FUN3D v12.4 Training

Session 2: Compilation and Installation

Bill Jones

Learning Goals

- What this will teach you
  - How to configure and compile the FUN3D suite
  - Configuration options
    - Enable/Disable capabilities
    - Specify the location of 3rd party libraries and tools
  - How *we* do it
- What you will not learn
  - How to build/install 3rd party libraries and tools
  - How to configure your system to compile Fortran 90/MPI code
- What should you already know
  - How to navigate through a *NIX shell
    - **`mkdir`**
    - **`cd`**
    - Absolute/relative paths

Setting

- Background
  - FUN3D uses the *de facto* industry standard build environment provided by GNU Autotools
  - Build of the FUN3D distribution does **not** require Autotools on your system
  - Provides localization through options to a configuration script
- Compatibility
  - Requires a Bourne Shell derivative (*NIX, OS X, MinGW, etc.)
  - Requires GNU `**make**`
  - Requires a functioning Fortran 95 compliant compiler (some optional capabilities rely on Fortran 2003 additions)
  - May not work with *non-standard* installation of 3rd party libraries
  - DiRTLib and SUGGAR++ assumptions for overset support
    - Required library names: **`libp3d.a`**, **`libdirt.a`**, **`libdirt_mpich.a`**, **`libsuggar.a`**, and **`libsuggar_mpi.a`**
  - Developers will need GNU Autotools installed

Nuts and Bolts (1 of 4)

- Two step process
  - `configure` selects capabilities and localizes to system
  - `make` creates executables
- Distribution contains a `configure` script
  - Familiar to Linux users/administrators who have built open source packages
  - Must **NOT** be edited by hand
  - Custom localization through command line options
- The `configure` script creates **Makefiles**
  - **Makefiles** are customized/localized for a specific *configuration*
  - Not practical for human consumption
  - Must **NOT** be edited by hand
    - All localization is managed through the `configure` script
  - Checks various details required by compilation
    - Fails fast (prior to compilation of FUN3D) if problems are detected with the configuration options (no compiler, missing libraries, etc.)

Nuts and Bolts (2 of 4)

- `configure --help` will show a list of all options
  - Command line options
  - Environment variables
  - Order independent (uses last value if specified multiple times)
- FUN3D optional Features of general interest

```
--disable-FEATURE          do not include FEATURE
                           (same as --enable-FEATURE=no)
--enable-FEATURE[=ARG]     include FEATURE [ARG=yes]

--enable-design            build Adjoint design tools [no]
--enable-hefss             build with High Energy Physics [no]
--enable-ftune             tailor Fortran compiler options for FUN3D [yes]
```

Nuts and Bolts (3 of 4)

- FUN3D optional Packages of general interest

```
--with-PACKAGE[=ARG]          use PACKAGE [ARG=yes]
--without-PACKAGE             do not use PACKAGE (same as --with-PACKAGE=no)

--with-mpi[=ARG]              Path to MPI library (installation root)
--with-mpibin[=ARG]           MPI binary directory      (relative, absolute, without)
--with-mpif90[=ARG]           MPI Fortran compiler wrapper  (relative, absolute, without)
--with-mpicc[=ARG]            MPI C compiler wrapper    (relative, absolute, without)
--with-mpiexec[=ARG]          MPI execution startup script  (relative, absolute, without)
--with-mpibin[=ARG]           MPI bin directory     (relative, absolute, without)
--with-mpiinc[=ARG]           Path to "mpif.h"      (relative, absolute, without)
--with-parmetis[=ARG]         ParMetis install path     (contains lib/libparmetis.a)
--with-dirtlib[=ARG]          use DiRTlib overset library   (contains lib/libdirt.a)
--with-suggar[=ARG]           use SUGGAR overset library    (contains lib/libsuggar.a)
--with-tecio[=ARG]            Tecplot I/O library install path  (contains lib/libtecio.a)
--with-refine[=ARG]           use refine adaptation package (installation root)
--with-refineFAKEGeom[=ARG]   to specify refine FAKEGeom libs [-lFAUXGeom]
--with-knife[=ARG]            use Knife cut cell package    (installation root)
--with-CGNS[=ARG]             CGNS library path     (installation root)
--with-PORT[=ARG]             use PORT optimization library (contains lib/libport.a)
--with-KSOPT[=ARG]            use KSOPT optimization library    (contains lib/libksopt.a)
--with-SNOPT[=ARG]            use SNOPT optimization library    (contains lib/libsnopt.a)
```

Nuts and Bolts (4 of 4)

- FUN3D environment variables of general interest

FC	**Fortran compiler command** (overridden by `--with-mpif90`)
FCFLAGS	**Fortran compiler flags** (adds to default unless `--disable-ftune`)
LDFLAGS	**linker flags, e.g. -L<lib dir> if you have libraries in a nonstandard directory <lib dir>**
CC	**C compiler command**
CFLAGS	**C compiler flags**
CPPFLAGS C/C++	**preprocessor flags, e.g. -I<include dir> if you have headers in a nonstandard directory <include dir>**
CPP	**C preprocessor**

- `make` is used to build the executables
  - Will reside in respective directories (e.g. **nodet** is in **FUN3D_90**)

Basic Operation

- Construct the vanilla **serial** executable
- Unpack your FUN3D distribution
  - Creates a directory "fun3d-12.4-69663"
- Enter the FUN3D distribution directory
- Run the `configure` script and build executables with `make`

```
$ mkdir serial
$ cd serial
$ ../configure
$ make
```

- Note that this will search for a supported compiler in your path
  - Chooses the first one found based on pre-defined order
  - Override this with the `FC=mycompiler` option
    - MPI configurations will use the `--with-mpif90` wrapper if given

Did It Work? Expected Output

```
…
Configuration (FUN3D):
      Source code location: ..
      Version:              12.4-69663
      Fortran Compiler:     ifort
      Fortran basis:        ifort
      Fortran flags:        -O2 -ip -align
                            -fno-alias -vec-report0
      C Compiler:           gcc
      C flags:              -g -O2
      Linker flags:         -lm
      Dependencies:         Normal

build:

      High Energy Physics:  no
      Cmplx Variable Tools: no
      Python bindings:      no
      PDF documentation:    yes

bindings:

      Libcore:              internal
      refine:               subpackage
      CAPRI support:        no
```

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```
      knife:                subpackage
      MPI support:          no
      CUDA support:         no
      Zoltan:               no
      ParMETIS:             no
      Tecplot I/O:          no
      6DOF libraries:       no
      DiRTlib support:      no
      SUGGAR support:       no
      DYMORE support:       no
      RCAS_SDX support:     no
      CGNS support:         no
      PORT support:         no
      NPSOL support:        no
      DOT support:          no
      KSOPT support:        no
      SNOPT support:        no
      SMEMRD support:       version 1.3.1
      SSDC support:         no
      SBOOM support:        no
      VisIt support:        no
```

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- Executables created relative to the *serial* sub-directory
  - **`FUN3D_90/nodet`**, **`Adjoint/dual`**, **`Design/opt_driver`**

Extended Operation

(How we do it)

- Create a **parallel** version of the code
- Build in a separate *configuration* subdirectory
  - Stores object code and executables only
  - Does not *pollute* the source tree with object code
  - Multiple configurations utilize the same source

```
$ mkdir mpi
$ cd mpi
$ ../configure --with-mpi=/path/to/mpi \
               --with-parmetis=/path/to/parmetis
$ make
```

Did It Work? Expected Output

```
…
Configuration (FUN3D):
      Source code location: ..
      Version:              12.4-69663
      Fortran Compiler:     /path/to/mpi/bin/mpif90
      Fortran basis:        ifort
      Fortran flags:        -O2 -ip -align
                            -fno-alias -vec-report0
      C Compiler:           /path/to/mpi/bin/mpicc
      C flags:              -g -O2
      Linker flags:         -lm
      Dependencies:         Normal

build:

      High Energy Physics:  no
      Cmplx Variable Tools: no
      Python bindings:      no
      PDF documentation:    yes

bindings:

      Libcore:              internal
      refine:               subpackage
      CAPRI support:        no
```

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```
      knife:                subpackage
      MPI support:          yes
      CUDA support:         no
      Zoltan:               no
      ParMETIS:             /path/to/parmetis
      Tecplot I/O:          no
      6DOF libraries:       no
      DiRTlib support:      no
      SUGGAR support:       no
      DYMORE support:       no
      RCAS_SDX support:     no
      CGNS support:         no
      PORT support:         no
      NPSOL support:        no
      DOT support:          no
      KSOPT support:        no
      SNOPT support:        no
      SMEMRD support:       version 1.3.1
      SSDC support:         no
      SBOOM support:        no
      VisIt support:        no
```

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- Executables created under the *configuration* directory
  - **`FUN3D_90/nodet_mpi`**, **`Adjoint/dual_mpi`**, **`Design/opt_driver`**

Troubleshooting/FAQ (1of 3)

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- Problems
  - "checking for Fortran compiler default output file name... configure: error: Fortran compiler cannot create executables See `config.log` for more details."
    - Make sure that Fortran compiler works in your environment
      - Adjust PATH, load appropriate GNU modules, MPI installation, etc.
  - Limited check of **`configure`** options
    - Bad "**`--enable-*`**" and "**`--with-*`**" options silently ignored
  - Option values containing spaces must be quoted from shell
    - e.g. **`FCFLAGS="-g -O2 -m32 -fno-common"`**
  - Do **NOT** configure in top level distribution directory and **then** try to make individual configuration directories
    - **`make distclean`** to clean a previous configuration of the source
  - Look/send "**`config.log`**" file
    - Also includes configuration options at the top (less quoted values w/ spaces)

Troubleshooting/FAQ (2 of 3)

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- Can I…
  - Override the default compiler options?
    - Yes, **`--disable-ftune FCFLAGS="-what-ever-you-want"`**
      - Remember some compilers always need certain options
  - Explicitly specify my compiler?
    - You can, with **`FC=compiler`**, <u>but</u> this will be overridden if using **`` `--with-mpif90` ``**
  - Fix anything by manually editing the **`` `configure` ``** script or **`Makefiles`**?
    - **<u>NO</u>**! and we cannot support any such action
    - Anything that you can safely change is governed by a configure option
  - Install the executables in a central location?
    - Yes, **`` `make install` ``** will install executables, etc. under the location given by the **`"--prefix=/your/path"`** option to **`` `configure` ``**

Troubleshooting/FAQ (3 of 3)

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- What if I…
  - Have a proprietary MPI installation?
    - Some HPC resources have proprietary MPI installations using non-standard paths and names
    - Use "**--with-mpibin**", "**--with-mpiinc**", "**--with-mpif90**", and "**--with-mpiexec**" along with their "**—without-***" counterparts as needed to specify the binary and include paths as well as the name for the `**mpif90**` compiler wrapper and, if needed, the `**mpiexec**` script
    - Paths can be absolute or relative to the **--with-mpi** and **--with-mpibin** values

```
$ ./configure --with-mpi=/path/to/mpi
    --with-mpif90=my_mpif90 --without-mpiexec …
```

  - My MPI executables will not run
    - Check the consistency of your MPI compilation/runtime installations
    - The MPI installation used for compilation is available as **MPI Prefix:** from

```
$ /path/to/nodet/nodet_mpi --version
```

What We Covered

- How to configure and compile the FUN3D suite
  - Execute `` `configure` `` to localize a configuration
  - Build the executables with `` `make` ``
- Configuration options
  - Enable/Disable Features
  - With/Without Packages (3rd party libraries and tools)
  - Custom environment variables
- Use separate *configuration* subdirectories
  - Keeps source and object code separate
  - Allows multiple *configurations* under one source
  - Invoke as `` `../configure …` ``