



EFDA

EUROPEAN FUSION DEVELOPMENT AGREEMENT

Task Force
INTEGRATED TOKAMAK MODELLING

*ITM Training
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Introduction to the ITM AMNS system

Presented by D. Coster

TF Leader : G. Falchetto,
Deputies: R. Coelho, D. Coster

EFDA CSU Contact Person: D. Kalupin

- background on the ITM AMNS system
- an example
- more information about the AMNS system

- The AMNS system aims to
 - provide an easy-to-use interface to AMNS data
 - provide validated AMNS data
 - capture the provenance of the data used in each simulation

- The AMNS system consists of
 - programs for importing AMNS data into the ITM system
 - AMNS data stored in the ITM database under the machine name “amns”
 - fortran callable library to access the AMNS data
 - c callable library to access the AMNS data
 - example programs to demonstrate the use of AMNS data
 - documentation on the ITM web pages about the AMNS system

- The example program is designed to calculate the density of charge states that represents an equilibrium between ionization and recombination

$$\frac{\partial n_i}{\partial t} = S_{i-1}n_{i-1} - R_i n_i - S_i n_i + R_{i+1}n_{i+1}$$

$$\begin{pmatrix} \cdot & \cdot & 0 & 0 & 0 & 0 & 0 \\ \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & S_{i-1} & -(S_i + R_i) & R_{i+1} & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ 0 & 0 & 0 & 0 & 0 & \cdot & \cdot \end{pmatrix} \begin{pmatrix} n_0 \\ \cdot \\ n_{i-1} \\ n_i \\ n_{i+1} \\ \cdot \\ n_z \end{pmatrix} = \begin{pmatrix} 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \end{pmatrix}$$

- The program needs the ionization and recombination rates
- Will output these as a function of T_e for a fixed chosen density
- Documentation of AMNS system
 - AMNS User Interface

- Check out the code
 - svn co <http://gforge.efda-itm.eu/svn/amnsproto/tags/examples/coronal>
- Move to the new directory
 - cd coronal
- Compile and run the code
 - make
- Plot the results
 - obj/plot_densities
 - obj/plot_radiation

- From the Makefile
 - `SYS=amd64_pgi`
 - `-include obj/SYS`
 - `OBJECTCODE=obj/${SYS}`
 - `include config/${SYS}`
- default to “amd64_pgi” but override this with the contents of “obj/SYS” if this exists
- set the object code location
- pick up the compiler configuration files

- From the Makefile
 - VPATH=src
 - MOD=mod
- set the search path for source and the suffix for modules

- From the Makefile
 - AMNSLIBDIR = /afs/efda-itm.eu/user/a/
amnsdata/public/library/4.09a/lib/\${SYS}
- set the location for picking up the AMNS library

- From the Makefile
 - run_coronal: \${OBJECTCODE}/coronal
 - @time \$^ \${ARGS}
- set the default in the makefile to run the example
- this has a dependency on the compiled code
- it also allows for passing arguments to the code

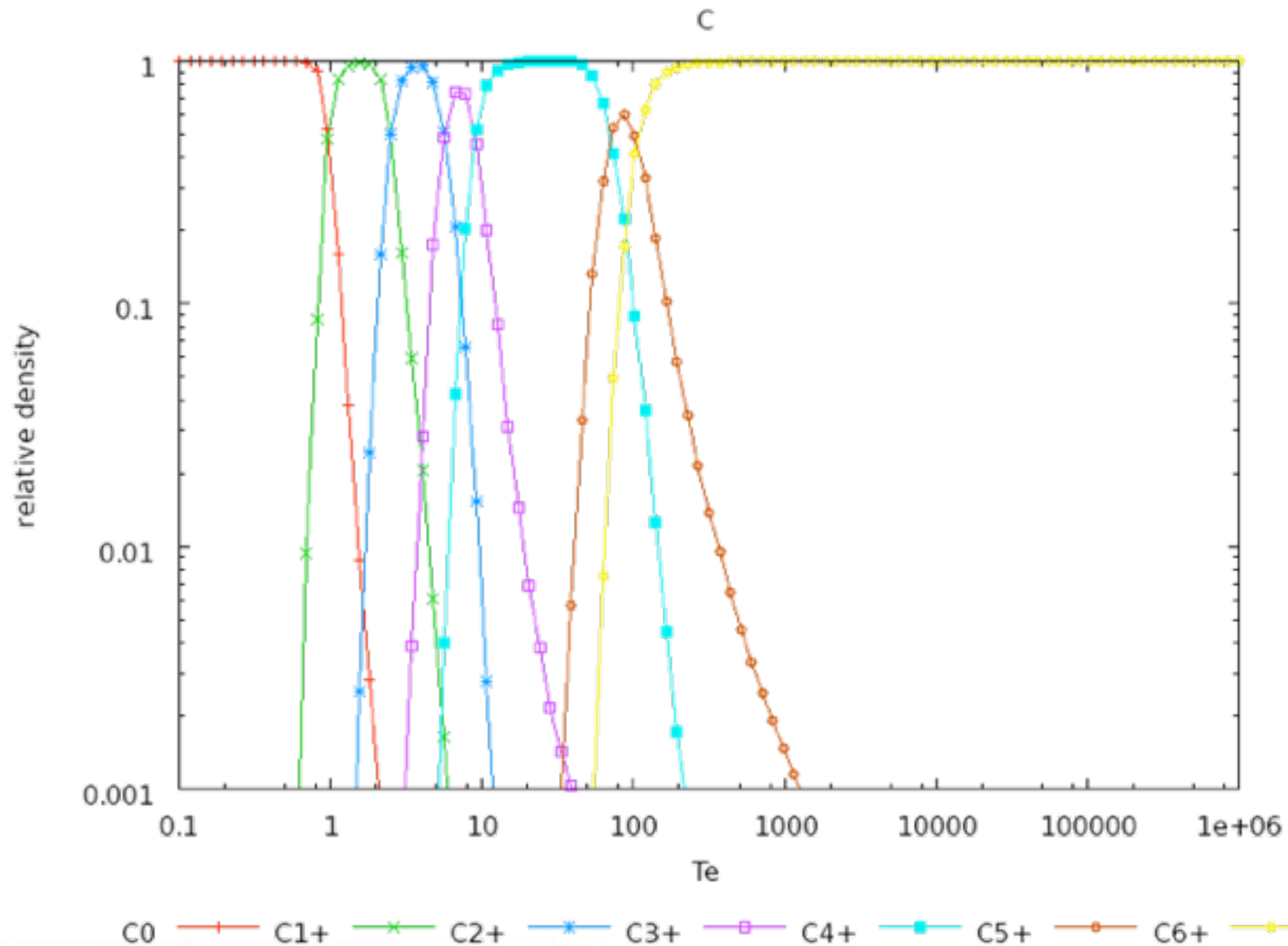
- From the Makefile
 - `${OBJECTCODE}/coronal : ${OBJECTCODE}/coronal.o`
 - `${FC} ${FCOPTS} -o $@ $^ -L${AMNSLIBDIR} -lamns ${LIBS} ${UAL_lib}`
- how to link the code

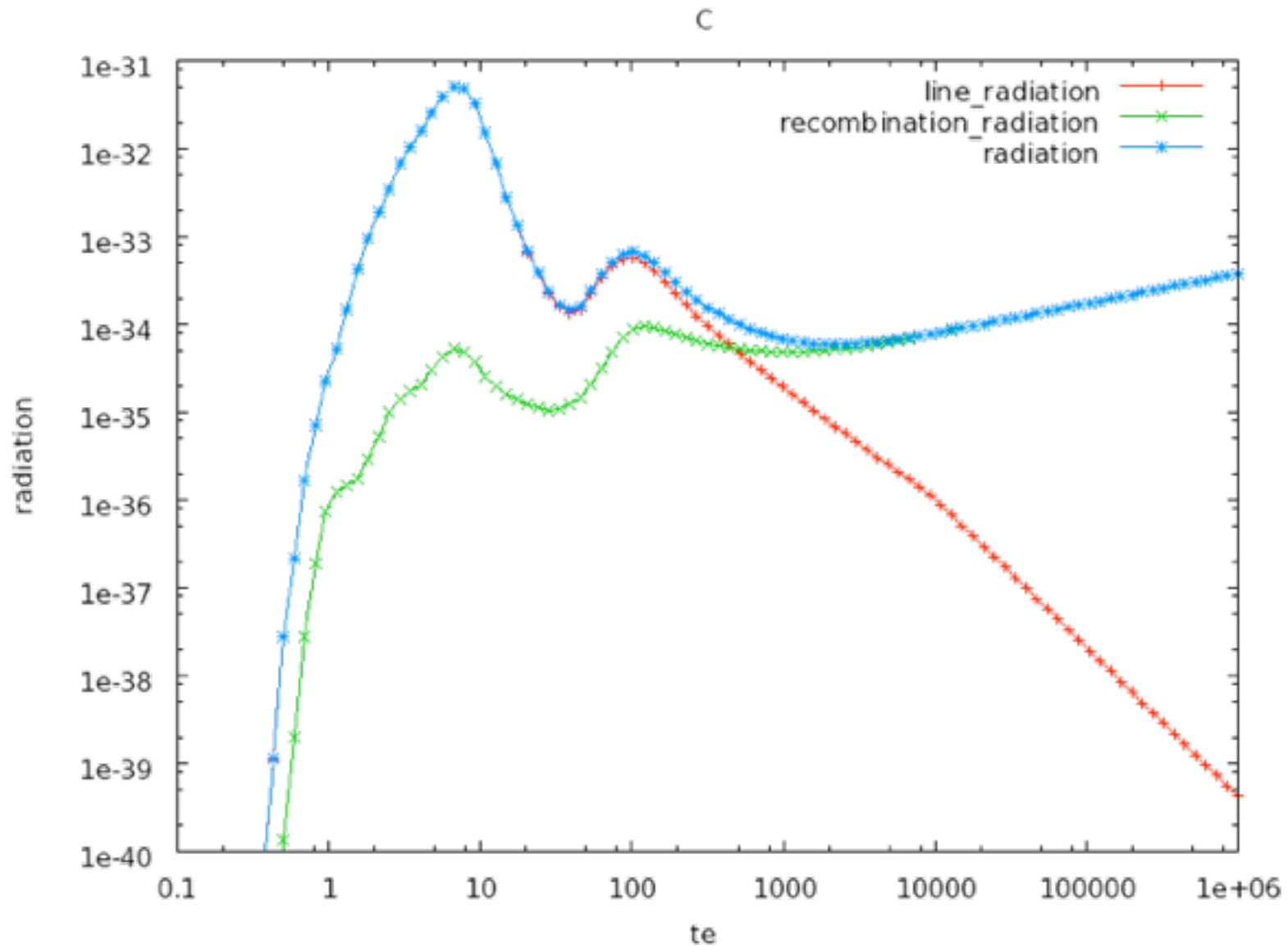
- From the Makefile
 - `${OBJECTCODE}/%.o : %.F90`
 - `${FC} ${FCOPTS} -I${OBJECTCODE} $`
`{UAL_inc} -I${AMNSLIBDIR} -c $< -o $@`
 - `-mv *.${MOD} ${OBJECTCODE}/`
- how to compile the code

- Gives some command lines for running and plotting a variety of species
 - make ARGS="1 2"
 - obj/plot_densities
 - obj/plot_cps D.ps
 - obj/plot_radiation
 - obj/plot_cps D_rad.ps

 - make ARGS="6 12.01"
 - obj/plot_densities
 - obj/plot_cps C.ps
 - obj/plot_radiation
 - obj/plot_cps C_rad.ps

 - ...





- Data is stored in CPOs with the tokamak name “amns”
- The shot number is $M*100+A$ for reactions that depend on the isotope, and A for reactions that don't (M =mass; A =nuclear charge)
- The run number represents the version number of data for a particular run number
- An MySQL database is used to store which versions of the data have been stored
- If the user has no AMNS data, that from “amnsdata” is used instead