

ITER example classifying AE-EP transport scenarios: workflow L1 for standard 15 MA scenario

[Polevoi, 2002]



 degree of (non-linear) Alfven eigenmode (AE) resonance overlap will determine the nature of EP transport [Berk 1995] in ITER and DEMO

example: I 5MA ITER scenario: linear TAE-α resonances depend strongly on q0: strongly overlapping (I), intermediate (II) and scarce

(III) TAEs spectra can exist here: only linearly unstable modes are shown

(N.B: for small particle orbits, $P\phi$ and s 1.15 are similar, thus the radial mode overlap is a proxy for resonance overlap)



on-axis safety factor

IPP

automated HAGIS/LIGKA workflow L2/L3/L4 for ITER 15 MA including FLR/FOW

including some global information (kr,k \perp) in local model:



10-100 times faster than global solver: analytical finite orbit width version of LIGKA based on analytical theories [Zonca 1996/98]; benchmarked in relevant limit [Lauber 2018]:

reasonable agreement for intermediate and high mode numbers

IPP

automated HAGIS/LIGKA workflow HI including FLR/FOW



runtime: minutes - few hours for whole range of mode numbers global effects crucial, non-perturbative features observed; no constant between A~(γ/w)² IPP

automated HAGIS/LIGKA workflow H2 including FLR/FOW



automated HAGIS/LIGKA workflow H2 including FLR/FOW



above simulations are 'worst case':

reduced amplitudes found when zonal flow dynamics is included (forced excitation)! [Todo, 2011, Biancalani, 2016, Vlad 2017]

recently: similar behaviour found using global ORB5 [T. Hayward-Schneider, 2019] (in contrast to other studies [M. Fiztgerald, 2015] not observing this transition)



the LIGKA model equations

$$f = h + H_1 \frac{\partial F_0}{\partial E} - [e \frac{\partial F_0}{\partial E} - \frac{c \nabla F_0}{i \omega B} \cdot (\mathbf{b} \times \nabla)] J_0 \psi,$$

comparison analytical theory - numerical integration; can be calculated including FLR/FOW effects (100 times faster!)



similar dependence for analytical expression vs numerical result
D and T damping differ - exponential dependence of ion LD!
numerical damping is typically smaller than analytical expression (assumption v=v//)

 $v_{th,e}/v_{A0} \sim 13$ $v_{th,D}/v_{A0} \sim 0.2$ $v_{th,T}/v_{A0} \sim 0.16$ $v_{th,He}/v_{A0} \sim 0.14$ $v_{th,Be}/v_{A0} \sim 0.09$