



EFDA

EUROPEAN FUSION DEVELOPMENT AGREEMENT

Task Force
INTEGRATED TOKAMAK MODELLING

Kepler for beginners
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Tutorial/demonstration: Kepler for beginners

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- **Goal : build and execute a Kepler workflow**
- **Pre-requisites :**
 - Create a private database in your account
 - for a given machine (tokamak name) and data structure version :
 - `/afs/efda-itm.eu/project/switm/scripts/create_user_itm_dir TokamakName DataVersion`
 - Example : create a tree for tokamak name test (allowed for testing purposes)
`/afs/efda-itm.eu/project/switm/scripts/create_user_itm_dir test 4.08b`
 - Copy KEPLER in your directory :
 - If there is a previous release of Kepler in your directory :
 - `mv your_kepler your_kepler_save`
 - remove the KEPLER cache : `rm -rf ~/.kepler`
 - get KEPLER :
`cd ~; tar xvf /afs/efda-itm.eu/project/switm/kepler/xxxx/kepler.tar`
where `xxxx` is the current release of the UAL
This command creates in your home directory a subdirectory named `kepler`
 - Set the environment variables :
 - put in the configuration file `.cshrc`:
`source /afs/efda-itm.eu/project/switm/scripts/ITMv1 kepler TokamakName DataVersion`
NB : if this command is not included in the `.cshrc` file, it is mandatory to type it each time you open a new window

FORTRAN : Create a subroutine

- Identify the CPO in and out
- Create a subroutine with CPOs as argument
 - This example can be found ~signoret/public/workflow_examples/fortran

```

subroutine coreprof2mhd(coreprofin,mhdout)

use eulTM_schemas
implicit none
integer,parameter :: DP=kind(1.0D0)
!
! Always describe cpo as array
! In case of time slice, the size of the input cpo is 1
!
type (type_coreprof),pointer :: coreprofin(:)
type (type_mhd),pointer :: mhdout(:)
integer :: i,j,k,l
!!!!!!!!!!!!!!!!!!!! Physics calculations
! The output CPO must be allocated with its number of time slices
! (1 for a single time slice physics module)
allocate(mhdout(size(coreprofin)))
! Fill in the output CPO (Physical data)
do i=1,size(coreprofin)
  ! Time : copy from input CPO
  mhdout(i)%time = coreprofin(i)%time ! THE TIME FIELD MUST BE FILLED
(MANDATORY)
  ! Psi : copy from input CPO
  allocate(mhdout(i)%psi(size(coreprofin(i)%psi%value)))
  mhdout(i)%psi = coreprofin(i)%psi%value
! Example of a physics quantity that would have been calculated by the module
allocate(mhdout(i)%frequency(3))
mhdout(i)%frequency(1) = 1.1D0
mhdout(i)%frequency(2) = 1.2D0
mhdout(i)%frequency(3) = 1.3D0
.....

allocate(mhdout(i)%disp_perp(3,2,2));
do j=1,3
  do k=1,2
    do l=1,2
      mhdout(i)%disp_perp(j,k,l)=i*1.0+k*(-1.2)+l;
    enddo
  enddo
enddo
return
end subroutine

```

FORTRAN : testbed

- To make sure the code handles the CPOs correctly, run it in a “testbed”
 - create a main program which :
 - implements UAL calls to open, read, write, close the shot in the database
 - Calls the subroutine

```

program test
use euITM_schemas
use euITM_routines
implicit none
integer,parameter :: DP=kind(1.0D0)

interface
subroutine coreprof2mhd(coreprofin,mhdout)
use euitm_schemas
type (type_coreprof),pointer :: coreprofin(:) ← Declare the subroutine
type (type_mhd),pointer :: mhdout(:)
end subroutine
end interface

type (type_coreprof),pointer :: coreprofin(:)
type (type_mhd),pointer :: mhdout(:)

integer :: idxin, idxout, shot, runin, runout, refshot, refrun
character(len=5)::treename

shot = 1983
runin = 1
runout = 2
refshot = 0 ! Dummy, not used
refrun =0 ! Dummy, not used
treename = 'euitm' ! Mandatory, do not change
  
```

Example :

~signoret/public/workflow_examples/fortran/standalone.f90

```

write(*,*) 'Open shot in MDS !' (continued)
call euitm_open(treename,shot,runin,idxin)

write(*,*) 'Reading the input CPO :'
call euitm_get(idxin,"coreprof",coreprofin)

write(*,*) 'Calling the coreprof2mhd subroutine :'
call coreprof2mhd(coreprofin,mhdout)

write(*,*) 'Creating output run :'
! This is the UAL function that creates a shot in the ITM MDS+ tree
call euitm_create(treename,shot,runout,refshot,refrun,idxout)

write(*,*) 'Put result'
call euitm_put(idxout,"mhd",mhdout)

write(*,*) 'Closing Database :'
call euitm_close(idxin,treename,shot,runin)
call euitm_close(idxout,treename,shot,runout)

write(*,*) 'Deallocate CPOs :'
call euitm_deallocate(coreprofin)
call euitm_deallocate(mhdout)
end
  
```

- Prepare a makefile to compile both the subroutine as a library and the standalone program

Example :

~signoret/public/workflow_examples/fortran/Makefile

```

F77=pgf90
F90=pgf90
CC=gcc
COPTS = -r8 -fPIC -Mnosecond_underscore -g
LIBS = -L$(UAL)/lib -IUALFORTRANInterface_pgi
INCLUDES = -I$(UAL)/include/amd64_pgi

all: coreprof2mhd.o standalone libcpo2cpof

libcpo2cpof: coreprof2mhd.o
    ar -rvs libcpo2cpof.a coreprof2mhd.o

standalone: standalone.f90
    $(F90) $(COPTS) -o $@ $^ coreprof2mhd.o ${INCLUDES} $(LIBS)

coreprof2mhd.o: coreprof2mhd.f90
    $(F90) $(COPTS) -c -o $@ $^ ${INCLUDES} $(LIBS)

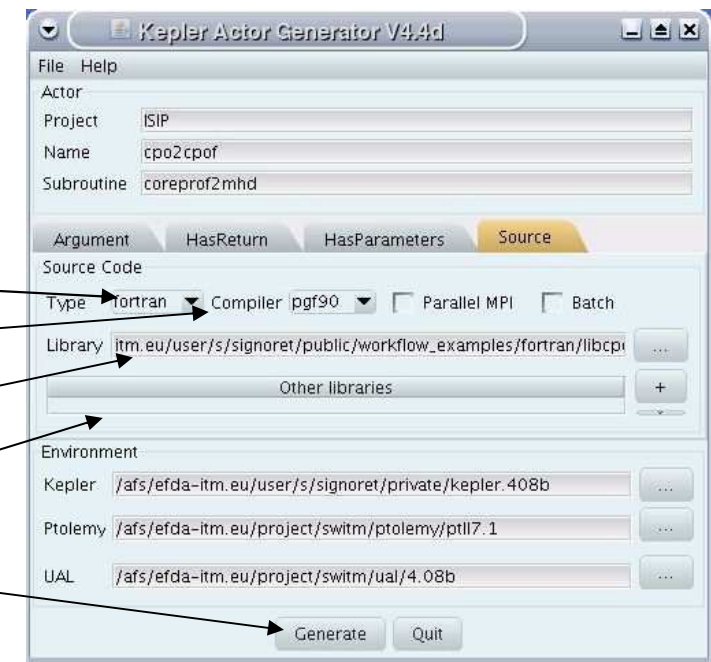
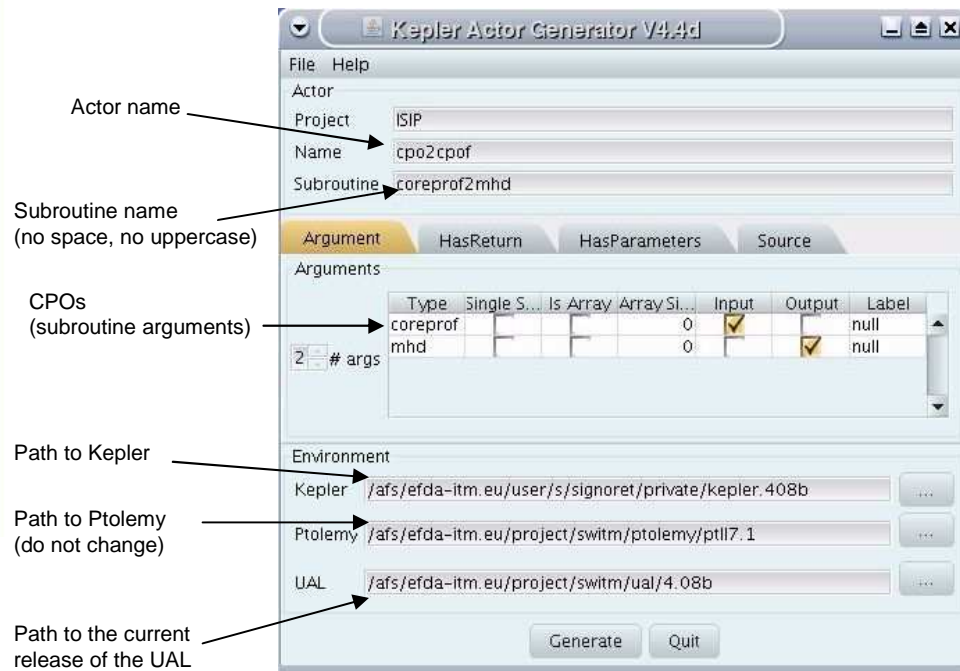
clean:
    rm -f *.o standalone *.a
  
```

-fPIC flag mandatory to make the library
 -g flag for debugging
 \$UAL = path to the current version of the UAL
 Building the library
 Building the standalone program

- run the standalone program to test the physics code

Run fc2k

- Launch : fc2k &



Automatic generation of the **actor**
(code, component, ...)

language

Compiler

Actor library

Additional libraries

Example : in fc2k load the file
~signoret/public/workflow_examples/fortran/cpo2cpof_fc2k.xml

Workflow: introduction to KEPLER

Run control buttons

Scheduler: time behavior

Catalogues

Many actors:
 database, math,
 display, web
 service, grid
 service, R-
 expression, ...

Many directors: DE,
 CT, SDF, PN,
 DDF and more
 in Ptolemy II
 (FSM, Giotto,
 DDE)

This model shows the solution to the classic Lotka-Volterra predator prey dynamics model. It uses the Continuous Time domain to solve two coupled differential equations, one that models the predator population and one that models the prey population. The results are plotted as they are calculated showing both population change and a phase diagram of the dynamics.

Rich Williams, 2003, NCEAS

Design of a workflow (1)

Choose the actors

1. In the Catalogues space, expand 'ISIP', drag and drop to the workspace the actors ualinit, cpo2cpof, ualcollector
 ualinit reads the input shot in the database
 ualcollector writes the resulting cpo in the database
2. Search 'constant' actor : drag and drop 4 constant and 3 String constant actors to the workspace
3. Search 'display' actor; drag and drop it
4. Define the director : choose a SDF director

Customize the constant actor

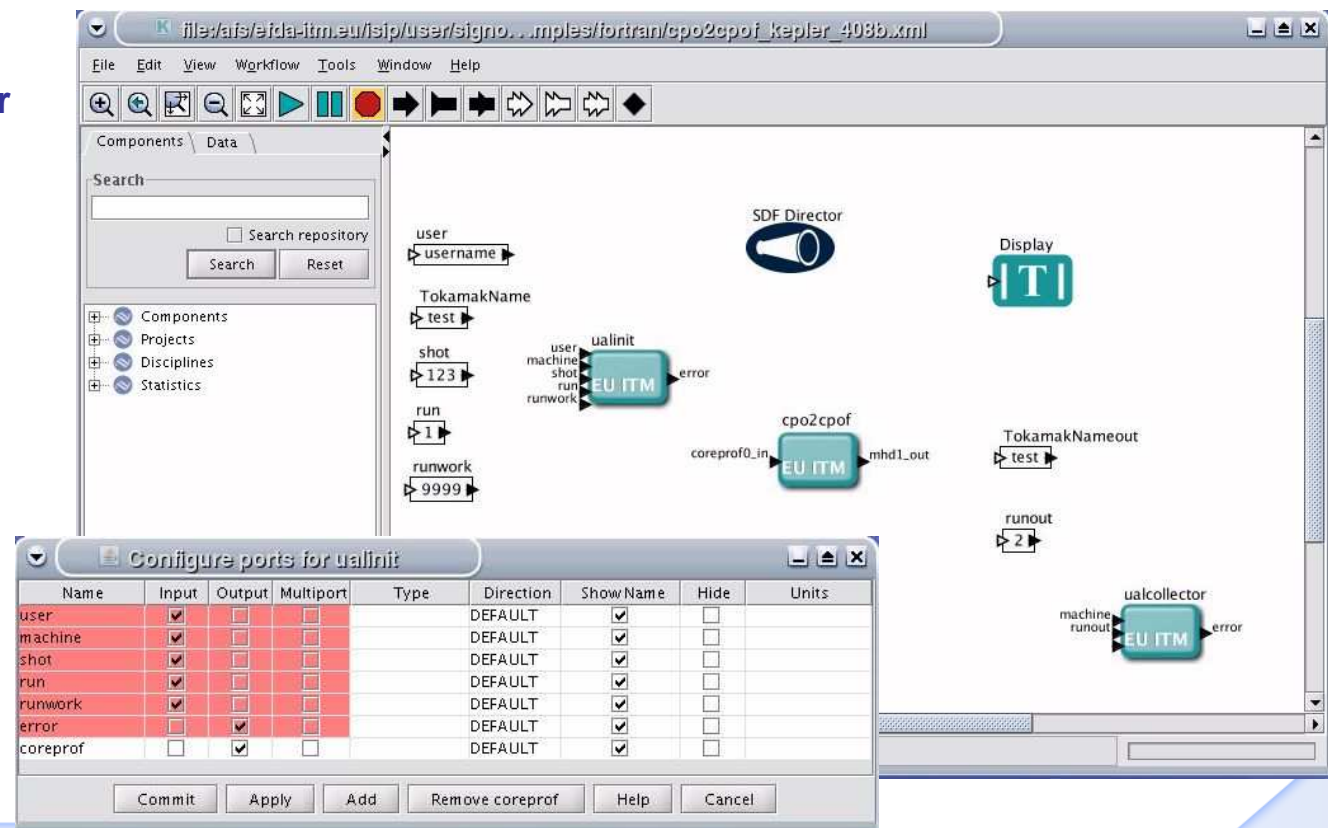
- Change the actors names
 right click on an actor, choose 'Customize Name')
- Set the values of the actors
 (right click on an actor, choose 'Configure Actor')

Customize ualinit

1. Add a port for each cpo

Connect the actors

Don't forget to save the workflow!



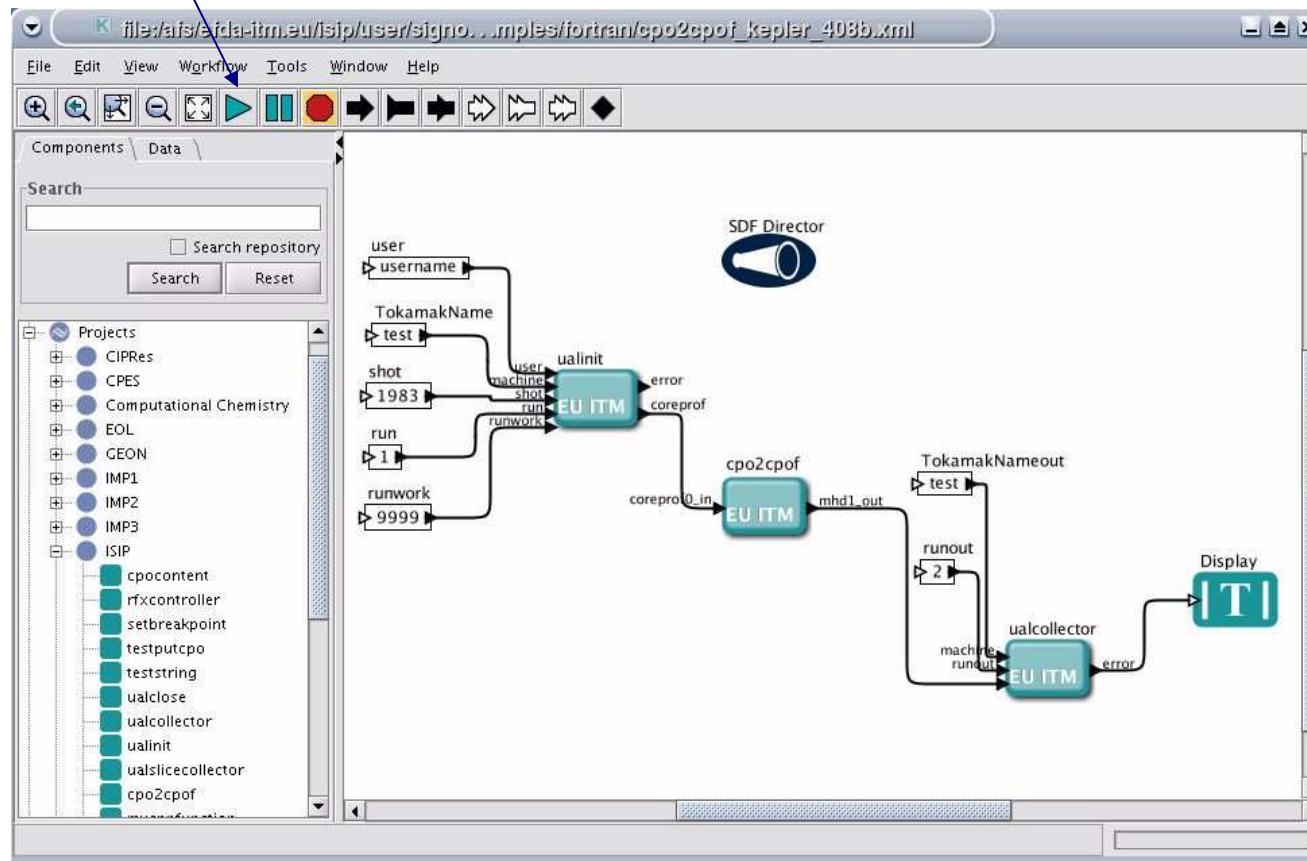
The screenshot shows the Kepler workflow editor interface. The main workspace contains several actors: 'user', 'TokamakName', 'shot', 'run', 'runwork', 'SDF Director', 'ualinit', 'cpo2cpof', 'Display', and 'ualcollector'. The workflow is connected as follows: 'user' (username) feeds into 'TokamakName' (test), which feeds into 'shot' (123). 'shot' feeds into 'run' (1), which feeds into 'runwork' (9999). 'runwork' feeds into 'ualinit' (EU ITM), which has an 'error' output. 'ualinit' feeds into 'cpo2cpof' (EU ITM), which has a 'coreprof0_in' input and an 'mhd1_out' output. 'cpo2cpof' feeds into 'Display' (T) and 'ualcollector' (EU ITM). 'ualcollector' has a 'machine runout' input and an 'error' output. The 'SDF Director' actor is also present but not connected.

The 'Configure ports for ualinit' dialog box is open, showing a table of ports:

Name	Input	Output	Multiport	Type	Direction	Show Name	Hide	Units
user	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>		DEFAULT	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
machine	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>		DEFAULT	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
shot	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>		DEFAULT	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
run	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>		DEFAULT	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
runwork	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>		DEFAULT	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
error	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>		DEFAULT	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
coreprof	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>		DEFAULT	<input checked="" type="checkbox"/>	<input type="checkbox"/>	

Run the simulation

1. **Optional : Select Tools → Animate at runtime** : set a delay in milliseconds to follow the execution of the workflow
2. **Push the button**



NB : the following examples are available for UAL 4.08b

- **Processing a single slice (FORTRAN)**

- Parameters for fc2k :

- ~signoret/public/workflow_examples/fortran/cposlice2cposlicef_fc2k.xml

- Worklow to open in Kepler :

- ~signoret/public/workflow_examples/fortran/cposlice2cposlicef_kepler_408b.xml

- **Processing a cpo (C++) :**

- Parameters for fc2k :

- ~signoret/public/workflow_examples/cpp/cpo2cpocpp_fc2k.xml

- Worklow to open in Kepler :

- ~signoret/public/workflow_examples/cpp/cpo2cpocpp_kepler_408b.xml

- **rmactor** :
 - Delete an actor from your Kepler installation
 - Usage : `rmactor actorname`
- **extract_actor**
 - Export all the files from your Kepler installation for a given actor and copy them in a tar file
 - Usage : `extract_actor actorname`
- **import_actor**
 - Import all the files from an actor tar file into your Kepler installation
 - Usage : `import_actor [options] [path/]actorname`
 - `import_actor -h` : display the detailed usage
- **NB** : users can exchange their own actors using `extract_actor` and `import_actor`

References

1. KEPLER:
 - <http://www.kepler-project.org/>
 - <http://users.sdsc.edu/~altintas/KeplerTutorial/>
2. ITM: <http://www.efda-taskforce-itm.org/>
3. Gateway: <http://www.efda-itm.eu>
4. Euforia: <http://www.euforia-project.eu/EUFORIA/>
5. Ptolemy II: <http://ptolemy.eecs.berkeley.edu/publications/>
6. See also :
 - UAL Tutorial
 - ITM Tools Tutorial