



EFDA

EUROPEAN FUSION DEVELOPMENT AGREEMENT

Task Force  
INTEGRATED TOKAMAK MODELLING

*ITM General Meeting 2010  
Lisbon, September 13-15, 2010*

# Hybrid MHD-Gyrokinetic codes for studying the mutual nonlinear interaction of shear Alfvén modes and energetic particles

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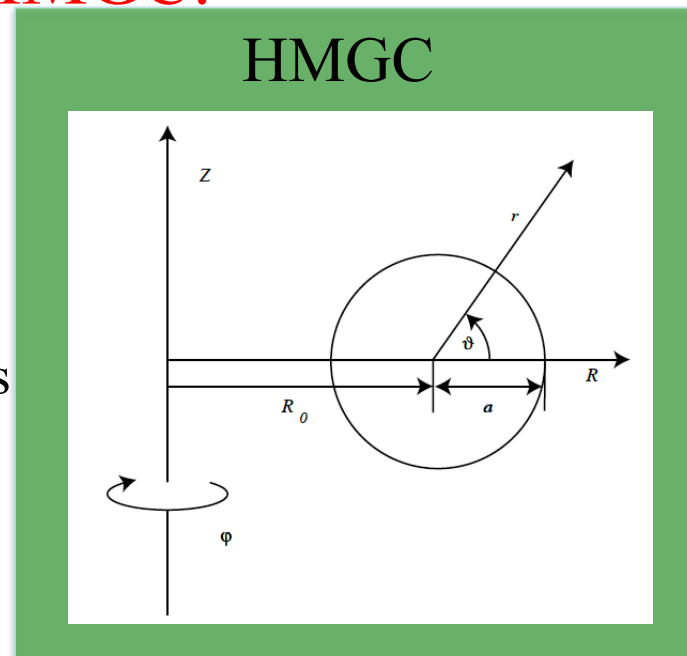
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## • WP10-ITM-IMP5-ACT5: HMGC

### – Status and recent developments of HMGC:

- Routinely runs on the Gateway as a stand-alone module
- Description of initial fast particle distribution function in the space of constant of motion
- The new version of HMGC can have two species of kinetic particles with different (anisotropic) initial distribution functions (e.g., slowing down (NBI, alphas) and bi-Maxwellian (ICRH))
- Added thermal ion compressibility and diamagnetic effects in addition to EP kinetic behaviours (eXtended HMGC, XHMGC)



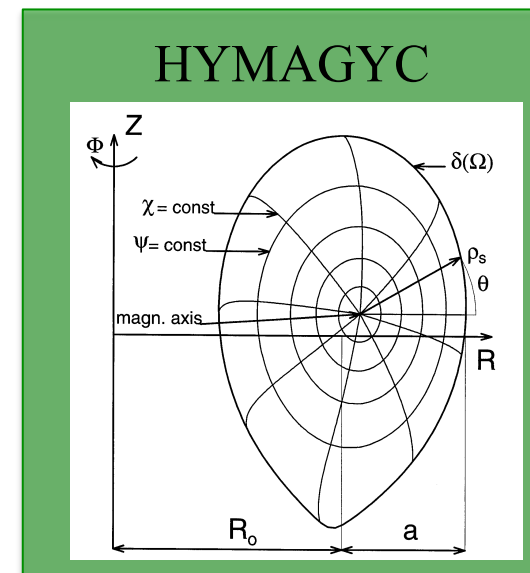
### – Deliverables before end 2010:

- A simple interface to equilibrium CPOs

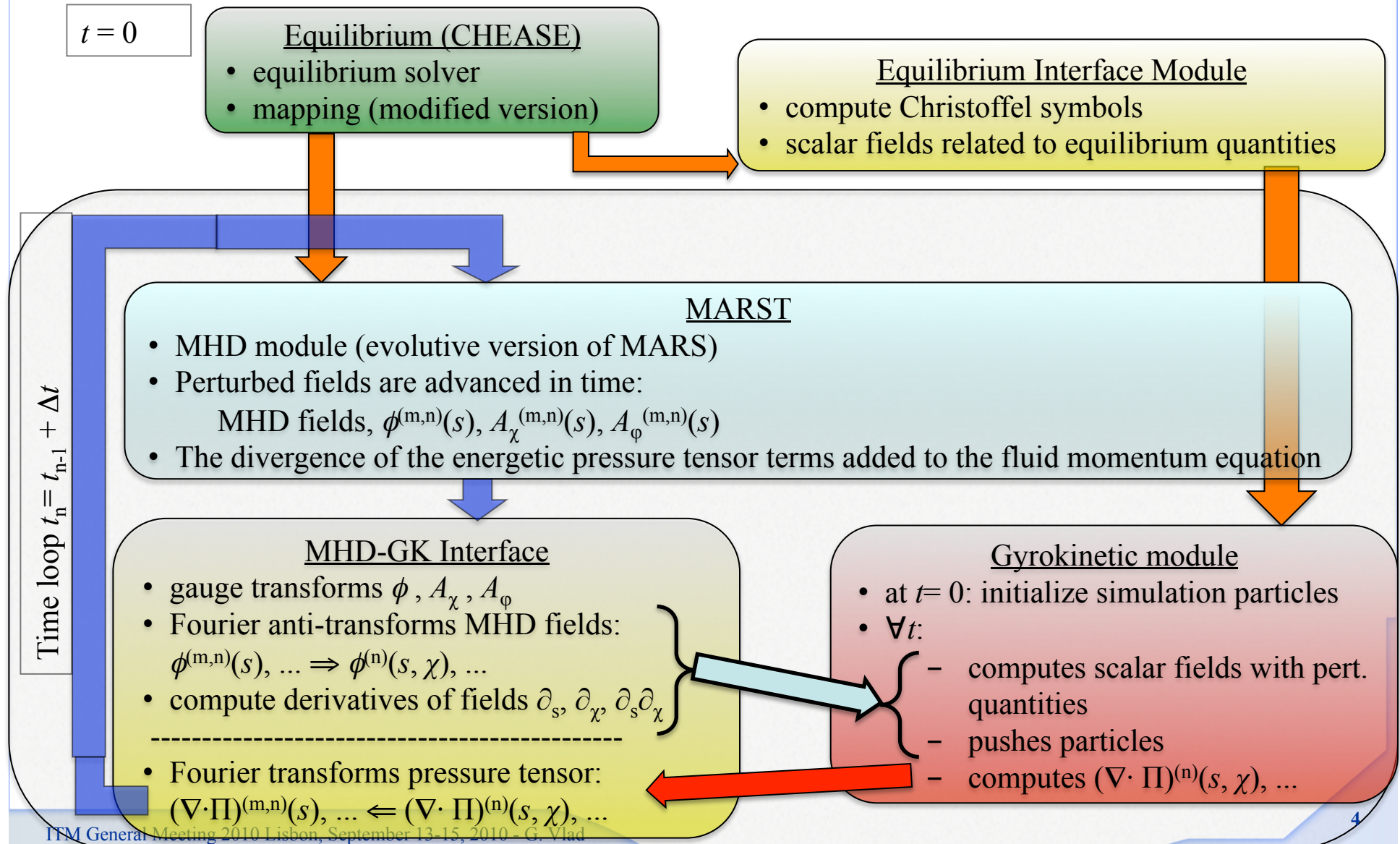
## • WP10-ITM-IMP5-ACT5: HYMAGYC

The new Frascati hybrid MHD-Gyrokinetic code, [HYMAGYC](#):

- Thermal (core) plasma:
  - described by full, resistive MHD linear equations
  - e.m. potentials required by Gyrokinetic module:  $\mathbf{A}$ ,  $\phi$
  - Fluid nonlinearities will not be retained
- Energetic-ion population:
  - particle gyrocenter-coordinates are evolved by solving gyrokinetic eqs. up to order  $O(\varepsilon^2)$  and  $O(\varepsilon\varepsilon_B)$
  - perturbed quantities satisfy the nonlinear gyrokinetic ordering of Frieman-Chen, Phys. Fluids (1982) 23, 502  
 $\omega/\Omega_E \approx k_{\parallel}\rho_E = O(\varepsilon)$ ,  $k_{\perp}\rho_E = O(1)$   
 (with  $\Omega_E$  the Larmor frequency and  $k_{\parallel}$  the component of the wave vector parallel to the magnetic field)
  - returns energetic particles pressure tensor  $\Pi^{ij}$  computed in terms of the particle distribution function in gyrocenter coordinates
- Flux coordinates system  $(s, \chi, \varphi)$



## • WP10-ITM-IMP5-ACT5: HYMAGYC



- WP10-ITM-IMP5-ACT5: HYMAGYC
  - parallelization with a OpenMP (inter-node) + MPI (intra-node) scheme
  - run as standalone program on the Gateway
  - MHD module fully integrated with equilibrium CPOs (thanks to IMP12 collaboration) *see later for details*
  - Fast particles module currently under testing
  - integration with fast particles CPOs to be done

- Inputs for MARST module from equilibrium CPOs (in order to substitute the standard input prepared by CHEASE)

– Quantities required by MARS/MARST module:

- integer and half-integer flux variable mesh:  $s \equiv \sqrt{\frac{|\psi - \psi_0|}{|\psi_{\text{edge}} - \psi_0|}} \in [0, 1]$

- Fourier components of the covariant metric tensor elements; CPOs contain the contravariant elements (thus, the matrix needs to be inverted):

$$g^{ss} = \left( \frac{\partial s}{\partial \psi} \right)^2 g^{\psi\psi} = \left( \frac{\partial s}{\partial \psi} \right)^2 |\nabla\psi|^2$$

$$g^{s\chi} = g^{\chi s} \left( \frac{\partial s}{\partial \psi} \right) g^{\psi\chi} = \left( \frac{\partial s}{\partial \psi} \right) \nabla\psi \cdot \nabla\chi$$

$$g^{\chi\chi} = |\nabla\chi|^2$$

$$g^{\phi\phi} = |\nabla\phi|^2 \equiv \frac{1}{R^2}$$

- $\chi$  poloidal like angular variable, such that Jacobian is:  $J_s = [(\nabla s \times \nabla\chi) \cdot \nabla\phi]^{-1}$
- Equilibrium magnetic field, current, pressure:  $B^\chi, B^\phi, j^\chi, j^\phi, p, dp/ds$
- A vacuum mesh (simple model for free boundary modes)

# Energetic particle activities

- Issues arisen during equilibrium CPOs interface (1)
  - 1D profile quantities `equilibrium%profiles_1d%...`  
 (e.g., `%q`, `%psi`, `%F_dia`, `%ffprime`, `%p`, `%pprime`):  
 data are “almost equispaced” in the variable  $s \propto \sqrt{\psi}$
  - Coordinate system 2D quantities:

<code>equilibrium%coord_sys%</code>	physical quantity	comment
<code>grid%dim1</code>	$\psi$	“almost equispaced” in $s$ , in principle different from 1D definition
<code>grid%dim2</code>	$\chi$	$\chi$ mesh $\chi = (ndim2 - 1)/ndim2 * 2\pi$
<code>grid_type</code>	description	<i>Straight field line</i> : the only choice for HELENA, which implies also $J = \frac{qR^2}{2\pi F_{dia}}$ <i>Non-straight field line</i> : choice available for CHEASE, $J_\psi = C(\psi)R^{NER} \nabla\psi ^{NEGP}$
<code>position%r</code> <code>position%z</code>	$R(\psi, \chi)$ $Z(\psi, \chi)$	coordinates in the poloidal cross section
<code>jacobian</code>	$J(\psi, \chi)$	Jacobian
<code>g_11</code> , <code>g_22</code> , <code>g_33</code> , <code>g_12</code>	$g^{ss}(\psi, \chi)$ , $g^{\chi\chi}(\psi, \chi)$ , $g^{\phi\phi}(\psi, \chi)$ , $g^{s\chi}(\psi, \chi)$	contravariant(!) metric tensor elements

# Energetic particle activities

- Issues arisen during equilibrium CPOs interface (2)
  - Regular mesh don not means “equidistant mesh”!
  - signs for  $I_p$  and  $B_0$  must be taken into account (note that usual MHD stability codes ignore those signs, whereas CPOs do not!)
  - Only one mesh is provided in the CPOs (the integer one, which includes the magnetic axis and the plasma boundary), thus interpolation is required: note that some of the metric quantities could diverge toward the origin, care should be taken in interpolating such quantities!
  - Also, interpolation of 1D quantities on the `dim1` mesh is required, in order to compute all the quantities required by MARS/MARST
  - Several issues on the Jacobian definition, choice of the coordinate system for the stability mesh, etc. have been considered and agreed between ITM (see WEB page [https://www.efda-itm.eu/~wwwimp3/TEST/ITM/html/itm\\_conventions.html](https://www.efda-itm.eu/~wwwimp3/TEST/ITM/html/itm_conventions.html))
  - Not jet available in equilibrium CPOs:
    - maps  $\psi(R, Z)$ ,  $\chi(R, Z)$  (ready from IMP12?)
    - Christoffel symbols (require derivatives of the metric tensor elements)



- Gateway MARS/MARST version
  - The exercise of modifying the MARS/MARST MHD module used by HYMAGYC in order to read the equilibrium quantities from the CPOs is almost completed
  - a C preprocessor produces the FORTRAN for standard use or the one required to interface the code with the CPOs:
    - transforms the module from a “main” to a “subroutine(equilibrium\_in)”
    - declares `type_equilibrium`, and `interface` quantities
    - the standard input routine is substituted by the ones which reads the CPOs and computes all the derived quantities (subroutine `input_gw(equilibrium_in)`, which, on turn, call the module `read_eq_CPOs_MARS.f90`)
  - A driver to run and test the Gateway version of the code reading an equilibrium CPOs (`driver_mars_fftw_gw_linux.f90`) has been written
  - Tested on CPOs produce by HELENA and CHEASE
  - Still to be done:
    - write the XML code specific parameters inputs
    - write the output results to the MHD stability CPOs