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# **Alpha channeling, edge turbulence, A&M reactions: a journey into the physics of the tokamak**

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Agenzia nazionale per le nuove tecnologie,  
l'energia e lo sviluppo economico sostenibile

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## Career:

- **2008**: PhD in Relativistic Astrophysics, University of Rome “Sapienza”
- **2008-2016**: some postdocs across Europe (Universities of Rome “Sapienza”, Erlangen-Nürnberg, Wrocław)
- **2017-2019**: postdoc at University of Rome “Tor Vergata”
- **2019-2021**: postdoc at PIIM laboratory, CNRS/Aix-Marseille University
- **2021-2023**: postdoc at Jülich Forschungszentrum
- **2023-?**: researcher at ENEA Frascati

# Summary

- Alpha Channelling  
EUROfusion grant at University of Roma Tor Vergata 2017-2019
- Simulations of Edge Plasma Oscillations  
Postdoc at PIIM Laboratory (CNRS, AMU) 2019-2021
- Collisional Radiative Models in EIRENE simulations  
Postdoc at Forschungszentrum Julich 2021-2023 (TSVV5)
- Turbulence simulations near the X-point  
ENEA (since June 2023)

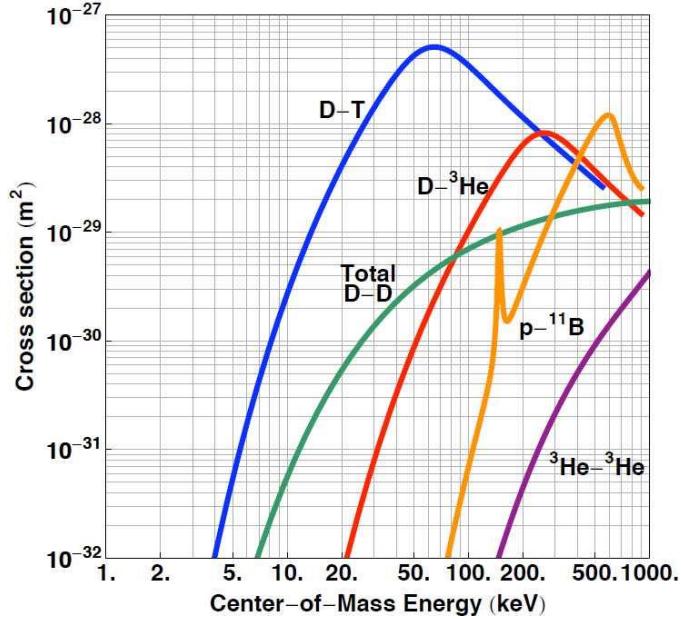


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# Alpha Channelling

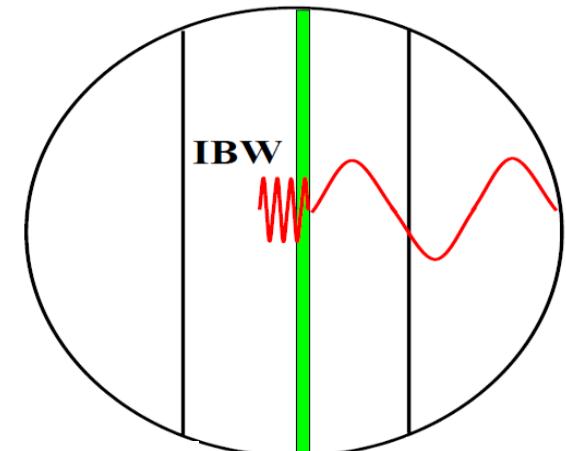
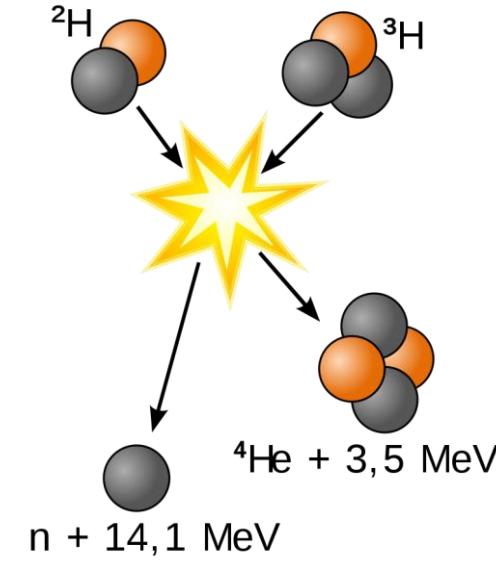
# Motivation: can we heat bulk ions directly through alpha particles?

## Fusion reactions:



Classically alpha particles mainly slow-down on electrons, but in a reactor the best performance is achieved in HOT ION mode.

The idea of alpha channelling is to use some waves (IBWs) that can extract the kinetic energy of alpha particles (through stochastic interaction) and release it to bulk ions (via cyclotron interaction).



# Stochastic cooling

$$\Delta\mu = \frac{ZeN}{m\omega} \Delta\varepsilon$$

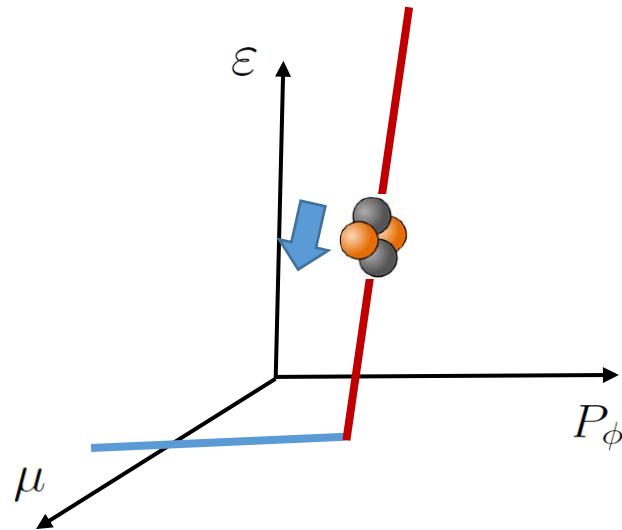
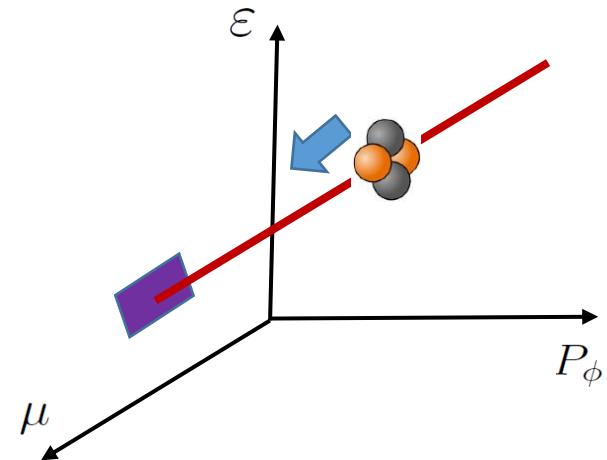
$$\Delta P_\phi = \frac{n_\phi}{\omega} \Delta\varepsilon$$

Original idea: drive most of alpha particles to the plasma boundary

$$\left. \begin{array}{l} \Delta P_\phi \sim \frac{Ze}{c} \psi^w \\ \Delta\varepsilon \sim \varepsilon_\alpha \end{array} \right\} \quad \xrightarrow{\hspace{1cm}} \quad n_\phi \sim 1000$$

too high toroidal  
wave number (not  
realizable, not even  
efficient)

[Fisch and Rax *PRL* 92, Snyder, Hermann and Fisch *JFE* 94]



Two waves scenario: more advantageous to use two waves:

\_ IBW to extract energy

\_ low-frequency wave (TAE?) to move particles to the wall

[Fisch and Hermann *NF* 95, Hermann and Fisch *PRL* 97]

# Goals of the EUROfusion research grant

Very limited evidences of alpha channelling have been found so far both in experiments and in numerical simulations. [Darrow et al *NF* 96, Clark and Fisch *PoP* 00, Wong et al. *PRL* 04] [Cook et al *PRL* 10, Cook et al *PPCF* 11, Ochs and Bertelli and Fisch *PoP* 15]

The goals of the project are:

- *the quantitative determination of the amount of fusion alpha power that can be transferred to IBW*
- *the determination of the conditions for minimal absorption by the electron and optimal transfer of the energy from IBW to the thermal ions and of the required injected power.*
- *the optimization of the characteristic of the Alfvénic turbulence to maximize the effect of alpha channelling*

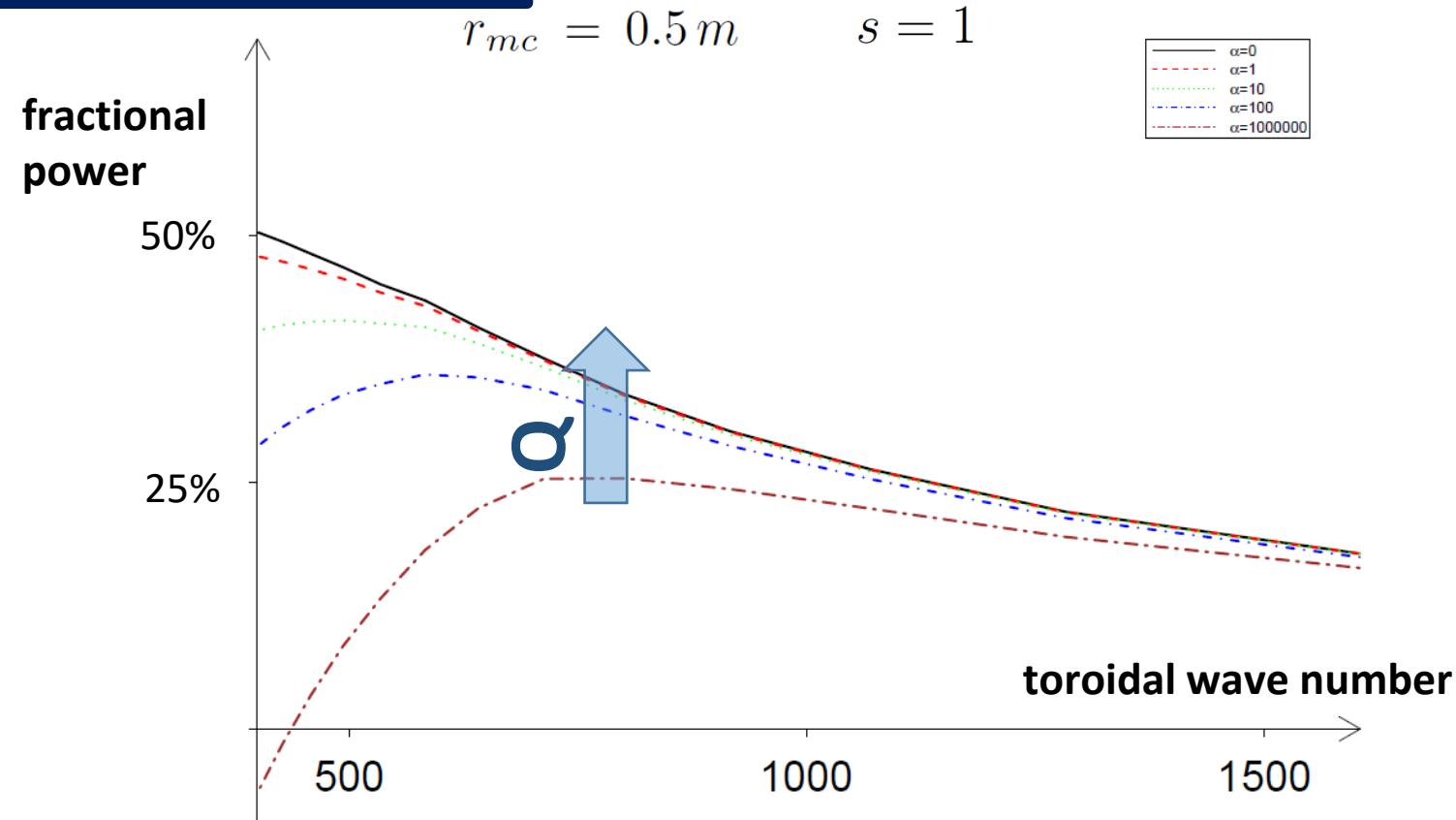
# Results: fractional power vs toroidal wave number

## Q: outward flux to model the low-frequency wave

By increasing the low-frequency wave flux the maximal power increases and moves towards low toroidal wave numbers.

**Only the scenario with two waves can provide efficient alpha channeling.**

In the optimal case, about **55% of alpha particle energy** can be extracted by IBWs.

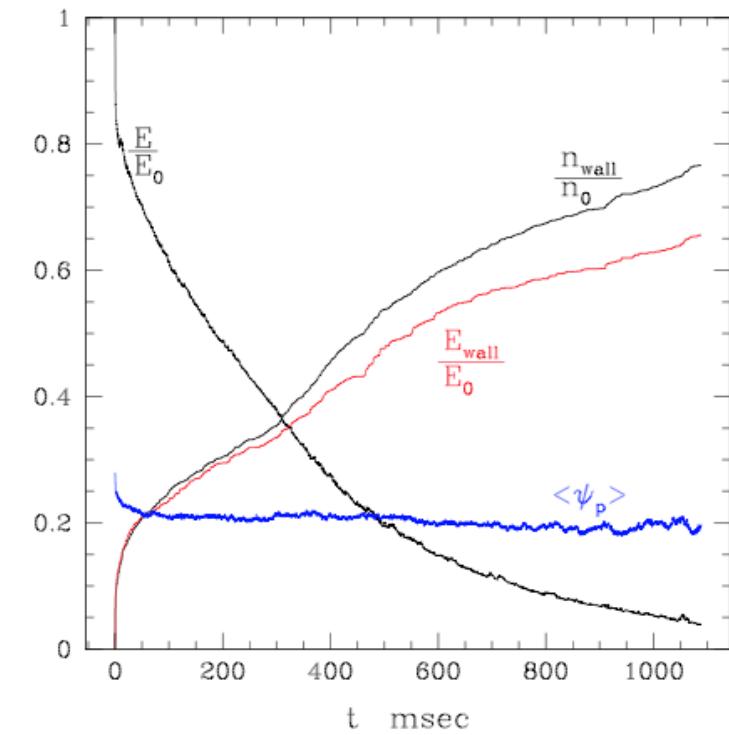


[Cianfrani and Romanelli *NF* 18]  
[Cianfrani and Romanelli *NF* 19]

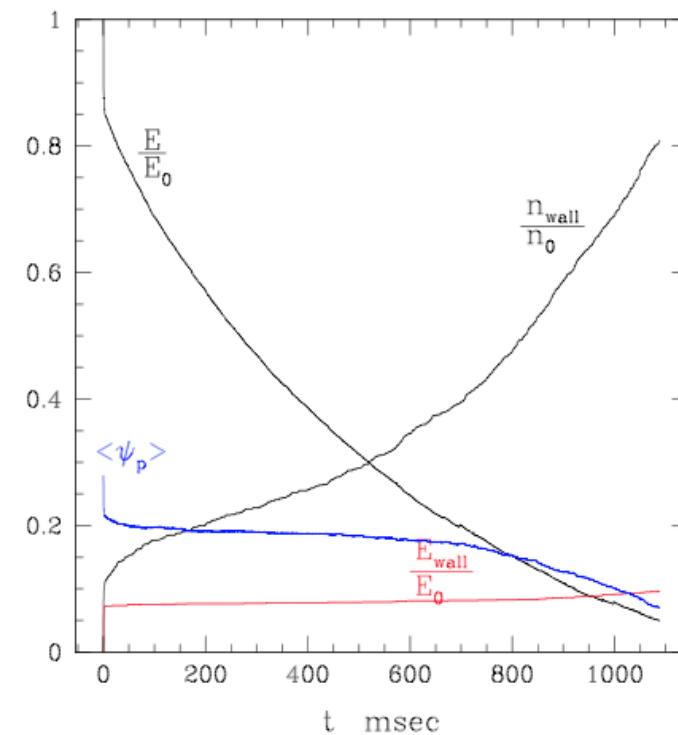
# ORBIT simulations

The results on the alpha particles cooling have been **confirmed** by the simulations performed via the ORBIT code.

[White, Romanelli, Cianfrani and Valeo PoP 21]



These simulations also outlines how **TAEs** provides **too high losses** of particle energy (over ten percent of the particle energy to be deposited on the walls). **high ash removal high energy losses**



Promising results with **micro-turbulence** playing the role of the second low-frequency wave.

**high ash removal low energy losses**

# Simulations of Edge Plasma Oscillations

**EMEDGE3D**: two-fluid code simulating the edge of a tokamak plasma (constant density, equal electron and ion temperatures and a constant toroidal external magnetic field).

Reduced MHD equations for electronic pressure, electrostatic potential, electromagnetic potential (drift ordering, including **DIAMAGNETIC VELOCITY**)

Charge Balance  $\partial_t \nabla_{\perp}^2 \phi + \{\phi, \nabla_{\perp}^2 \phi\} = -\nabla_{\parallel} \nabla_{\perp}^2 \psi - \omega_D \mathbf{G} p + \mu_{\perp} \nabla_{\perp}^4 \phi$

Energy Balance  $\partial_t p + \{\phi, p\} = -\Gamma \nabla_{\parallel} \nabla_{\perp}^2 \psi + \delta_c \mathbf{G}(\phi - C_{dia} p) + \chi_{\perp} \nabla_{\perp}^2 p + S$

Ohm's Law  $\beta_e \partial_t \psi = -\frac{1}{\epsilon_{\parallel}^2} \nabla_{\parallel}(\phi - C_{dia} p) + \eta \beta_e \nabla_{\perp}^2 \psi$

Curvature

ExB advection

Parallel current div

Viscosity

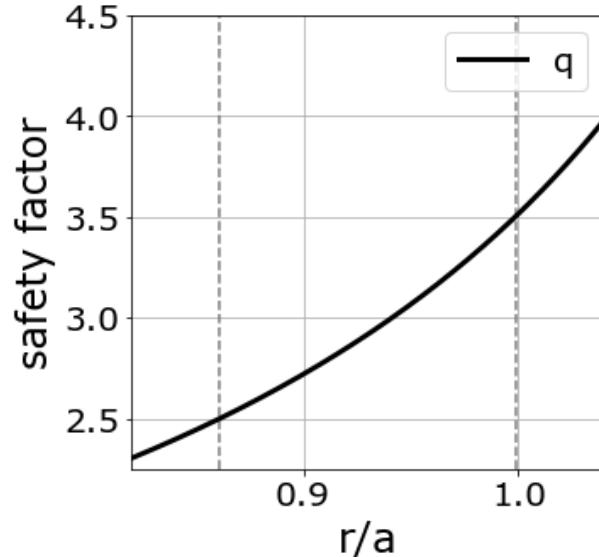
Diffusion

Resistivity

Source

## Medium-size tokamak

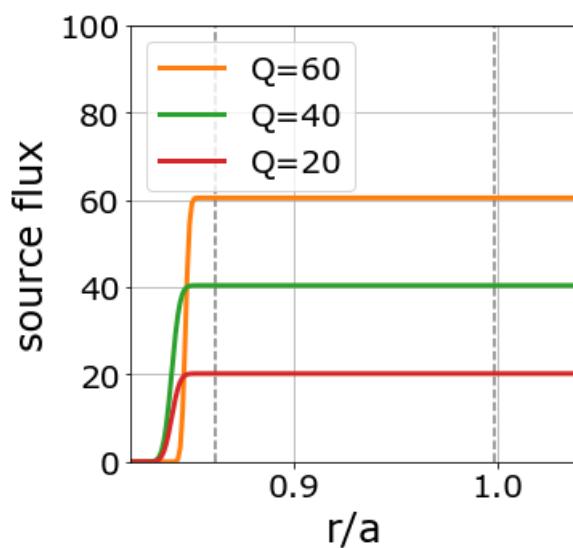
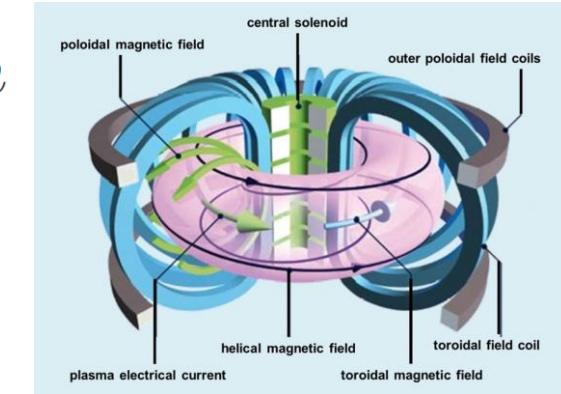
$$B_0 = 1T \quad R_0 = 1.75m \quad a = 0.45m$$



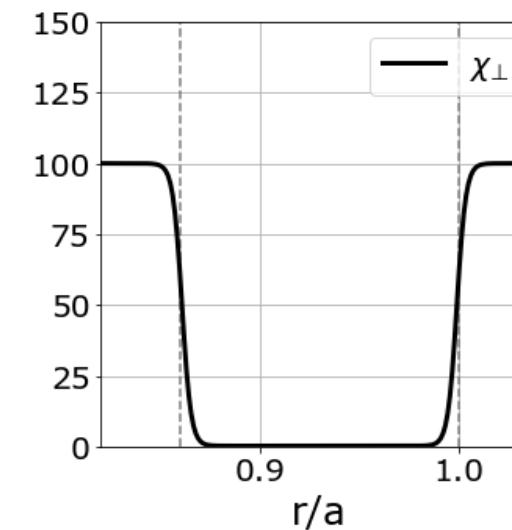
$$\begin{aligned} n &= 2.5 \times 10^{19} m^{-3} \\ T_i = T_e &= 50 eV \\ \beta_e &= 0.2\% \end{aligned}$$

typical values  
for a JET  
discharge close  
to LH transition

[Bourdelle, NF 2014]

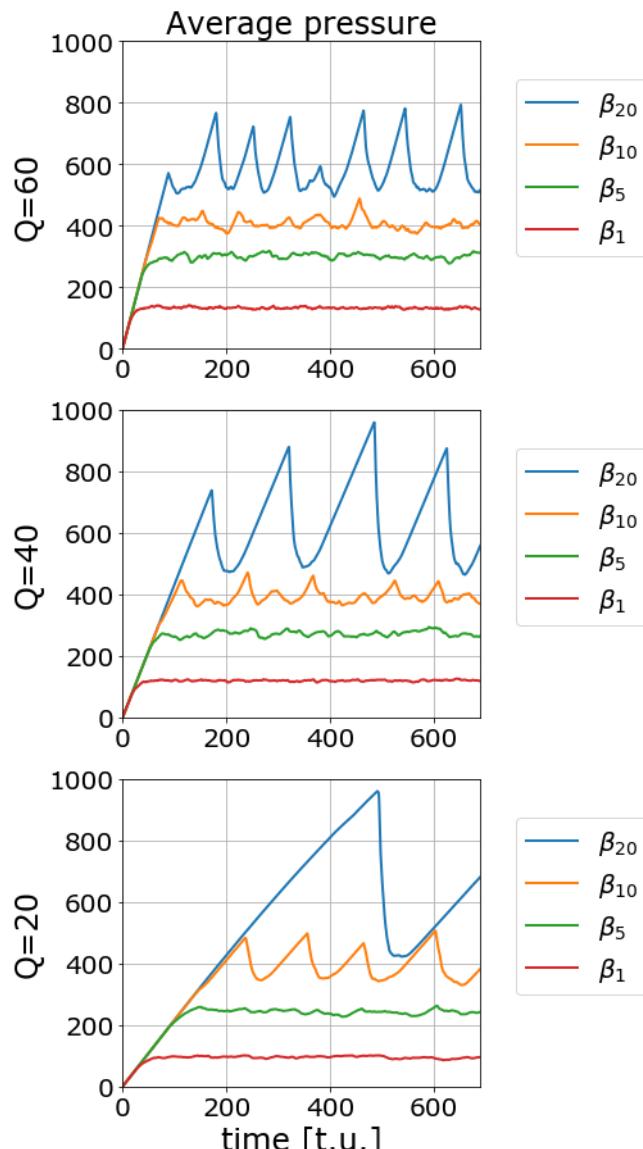


source in the  
inner buffer



dissipation is  
sharply increased  
in order to quench  
turbulence (high  
perpendicular  
diffusivity).

# Results: oscillations at low $\beta_e$

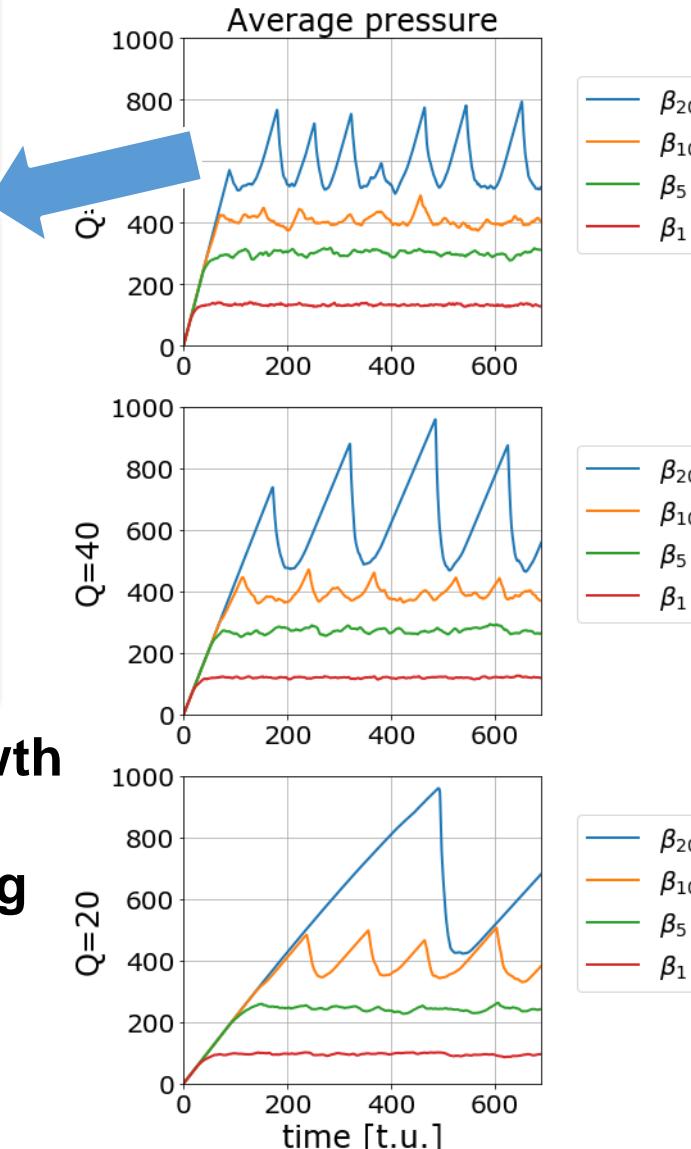
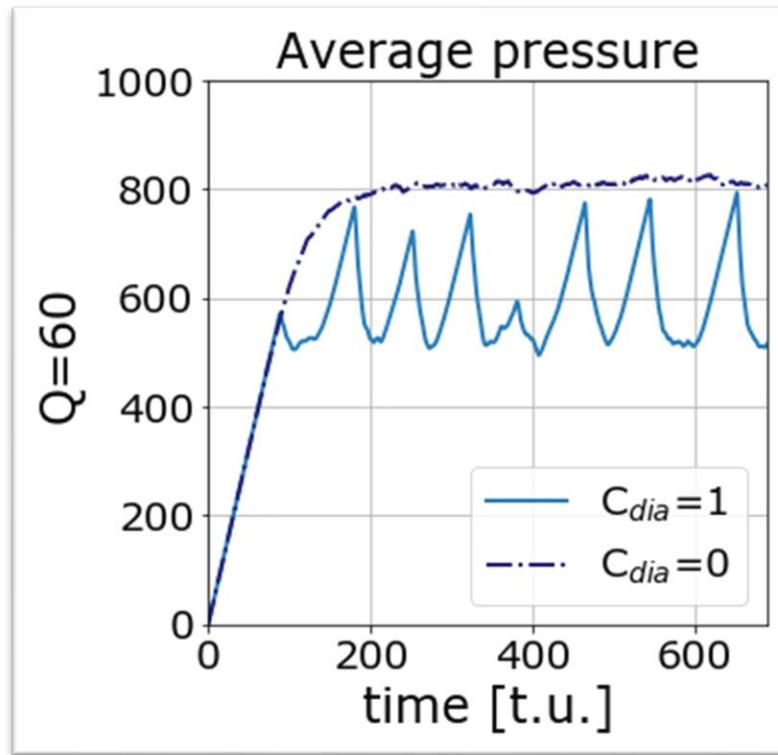


$$\beta_d = \beta_e/d$$

low  $\beta_e$   
↓  
low density

$100 \text{ t.u.} = 0.61 \text{ ms}$

# Results: oscillations at low $\beta_e$



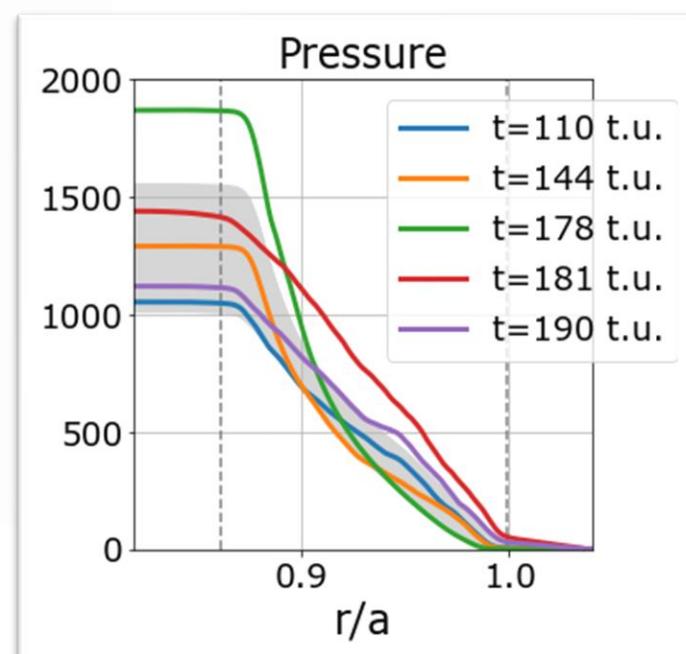
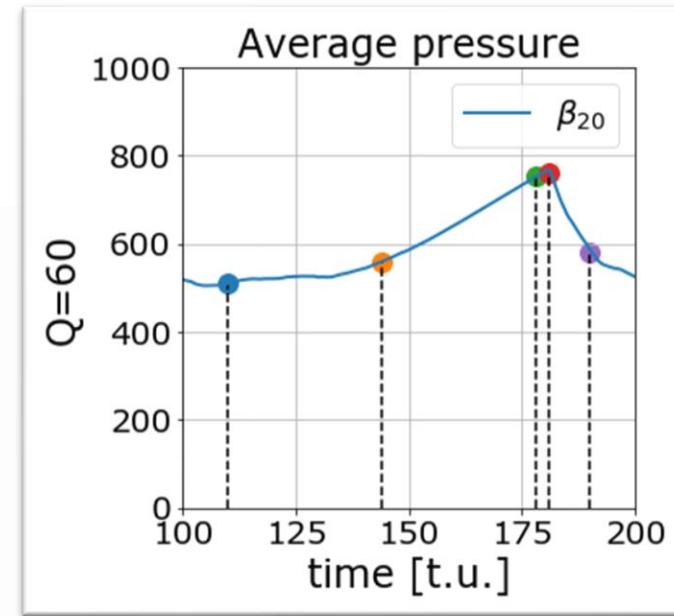
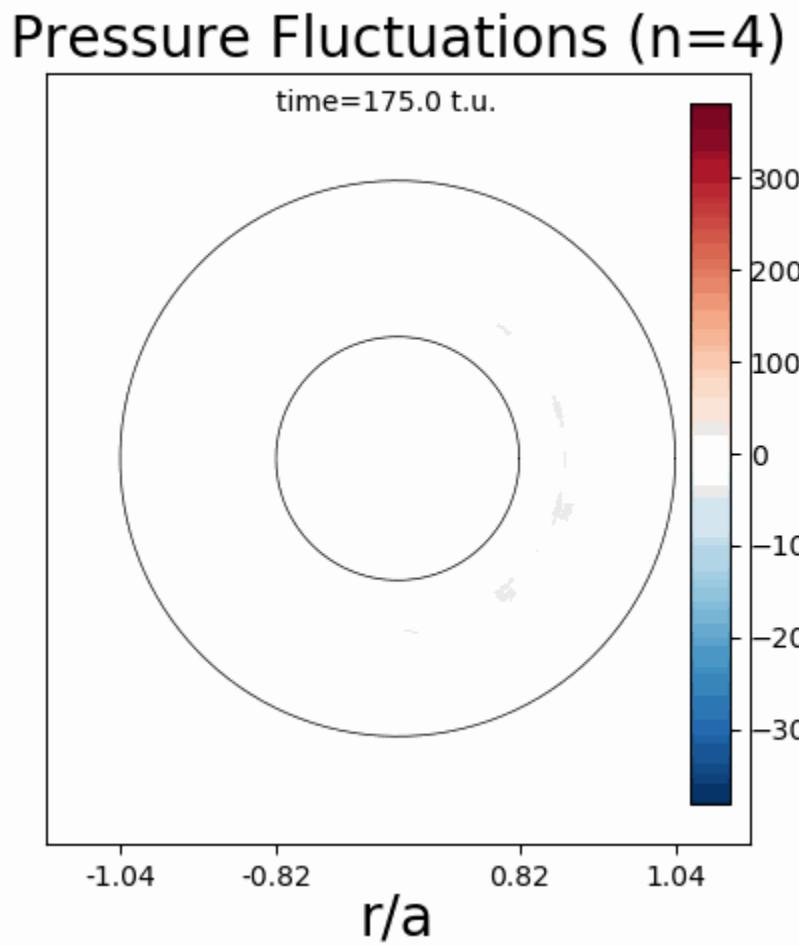
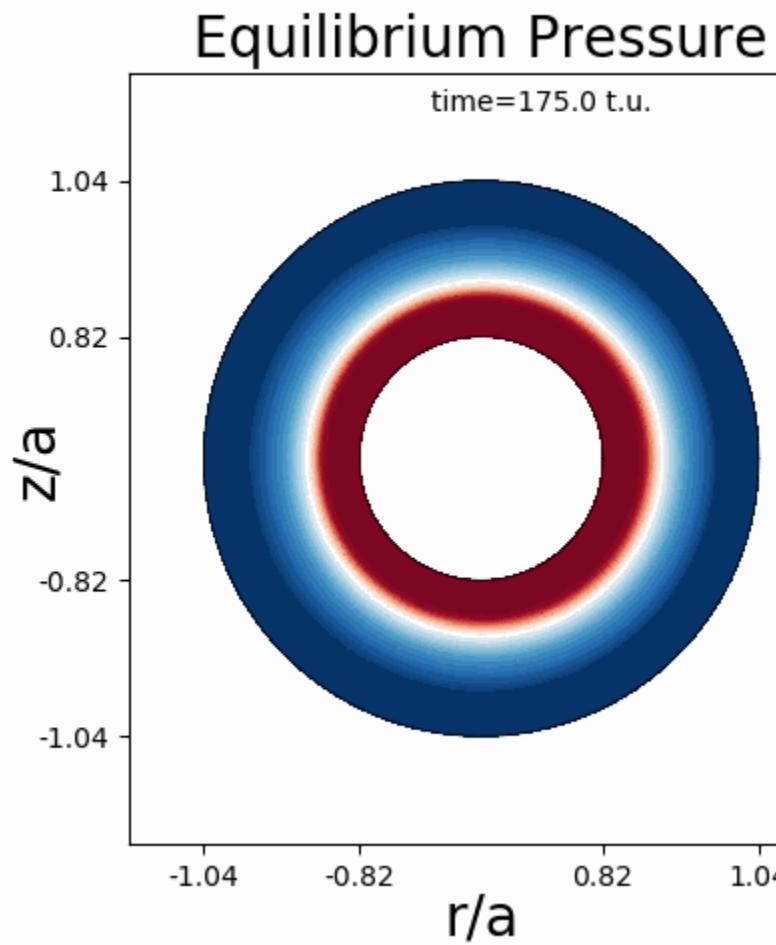
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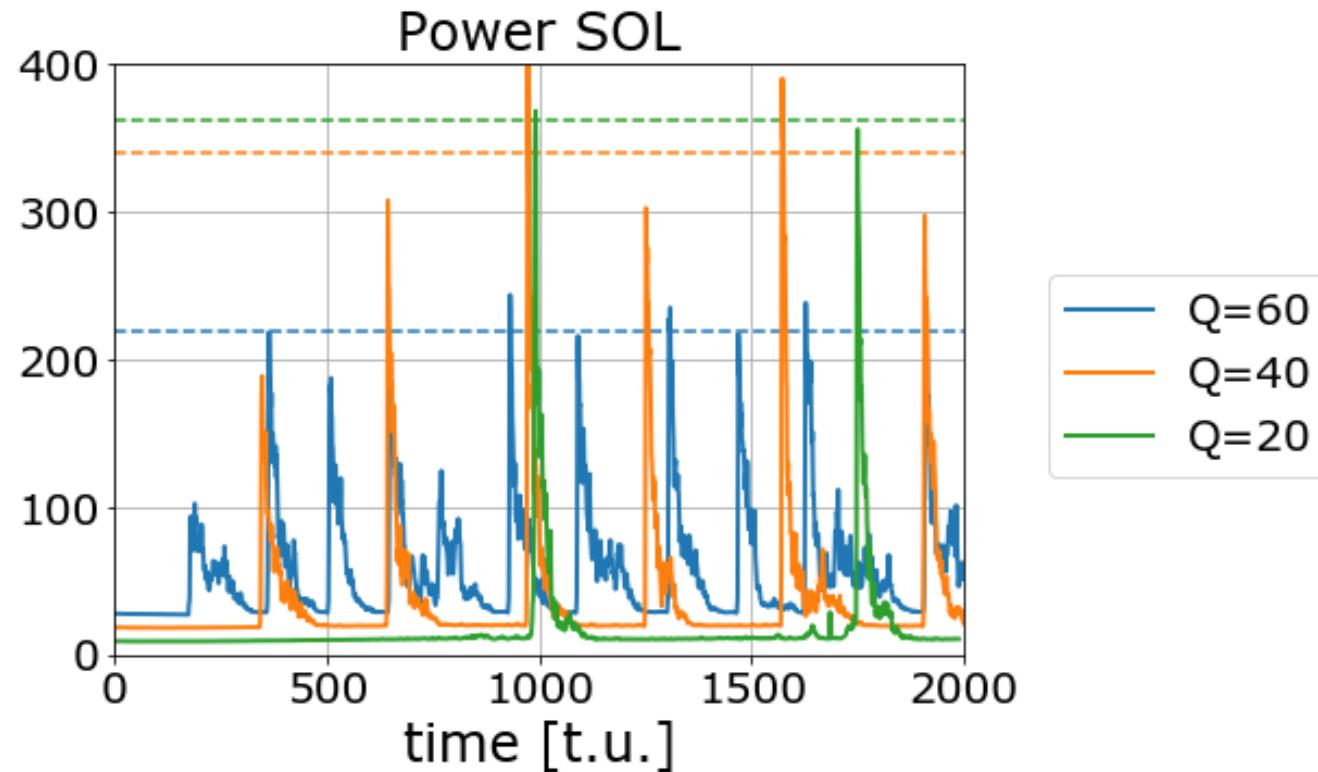
oscillations due to explosive growth  
of unstable modes previously  
stabilized by diamagnetic coupling  
(relaxation of a transport barrier).

$100 \text{ t.u.} = 0.61 \text{ ms}$

## 2D Animations



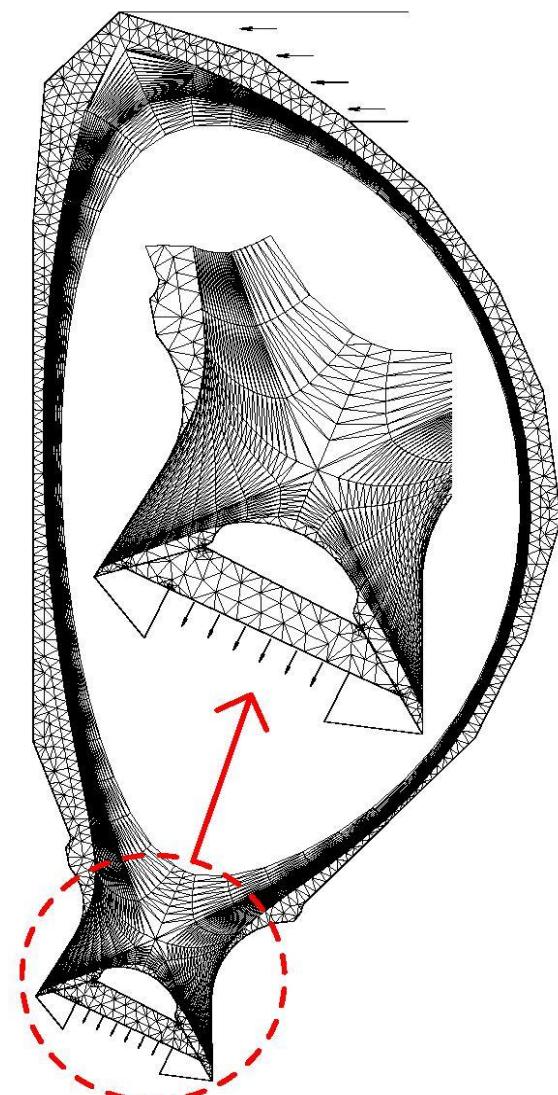
# Results: type-III ELM



the oscillation frequency decreases  
with increasing energy flux into SOL

# **Collisional Radiative Models in EIRENE simulations**

# What is EIRENE in a nutshell (e.g. in SOLPS)?..



## Macroscopic:

CFD codes  
(computational fluid  
dynamic):  
B2, Edge2D, EMC3,  
SOLEdge3X, etc...

## Plasma flow parameters

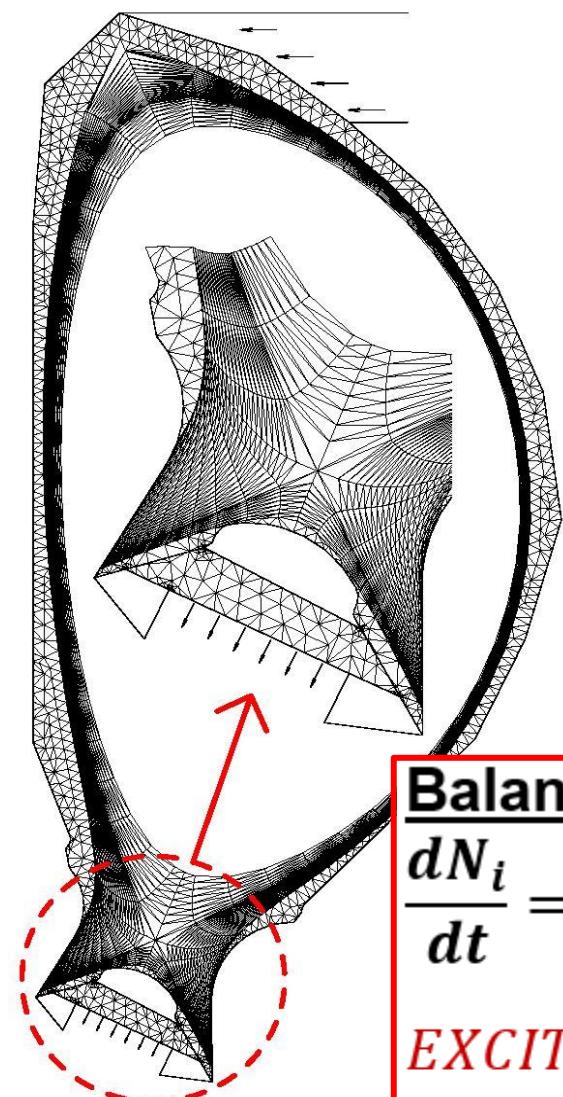
## Microscopic

**EIRENE-NGM:** a 3D  
Monte Carlo multi-  
species transport  
code incl. radiation  
transfer, kinetic  
or **F-K** hybrid.

## Source terms (Particle, Momentum, Energy)

CRMs for atomic and  
molecular neutrals H,  
 $H^*$ ,  $H_2$ ,  $H_2(v)$ ,  $H_2^*$ ,  $H^-$ ,  
 $H_2^+$ , ..., impurities.  
Ionisation, CX,  
recombination etc.

# What is EIRENE in a nutshell (e.g. in SOLPS)?..



## Macroscopic:

CFD codes (computational fluid dynamic): B2, Edge2D, EMC3, SOLEdge3X, etc...

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## Microscopic

**EIRENE-NGM:** a 3D Monte Carlo multi-species transport code incl. radiation transfer, kinetic or **F-K** hybrid.

## Source terms (Particle, Momentum, Energy)

CRMs for atomic and molecular neutrals H, H\*, H<sub>2</sub>, H<sub>2</sub>(v), H<sub>2</sub><sup>\*</sup>, H<sup>·</sup>, H<sub>2</sub><sup>+</sup>, ..., impurities ionisation, CX, recombination etc.

## Balance equations:

$$\frac{dN_i}{dt} = \sum_{j \neq i} A_{ji} N_j + n_e \cdot (EXCIT + IZ + CX + REC)$$

$$EXCIT = \sum_{j \neq i} \langle v \sigma_{ji} \rangle N_j$$

$$IZ = \sum_m \langle v \sigma_{mi} \rangle N_m^- + \sum_z \langle v \sigma_{zi} \rangle N_z^{2-} + \dots$$

$$REC = \sum_k \langle v \sigma_{ki} \rangle N_k^+ + \sum_l \langle v \sigma_{li} \rangle N_l^{2+} + \dots$$

CRM = list of states + transition data

Often used:

$\langle v \sigma_{ji} \rangle (T_e, n_e)$ - effective Maxwellian averaged rates

# PLOUTOS: what is inside?..

- 1) CRM solver (transport excluded, pure A&M side of the problem) with extensive features
- 2) Plotting, solver results visualisation
- 3) Flexible reaction tables for EIRENE Databases (AMJUEL, HYDHEL, H2VIBR) interconnected with the solver and plotting

## *TSVV5 Activities*

- *grouping of reactions*: easier comparison btw data from different sources
- *new column/features*: sparse information now available
- *default cases*: testing purposes and model for molecular decay
- *json output*: reusability for new solver runs and other codes (EIRENE)

Conclusions Eirene Code Camp 22

we want to promote .. [PLOUTOS] .. as a standard pre-processing tool for Eirene

# PLOUTOS: what is inside?..

- 1) CRM solver (transport excluded)
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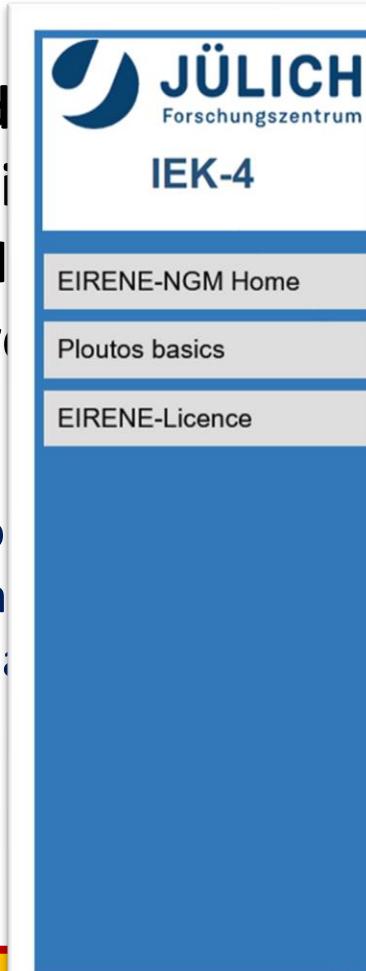
interconnected with the solver

## TSVV5 Activities

- **grouping of reactions:** easier comparison
- **new column/features:** sparse input
- **default cases:** testing purposes and examples
- **json output:** reusability for new applications

Conclusions Eirene Code Camp 22

we want to promote .. [PLOUTOS]



available in [eirene.de](http://eirene.de)  
(for registered users only).



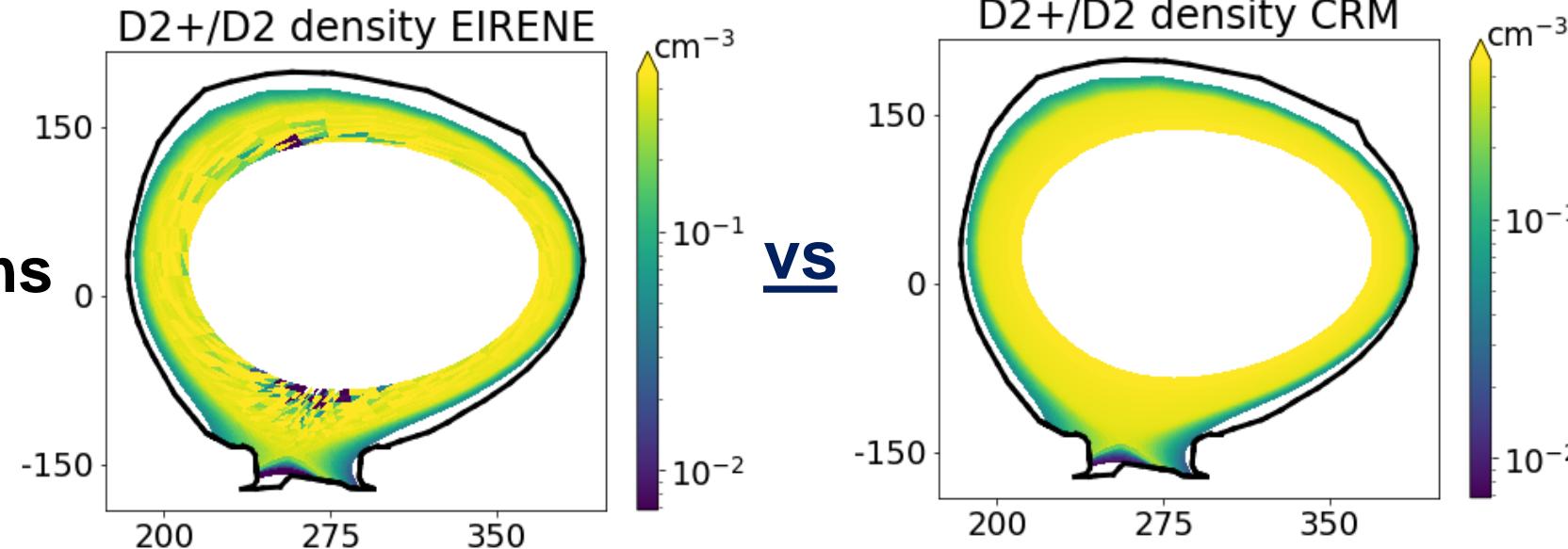
Statue of the baby Ploutos,  
personification of wealth.  
National Archeological Museum,  
Athens.

- Plot cross section and rates from AMJUEL, HYDHEL, H2VIBR.
- Assess data quality and origin.
- Solve stationary CRMs with spectral analysis and sensitivity analysis.
- Prepare A&M input file for EIRENE.

Enter Ploutos for H2-F.Cianfrani,2022-2023

# CRM vs EIRENE: TRANSPORT

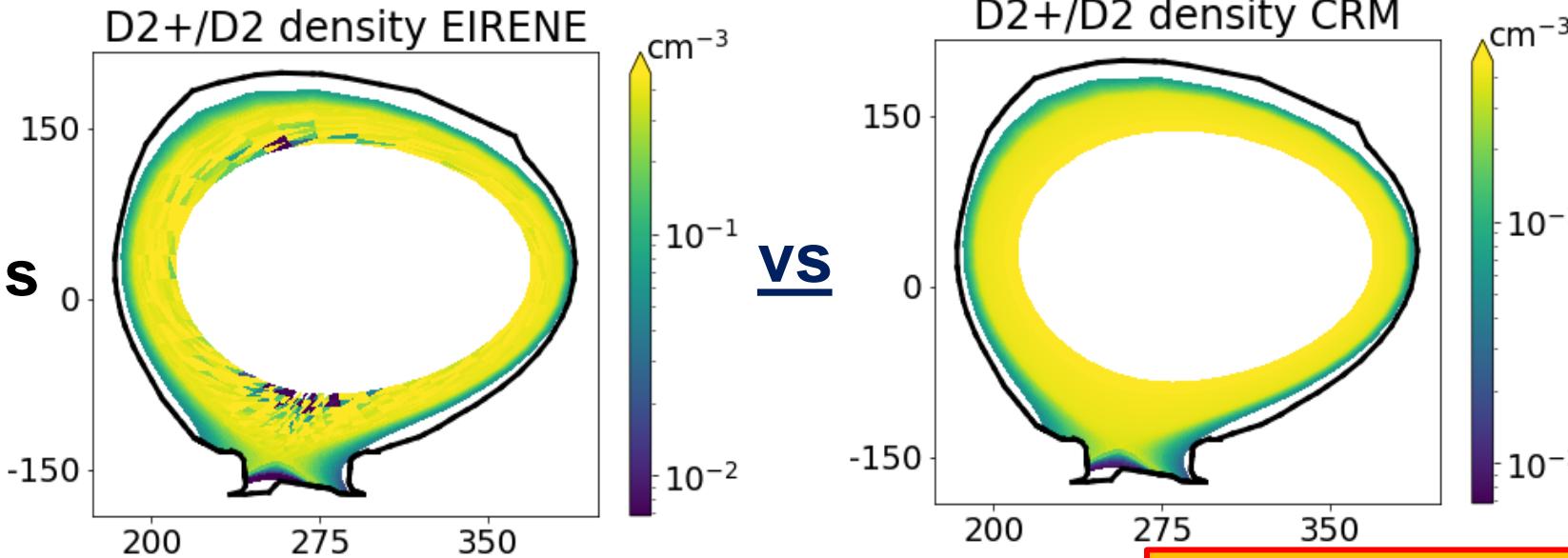
**EDGE2D-  
EIRENE  
simulations  
JET shot  
81472:**



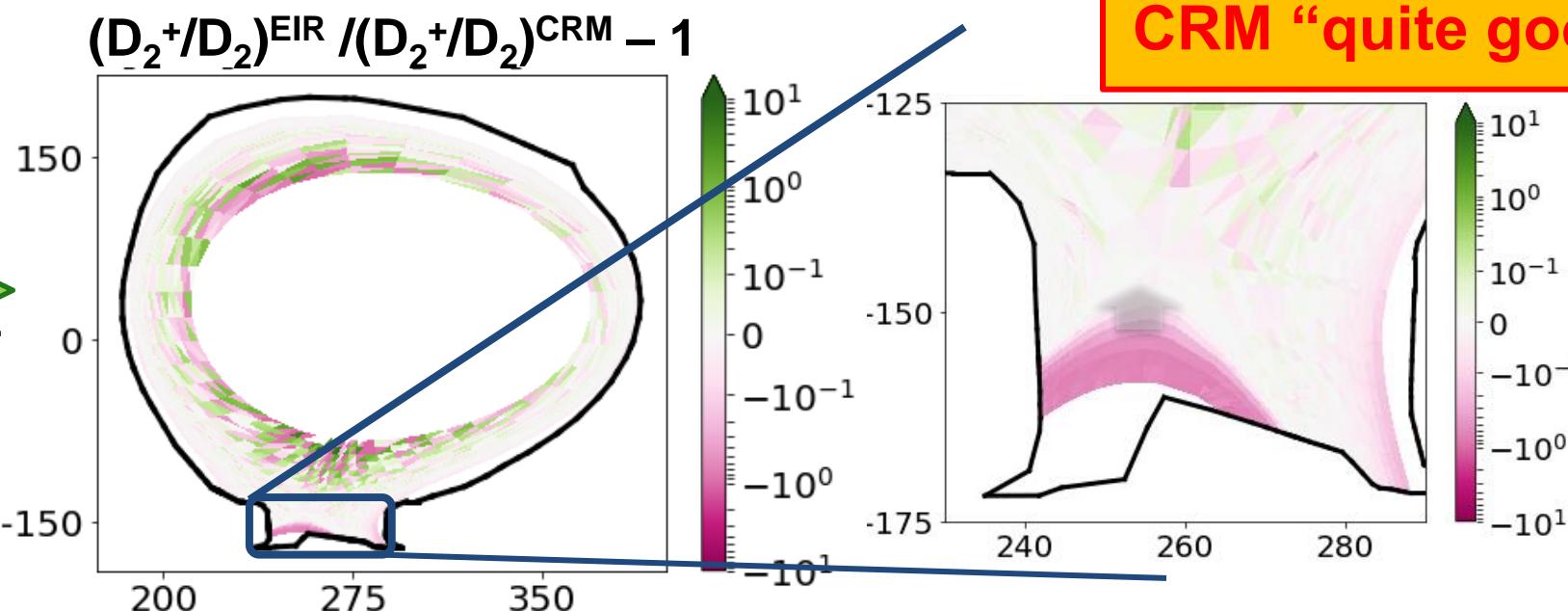
**CRM in each  
simulation cell**

# CRM vs EIRENE: TRANSPORT

EDGE2D-  
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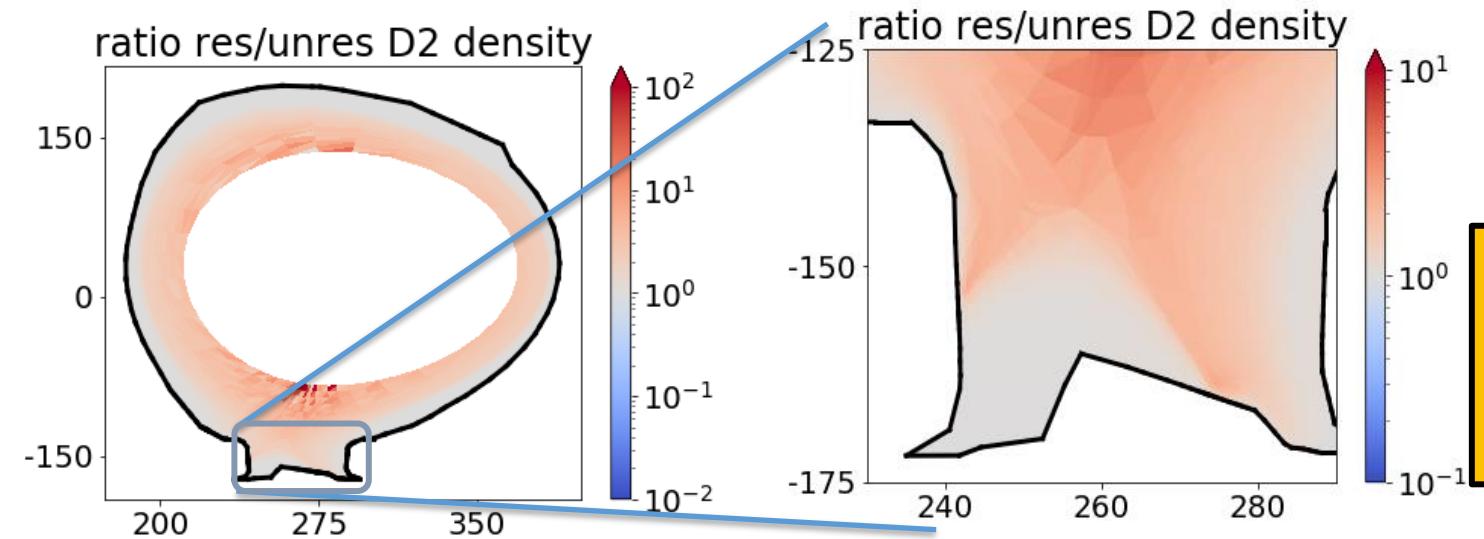


CRM “quite good” in SOL...

transport of  
molecules  
from private  
flux region

# Impact of vibrational resolution

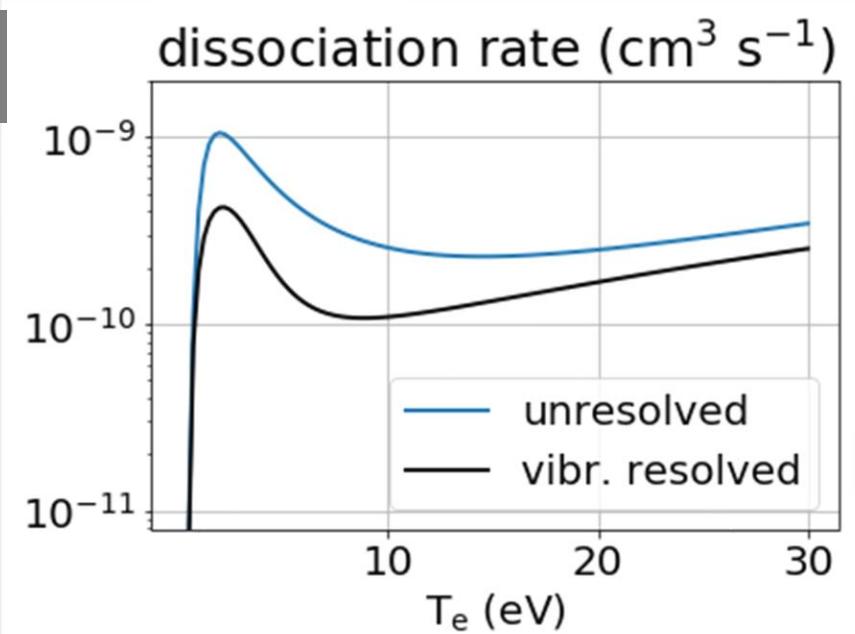
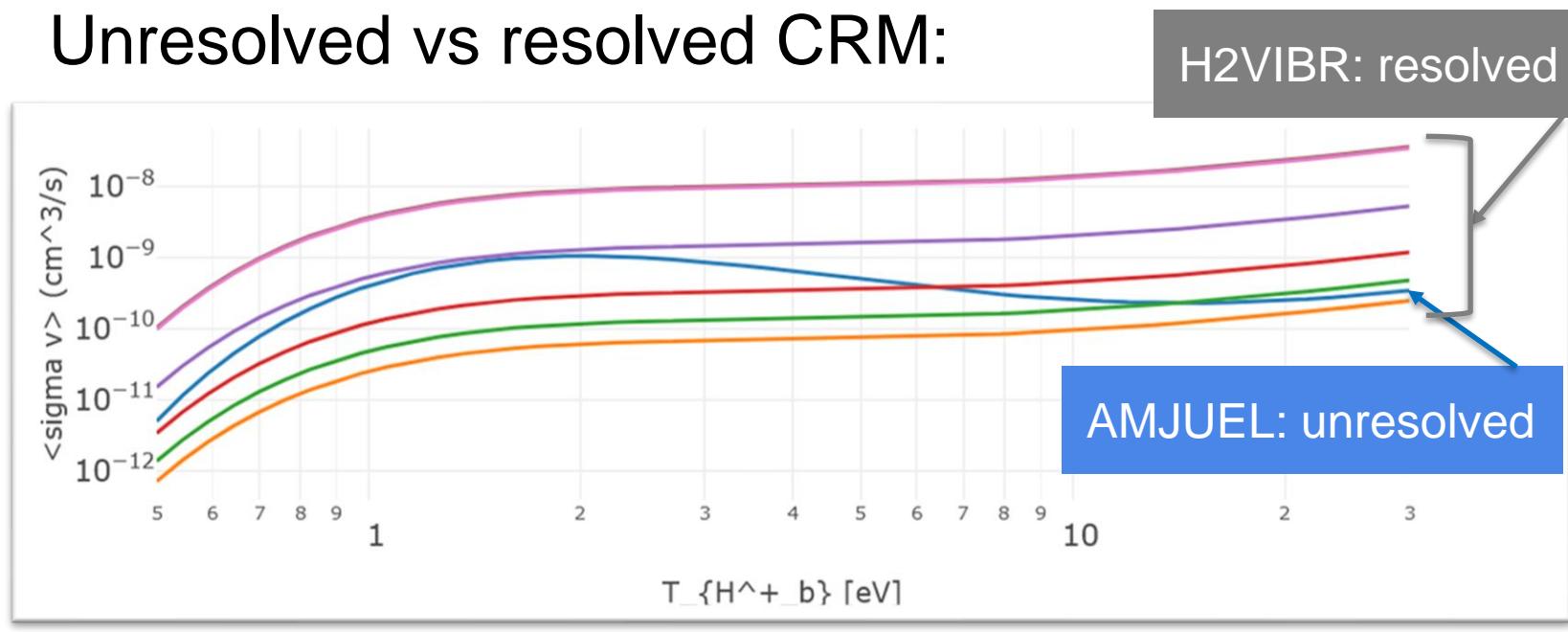
Unresolved vs resolved EIRENE simulations:



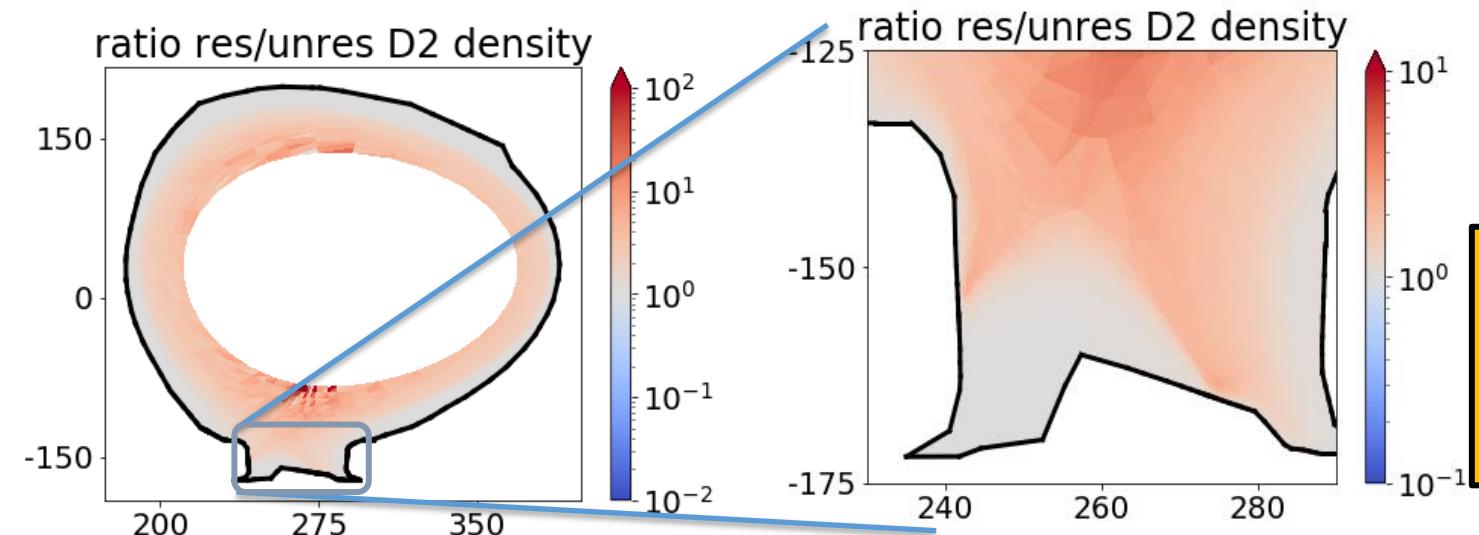
**Larger molecular density  
(factor  $\approx 2-3$ ) along the  
separatrix with vibr. resolution.**

# Impact of vibrational resolution

Unresolved vs resolved CRM:



Unresolved vs resolved EIRENE simulations:

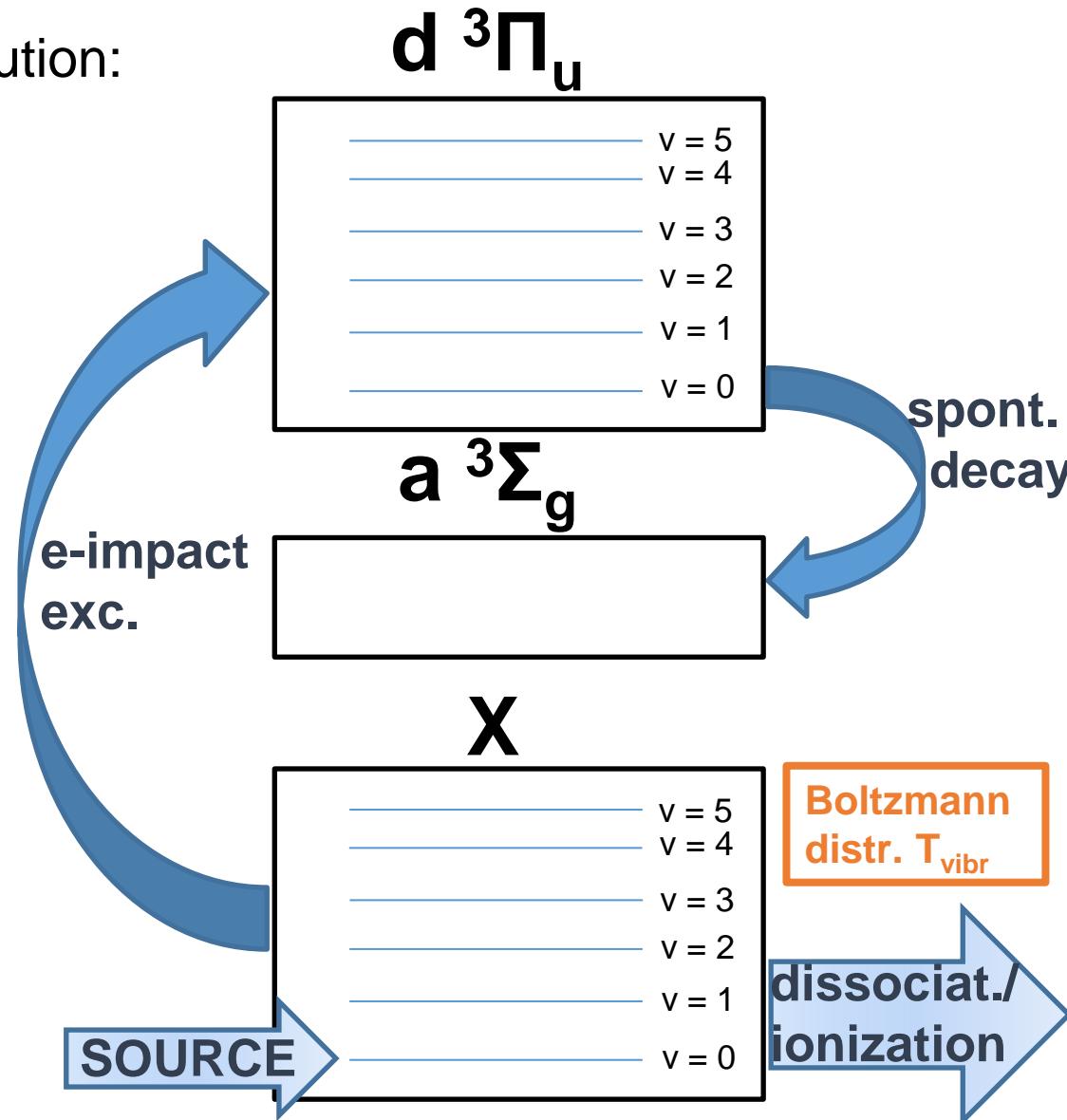


**reduced dissociation rate  
from vibr. resolved data.**

**Larger molecular density  
(factor  $\approx 2-3$ ) along the  
separatrix with vibr. resolution.**

# EXCITED MOLECULAR STATES:

Model for excited states on top of ground state distribution:

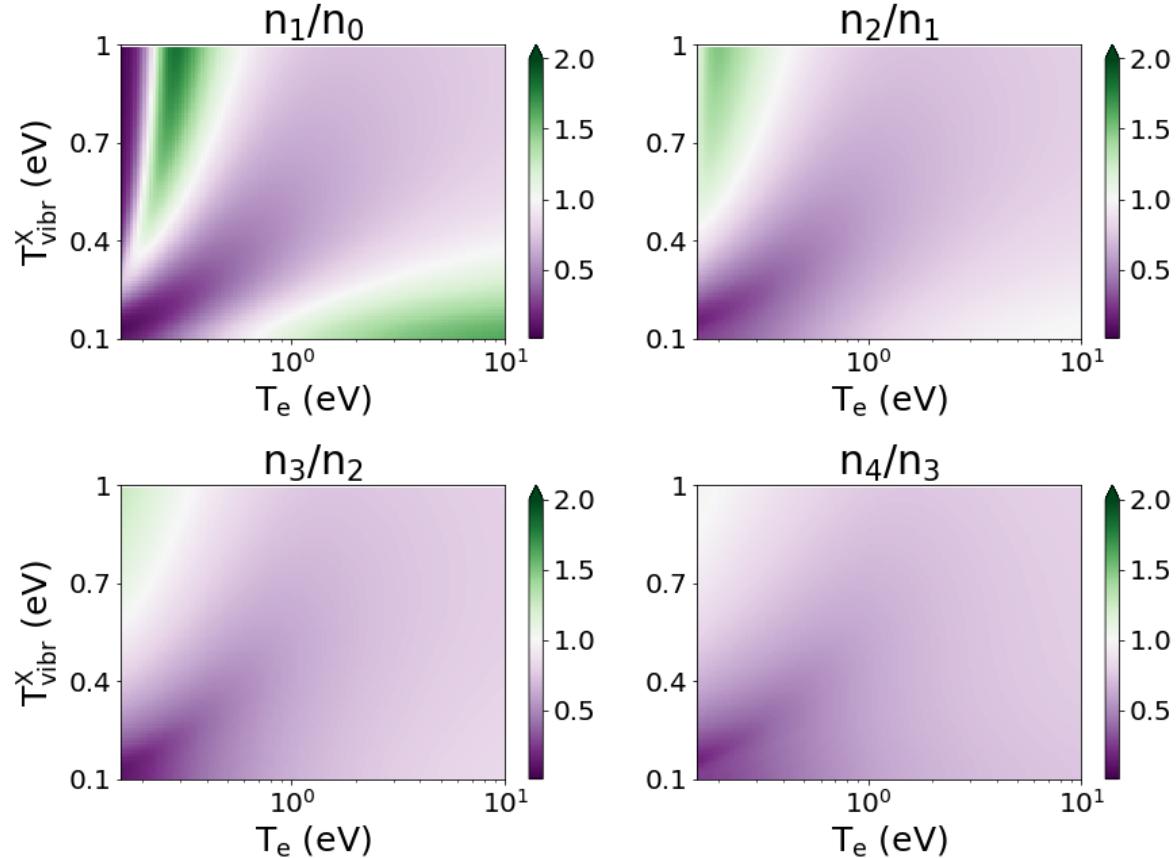


# EXCITED MOLECULAR STATES:

MCCC database

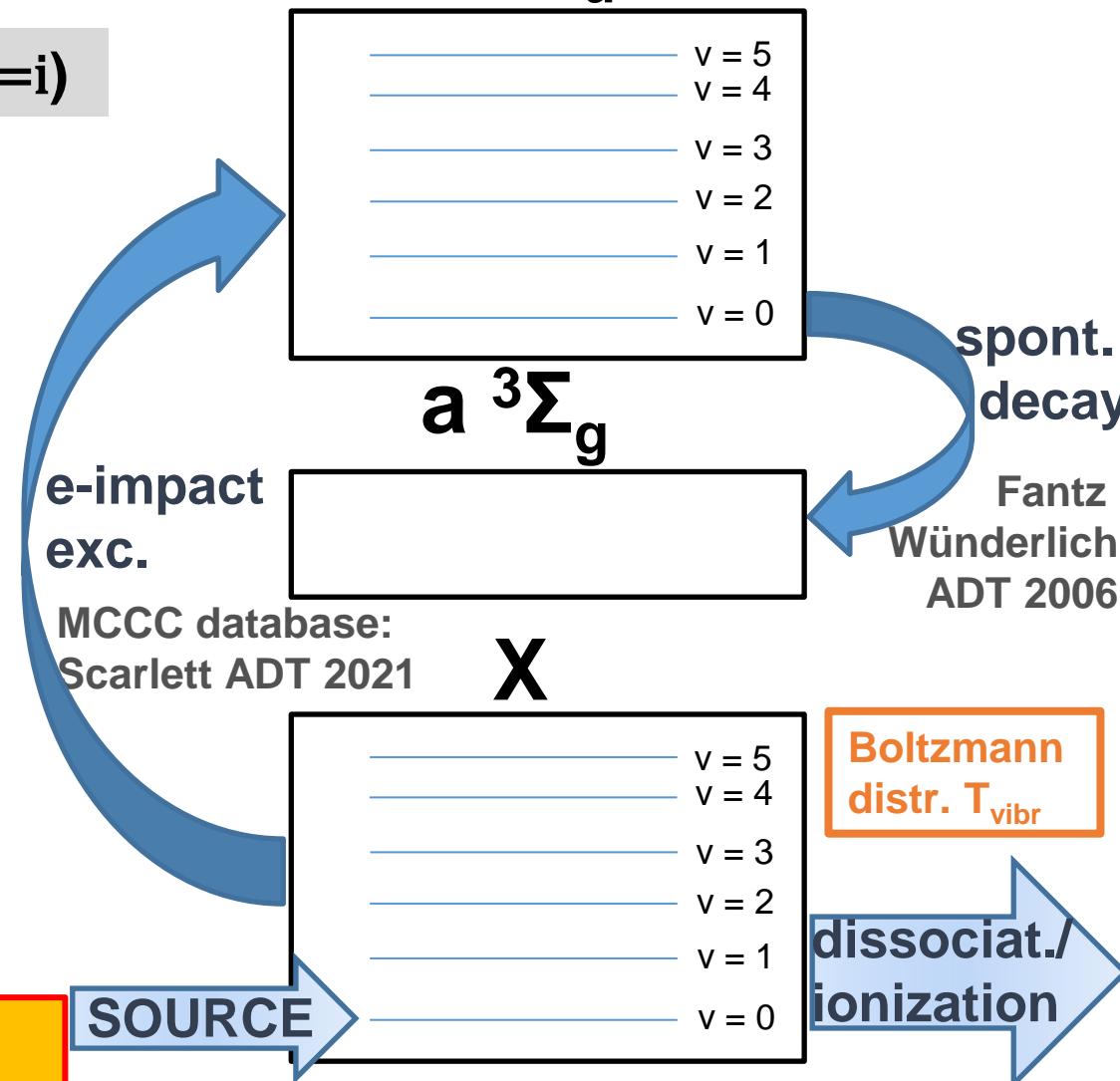
Model for excited states on top of ground state distribution:

- population coeff. excited state  $d\ 3\Pi_u$ :  $n_i = n(d, v=i)$



ratio of Fulcher band intensities as a proxy for ground state vibr. distribution

$d\ 3\Pi_u$



Boltzmann distr.  $T_{vibr}$

dissociat./ionization

## Actions:

- **ModCR:** CRM solver (CVODE) for integration with EIRENE
  - Ongoing
- **MCCC database:** inclusion into EIRENE
  - planned, not started yet



### ***MCCC DATABASE***

MCCC database - Molecular Convergent Close-Coupling  
e-molecule scattering cross sections  
<https://www.mccc-db.org/>

# Turbulence simulations near the X-point

**Motivation:** understand the physics of turbulence generation near the X-point through fast 3D simulations

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ASSUMPTIONS: constant temperature, ignoring curvature-parallel velocity-drifts-magnetic potential-diamagnetic terms, low frequency....

--> **SIMPLEST SYSTEM FOR 3D TURBULENCE SIMULATIONS**

$$\left\{ \begin{array}{l} \partial_t \Delta_{\perp} \phi + \{\phi, \Delta_{\perp} \phi\} = B \nabla_{\parallel}^2 (\tilde{p} - \phi) + \mu \Delta_{\perp}^2 \phi \\ \partial_t \tilde{p} + \{\phi, \tilde{p}\} = -\hat{k} \partial_y \phi + D \nabla_{\parallel}^2 (\tilde{p} - \phi) + \chi_{\perp} \Delta_{\perp} \tilde{p} \end{array} \right.$$

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Viscosity

Drift coupling

Diffusion

Advection:

$$\{f, g\} = \partial_x f \nabla_{\hat{y}} g - \partial_x g \nabla_{\hat{y}} f$$

Free energy source:

background pressure gradients

$$p = p_0 + \tilde{p}$$

$$-\partial_x p_0 = p_0/l_0 \propto \hat{k}$$

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**Advection:** nonlinear inverse energy cascade to large parallel scales (n=0)

**Drift coupling:** near the X-point it is toroidal, thus proportional to n.



Turbulence suppression? HW effective model still valid? Is shear able to sustain drift response? What's the role of other parameters?

## Geometry:

$$x = \frac{r - r_0}{L_{\perp}}, y = \frac{a\theta}{L_{\perp}}, z = \frac{R\varphi}{L_{\parallel}}$$

full toroidal  
local in poloidal plane

$$k_{\parallel} \gg k_{\perp}$$

□ **SLAB:** close X point

$$\nabla_{\parallel} = \partial_z \quad \Delta_{\perp} = \partial_x^2 + \partial_y^2$$

□ **TOKAMAK:**

$$\nabla_{\parallel} = \hat{b} \cdot \vec{\nabla} \quad \Delta_{\perp} = (\nabla - \nabla_{\parallel})^2$$

$$\hat{b} = \frac{\vec{B}}{B}$$

Laplacian and parallel derivative are non-linear in FS if B's are not constant (shear).

## Metodology: 4th order Runge-Kutta in Fourier space

- at each iteration pressure and vorticity are advanced
- electrostatic potential is computed inverting the Laplacian
  - **w/o shear:** inverse Laplacian is analytical in FS and can be easily inverted.
  - **with shear:** Laplacian and inverse are numerically pre-calculated before starting time iteration and stored as csr matrix.
- all non-linear terms are computed in real space with Fast Fourier Transforms (**FFT**) and pulled back in FS with inverse FFT (**pseudo-spectral**)
  - **w/o shear:** advection
  - **with shear:** advection + all terms with B's (drift,..)

## Parameters:

Braginsky parameters for typical DTT case:

$$a=68\text{cm} \quad R=2.14\text{m} \quad B=3 \times 10^4 \text{Gs} \quad n=5 \times 10^{13} \text{ cm}^{-3}$$

Background pressure gradients:  $l_0=40\text{cm}$

Two cases for diffusivity: Classical( $1.2\text{m}^2/\text{s}$ ) vs Neoclassical (x10)

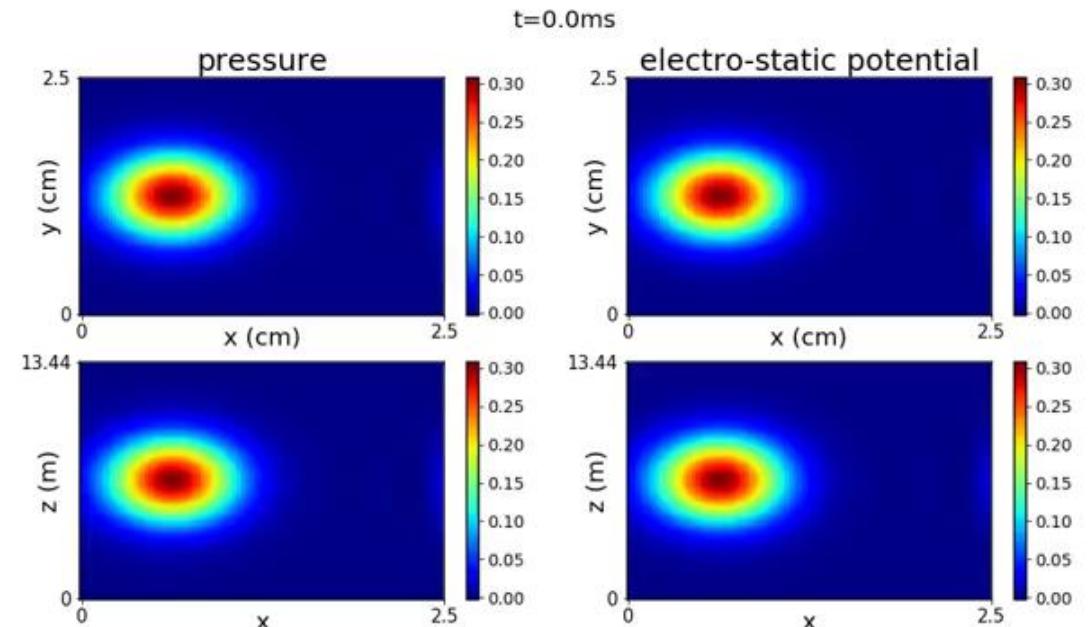
## Initial values:

3D wave packet centered inside the simulation domain.

## Boundary Conditions:

periodic boundary conditions

NEW: vanishing Dirichlet "radial" BC

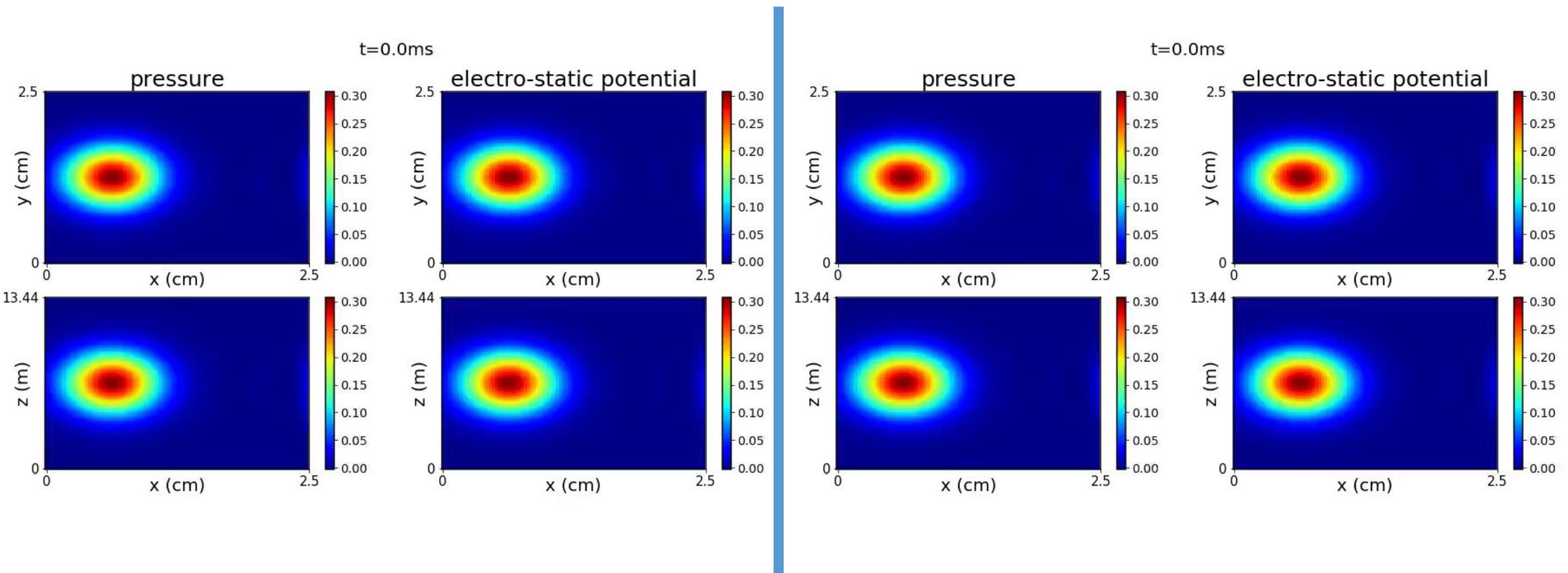


## Simulations: SLAB

classical diffusivity

vs

neoclassical diffusivity

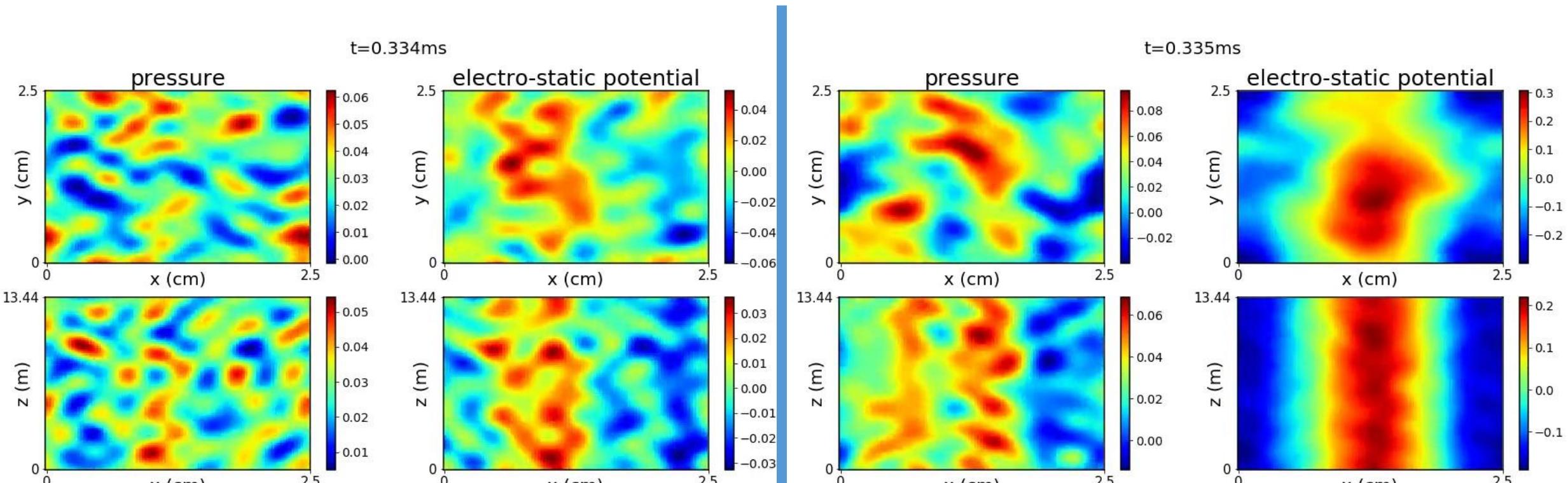


## Simulations: SLAB

classical diffusivity

vs

neoclassical diffusivity



Stronger turbulence  
large drift response  
wider toroidal spectrum

Turbulence reduction  
large diffusion  
inverse cascade

## Outlook:

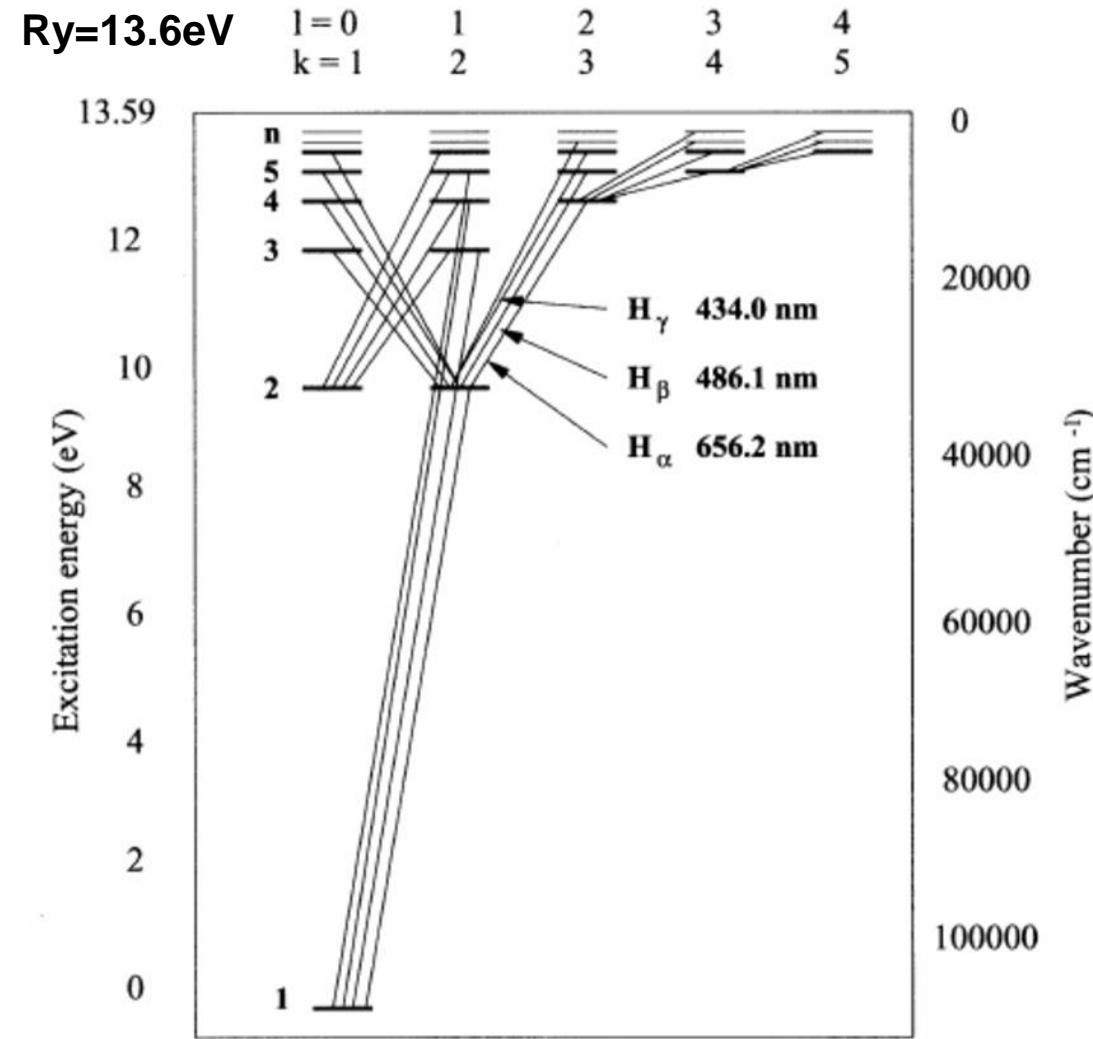
- comparison between different magnetic geometries → X point
- inclusion of more effects (curvature, electro-magnetic, gyrofluid corrections,..)
- consistent diffusivity: test particle motion on turbulent scenario
- global (poloidal) simulations
- predictions for DTT

**Fabio Moretti's talk..**

## **Additional material**

# Collisional-Radiative Models (CRM)

Grotian diagram for atomic H (D, T)



**Balance equations:**

$$\frac{dN_i}{dt} = \sum_{j \neq i} A_{ji} N_j + n_e \cdot (EXCIT + IZ + CX + REC)$$

$$EXCIT = \sum_{j \neq i} \langle v \sigma_{ji} \rangle N_j$$

$$IZ = \sum_m \langle v \sigma_{mi} \rangle N_m^- + \sum_z \langle v \sigma_{zi} \rangle N_z^{2-} + \dots$$

$$REC = \sum_k \langle v \sigma_{ki} \rangle N_k^+ + \sum_l \langle v \sigma_{li} \rangle N_l^{2+} + \dots$$

...

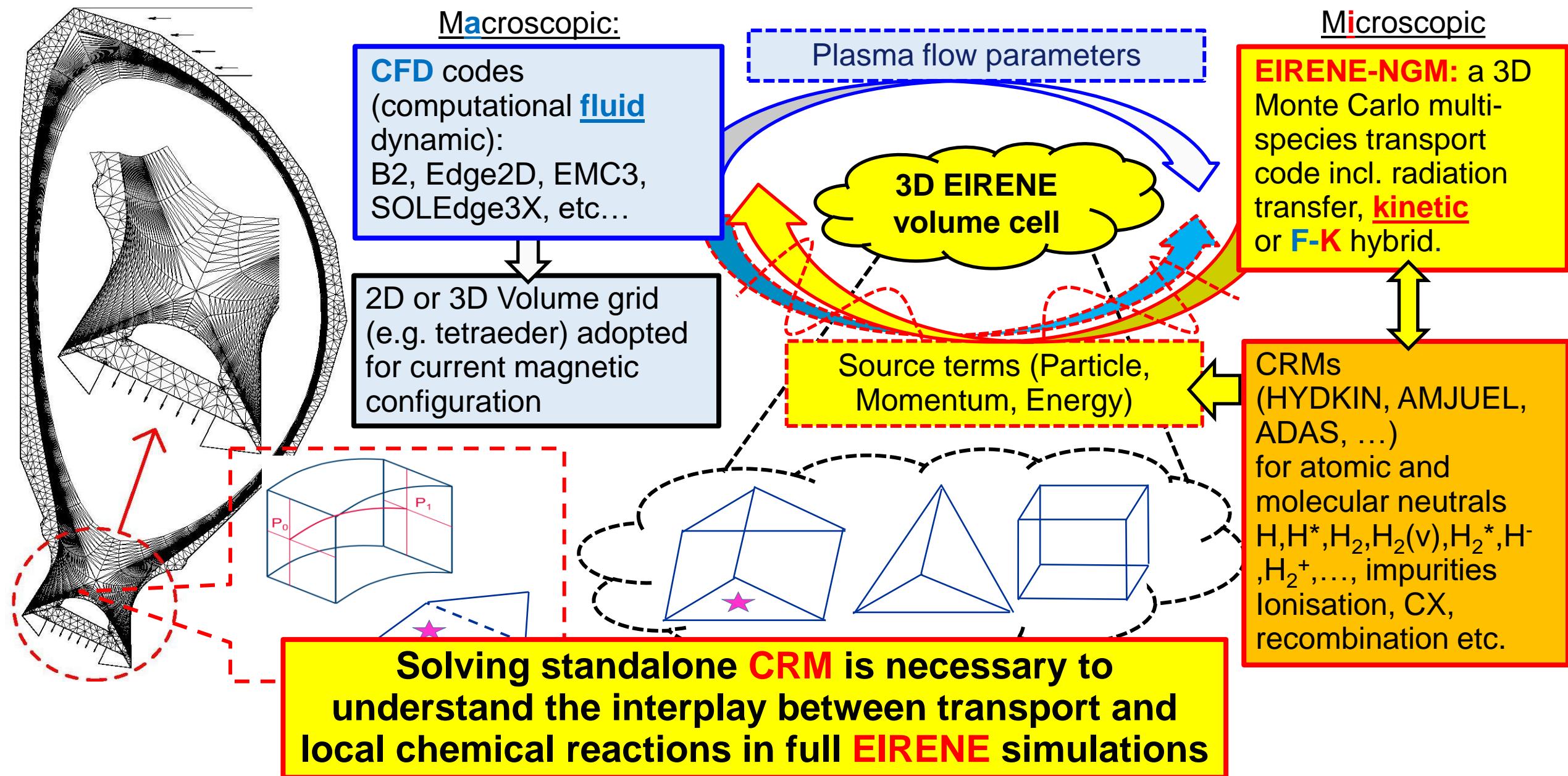
*j, k, l, m, z, ... states can be fine-superfine resolved or, opposite, bundled into few quasi-metastables (MS)*

CRM = list of states + transition data

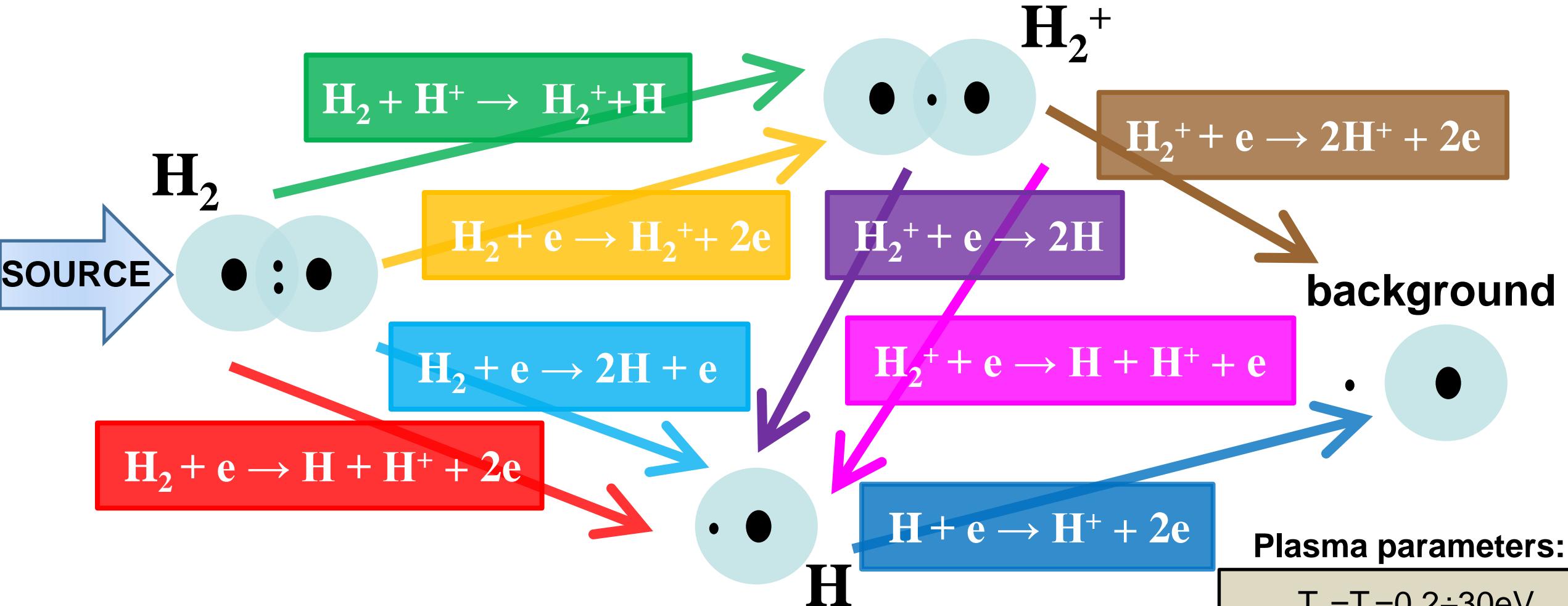
Often used:

$\langle v \sigma_{ji} \rangle (T_e, n_e)$  - effective Maxwellian averaged rates

# What is EIRENE in a nutshell (e.g. in SOLPS)?..



# CRM FOR MOLECULES



Databases:

AMJUEL, HYDHEL

$T_e = T_i = 0.2 \div 30 \text{ eV.}$

$n_e = n_i = 10^{19} \text{ m}^{-3}$

$n_e = n_i = 10^{21} \text{ m}^{-3}$

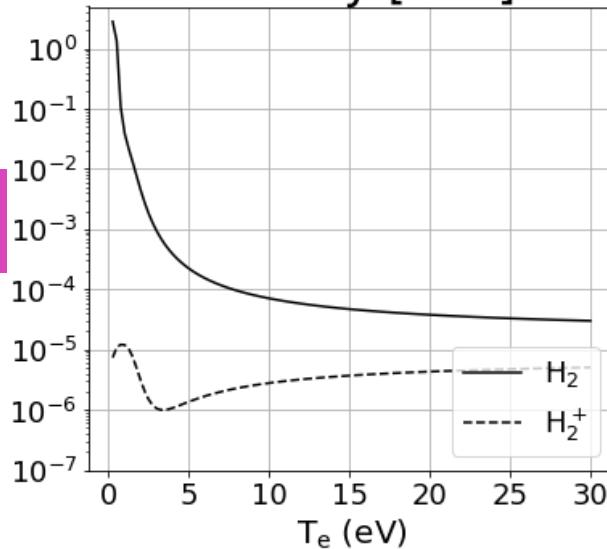
JET

EU-DEMO

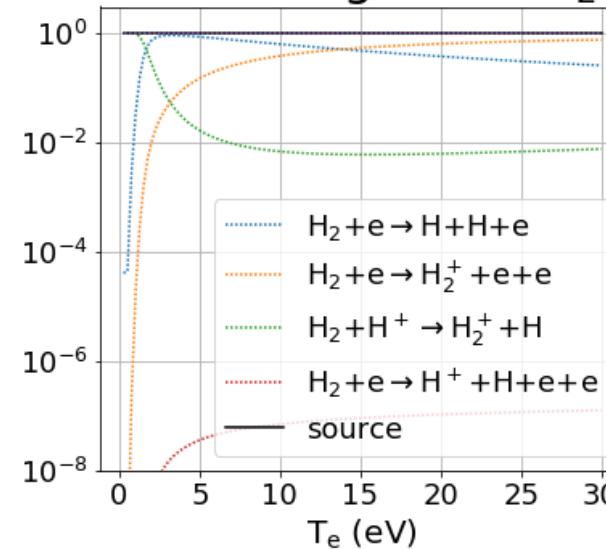
# CRM FOR MOLECULES: leading reactions with temperature

EU-DEMO

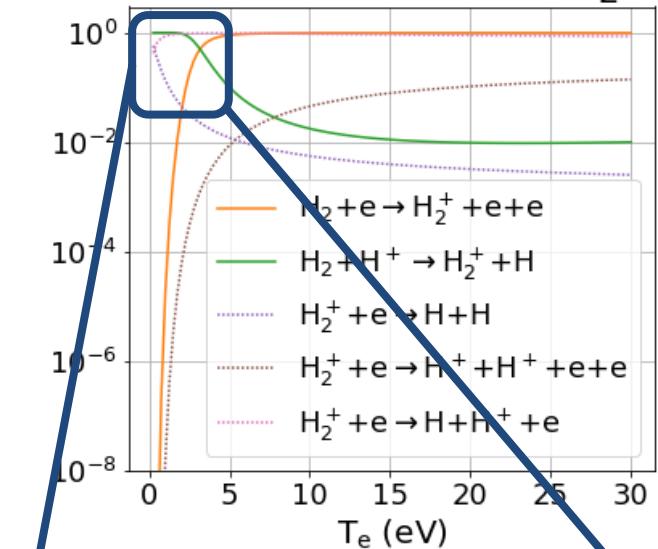
Density [a.u.]



Branching ratios  $\text{H}_2$



Branching ratios  $\text{H}_2^+$



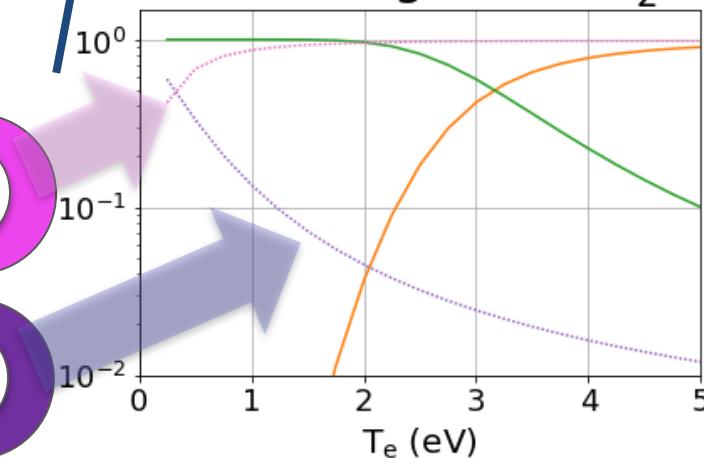
low temperature ( $T < 2$  eV) leading reaction chains:

MAR/MAD

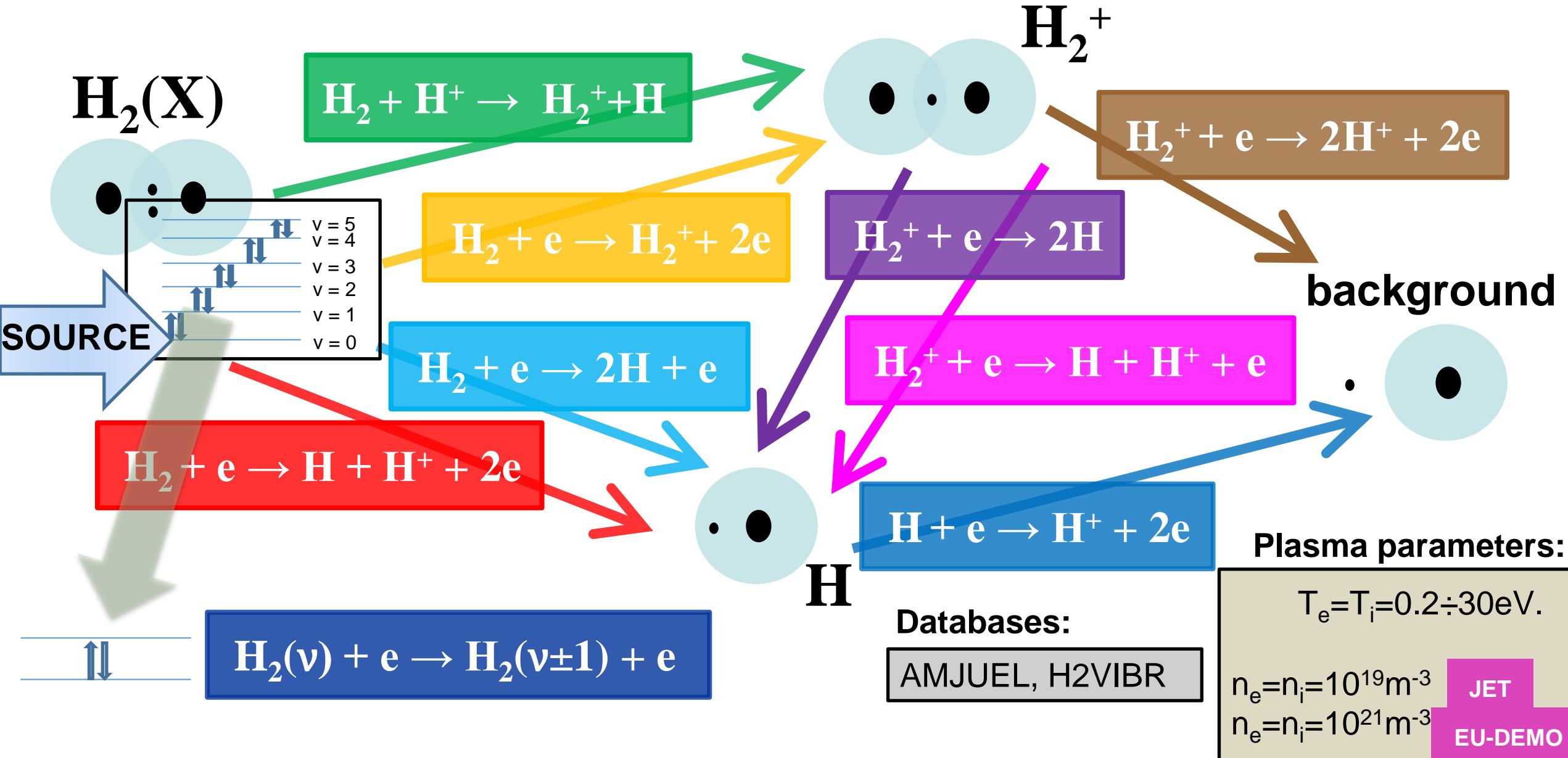
competition at very  
low temperature



Branching ratios  $\text{H}_2^+$

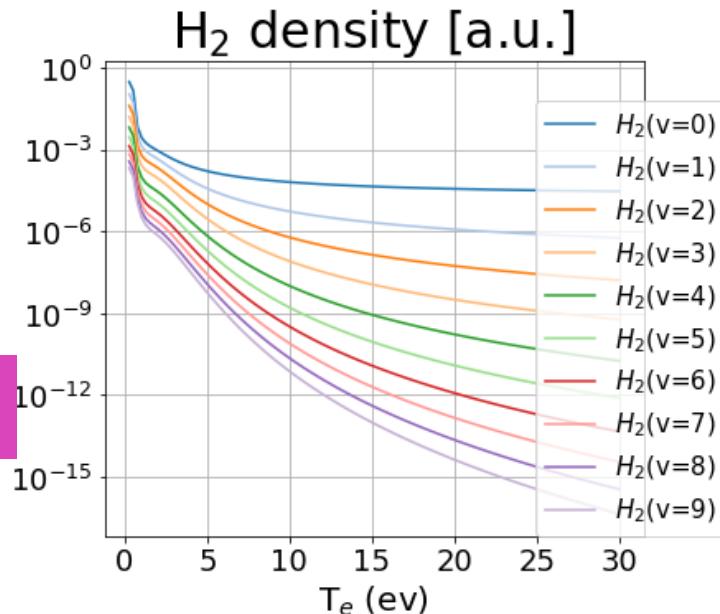


# CRM FOR MOLECULES (vibr. resolved)

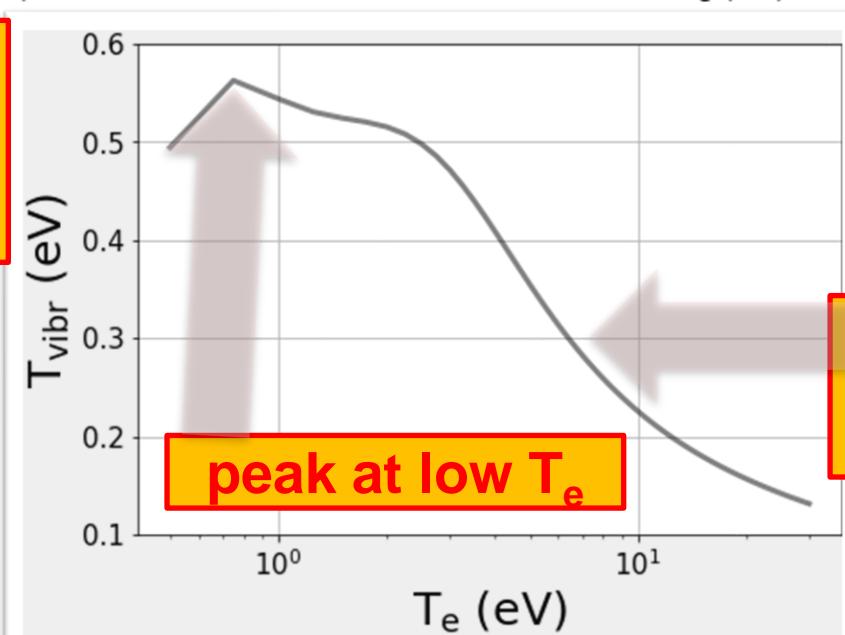
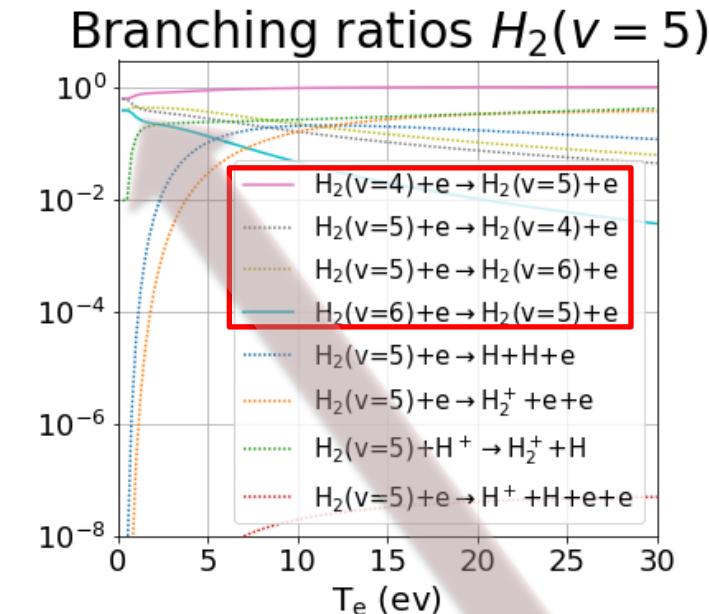
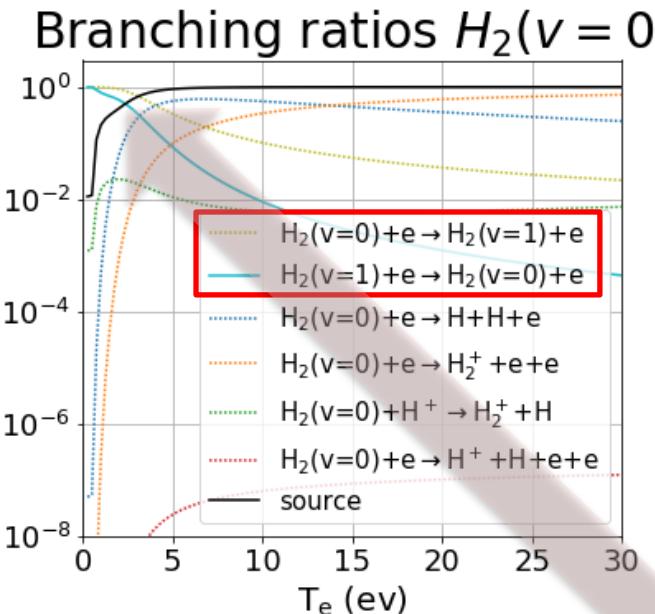


# CRM FOR MOLECULES: leading reactions with temperature

EU-DEMO



Maxwellian distribution (not obvious..)



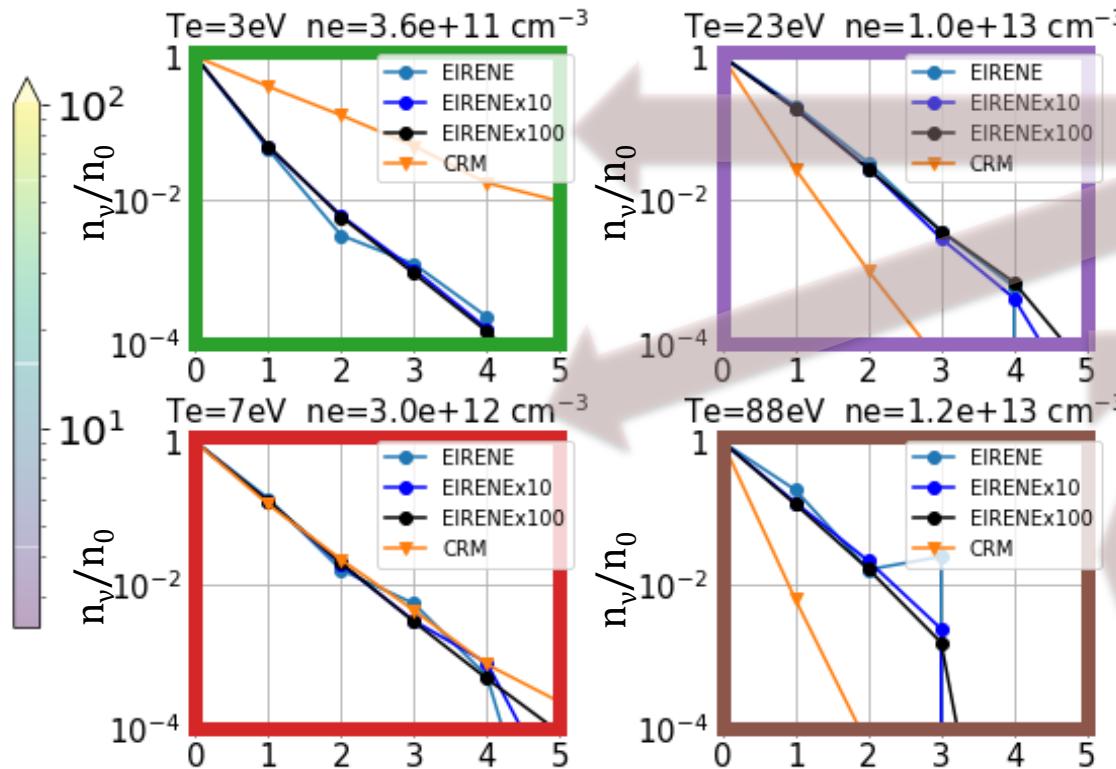
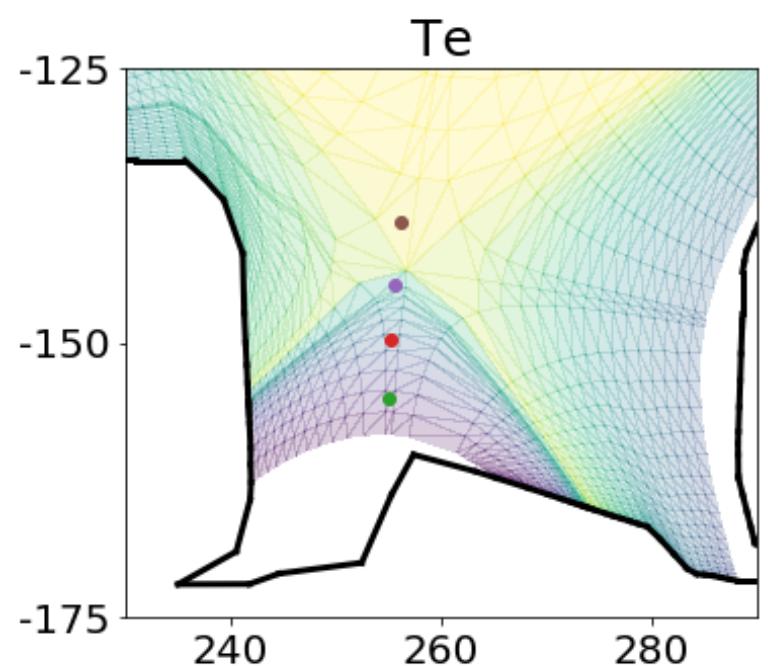
$T_{vibr}$  increases with decreasing  $T_e$

vibrational transitions are the main reaction channels at low temperatures

# CRM vs EIRENE (vibr. resolved):

Simulations: SOURCE  $v=0$

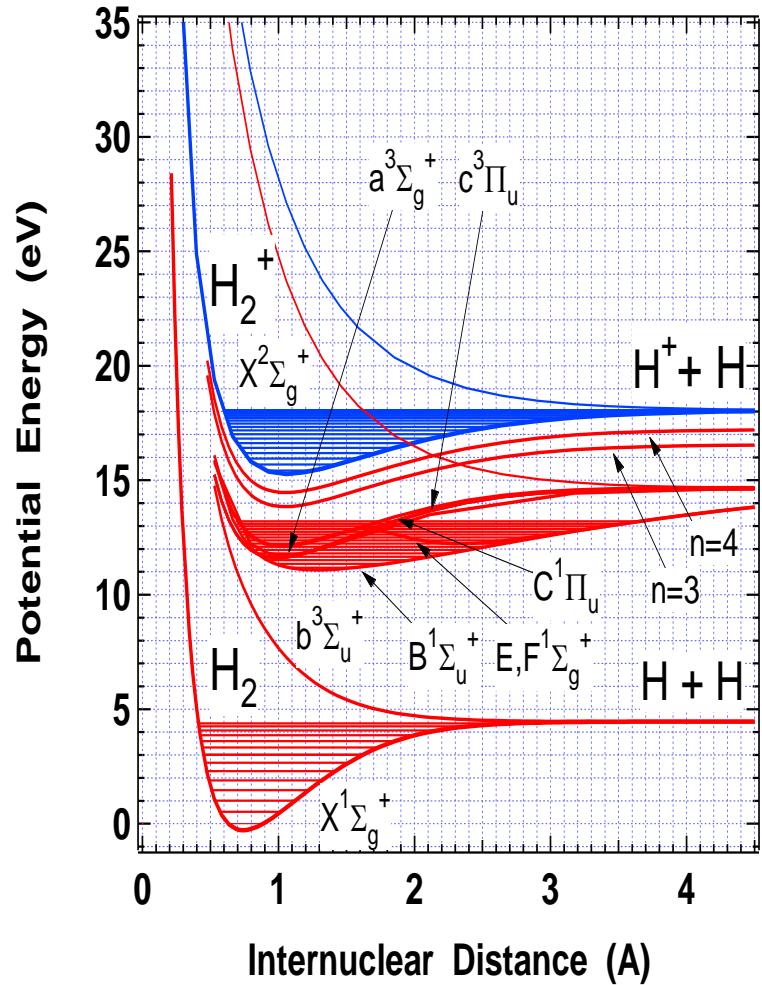
EIRENE: 793'300 histories  
EIRENEx10: 7'933'000 histories  
EIRENEx100: 79'330'000 histories



transport is important for vibrational distribution-  
full particle simulations (long and heavy)...

# EXCITED MOLECULAR STATES: MCCC DB

□ Preliminarily:

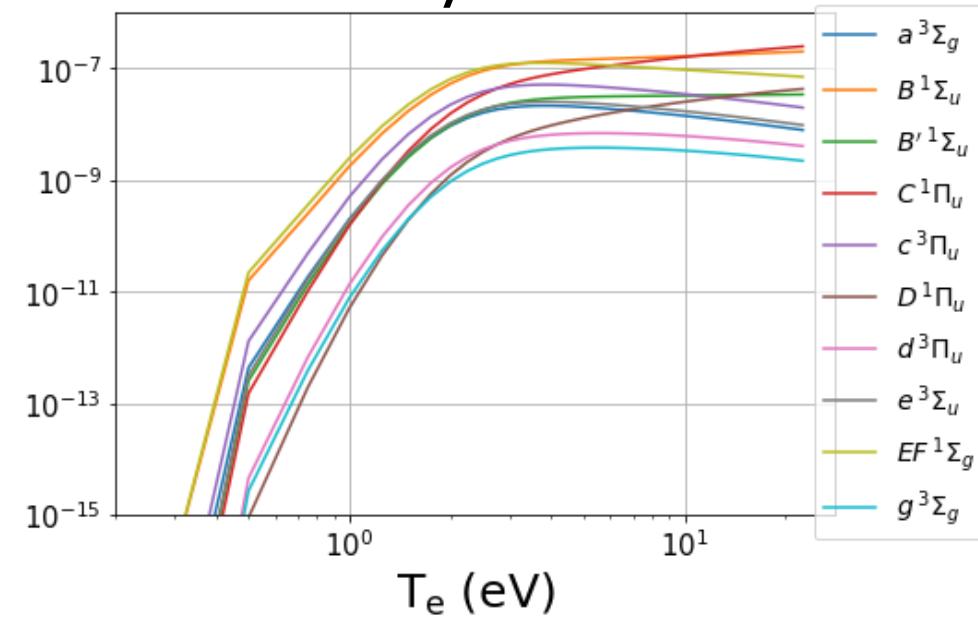
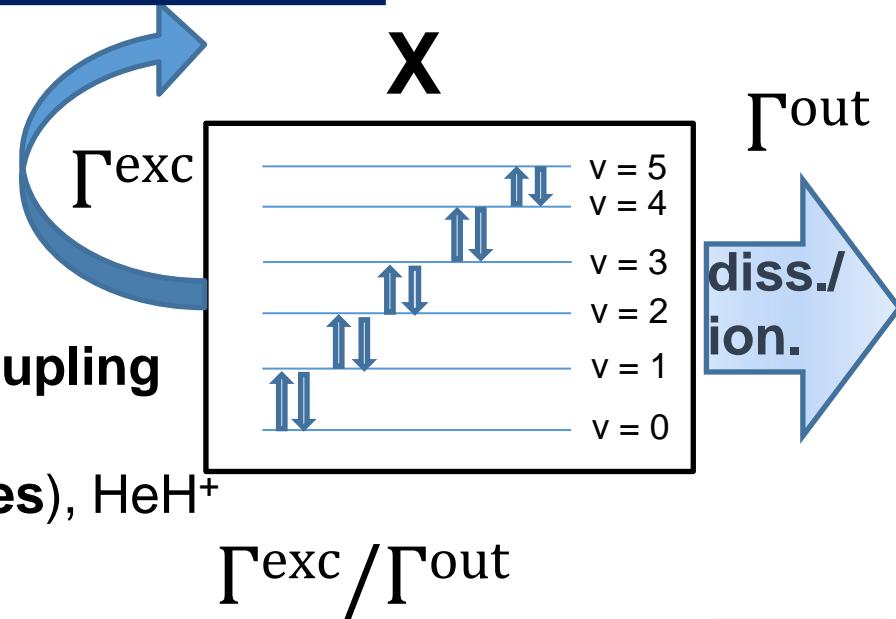


MCCC database:  
Scarlett ADT 2021

- Molecular Convergent Close-Coupling
- electron and positron collisions
- target:  $\text{H}_2$  (isotopes),  $\text{H}_2^+$  (isotopes),  $\text{HeH}^+$
- dissociation, ionizat., excitation + **vibrational resolution**
- output: tabular

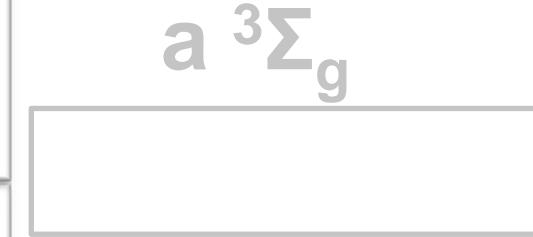
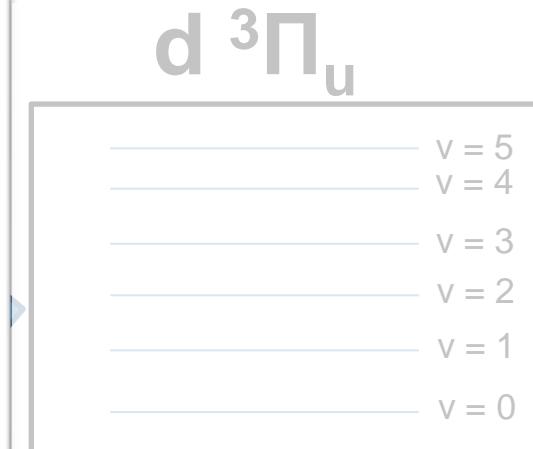
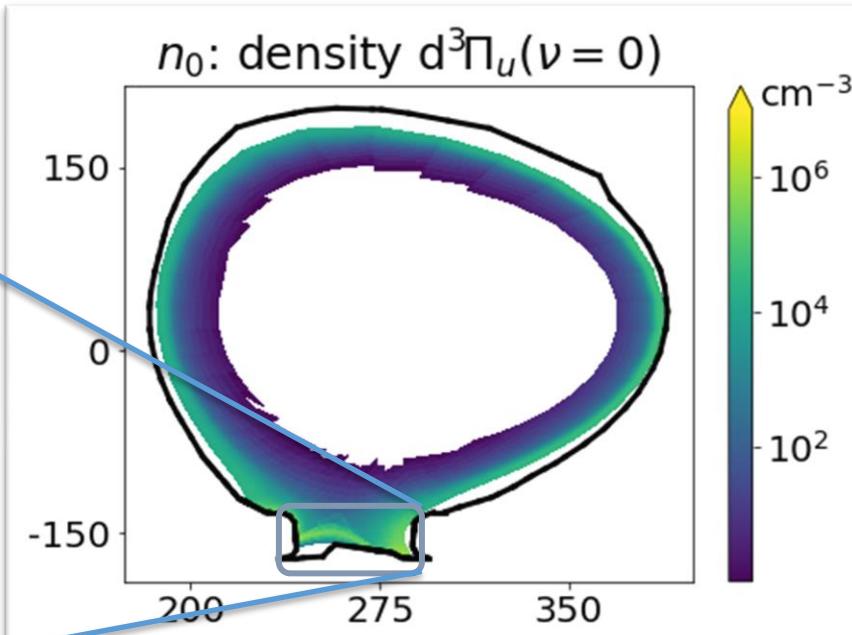
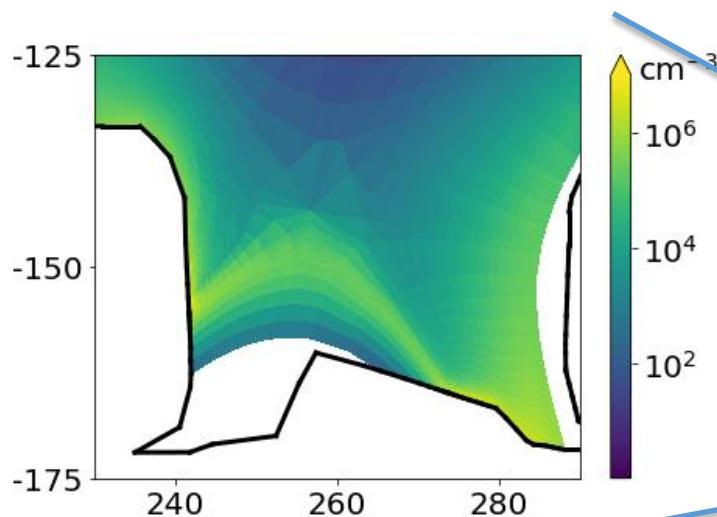
$\Gamma^{\text{exc}} \ll \Gamma^{\text{out}}$

Model for excited  
states on top of ground  
state distribution

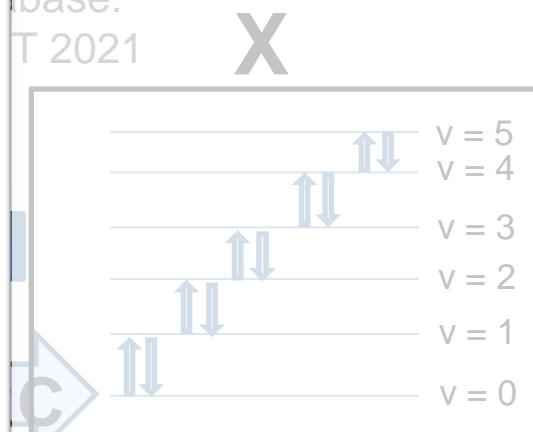
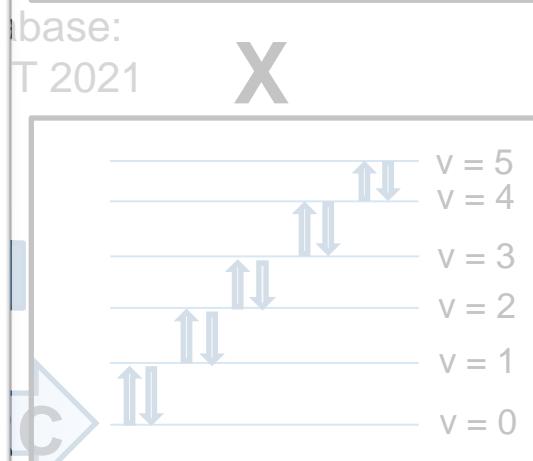
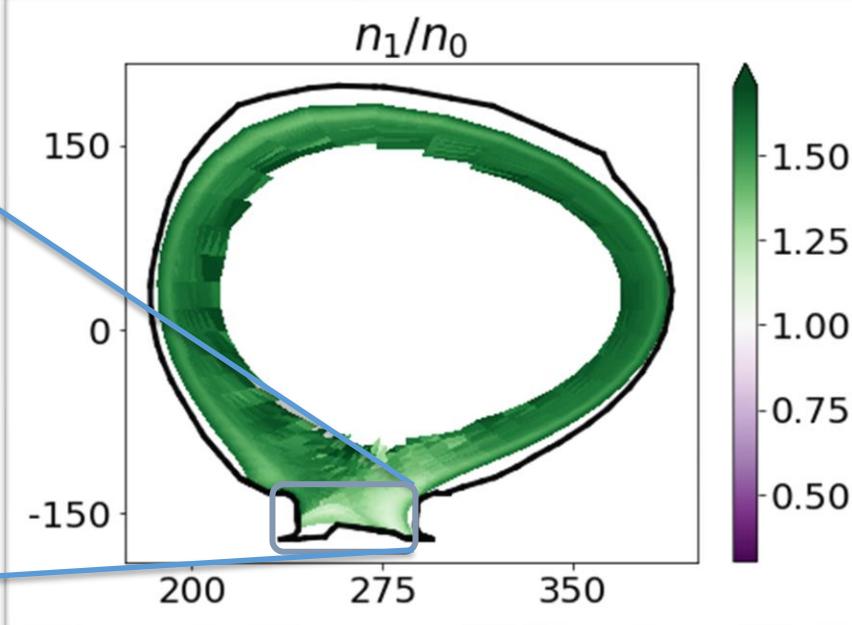
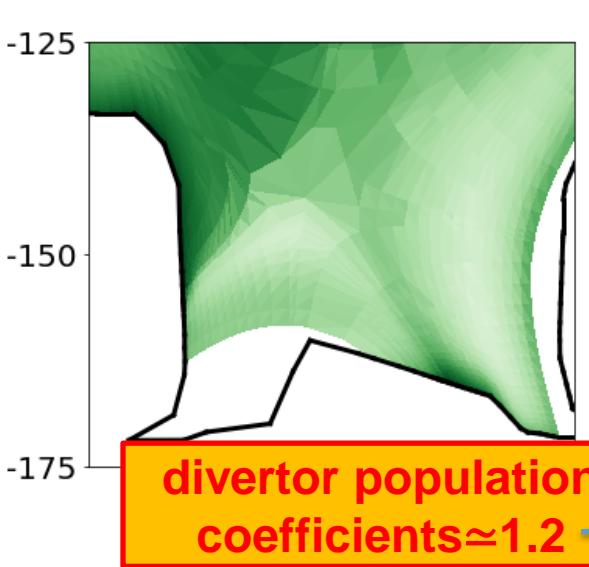


# EXCITED MOLECULAR STATES: EIRENE

excited state  $d\ 3\Pi_u$ :



spont. decay  
Fantz  
Wunderlich  
ADT 2006

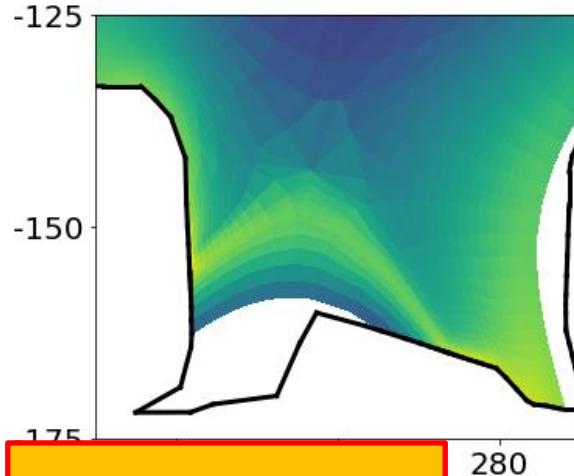


base:  
T 2021  
X  
C  
dissociat./  
ionization

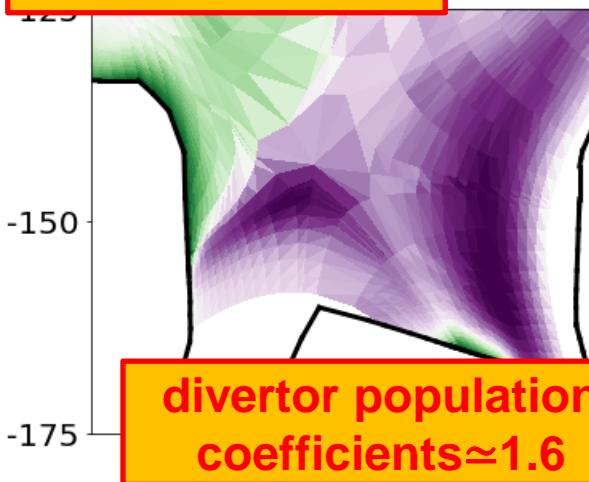
# EXCITED MOLECULAR STATES: EIRENE

□

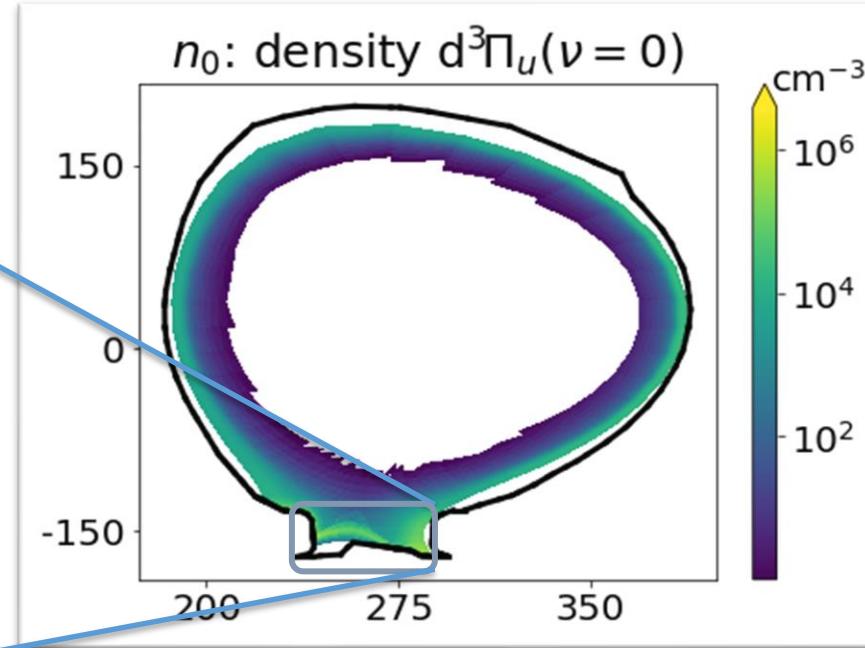
excited state  $d\ 3\Pi_u$ :



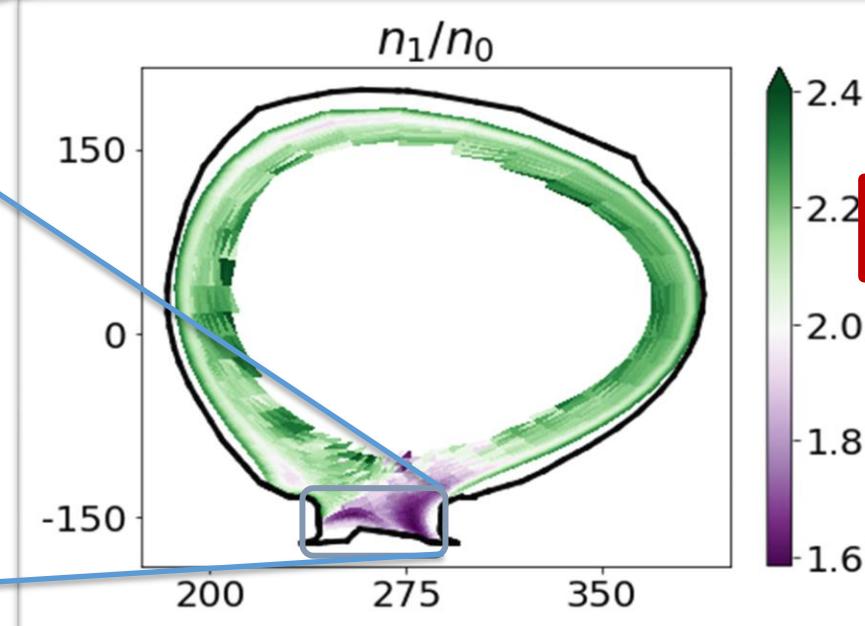
**DEUTERIUM**



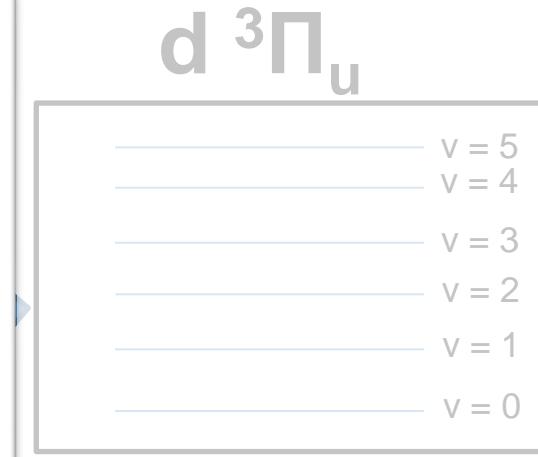
divertor population  
coefficients  $\approx 1.6$



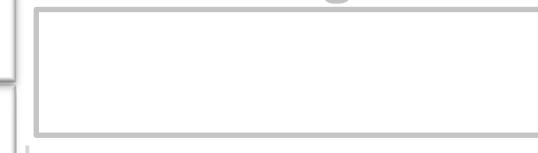
$n_0$



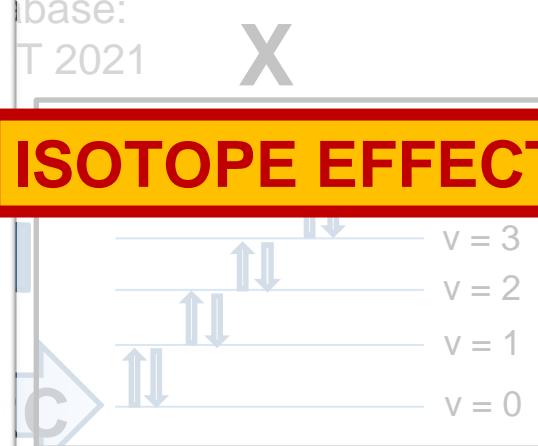
$n_1/n_0$



$d\ 3\Pi_u$



$a\ 3\Sigma_g$



**ISOTOPE EFFECT**

spont.  
decay  
Fantz  
Wunderlich  
ADT 2006

dissociat./  
ionization