

# **Running Global Model Parallel Experiments**



**Version 6.0**

**February 18<sup>th</sup>, 2015**

**NOAA/NWS/NCEP/EMC  
Global Climate and Weather Modeling Branch**

# Contents

1. Introduction .....	4
2. Operational Overview .....	5
2.1. Timeline of GFS and GDAS .....	5
2.2. Operational run steps .....	6
3. The Parallel Environment .....	7
4. Directories & Scripts .....	8
5. Data .....	11
5.1. Global Dump Archive .....	11
5.1.1. Location .....	11
5.1.2. Grouping .....	11
5.1.3. Files & Availability .....	11
5.2. Input/output files .....	12
5.2.1. Restart / initial conditions files .....	13
5.2.2. Observation files .....	14
5.2.3. Diagnostic files .....	15
5.3. Finding GDAS and GFS production run files .....	17
6. System Settings .....	19
6.1. Grid dimensions .....	19
6.2. Global Model Variables .....	A
7. Setting up an experiment .....	20
7.1. Important terms .....	20
7.2. Setting up your environment .....	20
7.3. Configuration file .....	21
7.4. Reconcile.sh .....	21
7.5. Rlist .....	22
7.6. Initial Conditions / Required Forcing Files .....	24
8. Submitting & running your experiment .....	25
8.1. Plotting output .....	26
8.2. Experiment troubleshooting .....	27
8.3. Tutorials .....	27
9. Parallels .....	28
10. Subversion & Trac .....	28
11. Related utilities .....	28
11.1. copygb .....	28
11.2. sfchr .....	29
11.3. sighdr .....	29
11.4. ss2gg .....	31
Appendix A: Global model variables.....	32

**Contacts:**

- Global Model POC - Kate Howard ([kate.howard@noaa.gov](mailto:kate.howard@noaa.gov))
- Global Branch Chief - Mark Iredell ([mark.iredell@noaa.gov](mailto:mark.iredell@noaa.gov))

**Version 6.0 Change Notes:**

- Added data section.
  - Moved data related information to new “Data” section, including dump archive.
  - Expanded dump archive section.
  - Updated data information based on Q1FY15 GFS/GDAS implementation.
- Other general updates from Q1FY15 GFS/GDAS implementation and recent machine changes.
- Updated tutorial information.

# What is the Global Forecast System?

The **Global Forecast System (GFS)** is a global numerical weather prediction system containing a global computer model and variational analysis run by the U.S. National Weather Service (NWS). The mathematical model is run four times a day, and produces forecasts for up to 16 days in advance, with decreased spatial resolution after 10 days. The model is a spectral model with a resolution of T1534 from 0 to 240 hours (0-10 days) and T574 from 240 to 384 hours (10-16 days). In the vertical, the model is divided into 64 layers and temporally, it produces forecast output every hour for the first 12 hours, every 3 hours out to 10 days, and every 12 hours after that.

## 1. Introduction

So you'd like to run a GFS experiment? This page will help get you going and provide what you need to know to run an experiment with the GFS. Before continuing, some information:

- This page is for users who can access the R&D machine (Zeus), WCOSS (Gyre/Tide), or the S4 system.
- This page assumes you are new to using the GFS model and running GFS experiments. If you are familiar with the GFS Parallel System, or are even a veteran of it, feel free to jump ahead to specific sections.
- If at any time you are confused and can't find the information that you need please feel free to email for help.

To join the global model mailing list:

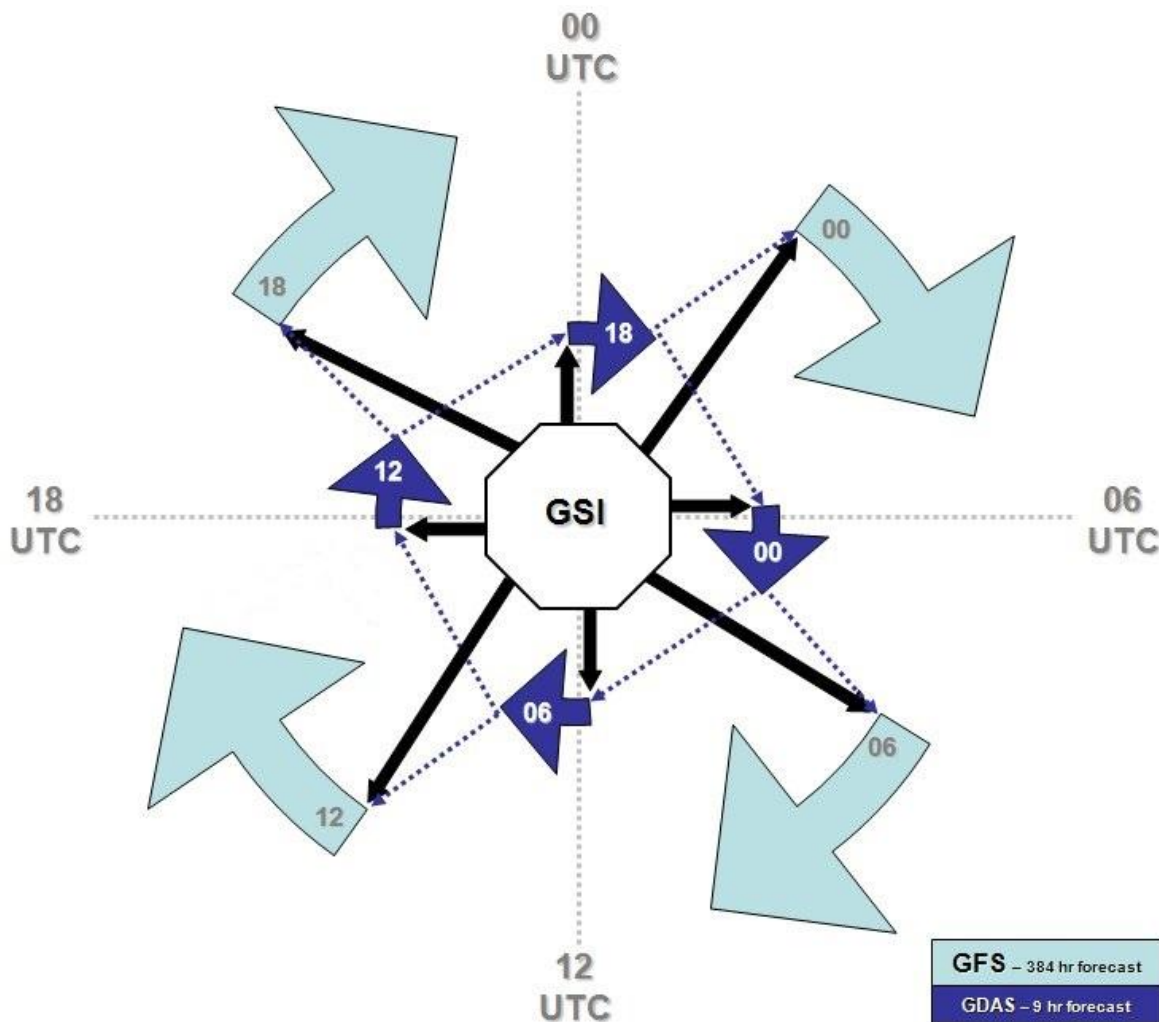
Global parallel announcements -  
<https://lstsrv.ncep.noaa.gov/mailman/listinfo/ncep.list.emc.glopara-announce>

## 2. Operational Overview

The Global Forecast System (GFS) is a three-dimensional hydrostatic global spectral model run operationally at NCEP. The **GFS** consists of two runs per six-hour cycle (00, 06, 12, and 18 UTC), the "early run" **gfs** and the "final run" **gdas**:

- **gfs/GFS** refers to the "early run". In real time, the early run, is initiated approximately 2 hours and 45 minutes after the cycle time. The early gfs run gets the full forecasts delivered in a reasonable amount of time.
- **gdas/GDAS** refers to the "final run", which is initiated approximately six hours after the cycle time.. The delayed gdas allows for the assimilation of later arriving data. The gdas run includes a short forecast (nine hours) to provide the first guess to both the gfs and gdas for the following cycle.

### 2.1 Timeline of GFS and GDAS



\*Times are approximate

## 2.2 Operational run steps

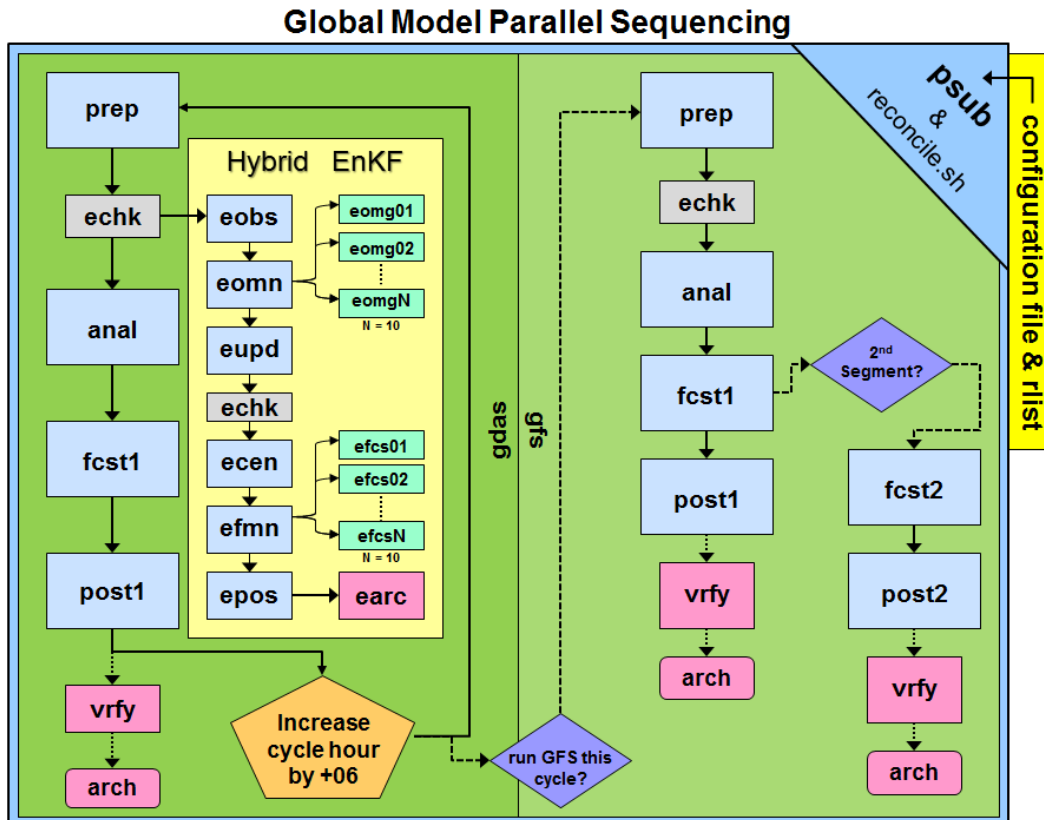
- **dump** - Gathers required (or useful) observed data and boundary condition fields (done during the operational GFS run); used in real-time runs, already completed for archived runs. Unless you are running your experiment in real-time, the dump steps have already been completed by the operational system (gdas and gfs) and the data is already waiting in a directory referred to as the dump archive.
- **storm relocation** - In the presence of tropical cyclones this step adjusts previous gdas forecasts if needed to serve as guess fields. For more info, see the relocation section of Dennis Keyser's Observational Data Dumping at NCEP document. The storm relocation step is included in the prep step (gfsprep/gdasprep) for experimental runs.
- **prep** - Prepares the data for use in the analysis (including quality control, bias corrections, and assignment of data errors) For more info, see Dennis Keyser's PREPBUFR PROCESSING AT NCEP document.
- **analysis** - Runs the data assimilation, currently Gridpoint Statistical Interpolation (GSI)
- **enkf** - Multiple jobs which run the hybrid ensemble Kalman filter–three-dimensional variational (3DVAR) analysis scheme
- **forecast** - From the resulting analysis field, runs the forecast model out to specified number of hours (9 for gdas, 384 for gfs)
- **post** - Converts resulting analysis and forecast fields to WMO grib for use by other models and external users.

Additional steps run in experimental mode are (pink boxes in flow diagram in next section):

- verification (gfs vrfy / gdas vrfy)
- archive (gfs arch / gdas arch) jobs

### 3. The Parallel Environment

**GFS** experiments employ the global model parallel sequencing (shown below). The system utilizes a collection of job scripts that perform the tasks for each step. A job script runs each step and initiates the next job in the sequence. Example: When the anal job finishes it submits the forecast job. When the forecast job finishes it submits the post job, etc.



Flow diagram of a typical experiment with Hybrid EnKF turned ON

As with the operational system, the **gdas** provides the guess fields for the **gfs**. The **gdas** runs for each cycle (00, 06, 12, and 18 UTC), however, to save time and space in experiments the **gfs** (right side of the diagram) is initially setup to run for only the 00 UTC cycle. (See the "run **GFS** this cycle?" portion of the diagram) The option to run the **GFS** for all four cycles is available (see `gfs_cyc` variable in configuration file).

As mentioned in section 2.2, an experimental run is different from operations in the following ways:

- Dump step is not run as it has already been completed during real-time production runs
- Addition steps in experimental mode:
  - verification (vrfy)
  - archive (arch)

## 4. Directories & Scripts

Copies of the GFS svn project trunk on various machines:

WCOSS: /global/save/emc.glopara/svn/gfs/trunk/para  
Zeus: /scratch2/portfolios/NCEPDEV/global/save/glopara/svn/gfs/trunk/para  
S4: /usr/local/jcsda/nwprod\_v2012  
SVN: <https://svnemc.ncep.noaa.gov/projects/gfs/trunk/para>

**NOTE:** It is not advised to run your experiments using the GFS trunk copies above, unless your experiment is really short or you check out a copy of the trunk yourself. The trunk is a moving target! It is best to run your experiments using a recent GFS tag.

**bin** - These scripts control the flow of an experiment

**pbeg** Runs when parallel jobs begin.  
**pcne** Counts non-existent files  
**pcon** Searches standard input (typically rlist) for given pattern (left of equal sign) and returns assigned value (right of equal sign).  
**pcop** Copies files from one directory to another.  
**pend** Runs when parallel jobs end.  
**perr** Runs when parallel jobs fail.  
**plog** Logs parallel jobs.  
**pmkr** Makes the rlist, the list of data flow for the experiment.  
**psub** Submits parallel jobs (check here for variables that determine resource usage, wall clock limit, etc).

**jobs** - These scripts, combined with variable definitions set in configuration, are similar in function to the wrapper scripts in /nwprod/jobs, and call the main driver scripts. E-scripts are part of the Hybrid EnKF.

**anal.sh** Runs the analysis. Default ex-script does the following:  
1) update surface guess file via global\_cycle to create surface analysis;  
2) runs the atmospheric analysis (global\_gsi);  
3) updates the angle dependent bias (satang file)

**arch.sh** Archives select files (online and hpss) and cleans up older data.

**copy.sh** Copies restart files. Used if restart files aren't in the run directory.

**dcop.sh** This script sometimes runs after dump.sh and retrieves data assimilation files.

**dump.sh** Retrieves dump files (not used in a typical parallel run).

**earc.sh** Archival script for Hybrid EnKF.  
1) Write select EnKF output to HPSS,  
2) Copy select files to online archive,  
3) Clean up EnKF temporary run directories,  
4) Remove "old" EnKF files from rotating directory.

**ecen.sh** Multiple functions:



- 1) Compute ensemble mean analysis from 80 analyses generated by eupd,
- 2) Perturb 80 ensemble analyses,
- 3) Compute ensemble mean for perturbed analyses,
- 4) Chgres T574L64 high resolution analysis (sanl/siganl) to ensemble resolution (T254L64),
- 5) Recenter perturbed ensemble analysis about high resolution analysis.

**echk.sh** Check script for Hybrid EnKF.  
 1) Checks on availability of ensemble guess files from previous cycle. (The high resolution (T574L64) GFS/GDAS hybrid analysis step needs the low resolution (T254L64) ensemble forecasts from the previous cycle);  
 2) Checks availability of the GDAS sanl (siganl) file (The low resolution (T254L64) ensemble analyses (output from eupd) are recentered about the high resolution (T574L64). This recentering can not be done until the high resolution GDAS analysis is complete.)

**efcs.sh** Run 9 hour forecast for each ensemble member. There are 80 ensemble members. Each efcs job sequentially processes 8 ensemble members, so there are 10 efcs jobs in total.

**efmn.sh** Driver (manager) for ensemble forecast jobs. Submits 10 efcs jobs and then monitors the progress by repeatedly checking status file. When all 10 efcs jobs are done (as indicated by status file) it submits epos.

**eobs.sh** Run GSI to select observations for all ensemble members to process. Data selection done using ensemble mean.

**eomg.sh** Compute innovations for ensemble members. Innovations computed by running GSI in observer mode. It is an 80 member ensemble so each eomg job sequentially processes 8 ensemble members.

**eomn.sh** Driver (manager) for ensemble innovations jobs. Submit 10 eomg jobs and then monitors the progress by repeatedly checking status file. When all 10 eomg jobs are done (as indicated by status file) it submits eupd.

**epos.sh** Compute ensemble mean surface and atmospheric mean ensemble files.

**eupd.sh** Perform EnKF update (i.e., generate ensemble member analyses).

**fcst.sh** Runs the forecast.

**prep.sh** Runs the data preprocessing prior to the analysis (storm relocation if needed and generation of prepbufr file).

**post.sh** Runs the post processor.

**vrify.sh** Runs the verification step.

**exp** - This directory typically contains config files for various experiments and some rlists.

Filenames with "config" in the name are configuration files for various experiments. Files ending in "rlist" are used to define mandatory and optional input and output files and files to be archived. For the most up-to-date configuration file that matches production see section 5.2.

**scripts** - Development versions of the main driver scripts. The production versions of these scripts are in /nwprod/scripts.

**ush** - Additional scripts pertinent to the model typically called from within the main driver scripts, also includes:

**reconcile.sh**      This script sets required, but unset variables to default values.

# 5. Data

## 5.1 Global Dump Archive

### 5.1.1 Location

An archive of global dump data is maintained in the following locations:

```
WCOS: /globaldump/YYYYMMDDCC
Zeus: /scratch2/portfolios/NCEPDEV/global/noscrub/dump/YYYYMMDDCC
S4: /usr/local/jcsda/dataset/global/YYYYMMDDCC
```

...where: YYYY = year, MM = month, DD = day, CC = cycle (00, 06, 12, or 18)

### 5.1.2 Grouping

The dump archive is divided into sub-directories:

- gdas[gfs] - main production dump data
- gdas[gfs]nr - non-restricted copies of restricted dump files
- gdas[gfs]x - experimental data, planned implementation
- gdas[gfs]y - experimental data, no planned implementation
- gdas[gfs]p - parallel dump data (short term)

Example of a typical 00z dump archive folder:

```
/global/save/emc.glopara/dump_archive[121]11 /globaldump/2014100100
total 512
drwxr-xr-x 2 emc.glopara global 131072 Oct 1 02:13 gdas
drwxr-xr-x 2 emc.glopara global 512 Oct 1 02:14 gdasnr
drwxr-xr-x 2 emc.glopara global 512 Oct 1 02:16 gdasx
drwxr-xr-x 2 emc.glopara global 512 Oct 1 02:16 gdasy
drwxr-xr-x 2 emc.glopara global 131072 Sep 30 23:07 gfs
drwxr-xr-x 2 emc.glopara global 512 Sep 30 23:08 gfsnr
drwxr-xr-x 2 emc.glopara global 512 Sep 30 23:09 gfsx
drwxr-xr-x 2 emc.glopara global 512 Sep 30 23:09 gfsy
```

### 5.1.3 Files & Availability

Time period covered by dump archive (as of 2/1/15): **2012010100 - present**

Not all data is available every day of the covered time period. The table in section 5.2.2 lists the files found within the global dump archive.

## 5.2 Input/output files

Many of the parallel files are in GRIB or BUFR formats, the WMO standard for gridded and ungridded meteorological data, respectively.

Other parallel files such as restart files are in flat binary format, and are not generally intended to be accessed by the general user.

Unfortunately but predictably, the global parallel follows a different file naming convention than the operational file naming convention. (The global parallel file naming convention started in 1990 and predates the operational file naming convention.)

The global parallel file naming convention is a file type followed by a period, the run (gdas or gfs), and the 10-digit current date \$CDATE in YYYYMMDDHH form:

FILETYPE.CDUMP.CDATE

(i.e. pgbf06.gfs.2008060400).

Some names may have a suffix, for instance if the file is compressed.

For the sake of users that are accustomed to working with production files or those who want to do comparisons, the equivalent production file name info is included here. Production file naming convention is the run followed by a period, the cycle name, followed by a period, and the file type. (i.e. gfs.t00z.pgrbf06). In the table below, only the file type is listed for production names.

The files are divided into the categories restart files, observation files, and diagnostic files. Some files may appear in more than one category. Some verification files in the diagnostics table do not include a run qualifier.

Guide to variables in sections 5.2.1, 5.2.2, and 5.2.3:

Variable	Description	Values
<b>\$CDUMP</b>	Dump type	gdas, gfs
<b>\$CDATE</b>	Cycle date	YYYYMMDDCC
<b>\$FF</b>	Forecast hour	00[000]-384
<b>\$FE</b>	Forecast hour (GDAS EnKF)	03, 06, 09
<b>\$MEM</b>	Hybrid EnKF member number	001-080
<b>\$GRP</b>	Hybrid EnKF member group number	01-10

## 5.2.1 Restart / Initial Condition (IC) Files

glopara filename	production base name (eg, gdas1.t00z.prepbuf)	file description	format
biascr.\$CDUMP.\$CDATE	abias	Information about sensor/instrument/satellite, channel, tlapmean, and bias predictor coefficients	text
biascr_pc.\$CDUMP.\$CDATE	abias_pc	Information about observation number and the estimates of analysis error variances for bias predictor coefficients.	text
bfg_\$CDATE_fhr\$FE_ensmean	<i>same as glopara filename</i>	Mean of ensemble surface forecasts at fhr\$FE	binary
bfg_\$CDATE_fhr\$FE_mem\$MEM	<i>same as glopara filename</i>	Surface forecast at fhr\$FE for member \$MEM starting from \$CDATE ICs	binary
pgbanl.\$CDUMP.\$CDATE	pgrbanl	pressure level data from analysis	GRIB2
pgbl\$FF.\$CDUMP.\$CDATE	pgrb2.2p50.f\$FF	2.5° pressure level data from forecast	GRIB2
pgbf\$FF.\$CDUMP.\$CDATE	pgrb2.1p00.f\$FF	1° pressure level data from forecast	GRIB2
pgbh\$FF.\$CDUMP.\$CDATE	pgrb2.0p50.f\$FF	0.5° pressure level data from forecast	GRIB2
pgbq\$FF.\$CDUMP.\$CDATE	pgrb2.0p25.f\$FF	0.25° pressure level data from forecast	GRIB2
pgbe\$FF.\$CDUMP.\$CDATE	TBD – not yet implemented	0.125° pressure level data from forecast	GRIB2
prepqc.\$CDUMP.\$CDATE	prepbuf	Conventional Observations with quality control	BUFR
radstat.\$CDUMP.\$CDATE	radstat	Radiance assimilation statistics	binary
sfcanl.\$CDUMP.\$CDATE	sfcanl	surface analysis	binary
sfcanl_\$CDATE_ensmean	<i>same as glopara filename</i>	mean of ensemble surface ICs valid at \$CDATE	binary
sfcanl_\$CDATE_mem\$MEM	<i>same as glopara filename</i>	Surface ICs for member \$MEM valid at \$CDATE; input to ensemble forecasts	binary
siganl.\$CDUMP.\$CDATE	sanl	atmospheric analysis (aka sigma file)	binary
sanl_\$CDATE_ensmean	<i>same as glopara filename</i>	Mean of ensemble atmospheric analyses generated by EnKF update code valid at \$CDATE	binary
sanl_\$CDATE_mem\$MEM	<i>same as glopara filename</i>	Atmospheric analyses generated by EnKF update code for member \$MEM valid at \$CDATE	binary
sfcf\$FF.\$CDUMP.\$CDATE	bf\$FF	surface boundary condition at forecast hour \$FF	binary
sfg_\$CDATE_fhr\$FE_ensmean	<i>same as glopara filename</i>	Mean of ensemble atmospheric forecasts at fhr\$FE	binary
sfg_\$CDATE_fhr\$FE_mem\$MEM	<i>same as glopara filename</i>	Atmospheric forecast at fhr\$FE for member \$MEM starting from \$CDATE ICs	binary
sfg_\$CDATE_fhr\$FEs_mem\$MEM	<i>same as glopara filename</i>	Spectrally smoothed atmospheric forecast at fhr\$FE for member \$MEM starting from \$CDATE ICs	binary
sig\$FF.\$CDUMP.\$CDATE	sf\$FF	atmospheric model data at forecast hour \$FF	binary
siganl_\$CDATE_mem\$MEM	<i>same as glopara filename</i>	Atmospheric ICs for member \$MEM valid at \$CDATE at END of ecen; input to ensemble forecasts	binary

## 5.2.2 Observation files

glopara filename (FILE.\$CDUMP.\$CDATE, unless otherwise noted)	production base name (eg.gdas1.t00z.engicegrb)	file description	format
1bamua	1bamua.tm00.bufr_d	AMSU-A NCEP-proc. br. temps	BUFR
1bhrs4	1bhrs4.tm00.bufr_d	HIRS-4 1b radiances	BUFR
1bmhs	1bmhs.tm00.bufr_d	MHS NCEP-processed br. temp	BUFR
adpsfc	adpsfc.tm00.bufr_d	Surface land	BUFR
adpupa	adpupa.tm00.bufr_d	Upper-air	BUFR
aircar	aircar.tm00.bufr_d	MDCRS ACARS Aircraft	BUFR
aircft	aircft.tm00.bufr_d	Aircraft	BUFR
airsev	airsev.tm00.bufr_d	AQUA-AIRS AIRS/AMSU-A/HSB proc. btemps- every FOV	BUFR
ascatt	ascatt.tm00.bufr_d	METOP-2 ASCAT products (not superobed)	BUFR
ascatw	ascatw.tm00.bufr_d	METOP 50 KM ASCAT scatterometer data (reprocessed by wave_dcodquikscat)	BUFR
atms	atms.tm00.bufr_d	NPP Adv. Tech. Microwave Sounder (ATMS) radiances	BUFR
avcsam	avcsam.tm00.bufr_d	A.M.(N17,M2) AVHRR GAC NCEP-proc clr & sea btemps	BUFR
avcspm	avcspm.tm00.bufr_d	P.M.(N18-19) AVHRR GAC NCEP-proc clr & sea btemps	BUFR
bathy	bathy.tm00.bufr_d	Bathothermal	BUFR
cris	cris.tm00.bufr_d	NPP Cross-track Infrared Sounder (CrIS) radiances	BUFR
esamua	esamua.tm00.bufr_d	NOAA 15-19 AMSU-A proc. bright. temps from RARS	BUFR
eshrs3	eshrs3.tm00.bufr_d	NOAA 15-19 HIRS-3/-4 proc bright. temps from RARS	BUFR
esmhs	esmhs.tm00.bufr_d	NOAA 18-19 MHS processed bright. temps from RARS	BUFR
geoimr	geoimr.tm00.bufr_d	GOES 11x17 fov imager clear radiances	BUFR
goesfv	goesfv.tm00.bufr_d	GOES 1x1 fov sounder radiances	BUFR
gome	gome.tm00.bufr_d	METOP-2 Global Ozone Monitoring Exp.-2 (GOME-2)	BUFR
gpsipw	gpsipw.tm00.bufr_d	GPS - Integrated Precipitable Water	BUFR
gpsro	gpsro.tm00.bufr_d	GPS radio occultation data	BUFR
icegrb	engicegrb	Sea Ice Analysis	GRIB
imssnow96.grib2	imssnow96.grib2	IMS NH snow and ice cover analysis. 96 <sup>th</sup> mesh (or 4km) resolution.	GRIB2
mls	mls.tm00.bufr_d	Aura Microwave Limb Sounder (MLS) ozone data	BUFR
mtiasi	mtiasi.tm00.bufr_d	METOP-2 IASI 1C radiance data (variable channels)	BUFR
NPR.SNWN.SP.S1200.MESH16.grb	<i>same as glopara file</i>	AFWA NH snow depth analysis. 16 <sup>th</sup> mesh (or 23 km) resolution.	GRIB1
NPR.SNWS.SP.S1200.MESH16.grb	<i>same as glopara file</i>	AFWA SH snow depth analysis. 16 <sup>th</sup> mesh (or 23 km) resolution.	GRIB1
obsinput_\$CDATE_ensmean	<i>same as glopara file</i>	Tarball containing \$CDATE data (observations) selected using ensemble means; generated by eobs	tarball
omi	omi.tm00.bufr_d	Aura Ozone Monitoring Instrument (OMI) data	BUFR
osbuvs8	osbuvs8.tm00.bufr_d	SBUV layer ozone product (Version 8)	BUFR
proflr	proflr.tm00.bufr_d	Wind Profiler	BUFR
rassda	rassda.tm00.bufr_d	Radio Acoustic Sounding System Temp Profiles	BUFR
rtgssthr.grb[grib2]	rtgssthr.grb[grib2]	Global 5-minute RTG SST analysis	GRIB[2]
satwnd	satwnd.tm00.bufr_d	Satellite-derived wind reports	BUFR

seaice.5min.[grb][grib2]	seaice.5min.[grb][grib2]	EMC global 5-minute ice concentration analysis	GRIB
seaice.5min.blend.grb	seaice.5min.blend.grb	Global blended sea ice concentration analysis at 5-minute resolution. A blend of the EMC 5-min ice analysis and the 4km IMS ice cover analysis.	GRIB1
sfcshp	sfcshp.tm00.bufr_d	Surface marine	BUFR
snogrb	snogrb	Global 0.5-degree Snow cover and snow liquid equivalent analysis	GRIB1
snogrb_t###.\$LONB.\$LATB	Generated in dump step using a blend of the 4km IMS snow cover and the 23 km AFWA snow depth.	Snow depth and snow cover analysis on spectral t### grid (# = resolution, i.e. 1534)	GRIB1
ssmisu	ssmisu.tm00.bufr_d	DMSP SSM/IS 1C radiance data (Unified Pre-Proc.)	BUFR
sstgrb	sstgrb	Global 1.0-degree Sea Surface Temperature Analysis	GRIB1
statup	updated.status.tm00.bufr_d	Summary	text
stat01	status.tm00.bufr_d	Bufr status	text
tcvttl	syndata.tcvitals.tm00	Tropical Storm Vitals	text
tesac	tesac.tm00.bufr_d	TESAC	BUFR
trkob	trkob.tm00.bufr_d	TRACKOB	BUFR
vadwnd	vadwnd.tm00.bufr_d	VAD (NEXRAD) wind	BUFR

For more information on dump data types (as seen in production) visit this site:  
<http://www.nco.ncep.noaa.gov/pmb/nwprod/realtime/index.bufrdump.shtml>

### 5.2.3 Diagnostic files

<b>glopara filename</b>	<b>production base name (eg,gdas1.t00z.gsistat)</b>	<b>file description</b>	<b>format</b>
adpsfc.anl.\$CDATE		Surface observation and analysis fit file	GrADS
adpsfc.fcs.\$CDATE		Surface observation and forecast fit file3	GrADS
adpupa.mand.anl.\$CDATE		Rawinsonde observation and analysis fit file	GrADS
adpupa.mand.fcs.\$CDATE		Rawinsonde observation and forecast fit file3	GrADS
gsistat.\$CDUMP.\$CDATE	gsistat	GSI (obs-ges), qc, and iteration statistics	text
gsistat_.\$CDATE_ensmean	<i>same as glopara file</i>	gsistat file for \$CDATE; based on data selection run (eobs) using ensemble mean background fields	text
gsistat_.\$CDATE_mem\$MEM	<i>same as glopara file</i>	gsistat file for member \$MEM for \$CDATE	text
radstat_.\$CDATE_ensmean	<i>same as glopara file</i>	Radiance diagnostic file with \$CDATE observations; generated by eobs (data selection using ensemble mean)3	binary
radstat_.\$CDATE_mem\$MEM	<i>same as glopara file</i>	Radiance diagnost file for member \$MEM with \$CDATE observations	binary
cnvstat.\$CDUMP.\$CDATE	cnvstat	Conventional observation assimilation statistics	binary
cnvstat_.\$CDATE_ensmean	<i>same as glopara file</i>	Conventional diagnostic file with \$CDATE observations; generated by eobs (data selection using ensemble mean)	binary
cnvstat_.\$CDATE_mem\$MEM	<i>same as glopara file</i>	Conventional diagnostic file for member \$MEM with \$CDATE observations	binary
enkfstat_.\$CDATE	<i>same as glopara file</i>	EnKF update code stdout for \$CDATE	text
ensstat_.\$CDATE_all	<i>same as glopara file</i>	Log file denoting completion of averaging of ensemble forecasts (epos step) for \$CDATE	text
fcsstat_.\$CDATE_all	<i>same as glopara file</i>	Log file for denoting completion of all \$CDATE ensemble forecasts	text
fcsstat_.\$CDATE_grp\$GRP	<i>same as glopara file</i>	Log file for completion of group \$GRP ensemble forecasts for \$CDATE	text
flxf\$FF.\$CDUMP.\$CDATE	fluxgrbf\$FF	Model fluxes at forecast hour \$FF	GRIB

logf\$FF.\$CDUMP.\$CDATE	logf\$FF	Model logfile at forecast hour \$F	text
omgstat_.\$CDATE_all	<i>same as glopara file</i>	Log file denoting completion of all \$CDATE ensemble innovation jobs	text
omgstat_.\$CDATE_grp\$GRP	<i>same as glopara file</i>	Log file for completion of group \$GRP ensemble innovation job for \$CDAT	text
oznstat.\$CDUMP.\$CDATE	oznstat	Ozone observation assimilation statistics	binary
oznstat_.\$CDATE_ensmean	<i>same as glopara file</i>	Ozone diagnostic file with \$CDATE observations; generated by eobs (data selection using ensemble mean)	binary
oznstat_.\$CDATE_mem\$MEM	<i>same as glopara file</i>	Ozone diagnost file for member \$MEM with \$CDATE observations3	binary
pertdates_.\$CDATE	pertdates_.\$CDATE	Dates from from pertubation database used in \$CDATE additive inflation step (ecen	text
pcpstat.\$CDUMP.\$CDATE	pscpstat	Precipitation assimilation statistics	binary
prepqa.gdas.\$CDATE		Observations with QC plus analysis	BUFR
prepqc.\$CDUMP.\$CDATE	prepbufr	Conventional Observations with QC	BUFR
prepqf.gdas.\$CDATE		Observations with QC plus forecast	BUFR
radstat.\$CDUMP.\$CDATE	radstat	Radiance assimilation statistics	binary
sfshp.anl.\$CDATE		Ship observation and analysis fit file3	GrADS
sfshp.fcs.\$CDATE		Ship observation and forecast fit file	GrADS
tcinform_relocate.\$CDUMP.\$CDATE		Storm relocation information	text
tcvitals_relocate.\$CDUMP.\$CDATE		tropical cyclone vitals	text



### 5.3 Finding GDAS and GFS production run files

Locations below use the following:

```
YYYY = 4-digit year of run date
MM   = 2-digit month of run date
DD   = 2-digit day of run date
CC   = run cycle (00, 06, 12 18) in UTC
```

NCO maintains files for the last 10 days in WCOSS directories:

```
/com/gfs/prod/gdas.YYYYMMDD
/com/gfs/prod/gfs.YYYYMMDD
/com/gfs/prod/enkf.YYYYMMDD/CC
```

...and the last two days in Zeus directories:

```
/scratch2/portfolios/NCEPDEV/rstprod/com/gfs/prod/gdas.YYYYMMDD
/scratch2/portfolios/NCEPDEV/rstprod/com/gfs/prod/gfs.YYYYMMDD
/scratch2/portfolios/NCEPDEV/rstprod/com/gfs/prod/enkf.YYYYMMDD/CC
```

Locations of production files on HPSS (tape archive):

```
/NCEPPROD/hpssprod/runhistory/rhYYYY/YYYYMM/YYYYMMDD/
```

Example for December 16<sup>th</sup>, 2013 00Z (2013121600):

```
/NCEPPROD/hpssprod/runhistory/rh2013/201312/20131216
```

```
gdas files:    com_gfs_prod_gdas.2013121600.tar      (contains fcst ICs)
gfs files:    com_gfs_prod_gfs.2013121600.anl.tar    (contains fcst ICs)
               com_gfs_prod_gfs.2013121600.pgrb2.tar
enkf files:    com_gfs_prod_enkf.20131216_00.anl.tar    (contains efmn ICs)
               com_gfs_prod_enkf.20131216_00.fcs.tar
               com_gfs_prod_enkf.20131216_00.fcs03.tar
               com_gfs_prod_enkf.20131216_00.fcs09.tar
               com_gfs_prod_enkf.20131216_00.omg.tar
```

Example pulling ICs off HPSS for a fully-cycled GFS run with the Hybrid EnKF starting at 2013121600 gdas:

```
hpsstar get
/NCEPPROD/hpssprod/runhistory/rh2013/201312/20131216/com_gfs_prod_gdas.201312
1600.tar ./gdas1.t00z.abias ./gdas1.t00z.satang ./gdas1.t00z.sanl
./gdas1.t00z.sfcanl
hpsstar get
/NCEPPROD/hpssprod/runhistory/rh2013/201312/20131216/com_gfs_prod_enkf.201312
16_00.anl.tar
```

See "Notes" on next page...

## NOTES:

Make sure to rename the `gdas.t00z.*` files you pull from the first tarball. Those files need to be in the parallel naming convention (see previous page).

The 2nd command will pull all of the contents of that EnKF tarball. This is MUCH faster than trying to list all 160 ICs you'd need to pull from that tarball. You'll get a few extra files you don't need but oh well.

## 6. System Settings

### 6.1 Grid dimensions

<b>SPECTRAL RESOLUTION</b>	<b>EULERIAN</b>		<b>SEMI-LAGRANGIAN</b>	
	<b>LONB</b>	<b>LATB</b>	<b>LONB</b>	<b>LATB</b>
<b>T62</b>	192	94	128	64
<b>T126</b>	384	190	256	128
<b>T170</b>	512	256	352	176
<b>T190</b>	576	288	384	192
<b>T254</b>	768	384	512	256
<b>T382</b>	1152	576	768	384
<b>T574</b>	1760	880	1152	576
<b>T878</b>	2304	1152	1760	880
<b>T1148</b>			2304	1152
<b>T1534</b>			3072	1536
<b>T2014</b>			4032	2016
<b>T2046</b>			4096	2048
<b>T3070</b>			6144	3072

### 6.2 Global Model Variables

To view the full list of global model variables please see Appendix A.

## 7. Setting up an experiment

Steps:

1. Is your environment setup correctly? If you're not sure, check out the "Setting up your environment" section below.
2. Do you have restricted data access? If not go to:  
[http://www.nco.ncep.noaa.gov/sib/restricted\\_data/restricted\\_data\\_sib/](http://www.nco.ncep.noaa.gov/sib/restricted_data/restricted_data_sib/)  
and submit a registration form to be added to group rstprod.
3. Important terms
4. Set up experiment configuration file
5. Set up rlist
6. Submit first job

Additional information in this section:

1. Plotting model output
2. Experiment troubleshooting
3. Related utilities

### 7.1 Important terms

- **configuration file** - List of variables to be used in experiment and their configuration/value. The user can change these variables for their experiment. Description of variables.
- **job** - A script, combined with variable definitions set in configuration, which is similar in function to the wrapper scripts in /nwprod/jobs, and which calls the main driver scripts. Each box in above diagram is a job.
- **reconcile.sh** - Similar to the configuration file, the reconcile.sh script sets required, but unset variables to default values.
- **rlist** - List of data to be used in experiment. Created in reconcile.sh (when the pmkr script is run) if it does not already exist at beginning of experiment. More information on setting up your own rlist see section 5.4.
- **rotating directory (ROTDIR)** - Typically your "noscrub" directory is where the data and files from your experiment will be stored. Example on Zeus: /scratch2/portfolios/NCEPDEV/global/noscrub/\$LOGNAME/pr\$PSLOT

### 7.2 Setting up your environment

For successful GFS model runs it is important that your supercomputer environment be setup correctly. If you are unsure of what PATHs need setting, modules loaded, etc. then take a peek at the following .profile and .bashrc/.cshrc files:

MACHINE	.profile	.bashrc
WCOSS	/u/Kate.Howard/.profile	/u/Kate.Howard/.bashrc
Zeus	/home/Kate.Howard/.profile	/home/Kate.Howard/.cshrc

## 7.3 Configuration file

The following files have settings that will produce results that match production results. Copy this file, or any other configuration file you wish to start working with, to your own space and modify it as needed for your experiment.

MACHINE	LOCATION	FILE NAME	WHAT
<b>WCOSS</b>	/global/save/emc.glopara/svn/gfs/tags/... TAG_OF_CHOICE... /para/exp/	para_config_T254	Q1FY15 at T254
		para_config_T574	^ same at T574
		para_config_T670_T254	^ same at T670
		para_config_T1534*	^ same at T1534*
<b>Zeus</b>	/scratch2/portfolios/NCEPDEV/global/save/ glopara/svn/gfs/tags/TAG_OF_CHOICE/para/exp	Same as WCOSS	
<b>S4</b>	/home/khoward/GFS_tutorial/prtest	para_config	May 2012 version

\* Mimics production

Make sure to check the following user specific configuration file variables, found near the top of the configuration file:

<b>ACCOUNT</b>	LoadLeveler account, i.e., GFS-MTN (see more examples below for ACCOUNT, CUE2RUN, and GROUP)
<b>ARCDIR</b>	Online archive directory (i.e. ROTDIR/archive/prPSLOT)
<b>ATARDIR</b>	HPSS tape archive directory (see configuration file for example)
<b>CUE2RUN</b>	LoadLeveler (or Moab) class for parallel jobs (i.e., dev) (see more examples of CUE2RUN below)
<b>EDATE</b>	Analysis/forecast cycle ending date (YYYYMMDDCC, where CC is the cycle)
<b>EDUMP</b>	Cycle ending dump (gdas or gfs)
<b>ESTEP</b>	Cycle ending step (prep, anal, fcst1, post1, etc.)
<b>EXPDIR</b>	Experiment directory under save, where your configuration file, rlist, runlog, and other experiment scripts sit.
<b>GROUP</b>	LoadLeveler group (i.e., g01) (see more examples of GROUP below)
<b>PSLOT</b>	Experiment ID (change this to something unique for your experiment)
<b>ROTDIR</b>	Rotating/working directory for model data and i/o (i.e. /global/noscrub/\$LOGNAME/pr\$PSLOT)

## 7.4 Reconcile.sh

If concerned, make sure to take a look at the current reconcile script to assure that any changes you made in the configuration file are not overwritten. The reconcile script runs after reading in the configuration file settings and sets default values for many variables that may or may not be defined in the configuration file. If there are any default choices in reconcile that are not ideal for your experiment make sure to set those variables in your configuration file, perhaps even at the end of the file after reconcile has been run.

## 7.5 Rlist

You can start with an existing rlist and modify it by hand as needed or grab the sample that exists in the exp subdirectory of the tag (or other release) you wish to run (RECOMMENDED):

SVN: <https://svnemc.ncep.noaa.gov/projects/gfs/trunk/para/exp/prsample1.gsi.rlist>

The sample rlist files already contain the append.rlist entries.

If the rlist file does not exist when a job is submitted, pmkr will generate one based on your experiment configuration. However, it is currently advised that you **do not** use pmkr to create an rlist, but rather, pick up the sample rlist.

If the variable \$ARCHIVE is set to YES (the default is NO), this file is then appended automatically to the rlist by reconcile.sh, but only when the rlist is generated on the fly by pmkr. So, eg, if you submit the first job, which creates an rlist and then you realize that your ARCx entries are missing, creating the append\_rlist after the fact won't help unless you remove the now existing rlist. If you delete the errant rlist (and set \$ARCHIVE to YES, the next job you submit will see that the rlist does not exist, create it using pmkr, then append the \$append\_rlist file.

Also, along those lines, you may find that pmkr does not account for some new or development files. You can list those needed entries in the file pointed to by variable \$ALIST. The difference between \$ALIST and \$append\_rlist is that the latter only gets appended if variable \$ARCHIVE is YES.

Got all that?? (Now you know why it is sometimes easier to start with an existing rlist).

Brief overview of an rlist format:

Sample entries:

```
# rotational input
**/anal/ROTI    =      biascr.$GDUMP.$GDATE
**/anal/ROTI    =      satang.$GDUMP.$GDATE
**/anal/ROTI    =      sfcf06.$GDUMP.$GDATE
**/anal/ROTI    =      prepqc.$CDUMP.$CDATE
# optional input
**/anal/OPTI    =      sfcf03.$GDUMP.$GDATE
**/anal/OPTI    =      sfcf04.$GDUMP.$GDATE
**/anal/OPTI    =      sfcf05.$GDUMP.$GDATE
**/anal/OPTI    =      sfcf07.$GDUMP.$GDATE
**/anal/OPTI    =      sfcf08.$GDUMP.$GDATE
```

The left hand side is set of 4 patterns separated by slashes.

The first pattern represents the cycle (full date)

The second pattern represents the dump.

The third pattern represents the job.

The fourth pattern is a string that defines whether a file is optional/required input/output, eg:

DMPI - dump input from current cycle

DMPG - dump input from previous cycle  
 DMPH - dump input from two cycles prior  
 ROTI - required input from the rotating directory  
 OPTI - optional input from the rotating directory  
 ROTO - required output to the rotating directory (if the file is not available, a flag is set and the next job is not triggered)  
 OPTO - optional output to the rotating directory (save it if available, no worries if it's not)  
 ARCR - files to archive in online archive (should be required, but depends on setup of arch.sh)  
 ARCO - files to archive in online archive  
 ARCA - files saved to "ARCA" HPSS archive  
 ARCB - files saved to "ARCB" HPSS archive (check arch.sh job for other HPSS options... current version allows for ARCA thru ARCF)  
 COPI - required restart and files to initiate experiment with copy.sh job (fcst input)  
 DMRI - prerequisite dump file for submit (used in psub, but not used in job scripts to copy data!)

The right hand side typically represents a file.

An asterisk on either side is a wild card. Eg:

```
*/*/arch/ARCR = pgbf06.$CDUMP.$CDATE
```

The above entry in your rlist means that for any cycle, or any dump, the archivejob will copy pgbf06.\$CDUMP.\$CDATE to the online archive.

If you change that to:

```
*/gfs/arch/ARCR = pgbf06.$CDUMP.$CDATE
```

only the the gfs pgbf06 files will be copied to the online archive.

If you changed it to:

```
*00/gfs/arch/ARCR = pgbf06.$CDUMP.$CDATE
```

only the 00Z gfs pgbf06 files will be copied to the online archive.

If you changed it to:

```
20080501*/gfs/arch/ARCR = pgbf06.$CDUMP.$CDATE
```

only the May 1, 2008 gfs pgbf06 files will be copied to the online archive. (Not a likely choice, but shown as an example)

Changing that first example to:

```
*/*/arch/ARCR = pgbf*.$CDUMP.$CDATE
```

tells the archive job to copy the the pgb file for any forecast hour (from the current \$CDUMP and \$CDATE) to the online archive.

A more complex set of wildcards can be useful for splitting up the HPSS archive to keep tar files manageable. Eg:

```
# all gdas sigma files go to ARCA HPSS archive
*/gdas/arch/ARCA = sigf*.$CDUMP.$CDATE
```

```
# gfs sigf00 thru sigf129 go to ARCB HPSS archive
*/gfs/arch/ARCB = sigf??. $CDUMP.$CDATE
*/gfs/arch/ARCB = sigf1[0-2]?.$CDUMP.$CDATE
```

```
# gfs sigf130 thru sigf999 go to ARCC HPSS archive
*/gfs/arch/ARCC = sigf1[3-9]?.$CDUMP.$CDATE
*/gfs/arch/ARCC = sigf[2-9]??. $CDUMP.$CDATE
```

## 7.6 Initial Conditions / Required Forcing Files

The following files are needed to run the GFS/GDAS:

	PARALLEL	PRODUCTION
<b>NON-CYCLING / FREE FORECAST</b>	sfc anl.\$CDUMP.\$CDATE	gdas1.tCCz.sfc anl
	sig anl.\$CDUMP.\$CDATE	gdas1.tCCz.sanl
<b>CYCLING w/o HYBRID ENKF</b>	*biascr.\$CDUMP.\$CDATE	gdas1.tCCz.abias
	*biascr_pc.\$CDUMP.\$CDATE	gdas1.tCCz.abias_pc
	*radstat.\$CDUMP.\$CDATE	gdas1.tCCz.radstat
	sfc anl.\$CDUMP.\$CDATE	gdas1.tCCz.sfc anl
	sig anl.\$CDUMP.\$CDATE	gdas1.tCCz.sanl
	*biascr.\$CDUMP.\$CDATE	gdas1.tCCz.abias
<b>CYCLING w/ HYBRID ENKF</b>	*biascr_pc.\$CDUMP.\$CDATE	gdas1.tCCz.abias_pc
	*radstat.\$CDUMP.\$CDATE	gdas1.tCCz.radstat
	sfc anl.\$CDUMP.\$CDATE	gdas1.tCCz.sfc anl
	sig anl.\$CDUMP.\$CDATE	gdas1.tCCz.sanl
	sig anl_\$CDATE_mem\$MEM	sig anl_\$CDATE_mem\$MEM
	sfc anl_\$CDATE_mem\$MEM	sfc anl_\$CDATE_mem\$MEM

Where CC is the cycle (00, 06, 12, or 18 Z) & \$MEM is the member number (001-080)

\* Prior to Q1FY15 implementation these ICs should be pulled from pre-implementation parallels

So where do I find initial conditions (ICs)? See section 5.3



## 8. Submitting & running your experiment

1. Create directory \$EXPDIR (defined in configuration file)
2. Place a configuration file and rlist into \$EXPDIR
3. Create directory \$ROTDIR (defined in configuration file)
4. Copy required initial condition / forcing files into \$ROTDIR
5. Make the necessary edits to your configuration file to match the kind of experiment you wish to run (see section 7.3). Make sure to rename your rlist to match your experiment PSLOT (i.e. pr\$PSLOT1.gsi.rlist).
6. Then, it's time to submit! On command line type:

**\$PSUB \$CONFIG \$CDATE \$CDUMP \$CSTEP**

Where:

**\$PSUB** = psub script with full location path. It is always recommended to use the psub script from within the tag (or other release) you plan to run. The psub script currently works on both WCOSS and Zeus.

**\$CONFIG** = name of configuration file (with full location path if not submitting from within your \$EXPDIR)

**\$CDATE** = YYYYMMDDCC, initial/starting year (YYYY), month (MM), day (DD), and cycle (CC) for model run

**\$CDUMP** = dump (gdas or gfs) to start run

**\$CSTEP** = initial model run step (see flow diagram above for options)

Example on WCOSS:

```
/global/save/emc.glopara/svn/gfs/trunk/para/bin/psub para_config_T1534 2015011412 gdas fcst1
```

Notes:

- If you wish to cycle AND run the Hybrid EnKF then you need to submit both the fcst1 and efmn steps at the beginning.
- If you do not wish to cycle OR you do not wish to run the Hybrid EnKF then start with just the gdas fcst1 step.
- If you just wish to run a GFS free-forecast, start with the gfs fcst1 step.
- If you have a submit script that you are comfortable with then please feel free to use that to submit your experiment instead of the psub command, which should already be built into the submit script.

Additional information about running an experiment:

- The script "psub" kicks off the experiment and each parallel sequenced job.
- Remember that since each job script starts the next job, you need to define ESTEP as the job that follows the step with which you wish to end on. For example: You want to finish when your final planned cycle completes...your ESTEP could be "prep", which is the first step of the next cycle. Typically EDUMP is gdas...which means that if gfs\_cyc > 0 the next gfs cycle may be submitted even though it is the cycle after the end of your experiment.

A handy way to follow the status of your experiment is to do a tail of your runlog in your \$EXPPDIR directory:

```
tail -f pr$PSLOT.runlog          (where $PSLOT is your experiment tag)
```

## 8.1 Plotting output

Everyone has a favorite plotting program but one great option is GrADS. To use GrADS you'll first need to create a control file from your GRIB output:

1. Create GrADS readable ctl file using grib2ctl script:

Find copy here: /u/Wesley.Ebisuzaki/bin/grib2ctl.pl (WCOSS)

To run: **GRIB2CTL** [options] **INPUT** > **OUTPUT.ctl**

**GRIB2CTL** = full path of grib2ctl.pl or simply grib2ctl.pl if it's already in your environment

**INPUT** = the full name and path of the GRIB file

**OUTPUT** = the name of the ctl file you wish to create

[options] = full list of options can be found if you type "grib2ctl.pl" and hit enter. If you are making a ctl file from a forecast file then it is suggested to use the -verf option.

2. Create index file using gribmap:

```
gribmap -i OUTPUT.ctl
```

You should now have .ctl and .idx files.

3. Open GrADS (**grads** or **gradsc**) and then open your ctl file (**open OUTPUT.ctl**)

For information on using GrADS go here: <http://www.iges.org/grads/gadoc/>

## 8.2 Experiment troubleshooting

Machine issues? Contact appropriate helpdesk:

WCOSS - [wcooss-helpdesk@noaa.gov](mailto:wcooss-helpdesk@noaa.gov)  
Zeus - [rdhpcs.zeus.help@noaa.gov](mailto:rdhpcs.zeus.help@noaa.gov)  
S4 - [s4.admin@ssec.wisc.edu](mailto:s4.admin@ssec.wisc.edu)

As model implementations occur, ensure that you are using up-to-date versions of scripts/code and configuration file for your experiment. For instance, don't use the newest production executables with older job scripts. Changes may have been made to the production versions that will impact your experiment but may not be obvious.

For problems with your experiment please contact Kate Howard: [kate.howard@noaa.gov](mailto:kate.howard@noaa.gov)

**Please make sure to provide the following information in the email:**

- Machine you are working on (WCOSS, Zeus, or S4)
- Configuration file name and location
- Any other specific information pertaining to your problem, i.e., dayfile name and/or location.

## 8.3 Tutorials

Tutorials are available for users, who are perhaps unfamiliar with both the machine and the GFS. See the following table for tutorial information:

<u>Machine</u>	<u>Location</u>
<b>WCOSS</b>	/global/save/emc.glopara/TUTORIAL
<b>Zeus</b>	/scratch2/portfolios/NCEPDEV/global/save/glopara/TUTORIAL

## 9. Parallels

View the Global Parallel Spreadsheet here:

<https://docs.google.com/a/noaa.gov/spreadsheet/ccc?key=0AoyO6L08rs23dE9HdFhqa25YdUVyNUVZWTVrY01EeWc#gid=0%7C>

## 10. Subversion & Trac

GFS Trac page - <https://svnemc.ncep.noaa.gov/trac/gfs>  
SVN project page - <https://svnemc.ncep.noaa.gov/projects/gfs/>

GSM Trac page - <https://svnemc.ncep.noaa.gov/trac/gsm>  
SVN project page - <https://svnemc.ncep.noaa.gov/projects/gsm/>

## 11. Related utilities

Information on some useful related utilities:

**copygb** copies all or part of one GRIB file to another GRIB file, interpolating if necessary

**sfchdr** global\_sfchdr prints information from the header of a surface file

**sighdr** global\_sighdr prints information from the header of a sigma file

**ss2gg** ss2gg converts a sigma file to a grads binary file and creates a corresponding descriptor (ctl) file

### 11.1 copygb

The command copygb copies all or part of one GRIB file to another GRIB file, interpolating if necessary.

copygb can be found at: /nwprod/util/exec/copygb

Documentation is in: /nwprod/util/sorc/copygb.fd/copygb.doc

The NCEP grids for the -g option are listed in:

<http://www.nco.ncep.noaa.gov/pmb/docs/on388/tableb.html>

Documentation for the interpolation options are covered in: /nwprod/lib/sorc/ip/iplib.doc (though some parts may be outdated).

If you want to dig into any "w3" subroutines referenced, they generally have good docblocks in their source code. The directory is /nwprod/lib/sorc/w3 and there's a web doc at [http://www.nco.ncep.noaa.gov/pmb/docs/libs/w3lib/ncep\\_w3lib.shtml](http://www.nco.ncep.noaa.gov/pmb/docs/libs/w3lib/ncep_w3lib.shtml)

## 11.2 sfchr

global\_sfchr prints information from the header of a surface file

global\_sfchr can be found at:

```
/nwprod/exec/global_sfchr
```

```
Usage: global_sfchr sfcfile <variable.list >value.list
       or global_sfchr sfcfile variable >value
       or global_sfchr sfcfile
```

Running sfchr with no additional arguments (other than the input file) as in the last example allows for keyboard input of multiple variables, one at a time, until the program is interrupted (eg, via CTRL-c).

Enter "?" (without the quotes) as standard input and the possible input values will be printed.

Description of those possible values follows:

```
filetype - description ("GFS/SFC")
fhour    - forecast hour
ifhr     - integral forecast hour as string
idate    - initial date (YYYYMMDDHH)
iyr      - initial year
imo      - initial month
idy      - initial day
ihr      - initial hour
vdate    - valid date (YYYYMMDDHH)
vyr      - valid year
vmo      - valid month
vdy      - valid day
vhr      - valid hour
latb     - number of latitudes
lonb     - number of longitudes
ivs      - version number
lsoil    - number of soil levels
irealf   - floating point flag (=1 for 4-byte ieee, =2 for 8-byte ieee)
lpl      - number of longitudes for each latitude
zsoil    - soil depths (in meters)
```

## 11.3 sighdr

global\_sighdr prints information from the header of a sigma file

global\_sighdr can be found at:

```
/nwprod/exec/global_sighdr
```

```
Usage: global_sighdr sigfile <variable.list >value.list
       or global_sighdr sigfile variable >value
```

The following is from the docblock of /nwprod/sorc/global\_sighdr.fd/sighdr.f

```
program sighdr
!$$$ main program documentation block
```

```
!
! Main program: sighdr          Print information from sigma header
!   Prgmnr: Iredell           Org: np23           Date: 1999-08-23
!
! Abstract: This program prints information from the sigma header.
!   The following parameters may be printed out:
!     filetype
!     fhour
!     ifhr
!     idate
!     iyr
!     imo
!     idy
!     ihr
!     vdate
!     vyr
!     vmo
!     vdy
!     vhr
!     si
!     sl
!     ak
!     bk
!     siglev
!     jcap
!     levs
!     itrun
!     iorder
!     irealf
!     igen
!     latf
!     lonf
!     latb
!     lonb
!     latr
!     lonr
!     ntrac
!     icen2
!     ienst
!     iensi
!     idpp
!     ids1
!     idvc
!     idvm
!     idvt
!     idrun
!     idusr
!     pdryini
!     ncldt
!     ixgr
!     nxgr
!     nxss
!     ivs
!     nvcoord
!     vcoord
!     cfvars
```

## 11.4 ss2gg

ss2gg converts a sigma file to a grads binary file and creates a corresponding descriptor (ctl) file

Original Author: Mark Iredell

Usage: ss2gg sigfile(s) gggfile ctlfile idrt imax jmax

where:

sigfile(s) = sigma file(s) to be converted to grads readable ieee files

gggfile = output file name

ctlfile = name of grads descriptor file (output)

idrt = output grid type  
0 = linear S->N  
4 = gaussian  
256 = linear N->S

imax = integer number of longitude points for output grid

jmax = integer number of latitude points for output grid

```
!           (IDRT=4 FOR GAUSSIAN GRID,  
!  
!           IDRT=0 FOR EQUALLY-SPACED GRID INCLUDING POLES.  
!   imax    - Integer even number of longitudes for output grid  
!   jmax    - Integer number of latitudes for output grid
```

# Appendix A – Global Model Variables

VARIABLE	GROUP	DESCRIPTION
ACCOUNT	GENERAL	LoadLeveler account, i.e. GFS-MTN
adiab	FCST	Debugging, true=run adiabatically
AERODIR	FCST	Directory, usually set to \$FIX_RAD, see \$FIX_RAD
AIRSBF	ANAL	Naming convention for AIRSBF data file
ALIST	GENERAL	Extra set of files to be added to rlist if ARCHIVE=YES; used only if rlist is being generated on the fly in this step; done in reconcile.sh
AM_EXEC	FCST	Atmospheric model executable
AM_FCS	FCST	See \$FCSTEXECTMP
AMSREBF	ANAL	AMSR/E bufr radiance dataset
ANALSH	ANAL	Analysis job script, usually "anal.sh"
ANALYSSSH	ANAL	Analysis driver script
ANAVINFO	ANAL	Text files containing information about the state, control, and meteorological variables used in the GSI analysis
ANGUPDATESH	ANGU	Angle update script
ANGUPDATEEXEC	ANGU	Angle update executable
ANISO_A_EN	ENKF	TRUE = use anisotropic localization of hybrid ensemble control variable a_en
anltype	ANAL	Analysis type (gfs or gdas) for verification (default=gfs)
Apercent	FCST	For idvc=3, 100: sigma-p, 0: pure-theta
append_rlist	GENERAL	Location of append_rlist (comment out if not using)
AQCX	PREP	Prep step executable
ARCA00GDAS	ARCH	Points to HPSS file name for ARCA files for 00Z cycle GDAS
ARCA00GFS	ARCH	Points to HPSS file name for ARCA files for 00Z cycle GFS
ARCA06GDAS	ARCH	Points to HPSS file name for ARCA files for 06Z cycle GDAS
ARCA06GFS	ARCH	Points to HPSS file name for ARCA files for 06Z cycle GFS
ARCA12GDAS	ARCH	Points to HPSS file name for ARCA files for 12Z cycle GDAS
ARCA12GFS	ARCH	Points to HPSS file name for ARCA files for 12Z cycle GFS
ARCA18GDAS	ARCH	Points to HPSS file name for ARCA files for 18Z cycle GDAS
ARCA18GFS	ARCH	Points to HPSS file name for ARCA files for 18Z cycle GFS
ARCB00GFS	ARCH	Points to HPSS file name for ARCB files for 00Z cycle GFS
ARCB06GFS	ARCH	Points to HPSS file name for ARCB files for 06Z cycle GFS
ARCB12GFS	ARCH	Points to HPSS file name for ARCB files for 12Z cycle GFS
ARCB18GFS	ARCH	Points to HPSS file name for ARCB files for 18Z cycle GFS
ARCC00GFS	ARCH	Points to HPSS file name for ARCC files for 00Z cycle GFS
ARCC06GFS	ARCH	Points to HPSS file name for ARCC files for 06Z cycle GFS
ARCC12GFS	ARCH	Points to HPSS file name for ARCC files for 12Z cycle GFS
ARCC18GFS	ARCH	Points to HPSS file name for ARCC files for 18Z cycle GFS
ARCDIR	ARCH	Location of online archive
ARCDIR1	ARCH	Online archive directory
ARCH_TO_HPSS	ARCH	Make hpss archive
ARCHCFRRSH	ARCH	Script location
ARCHCOPY	ARCH	If yes then copy select files (ARCR and ARCO in rlist) to online archive
ARCHDAY	ARCH	Days to delay online archive step
ARCHIVE	ARCH	Make online archive
ARCHSCP	ARCH	If yes & user glopara, scp all files for this cycle to alternate machine
ARCHSCP_TO	ARCH	Remote system to receive scp'd data (mist->dew, dew->mist)
ARCHSH	ARCH	Archive script
ASYM_GODAS	ANAL	For asymmetric godas (default=NO)
ATARDIR	ARCH	HPSS tape archive directory
ATARFILE	ARCH	HPSS tape archive tarball file name, \$ATARDIR/\$ADAY.tar
AVG_FCST	FCST	Time average forecast output files
AVRG_ALL	AVRG	To submit averaging and archiving scripts; this should be set to 'YES' - valid for reanalysis
AVRGALLSH	AVRG	Script location
B1AMUA	ANAL	Location and naming convention of B1AMUA data file
B1HRS4	ANAL	Location and naming convention of B1HRS4 data file
B1MHS	ANAL	Location and naming convention of B1MHS data file
BERROR	ANAL	Location and naming convention of BERROR files
beta1_inv	ENKF	1/beta1 = the weight given to static background error covariance
BUFRLIST	PREP	BUFR data types to use
C_EXEC	FCST	Coupler executable
CAT_FLX_TO_PGB	POST	Cat flx file to pgb files (only works for ncep post and IDRT=0)
cnorm	FCST	Assumes all cloud water is inside cloud (true), operation (false)



<b>CCPOST</b>	POST	To run concurrent post
<b>ccwf</b>	FCST	Cloud water function, ras, 1: high res, 2: T62
<b>CDATE</b>	GENERAL	Date of run cycle (YYYYMMDDCC), where CC is the forecast cycle, e.g. 00, 06, 12, 18
<b>CDATE_SKIP</b>	ANAL	LDAS modified sfc files not used before this date; must be >24 hours from the start
<b>CDFNL</b>	VRFY	SCORES verification against selected dump, pgban.gdas or pgbanl.gfs
<b>CDUMP</b>	GENERAL	Dump name (gfs or gdas)
<b>CDUMPPFCST</b>	PREP	Fits-to-obs against gdas or gfs prep
<b>CDUMPPREP</b>	PREP	Prep dump to be used in prepqfit
<b>CFSRDMP</b>	DUMP	Location of CFS/climate dump archive
<b>CFSRR_ARCH</b>	ARCH	Script location
<b>CFSRRPLOTSH</b>	AVRG	Script location
<b>CFSV2</b>	FCST	CFS switch, YES=run CFS version 2
<b>ch1</b>	FCST	Hours in gdas fcst1 & post1 job wall-clock-limit [hours:minutes:seconds] (see reconcile script)
<b>ch1</b>	POST	See ch1 (FCST)
<b>ch2</b>	FCST	Same as ch1 but for segment 2
<b>ch2</b>	POST	See ch2 (FCST)
<b>cha</b>	ANAL	Analysis wall time; hours in job wall-clock-limit [hours:minutes:seconds] (see reconcile script)
<b>CHG_LDAS</b>	ANAL	To bring in new vegtyp table to LDAS
<b>CHGRESEXEC</b>	GENERAL	Chgres executable location
<b>CHGRESH</b>	GENERAL	Chgres script location
<b>CHGRESTHREAD</b>	GENERAL	Number of threads for chgres (change resolution)
<b>CHGRESVARS</b>	GENERAL	Chgres variables
<b>CLDASSH</b>	ANAL	CLDAS script
<b>climate</b>	FCST	CFS variable, grib issue
<b>CLIMO_FIELDS_OPT</b>	FCST	Interpolate veg type, soil type, and slope type from inputgrid, all others from sfcsub.f, 3: to coldstart higher resolution run
<b>cm1</b>	FCST	Minutes in gdas fcst1 & post1 job wall-clock-limit [hours:minutes:seconds] (see reconcile script)
<b>cm1</b>	POST	See cm1 (FCST)
<b>cm2</b>	FCST	Same as cm1 but for segment 2
<b>cm2</b>	POST	See cm2 (FCST)
<b>ema</b>	ANAL	Analysis wall time; minutes in job wall-clock-limit [hours:minutes:seconds] (see reconcile script)
<b>cmapdl</b>	GENERAL	Cmap dump location in \$COMDMP
<b>cmbDysPrf4</b>	ANAL	GODAS executable
<b>cmbDysPrfs4</b>	ANAL	GODAS executable
<b>CO2_seasonal_cycle</b>	FCST	CO2 seasonal cycle; global_co2monthlycyc1976_YYYY.txt
<b>CO2DIR</b>	FCST	Directory with CO2 files
<b>COMCOP</b>	GENERAL	Location where copy.sh looks for production (or alternate) files
<b>COMDAY</b>	GENERAL	Directory to store experiment "dayfile" output (dayfile contains stdout & stderr), see \$ROTDIR
<b>COMDIR</b>	GENERAL	See \$TOPDIR
<b>COMDMP</b>	GENERAL	Location of key production (or alternate) files (observation data files, surface boundary files)
<b>COMDMPTMP</b>	GENERAL	Temporary version of \$COMDMP
<b>COMROTTMP</b>	GENERAL	If set, replaces config value of \$ROTDIR
<b>CONFIG</b>	GENERAL	Configuration file name
<b>cont_eq_opt1</b>	FCST	TRUE = when the advected and nonlinear fields of the mass-continuity equation are separated into two parts so that a different interpolation can be used for each part - following the EC approach. Only use with herm_x = herm_y = herm_z = lin_xy = false and lin_xyz = true. Additionally, opt1_3d_cubic = true, if quasi-tricubic interpolation is used for nonlinear terms
<b>CONVINFO</b>	ANAL	Location of convinfo.txt file, conventional data
<b>COPYGB</b>	GENERAL	Location of copygb utility
<b>COUP_FCST</b>	FCST	NO: AM model only, YES: coupled A-O forecast (default=NO)
<b>COUP_GDAS</b>	FCST	YES: run coupled GDAS
<b>COUP_GFS</b>	FCST	YES: run coupled GFS forecast
<b>CQCX</b>	PREP	Prep executable
<b>crtrh</b>	FCST	For Zhao microphysics, if zhao_mic is .false., then for Ferrier-Moorthi microphysics
<b>cs1</b>	FCST	Seconds in gdas fcst1 & post1 job wall-clock-limit [hours:minutes:seconds] (see reconcile script)
<b>cs1</b>	POST	See cs1 (FCST)
<b>cs2</b>	FCST	Same as cs1 but for segment 2
<b>cs2</b>	POST	See cs2 (FCST)
<b>csa</b>	ANAL	Analysis wall time; seconds in job wall-clock-limit [hours:minutes:seconds] (see reconcile script)
<b>CSTEP</b>	GENERAL	Step name (e.g. prep, anal, fcst2, post1, etc.)
<b>ctei_rm</b>	FCST	Cloud top entrainment instability criterion, mstrat=true
<b>CTL_ANL</b>	POST	Parameter file for grib output
<b>CTL_FCS</b>	POST	Parameter file for grib output
<b>CTL_FCS_D3D</b>	POST	Parameter file for grib output
<b>CUE2RUN</b>	COMP	User queue variable; LoadLeveler class for parallel jobs (i.e. dev)
<b>CUE2RUN1</b>	COMP	Similar to \$CUE2RUN but alternate queue

<b>CUE2RUN3</b>	COMP	Similar to \$CUE2RUN but alternate queue
<b>cWGsh</b>	ANAL	GODAS script
<b>CYCLESH</b>	GENERAL	Script location
<b>CYCLEXEC</b>	GENERAL	Executable location
<b>CYINC</b>	GENERAL	Variable used to decrement GDATE {06}
<b>DATATMP</b>	GENERAL	Working directory for current job
<b>DAYDIR</b>	GENERAL	See \$ROTDIR
<b>DELTIM</b>	FCST	Time step (seconds) for segment 1
<b>DELTIM2</b>	FCST	Time step (seconds) for segment 2
<b>DELTIM3</b>	FCST	Time step (seconds) for segment 3
<b>DELTIM_EFCS</b>	ENKF	Time step for ensemble forecast
<b>diagtable</b>	PREP	Ocean and ice diagnostic file
<b>diagtable_1dy</b>	PREP	Ocean and ice diagnostic file
<b>diagtable_1hr</b>	PREP	Ocean and ice diagnostic file
<b>diagtable_3hr</b>	PREP	Ocean and ice diagnostic file
<b>diagtable_6hr</b>	PREP	Ocean and ice diagnostic file
<b>diagtable_hrs</b>	PREP	Ocean and ice diagnostic file
<b>diagtable_long</b>	PREP	Ocean and ice diagnostic file
<b>dlqf</b>	FCST	Fraction of cloud water removed as parcel ascends
<b>DMPDIR</b>	DUMP	Dump directory location
<b>DMPEXP</b>	DUMP	Dump directory location, gdasy/gfsy
<b>DMPOPR</b>	DUMP	Dump directory location
<b>DO_RELOCATE</b>	PREP	Switch; to perform relocation or not
<b>DO2ANL</b>	ANAL	Do second analysis run, depends on value of CDFNL
<b>DODUMP</b>	DUMP	For running in real-time, whether or not to run the dump step
<b>DOENKF</b>	ENKF	YES = turns on EnKF script processing
<b>DOHYBVAR</b>	ENKF	YES = tells analysis step to use ensemble background error products from previous cycle
<b>DSDUMP</b>	DUMP	CFS dump directory
<b>dt_aocpl</b>	FCST	Coupler timestep
<b>dt_cpld</b>	FCST	Coupled timestep
<b>dt_ocean</b>	FCST	Ocean timestep
<b>dt_rstrt</b>	FCST	OM restart writing interval/timestep (small)
<b>dt_rstrt_long</b>	FCST	OM restart writing interval/timestep (long)
<b>Dumphsh</b>	DUMP	Dump script location and name
<b>EDATE</b>	GENERAL	Analysis/forecast cycle end date - must be >CDATE; analysis/forecast cycle ending date (YYYYMMDDCC, where CC is the cycle)
<b>EDUMP</b>	GENERAL	Cycle ending dump (gdas or gfs)
<b>EMISDIR</b>	FCST	Directory, usually set to \$FIX_RAD, see \$FIX_RAD
<b>ENS_NUM_ANAL</b>	ENKF	Number of ensemble members
<b>ENS_NUM_ENKF</b>	ENKF	Number of ensemble members
<b>ENTHALPY</b>	FCST	Control the chgres and nceppost (default=NO)
<b>ESTEP</b>	GENERAL	Cycle ending step; stop experiment when this step is reached for \$EDATE; this step is not run
<b>EXEC_AMD</b>	FCST	Atmospheric model directory
<b>EXEC_CD</b>	FCST	Coupler directory
<b>EXEC_OMD</b>	FCST	Ocean model directory
<b>EXECcfs</b>	FCST	CFS executable directory location
<b>EXECDIR</b>	GENERAL	Executable directory (typically underneath HOMEDIR)
<b>execdir_godasprep</b>	PREP	GODAS prep executable directory, see \$EXECDIR
<b>EXECICE</b>	FCST	Sea ice executable directory, see \$EXECDIR
<b>EXPDIR</b>	GENERAL	Experiment directory under /save, where your configuration file, rlist, runlog, and other experiment scripts reside
<b>FAISS</b>	FCST	Scale in days to relax to sea ice to climatology
<b>fbak2</b>	FCST	Back up time for 2nd segment
<b>fbak3</b>	FCST	Back up time for 3rd segment
<b>FCSTEXECDIR</b>	FCST	Location of forecast executable directory (usually set to \$EXECDIR)
<b>FCSTEXECTMP</b>	FCST	Location and name of forecast executable
<b>FCSTSH</b>	FCST	Forecast script name and location
<b>FCSTVARS</b>	FCST	Group of select forecast variables and their values
<b>fcyc</b>	FCST	Surface cycle calling interval
<b>fdfi_1</b>	FCST	Digital filter time for AM 1st segment (default=3)
<b>fdfi_2</b>	FCST	Run digital filter for 2nd segment (default=0)
<b>fdump</b>	VRFY	Verifying forecasts from gfs: GFS analysis or gdas: GDAS analysis
<b>FH_END_POST</b>	POST	Implying use FHMAX (default=99999)
<b>FH_STRT_POST</b>	POST	Implying to use FHINI or from file \$ROTDIR/FHREST.\$CDUMP.\$CDATE.\$nknid (default=99999)
<b>FHCYC</b>	FCST	Cycling frequency in hours

<b>FHDFI</b>	FCST	Initialization window in hours (if =0, no digital filter; if =3, window is +/- 3hrs)
<b>FHGOC3D</b>	FCST	Hour up to which data is needed to force offline GOCART to write out data
<b>FHINI</b>	FCST	Initial forecast hour
<b>FHLWR</b>	FCST	LW radiation calling interval (hrs); longwave frequency in hours
<b>FHMAX</b>	FCST	Maximum forecast hour
<b>FHMAX_HF</b>	FCST	High-frequency output maximum hours; for hurricane track, gfs fcst only for 126-hr is needed
<b>FHOUT</b>	FCST	Output frequency in hours
<b>FHOUT_HF</b>	FCST	High frequency output interval in hours; for hurricane track, gfs fcst only for 126-hr is needed
<b>FHRES</b>	FCST	Restart frequency in hours
<b>FHROT</b>	FCST	Forecast hour to Read One Time level
<b>FHSTRT</b>	FCST	To restart a forecast from a selected hour, default=9999999
<b>FHSWR</b>	FCST	SW radiation calling interval (hrs); frequency of solar radiation and convective cloud (hours)
<b>FHZER</b>	FCST	Zeroing frequency in hours
<b>FIT_DIR</b>	VERFY	Directory for SAVEFITS output
<b>FIX_LIS</b>	PREP	Location of land model fix files
<b>FIX_OCN</b>	PREP	Location of ocean model fix files
<b>FIX_OM</b>	PREP	See \$FIX_OCN
<b>FIX_RAD</b>	PREP	Fix directory, usually set to \$FIXGLOBAL
<b>FIXDIR</b>	PREP	Fix file directory
<b>FIXGLOBAL</b>	PREP	Atmospheric model fix file directory
<b>flgmin</b>	FCST	Minimum large ice fraction
<b>fmax1</b>	FCST	Maximum forecast hour in 1st segment (default=192 hrs)
<b>fmax2</b>	FCST	Maximum forecast hour in 2nd segment (default=384 hrs)
<b>fmax3</b>	FCST	Maximum forecast hour in 3rd segment (default=540 hrs)
<b>FNAISC</b>	FCST	CFS monthly ice data file
<b>FNMASK</b>	FCST	Global slmask data file, also see \$SLMASK
<b>FNOROG</b>	FCST	Global orography data file
<b>FNTSFC</b>	FCST	CFS oi2sst data file
<b>FNVEGC</b>	FCST	CFS vegfrac data file
<b>FNVETC</b>	FCST	Global vegetable type grib file
<b>FORECASTSH</b>	FCST	Forecast script name and location
<b>fout_a</b>	FCST	GDAS forecast output frequency (default=3); used when gdas_fh is not defined (i.e. no long gdas fcst)
<b>fout1</b>	FCST	GFS sig, sfc, flx output frequency for 1st segment (default=3 hr)
<b>fout2</b>	FCST	GFS sig, sfc, flx output frequency for 2nd segment (default=3 hr)
<b>fout3</b>	FCST	GFS sig, sfc, flx output frequency for 3rd segment (default=3 hr)
<b>foutpgb1</b>	POST	NCEPPOST pgb frequency for 1st segment (default=fout1)
<b>foutpgb2</b>	POST	NCEPPOST pgb frequency for 2nd segment (default=fout1)
<b>foutpgb3</b>	POST	NCEPPOST pgb frequency for 3rd segment (default=fout1)
<b>fres1</b>	FCST	Interval for restart write, 1st segment (default=24 hr)
<b>fres2</b>	FCST	Interval for restart write, 2nd segment (default=24 hr)
<b>fres3</b>	FCST	Interval to write restart for 3rd segment (default=fres2)
<b>fseg</b>	FCST	Number of AM forecast segments; maximum=3 (default=1)
<b>FSNOL</b>	FCST	Scale in days to relax to snow to climatology
<b>FTSFS</b>	FCST	Scale in days to relax to SST anomaly to zero
<b>fzer1</b>	FCST	GFS output zeroing interval for 1st segment (default=6 hr)
<b>fzer2</b>	FCST	GFS output zeroing interval for 2nd segment (default=6 hr)
<b>fzer3</b>	FCST	GFS output zeroing interval for 3rd segment (default=6 hr)
<b>G3DPSH</b>	ANAL	G3DP script name and location
<b>gdas_cyc</b>	FCST	Number of GDAS cycles
<b>gdas_fh</b>	FCST	Default=999, i.e. no long fcst in GDAS step when <999, that would be the interval at which seasonal or longer from gdas initial conditions are made; for example, if gdas_fh=6 runs are made
<b>GDAS_GP</b>	POST	YES: use old post (global_postgp.sh), NO: ncep post
<b>GDUMP</b>	GENERAL	Dump to use for guess files (defaults to \$CDFNL, which defaults to "gdas")
<b>generate_ens</b>	ENKF	TRUE = generate internal ensemble based on existing background error
<b>GENPSICHI</b>	POST	Generate psi (streamfunction) and chi (velocity potential)
<b>GENPSICHIEXE</b>	POST	Executable for GENPSICHI
<b>gfs_cyc</b>	FCST	GFS cycles (00, 06, 12, and 18Z) (default=1 - (00Z) cycle)
<b>GFSDUMP</b>	DUMP	GFS dump subdirectory name and location, usually "\$DMPDIR/dump"
<b>gg_tracers</b>	FCST	Semilag option
<b>GLDASCYCHR</b>	FCST	GLDAS cycling frequency
<b>GODAS_DATA_DELAY</b>	ANAL	Delay for ocean data in days
<b>GODAS_WNDO</b>	ANAL	Data window for asymmetric godas
<b>GODASEXEC</b>	ANAL	GODAS executable
<b>GODASSH</b>	ANAL	GODAS script
<b>GRID_IDD</b>	FCST	3D output options

<b>GRID11FCST00gdas</b>	FCST	Grib identifier for 00z GDAS forecast output
<b>GRID11FCST06gdas</b>	FCST	Grib identifier for 06z GDAS forecast output
<b>GRID11FCST12gdas</b>	FCST	Grib identifier for 12z GDAS forecast output
<b>GRID11FCST18gdas</b>	FCST	Grib identifier for 18z GDAS forecast output
<b>grid25_1</b>	POST	Define this to interpolate pgb file to 2.5 x 2.5
<b>grid25_2</b>	POST	Same as grid25_1 but for segment 2 of post
<b>grid62_1</b>	POST	Define this to interpolate fix file to T62 grid
<b>GROUP</b>	GENERAL	LoadLeveler group (i.e. g01)
<b>group_name</b>	GENERAL	Similar to \$GROUP
<b>GSIDIR</b>	ANAL	GSI HOMEDIR, usually equals \$HOMEDIR
<b>GSIEXEC</b>	ANAL	GSI executable name and location
<b>GSIFIXDIR</b>	ANAL	Location of GSI fix files
<b>HOMEcfs</b>	FCST	CFS HOMEDIR, usually equals \$HOMEDIR
<b>HOMEDIR</b>	GENERAL	Home directory for parallel scripts
<b>HORZ_DIR</b>	VERFY	Directory for SAVEFITS output
<b>HPSSTAR</b>	ARCH	Location of hpsstar utility (creates, retrieves, and manages tarfiles on HPSS)
<b>HRKDAY</b>	GENERAL	Hours to keep dayfiles in ROTDIR
<b>HRKOCN_ANL</b>	GENERAL	Hours to keep ocean analysis file
<b>HRKOCN_GRB</b>	GENERAL	Hours to keep ocean grib output file
<b>HRKRES</b>	GENERAL	Hours to keep restart files
<b>HRKROT</b>	GENERAL	Hours to keep rotating archive
<b>HRKSIG</b>	GENERAL	Hours to keep sigma and sfc fcst files in directory \$ROTDIR
<b>HRKSIGG</b>	GENERAL	Hours to keep sigma files from analysis in directory ROTDIR
<b>HRKTMP</b>	GENERAL	Hours to keep tmpdir
<b>HRKVFY</b>	GENERAL	Hours to keep verification files in directory ROTDIR
<b>HYBRID</b>	FCST	Switch to run hybrid
<b>HYBRID_ENSEMBLE</b>	ENKF	GSI namelist for hybrid ensemble variables
<b>IAER</b>	FCST	111: with stratospheric aerosol, tropospheric aerosol LW, tropospheric aerosol SW
<b>ialb</b>	FCST	For original albedo, 0: climatology SW albedo based on surface vegetation types, 1: MODIS based land surface albedo
<b>ICO2</b>	FCST	0: fixed CO2 constant, 1: time varying global mean CO2, 2: changing CO2
<b>ictm</b>	FCST	CO2 option for radiation, YYYY#
<b>IDRT_NP</b>	POST	Master pgb from global_nceppost.sh, 4: gaussian, 0: linear
<b>IDSL</b>	FCST	Integer new type of sigma structure, 1: Phillips approach, 2: Henry, plain average
<b>idvc_a</b>	FCST	AM vertical coordinate for analysis, 2: sigma-p (Sela), 3: generalized (Juang)
<b>idvc_f</b>	FCST	For hybrid model forecast (2: Joe Sela, 3: Henry Juang)
<b>IDVM</b>	FCST	Integer new vertical mass variable ID
<b>idvt</b>	FCST	Integer new tracer variable ID; first number: # of cloud species, second number: location of ozone in tracer
<b>IEMS</b>	FCST	0: blackbody ground emission, 1: climatology on one-deg map
<b>IGEN</b>	FCST	Integer output generating code (See ON388 Table A), grib output identifier, GFS=82, CFS=197
<b>IGEN_ANL</b>	FCST	Same as IGEN but for analysis
<b>IGEN_FCST</b>	FCST	Same as IGEN but for forecast
<b>IGEN_OCNP</b>	FCST	Same as IGEN but for ocean analysis
<b>inch_1</b>	FCST	Interval of coupled run (default=360)
<b>inch_2</b>	FCST	Coupled model interval of increment hour look (segment 2)
<b>io_1</b>	FCST	Forecast pgb output lon resolution, 1st segment
<b>io_2</b>	FCST	Forecast pgb output lon resolution, 2nd segment
<b>io_3</b>	FCST	Forecast pgb output lon resolution, 3rd segment
<b>io_a</b>	ANAL	Analysis pgb output lon and lat resolution
<b>io_save</b>	ARCH	Longitude dimension for online archive pgb files (defaults to 144... only applies if lower res than posted pgb files)
<b>IOVR_LW</b>	FCST	0: random cloud overlap for LW, 1: maximum/random cloud overlap for LW
<b>IOVR_SW</b>	FCST	0: random cloud overlap for SW, 1: maximum/random cloud overlap for SW
<b>ISOL</b>	FCST	0: fixed solar constant, 1: changing solar constant
<b>ISUBC_LW</b>	FCST	0: standard LW clouds (no MCICA), 1: prescribed MCICA seeds, 2: random MCICA seeds
<b>ISUBC_SW</b>	FCST	0: standard SW clouds (no MCICA), 1: prescribed MCICA seeds, 2: random MCICA seeds
<b>iter_one_no_interp</b>	FCST	TRUE = omits the trilinear interpolation for the first iteration of the departure-point calculations
<b>IYS</b>	FCST	Sigma file format (options 198410, 200509 defined in /nwprod/sorc/global_fcst.fd/sigio_module.f)
<b>ivssfc</b>	FCST	Surface file version
<b>ivssig</b>	FCST	Sigma file version
<b>JCAP</b>	FCST	Wave number (0-192 hr), atmospheric model resolution (spectral truncation), eg. JCAP=382
<b>JCAP_A</b>	FCST	See \$JCAP
<b>JCAP_TMP</b>	FCST	See \$JCAP
<b>JCAP_ENKF</b>	ENKF	Spectral resolution for Hybrid EnKF; similar to JCAP
<b>JCAP_ENS</b>	ENKF	\$JCAP_ENKF; Project T254 ensemble into linear grid (512x256)

<b>JCAP2</b>	FCST	Wave number (192-384 hr) for 2nd segment, see \$JCAP
<b>JCAP3</b>	FCST	Wave number (384-540 hr) for 3rd segment, see \$JCAP
<b>jo_1</b>	FCST	Forecast pgb output lat resolution, 1st segment
<b>jo_2</b>	FCST	Forecast pgb output lat resolution, 2nd segment
<b>jo_3</b>	FCST	Forecast pgb output lat resolution, 3rd segment
<b>jo_a</b>	FCST	Analysis pgb output lon and lat resolution
<b>jo_save</b>	FCST	Lat dimension for online archive pgb files (defaults to 72... only applies if lower res than posted pgb files)
<b>JOBSDIR</b>	GENERAL	Job script directory (typically underneath HOMEDIR)
<b>JUST_AVG</b>	AVRG	Default=NO
<b>JUST_POST</b>	POST	Terminate jobs after finishing post
<b>JUST_TSER</b>	POST	Extract just time-series by running post
<b>km_mom4</b>	POST	Number of MOM4 levels
<b>ko_1</b>	FCST	Forecast pgb output lev resolution, 1st segment
<b>ko_2</b>	FCST	Forecast pgb output lev resolution, 2nd segment
<b>ko_3</b>	FCST	Forecast pgb output lev resolution, 3rd segment
<b>ko_a</b>	ANAL	Analysis pgb output lev resolution
<b>kto_1</b>	FCST	Forecast IPV (isentropic potential vorticity) output resolution, if kto is set to 0, then no IPV output
<b>kto_2</b>	FCST	Vertical levels for segment 2, post step
<b>kto_3</b>	FCST	Same as kto_2 but for segment 3
<b>l_hyb_ens</b>	ENKF	TRUE = turn on hybrid ensemble option
<b>LANLSH</b>	ANAL	Land analysis LANLSH name and location
<b>LATA</b>	ANAL	Grid used by hurricane relocation, analysis_grid lat dimension (typically linear gaussian grid)
<b>LATA_ENKF</b>	ENKF	ensemble analysis_grid lat dimension (typically linear gaussian grid)
<b>LATB</b>	FCST	Model grid lat dimension (aka quadratic grid)
<b>LATB_D3D</b>	FCST	3D diagnostic output grid parameter
<b>LATB_ENKF</b>	ENKF	ensemble forecast_grid lat dimension (aka quadratic grid)
<b>LATB2</b>	FCST	Same as \$LATB but for segment 2
<b>LATB3</b>	FCST	Same as \$LATB but for segment 3
<b>LATCH</b>	FCST	Integer number of latitudes to process at one time in global_chgres; defaults to 8 in the code; defaults to 48 in branch parallel scripts; set to 8 in configuration file if you must match production when moving from the 1st to 2nd fcst segment; otherwise, go with the branch parallel script default of 48 to save resources (check current version of global_chgres.fd/chgres.f to confirm the code default; check fcst.sh and reconcile for script default)
<b>ld3d_1</b>	FCST	Write out 3D diagnostics, .false.: no 3D diagnostics
<b>ld3d_2</b>	FCST	3D diagnostic for segment 2
<b>ld3d_3</b>	FCST	3D diagnostic for segment 3
<b>ldas_cyc</b>	ANAL	0: no ldas cycles (default=0)
<b>LDIAG3D</b>	FCST	Switch for 3D diagnostics (default=false)
<b>LEVS</b>	FCST	Number of atmospheric model vertical levels
<b>LEVS_ENKF</b>	ENKF	Number of levels in Hybrid EnKF forecasts; similar to LEVS
<b>lg3d_1</b>	FCST	GOCART option segment 1 (default=false)
<b>lg3d_2</b>	FCST	GOCART option segment 2 (default=false)
<b>lin_xy</b>	FCST	TRUE = when the advected and nonlinear fields of the mass-continuity equation are separated into two parts so that a different interpolation can be used for each part. Only use with herm_x = herm_y = herm_z = cont_eq_opt1= false, and lin_xyz = true.
<b>lingg_a</b>	FCST	Semilag option
<b>lingg_b</b>	FCST	Semilag option
<b>LINKFILESH</b>	GENERAL	Link file script
<b>liope</b>	FCST	Atmospheric variable for io pes (default=.true.)
<b>LISEXEC</b>	ANAL	GLDAS (aka LIS) executable
<b>LISSH</b>	ANAL	GLDAS (aka LIS) script
<b>LONA</b>	FCST	Grid used by hurricane relocation, analysis_grid lon dimension (typically linear gaussian grid)
<b>LONA_ENKF</b>	ENKF	ensemble analysis_grid lon dimension (typically linear gaussian grid)
<b>LONB</b>	FCST	Model grid lon dimension (aka quadratic grid)
<b>LONB_D3D</b>	FCST	3D diagnostic output grid parameter
<b>LONB_ENKF</b>	ENKF	ensemble forecast_grid lon dimension (aka quadratic grid)
<b>LONB2</b>	FCST	Same as \$LONB but for segment 2
<b>LONB3</b>	FCST	Same as \$LONB but for segment 3
<b>LONSPERLAT</b>	FCST	Forecast step, global_lonsperlat text file
<b>lsm</b>	FCST	Land surface model, 1: NOAH land model, 0: OSU land model
<b>LSOIL</b>	FCST	Number of soil layers
<b>MAKEPREPBUFRSH</b>	PREP	Makeprepbufr script, created prepbufr
<b>mdlist</b>	VRFY	Exps (up to 10) to compare in maps
<b>MEANDIR</b>	AVRG	Directory for monthly means
<b>MFCST00GFS</b>	GENERAL	Starting number for dayfile iterations

<b>mkEvNc4r</b>	ANAL	GODAS executable
<b>MODIS_ALB</b>	FCST	To use MODIS based albedo product
<b>MON_AVG</b>	AVRG	CFS option, monthly averages for long integrations, starts 00z first day of month
<b>MP_PULSE</b>	COMP	IBM computing resource variable
<b>mppnccombine</b>	FCST	Location and name of cfs_mppnccombine executable
<b>mstrat</b>	FCST	Switch to turn on/off Moorthi stratus scheme
<b>MTNDIR</b>	FCST	See \$FIXGLOBAL
<b>MTNVAR</b>	FCST	The global_mtnvar fortran code
<b>NARRSNO</b>	ANAL	How snow assimilation is performed, North American Reanalysis
<b>NCEPOST</b>	POST	Switch to use NCEP post (default=YES)
<b>NCP</b>	GENERAL	Location of ncp utility
<b>ncw</b>	FCST	For Ferrier microphysics
<b>n_ens</b>	ENKF	number of ensemble members
<b>NEW_DAYFILE</b>	GENERAL	To create new dayfile for every rerun
<b>newoz_nrl</b>	FCST	YES: use NRL ozone production and loss coefficients (default=YES)
<b>NGPTC</b>	FCST	For operational GFS, not reproducible with different NGPTC; number of horizontal points computed in the same call inside radiation and physics (defaults to JCAP/10)
<b>nknd_fcst</b>	FCST	For hindcasts from segment 2 only
<b>NLAT_A</b>	ANAL	Analysis grid parameter, JCAP > 574
<b>NLAT_ENS</b>	ENKF	`expr \$NLAT_A_ENKF + 2`; Project T254 ensemble into linear grid (512x256)
<b>NLON_A</b>	ANAL	Analysis grid parameter, JCAP > 574
<b>NLON_ENS</b>	ENKF	\$NLON_A_ENKF; Project T254 ensemble into linear grid (512x256)
<b>NMEM_ENS</b>	ENKF	\$ENS_NUM_ENKF; Project T254 ensemble into linear grid (512x256)
<b>NOANAL</b>	ANAL	NO: run analysis and forecast, YES: no analysis (default=NO)
<b>NOFCST</b>	FCST	NO: run analysis and forecast, YES: no forecast (default=NO)
<b>npe_node_a</b>	ANAL	Number of PEs/node for atmospheric analysis with GSI
<b>npe_node_ang</b>	ANGU	Number of PEs/node for global_angupdate
<b>npe_node_av</b>	AVRG	Number of PEs/node for avrg
<b>npe_node_f</b>	FCST	Number of PEs/node for AM forecast
<b>npe_node_o</b>	ANAL	Number of PEs/node for ocean analysis
<b>npe_node_po</b>	POST	Number of PEs/node for post step (default=16)
<b>npe_node_pr</b>	PREP	Number of PEs/node for prep step (default=32 for dew/mist/haze)
<b>nproco_1</b>	FCST	Number of processors for ocean model 1st segment
<b>nproco_2</b>	FCST	Number of processors for ocean model 2nd segment
<b>nproco_3</b>	FCST	Number of processors for ocean model 3rd segment
<b>NRLACQC</b>	PREP	NRL aircraft QC, if="YES" will quality control all aircraft data
<b>nsout</b>	FCST	Outputs every AM time step when =1 (default=0)
<b>NSST_ACTIVE</b>	FCST	NST_FCST, 0: AM only, no NST model, 1: uncoupled, non-interacting, 2: coupled, interacting
<b>nth_f1</b>	FCST	Threads for AM 1st segment
<b>nth_f2</b>	FCST	Threads for AM 2nd segment
<b>nth_f3</b>	FCST	Threads for AM 3rd segment
<b>NTHREADS_GSI</b>	ANAL	Number of threads for anal
<b>NTHSTACK</b>	FCST	Stacks for fcst step (default=128000000)
<b>NTHSTACK_GSI</b>	ANAL	Stack size for anal (default=128000000)
<b>NUMPROCANAL</b>	ANAL	Number of tasks for GDAS anal
<b>NUMPROCANALGDAS</b>	ANAL	Number of tasks for GDAS anal
<b>NUMPROCANALGFS</b>	ANAL	Number of tasks for GFS anal
<b>NUMPROCAVRGGDAS</b>	ANAL	Number of PEs for GDAS average
<b>NUMPROCAVRGGFS</b>	ANAL	Number of PEs for GFS average
<b>NWPROD</b>	GENERAL	Option to point executable to nwprod versions
<b>O3CLIM</b>	FCST	Location and name of global_o3clim text file
<b>O3FORC</b>	FCST	Location and name of global_o3prdlos fortran code
<b>OANLSH</b>	ANAL	Ocean analysis script
<b>OBSQC</b>	ENKF	GSI namelist for observation quality control variables
<b>OCN2GRIBEXEC</b>	POST	Ocean to grib executable
<b>OCNMEANDIR</b>	AVRG	Directory for ocn monthly means
<b>ocnp_delay_1</b>	POST	OM post delay time
<b>ocnp_delay_2</b>	POST	OM post delay time
<b>OCNPSH</b>	POST	Ocean post script
<b>OIQCT</b>	PREP	Prep step prepobs_oiqc.obers file
<b>oisst_clim</b>	ANAL	Ocean analysis fix field
<b>OM_EXEC</b>	FCST	Ocean model executable
<b>omres_1</b>	FCST	Ocean 1st segment model resolution (0.5 x 0.25) and number of processors
<b>omres_2</b>	FCST	Ocean 2nd segment model resolution (0.5 x 0.25) and number of processors
<b>omres_3</b>	FCST	Ocean 3rd segment model resolution (0.5 x 0.25) and number of processors

<b>OPANAL_06</b>	ANAL	For old ICs without LANDICE, only applicable for starting from existing analysis
<b>OPREPSH</b>	PREP	Ocean analysis prep script
<b>opt1_3d_qcubic</b>	FCST	See cont_eq_opt1 variable for more information
<b>OROGRAPHY</b>	FCST	Global orography grib file
<b>OUT_VIRTTEMP</b>	FCST	Output into virtual temperature (true)
<b>OUTTYP_GP</b>	POST	1: gfsio, 2: sigio, 0: both
<b>OUTTYP_NP</b>	POST	1: gfsio, 2: sigio, 0: both
<b>OVERPARMEEXEC</b>	POST	CFS overparm grib executable
<b>oz_univ_static</b>	ENKF	TRUE = decouple ozone from other variables and defaults to static B (ozone only)
<b>OZINFO</b>	ANAL	Ozone info file
<b>PARATRKR</b>	TRAK	Script location
<b>PARM_GODAS</b>	PREP	GODAS parm file
<b>PARM_OM</b>	PREP	Ocean model parm files
<b>PARM_PREP</b>	PREP	Prep step parm files
<b>PCONFIGS</b>	GENERAL	For running in real-time, configuration file
<b>PCPINFO</b>	ANAL	PCP info files
<b>PEND</b>	GENERAL	Location of pend script
<b>pfac</b>	FCST	Forecasting computing variable
<b>pgb_typ4prep</b>	PREP	Type of pgb file for prep step (default=pgbf)
<b>pgbf_gdas</b>	POST	GDAS pgbf file resolution, 4: 0.5 x 0.5 degree, 3: 1 x 1 degree
<b>PMKR</b>	GENERAL	Needed for parallel scripts
<b>polist_37</b>	POST	Output pgb (pressure grib) file levels
<b>polist_47</b>	POST	Output pgb (pressure grib) file levels
<b>post_delay_1</b>	POST	AM post delay time
<b>post_delay_2</b>	POST	AM post delay time
<b>POST_SHARED</b>	POST	Share nodes (default=YES)
<b>POSTGPEXEC_GP</b>	POST	Post executable, for enthalpy version
<b>POSTGPEXEC_NP</b>	POST	Post executable, ncep post
<b>POSTGPSH_GP</b>	POST	\$POSTGPEXEC_GP script
<b>POSTGPSH_NP</b>	POST	\$POSTGPEXEC_NP script
<b>POSTGPVARSNP</b>	POST	Similar to FCSTVARS but for post variables
<b>POSTSH</b>	POST	Post script
<b>POSTSPL</b>	POST	Special CFSRR analysis file created for CPC diagnostics
<b>PRECIP_DATA_DELAY</b>	ANAL	Delay for precip data in hours (for global lanl)
<b>PREPDIR</b>	PREP	Location of prep files/codes/scripts, usually \$HOMEDIR
<b>PREPFXDIR</b>	PREP	Location of prep fix files
<b>PREPQFITSH</b>	PREP	Name and location of a prep script
<b>PREPSH</b>	PREP	Name and location of main prep script
<b>PREX</b>	PREP	Prevents executable
<b>PROCESS_TROPCY</b>	PREP	Switch, if YES: run QCTROPYSH script (default ush/syndat_qctropcy.sh)
<b>PRPC</b>	PREP	Prep parm file
<b>PRPT</b>	PREP	Prep bufr table
<b>PRPX</b>	PREP	Prepdata executable
<b>PRVT</b>	PREP	Global error table for prep
<b>PSLOT</b>	GENERAL	Experiment ID
<b>PSTX</b>	PREP	Prep step, global_postevents executable
<b>PSUB</b>	GENERAL	Location of psub script
<b>q2run_1</b>	FCST	Additional queue for fcst segment 1
<b>q2run_2</b>	FCST	Additional queue for fcst segment 2
<b>QCAX</b>	PREP	Prep step, prepobs_acarsqc executable
<b>r2ts_clim</b>	ANAL	Ocean analysis fix field
<b>ras</b>	FCST	Convection parameter, relaxed
<b>readfi_exec</b>	FCST	CFS sea ice executable
<b>readin_localization</b>	ENKF	TRUE = read external localization information file
<b>readsst_exec</b>	FCST	CFS sea ice executable
<b>RECONCILE</b>	GENERAL	Location of reconcile script
<b>REDO_POST</b>	POST	Default=NO
<b>regrid_exec</b>	FCST	CFS sea ice executable
<b>RELOCATESH</b>	PREP	Name and location of relocation script
<b>RELOX</b>	PREP	Name and location of relocation executable
<b>RESDIR</b>	GENERAL	Restart directory
<b>RESUBMIT</b>	GENERAL	To resubmit a failed job (default=NO)
<b>RLIST</b>	GENERAL	List that controls input and output of files for each step
<b>RM_G3DOUT</b>	FCST	For GOCART related special output
<b>RM_ORIG_G3D</b>	FCST	For GOCART related special output

<b>ROTDIR</b>	GENERAL	Experiment rotating/working directory, for large data and output files
<b>RTMAERO</b>	ANAL	Location of CRTM aerosol coefficient bin file
<b>RTMCLDS</b>	ANAL	Location of CRTM cloud coefficient bin file
<b>RTMEMIS</b>	ANAL	Location of CRTM emissivity coefficient bin file
<b>RTMFX</b>	ANAL	Location of CRTM fix file(s)
<b>RUN_ENTHALPY</b>	FCST	Control the forecast model (default=NO)
<b>RUN_OPREP</b>	PREP	YES: run ocean prep to get tmp.prf and sal.prf
<b>RUN_PLOT_SCRIPT</b>	AVRG	Script location
<b>RUN_RTDUMP</b>	ANAL	YES: archived tmp.prf and sal.prf used
<b>rundir</b>	GENERAL	Verification run directory
<b>RUNLOG</b>	GENERAL	The experiment runlog
<b>SALTSFCRESTORE</b>	ANAL	GODAS script
<b>SATANGL</b>	ANAL	Name and location of satangbias file
<b>SATINFO</b>	ANAL	Name and location of satinfo file
<b>SAVEFITS</b>	VRFY	Fit to obs scores
<b>SBUVBF</b>	ANAL	Location and naming convention of osbuvs8 data file
<b>SCRDIR</b>	GENERAL	Scripts directory (typically underneath \$HOMEDIR)
<b>scrubtyp</b>	GENERAL	Scrub or noscrub
<b>semilag</b>	FCST	Semilag option
<b>SEND2WEB</b>	VRFY	Whether or not to send maps to webhost
<b>s_env_h</b>	ENKF	homogeneous isotropic horizontal ensemble localization scale (km)
<b>s_env_v</b>	ENKF	vertical localization scale (grid units for now)
<b>SET_FIX_FLDS</b>	COPY	Only useful wit copy.sh; create orographic and MODIS albedo related fix fields if they don't exist
<b>settls_dep3dg</b>	FCST	Set settls_dep3ds and settls_dep3dg to true for the SETTLS departure-point calculation
<b>settls_dep3ds</b>	FCST	Set settls_dep3ds and settls_dep3dg to true for the SETTLS departure-point calculation
<b>SETUP</b>	ANAL	GSI setup namelist
<b>SHDIR</b>	GENERAL	Similar to SCRDIR, just a directory setting
<b>sice_rstrt_exec</b>	FCST	Sea ice executable
<b>SICEUPDATESH</b>	FCST	Sea ice update script
<b>SIGGESENV</b>	ENKF	template for ensemble member sigma guess files
<b>SLMASK</b>	FCST	Global slmask data file, also see \$FNMASK
<b>snoid</b>	ANAL	Snow id (default=snod)
<b>SNOWNC</b>	ANAL	NetCDF snow file
<b>SSMITBF</b>	ANAL	SSM/I bufr radiance dataset
<b>sst_ice_clim</b>	ANAL	Fix fields for ocean analysis
<b>SSTICECLIM</b>	ANAL	Ocean analysis fix field
<b>SUB</b>	GENERAL	Location of sub script
<b>SYNDATA</b>	PREP	Switch (default=YES)
<b>SYNDX</b>	PREP	Syndat file, prep step
<b>tasks</b>	FCST	Number of tasks for 1st segment of forecast
<b>tasks2</b>	FCST	Number of tasks for 2nd segment of forecast
<b>tasks3</b>	FCST	Number of tasks for 3rd segment of forecast
<b>tasksp_1</b>	POST	Number of PEs for 1st segment of post
<b>tasksp_2</b>	POST	Number of PEs for 2nd segment of post
<b>tasksp_3</b>	POST	Number of PEs for 3rd segment of post
<b>thlist_16</b>	POST	Output theta levels
<b>time_extrap_etadot</b>	FCST	TRUE = with settls_dep3ds and settls_dep3dg =false, when a second-order accuracy of the vertical displacements are desired
<b>TIMEAVGEXEC</b>	AVRG	Executable location
<b>TIMEDIR</b>	GENERAL	Directory for time series of selected variables
<b>TIMELIMANAL</b>	ANAL	Wall clock time for AM analysis
<b>TIMELIMAVRG</b>	AVRG	CPU limit (hhmmss) for averaging
<b>TIMELIMPOST00GDAS</b>	POST	CPU limit for 00z GDAS post
<b>TIMELIMPOST00GFS</b>	POST	CPU limit for 00z GFS post
<b>TIMELIMPOST06GFS</b>	POST	CPU limit for 06z GFS post
<b>TIMELIMPOST12GFS</b>	POST	CPU limit for 12z GFS post
<b>TIMELIMPOST18GFS</b>	POST	CPU limit for 18z GFS post
<b>TIMEMEANEXEC</b>	AVRG	Executable location
<b>TOPDIR</b>	GENERAL	Top directory, defaults to '/global' on CCS or '/mtb' on Vapor if not defined
<b>TOPDRA</b>	GENERAL	Top directory, defaults to '/global' on CCS or '/mtb' on Vapor if not defined
<b>TOPDRC</b>	GENERAL	Top directory, defaults to '/global' on CCS or '/mtb' on Vapor if not defined



<b>TOPDRG</b>	GENERAL	Top directory, defaults to '/global' on CCS or '/mtb' on Vapor if not defined
<b>TRACKERSH</b>	TRAK	Tracker script location
<b>TSER_FCST</b>	FCST	Extract time-series of selected output variables
<b>USE_RESTART</b>	GENERAL	Use restart file under ROTDIR/RESTART if run is interrupted
<b>USHAQC</b>	PREP	See \$USHDIR
<b>USHCQC</b>	PREP	See \$USHDIR
<b>USHDIR</b>	GENERAL	Ush directory (typically underneath HOMEDIR)
<b>USHGETGES</b>	PREP	Directory location of getges.sh script
<b>USHICE</b>	PREP	See \$USHDIR
<b>USHNQC</b>	PREP	See \$USHDIR
<b>USHOIQC</b>	PREP	See \$USHDIR
<b>USHPQC</b>	PREP	See \$USHDIR
<b>USHPREV</b>	PREP	See \$USHDIR
<b>USHQCA</b>	PREP	See \$USHDIR
<b>USHSYND</b>	PREP	Directory, usually "\$PREPDIR/ush"
<b>USHVQC</b>	PREP	See \$USHDIR
<b>usrdir</b>	GENERAL	See \$LOGNAME
<b>uv_hyb_ens</b>	ENKF	TRUE = ensemble perturbation wind variables are u,v; FALSE = ensemble perturbation wind variables are stream function and velocity potential
<b>VBACKUP_PRCP</b>	VERFY	Hours to delay precip verification
<b>VDUMP</b>	VERFY	Verifying dump
<b>vlength</b>	VERFY	Verification length in hours (default=384)
<b>VERFY_ALL_SEG</b>	VERFY	NO: submit vrfy only once at the end of all segments, YES: submit for all segments (default=YES)
<b>vrfy_delay_1</b>	VERFY	AM verification delay time (in hhmm) for segment 1
<b>vrfy_delay_2</b>	VERFY	AM verification delay time for segment 2
<b>VERFYPRCP</b>	VERFY	Precip threat scores
<b>VERFYSCOR</b>	VERFY	Anomaly correlations, etc.
<b>VERFYTRAK</b>	VERFY & TRAK	Hurricane tracks
<b>VSDB_START_DATE</b>	VERFY	Starting date for vsdb maps
<b>VSDB_STEP1</b>	VERFY	Compute stats in vsdb format (default=NO)
<b>VSDB_STEP2</b>	VERFY	Make vsdb-based maps (default=NO)
<b>vsdbhome</b>	VERFY	Script home (default=\$HOMEDIR/vsdb)
<b>vsdbsave</b>	VERFY	Place to save vsdb database
<b>VSDBSH</b>	VERFY	Default=\$vsdbhome/vsdbjob.sh
<b>WEBDIR</b>	VERFY	Directory on web server (rzdm) for verification output
<b>webhost</b>	VERFY	Webhost (rzdm) computer
<b>webhostid</b>	VERFY	Webhost (rzdm) user name
<b>yzdir</b>	VERFY	Additional verification directory, based on personal directory of Yuejian Zhu
<b>zflxtvd</b>	FCST	Vertical advection scheme
<b>zhao_mic</b>	FCST	TRUE: Zhao microphysics option, FALSE: Ferrier microphysics