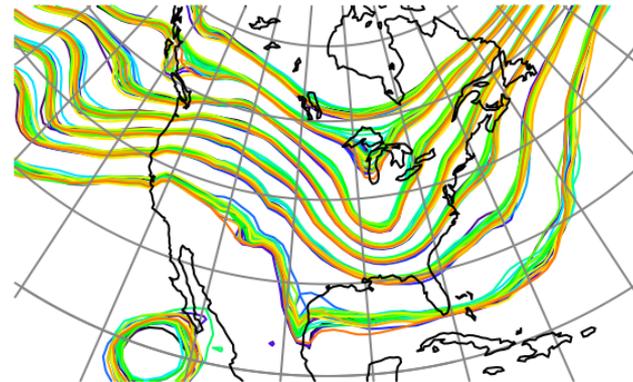


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## DART Tutorial Section 11: Creating DART Executables



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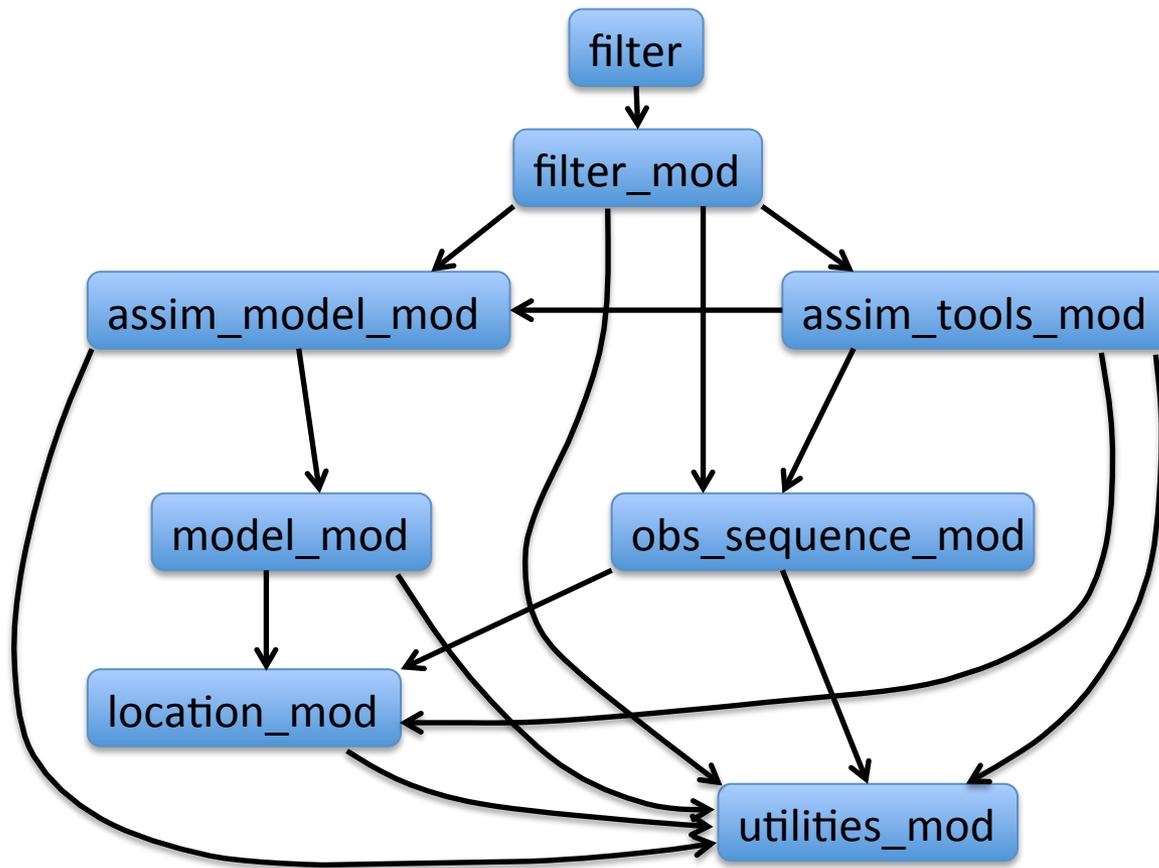


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# Fortran 90 'Use' Trees

DART requires use of F90 *use module\_mod*, only:  
No other mechanism for use of 'external' routine.



Program viewed as  
Directed Acyclic Graph  
(tree with shared leaves)

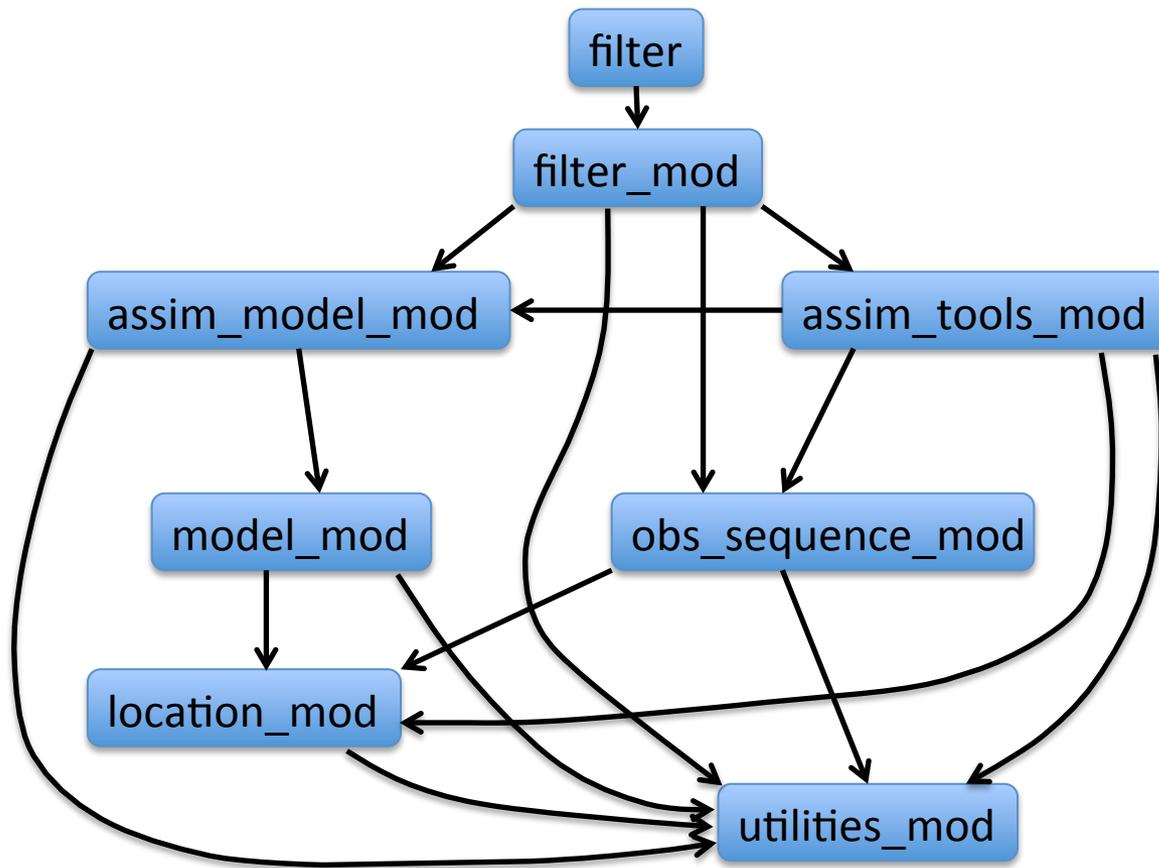
No cycles!

Partial tree for program  
filter is shown.

Vector represents use of  
target module.

# Generating a Makefile with mkmf

DART requires use of F90 *use module\_mod*, only:



Use ***mkmf*** Perl script to make a *Makefile*

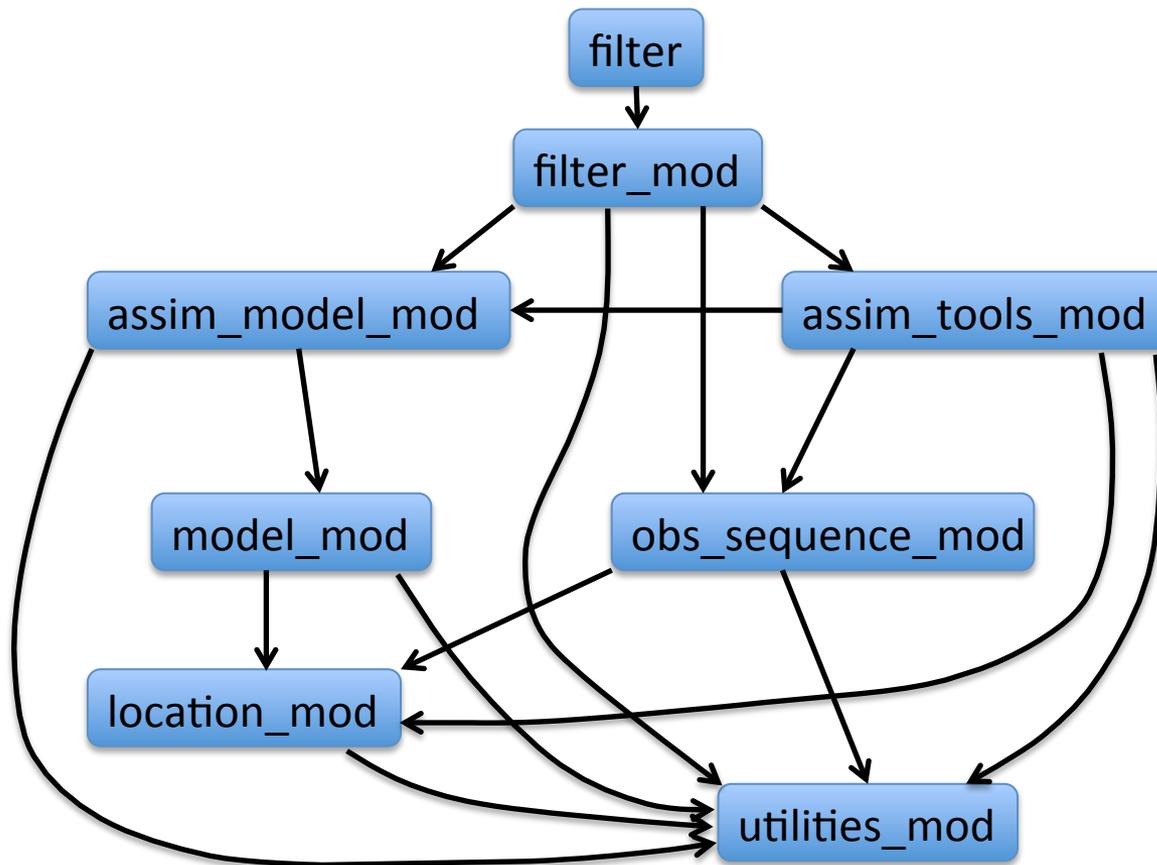
***mkmf*** requires a list of files to search for the main program and modules that are used.

This is called a *path\_names* file.

See *path\_names\_filter* in *models/lorenz\_63/work*.

# Generating a Makefile with mkmf

DART requires use of F90 *use module\_mod, only:*



**mkmf** searches files in `path_names` for one that contains *program filter*.

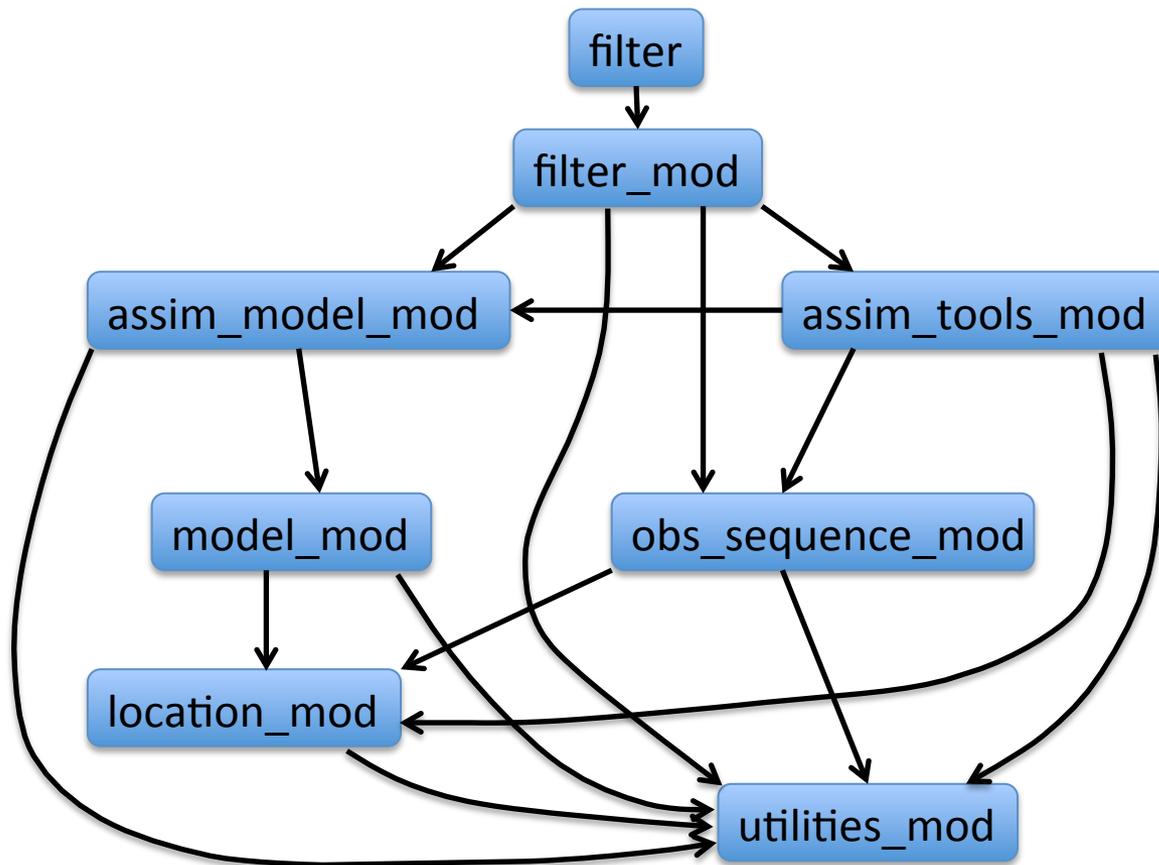
Finds first *use only* in filter.

Searches `path_names` files for this module recursively.

Builds a dependency graph like one at left.

# Generating a Makefile with mkmf

DART requires use of F90 *use module\_mod*, only:



From dependency graph, ***mkmf*** generates a standard *Makefile*.

Also creates a default namelist file, *input.nml.filter\_default*.

Examine to see namelists required, default values.

Enter ***make*** to create filter executable.

# mkmf Details

Each DART program has *mkmf\_* and *path\_names\_* files. Can see a selection of these in *models/lorenz\_63/work*.

Let's look at *mkmf\_perfect\_model\_obs* first.

File *mkmf\_perfect\_model\_obs* starts with scripting that controls whether code is built with mpi for parallel runs.

Key line starts with “*../../../build*” and contains

(default values in parentheses):

1. Relative location of mkmf program (*../../../build\_templates/mkmf*)
2. Name of executable program to create (*-p perfect\_model\_obs*)
3. Compiler options file (*-t ../../../build\_templates/mkmf.template*)
4. Relative base location for file search (*-a "../../.."*),
5. Name of the path\_names file (*path\_names\_perfect\_model\_obs*).

# mkmf Details

A variety of mkmf templates for different machines and compilers are available in the directory *build\_templates* (You can also see the Perl script *mkmf* there).

*mkmf* templates specify:

- What compiler to use,
- Where to find the netCDF libraries,
- Compile and link command line options.

Templates usually include comments about useful compiler and linker options for both debugging and production.

To change your template:

- Copy the appropriate *mkmf.template.xxx* file to *mkmf.template* and make any needed changes.

# mkmf Details

Two versions of the MPI module are available; one which really calls MPI and one with dummy routines.

The small models default to building a serial program.

The large models default to building an MPI parallel version.

To run *mkmf* for filter, execute shell script *mkmf\_filter*:

***./mkmf\_filter***    OR    ***csh mkmf\_filter***

# DART Modular Philosophy

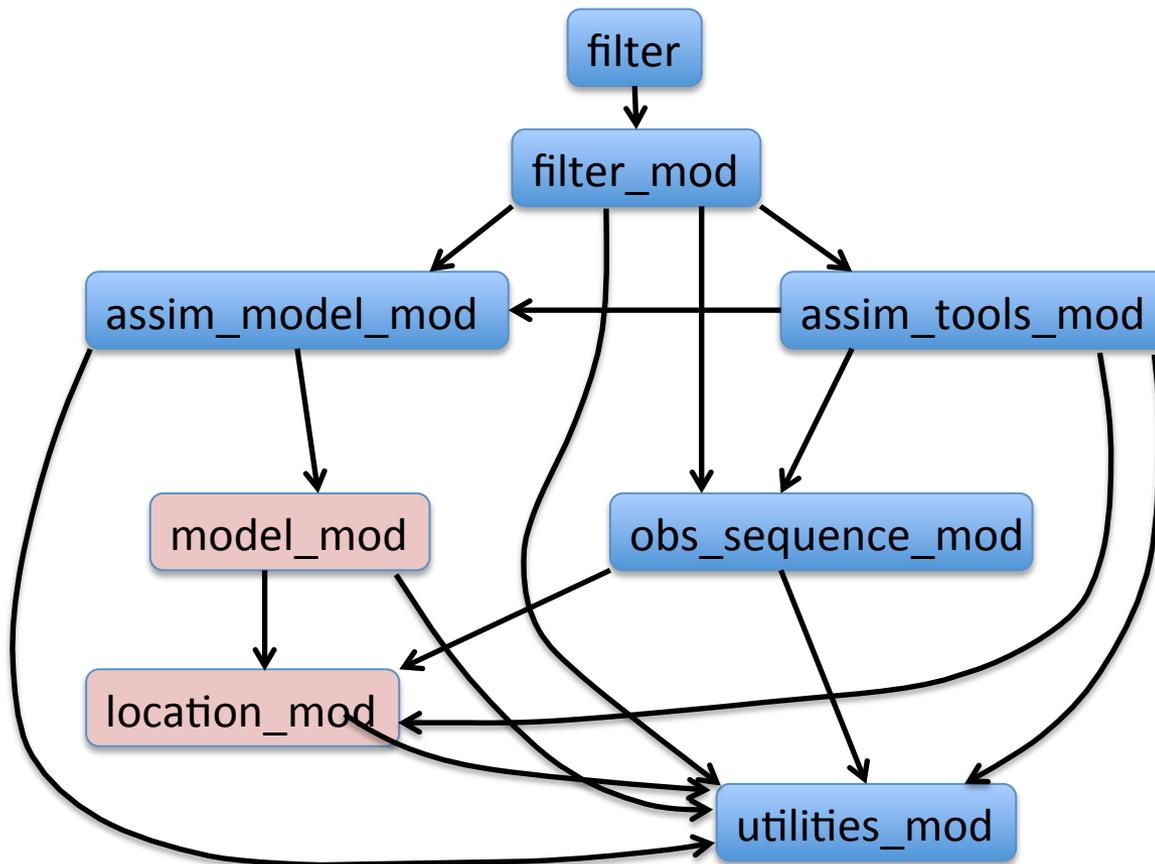
Can swap modules with same names and public interfaces.

Changing the following paths in *path\_names\_filter*:

models/**lorenz\_63**/model\_mod.f90 ➡ models/**cam-fv**/model\_mod.f90

assimilation\_code/location/**oned**/location\_mod.f90 ➡

assimilation\_code/location/**threed\_sphere**/location\_mod.f90



switches from Lorenz 63 to CAM-FV GCM!

Modules with multiple implementations have second directory level (see paths above).

Compare *path\_names* files in *models/lorenz\_63* with *models/cam-fv*

# Exercise: Compiling lorenz\_63 filter program

models/lorenz\_63/work/

1. Go to *models/lorenz\_63/work*
2. Remove all files with *.o* and *.mod* extensions.
3. Generate a *Makefile* and *input.nml.filter\_default* by:  
***ssh mkmf\_filter***
4. Generate program filter:  
Enter ***make***

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Multivariate assimilation.
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