi : integer, the desired number of mpi processes
```

```
#
```

```
# cpu_node: integer, the desired number of mpi processes per shared memory node
```

```
#
```

```
# task_omp: numero de processos openmp por processos mpi
```

```
#
```

```
# name : Name of the Process (job)
```

```
#
```

```
# TRC : Spectral resolution(Number of wave)
```

```
#
```

```
# LV : Vertical resolution(Number of layer in atmosphere)
```

```
#
```

```
# LABELI : initial data YYYYMMDDHH
```

```
#
```

```
# LABELW : restart data YYYYMMDDHH (if COLD ---> LABELW=LABELF)
```

```
#
```

```
# LABELF : Final data YYYYMMDDHH
#
# NMC : prefix of files NMC
#
# START : type of pos-processing (COLD WARM)
#
# FIRST : ---
#
#*****#
```