

Numerical Electron Degradation in Kilonovae

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Stockholms
universitet



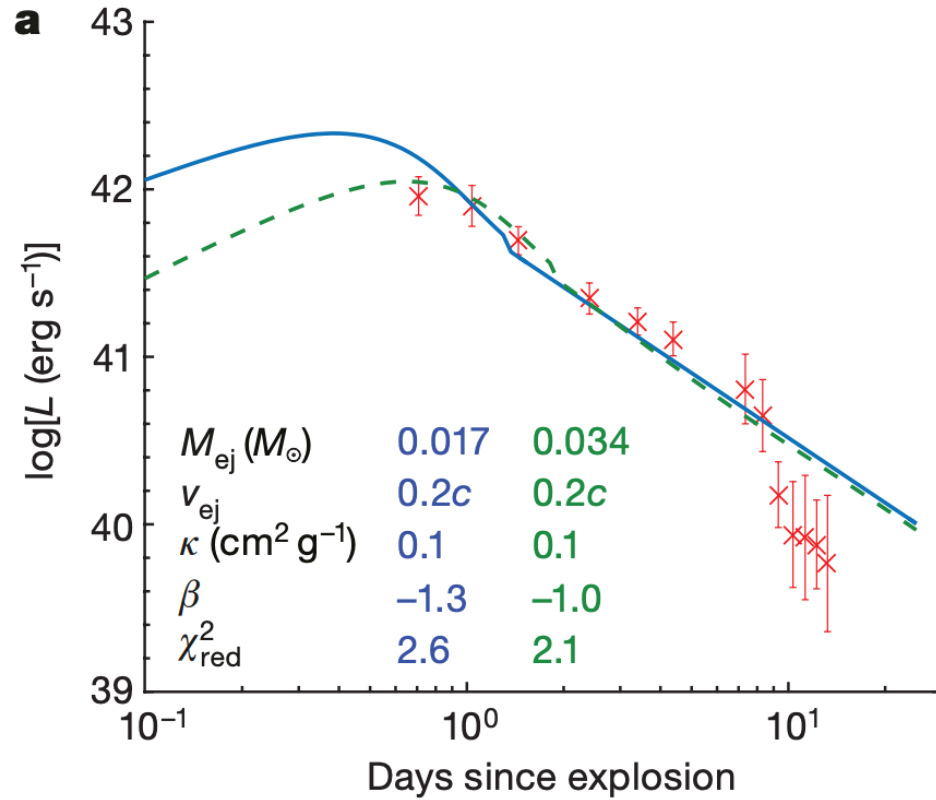
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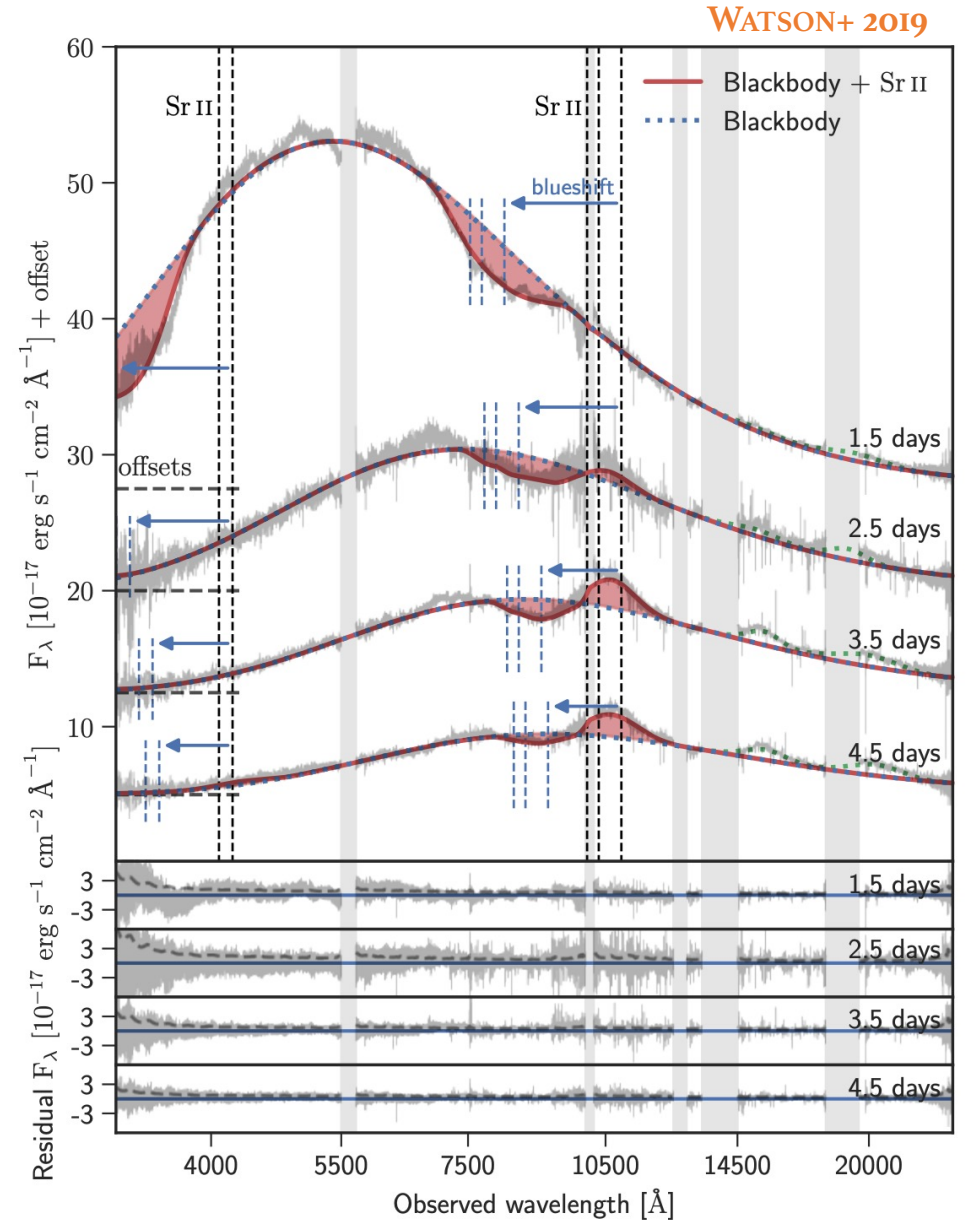
Outline

- Time dependent nebular phase and the non-thermal electrons
- Ionization cross-sections
- Numerical non-thermal electron degradation

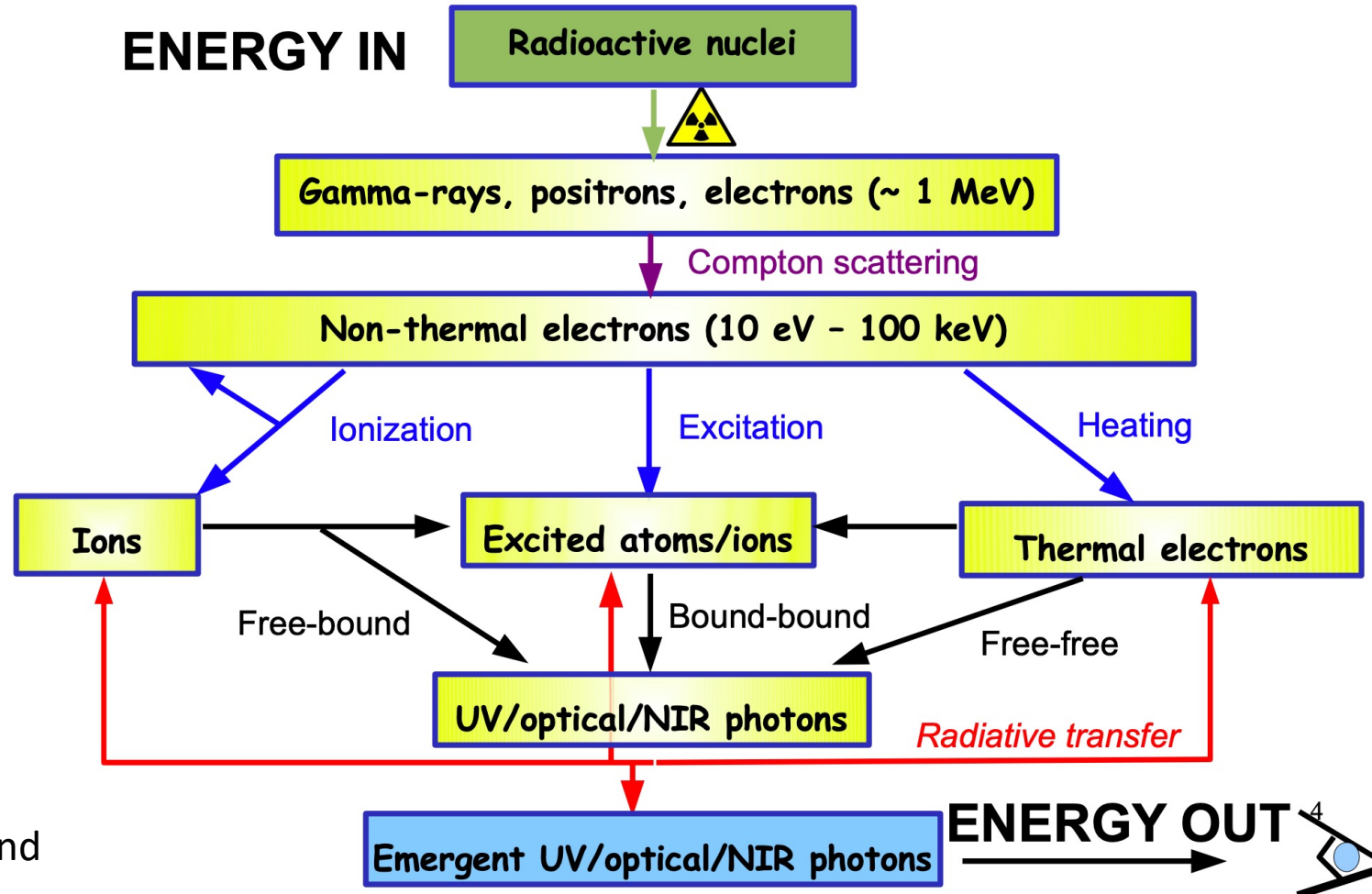
The nebular phase



SMARTT+ 2017



The nebular phase



Credits: A. Jerkstrand

Time-dependence

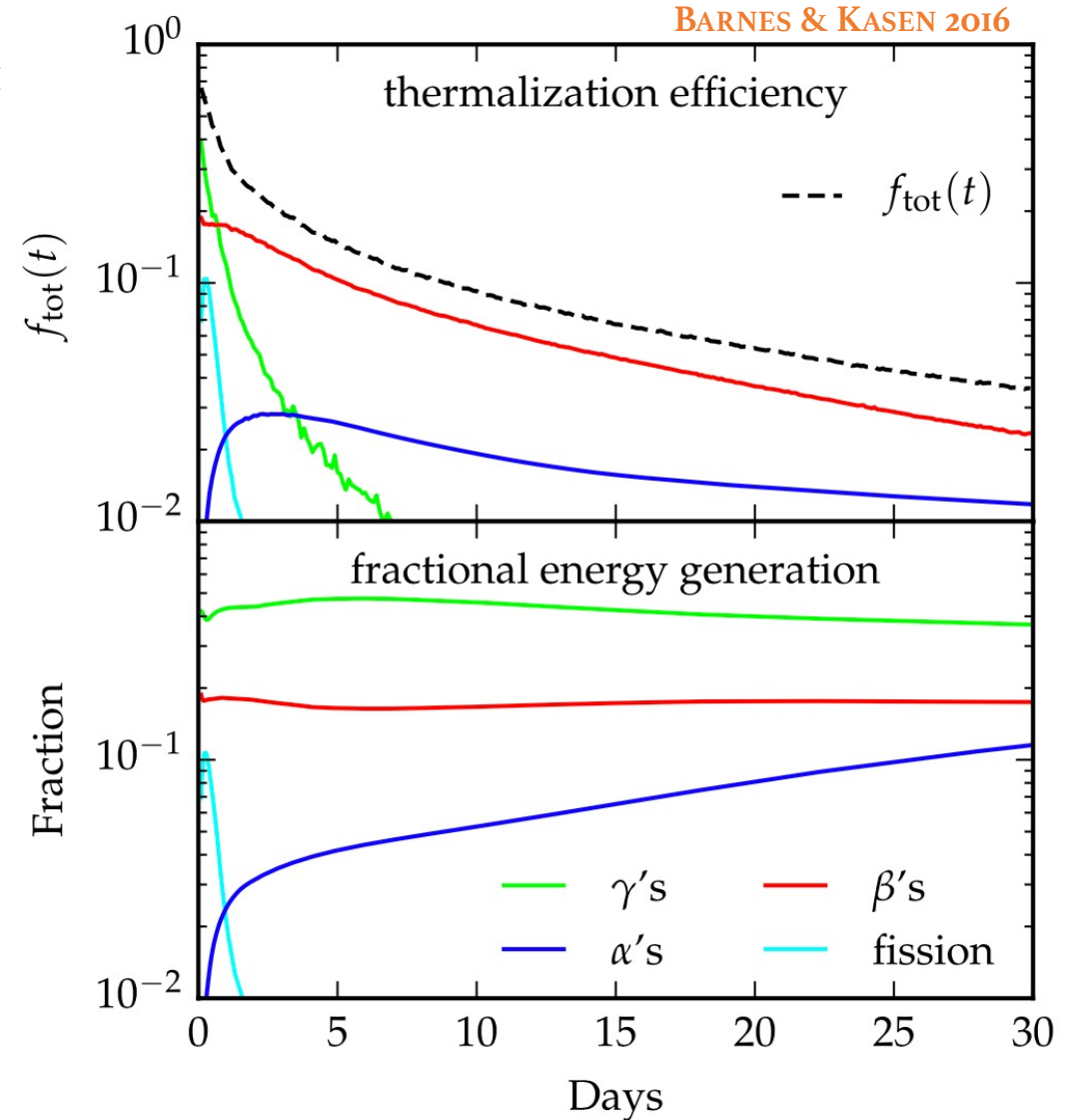
BARNES & KASEN 2016, KASEN & KASEN 2019, BARNES+2021

Time-dependent thermalisation efficiency

$$f_p(t) = \frac{\dot{E}_{\text{th}}}{\dot{E}_{\text{rad}}} = \frac{\ln \left[1 + 2 \left(\frac{t}{t_{\text{ineff,p}}} \right)^2 \right]}{2 \left(\frac{t}{t_{\text{ineff,p}}} \right)^2}.$$

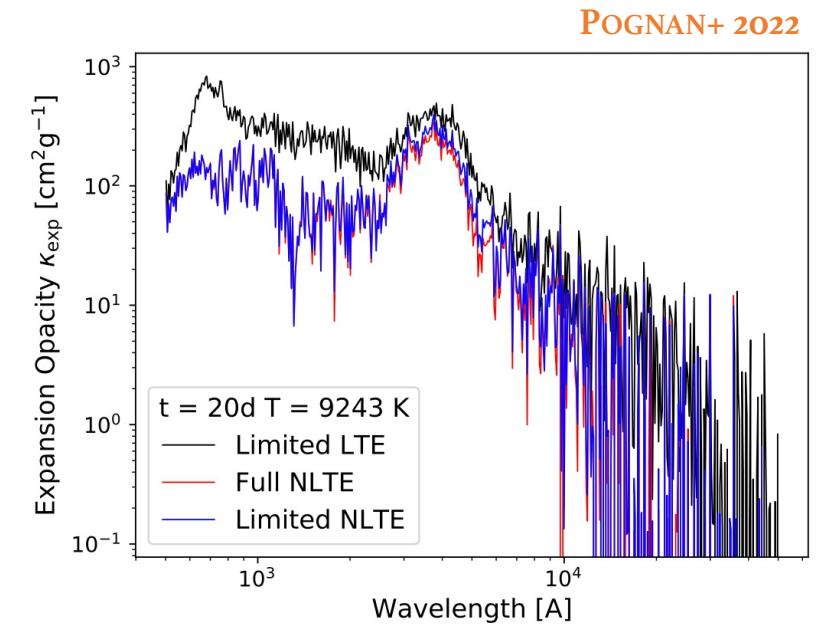
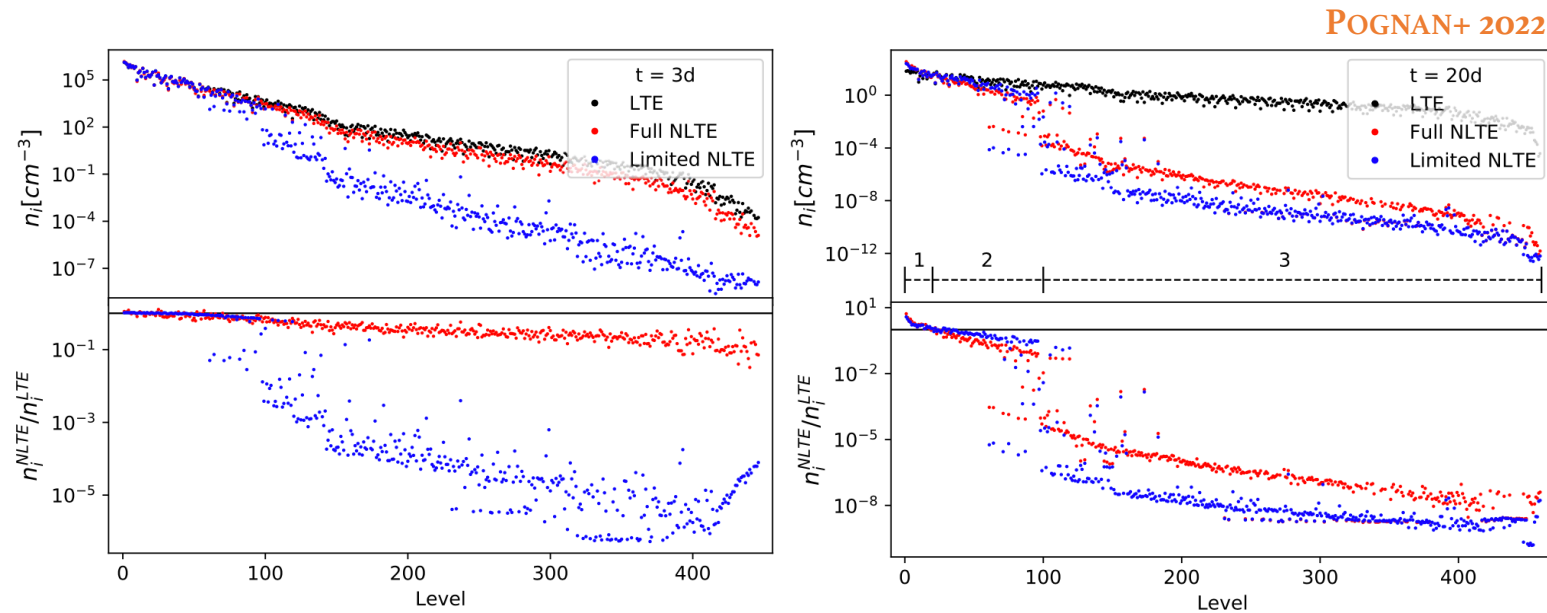
$$f_{\text{tot}}(t) = 0.36 \left[\exp(-at) + \frac{\ln(1 + 2bt^d)}{2bt^d} \right],$$

LTE: total heating fraction + Saha-Boltzmann



Local vs Non-local Thermodynamical Equilibrium

LTE IS INSUFFICIENT AFTER 10 DAYS



Time-dependence

BARNES & KASEN 2016, KASEN & BARNES 2019, BARNES+2021

Time-dependent thermalisation efficiency

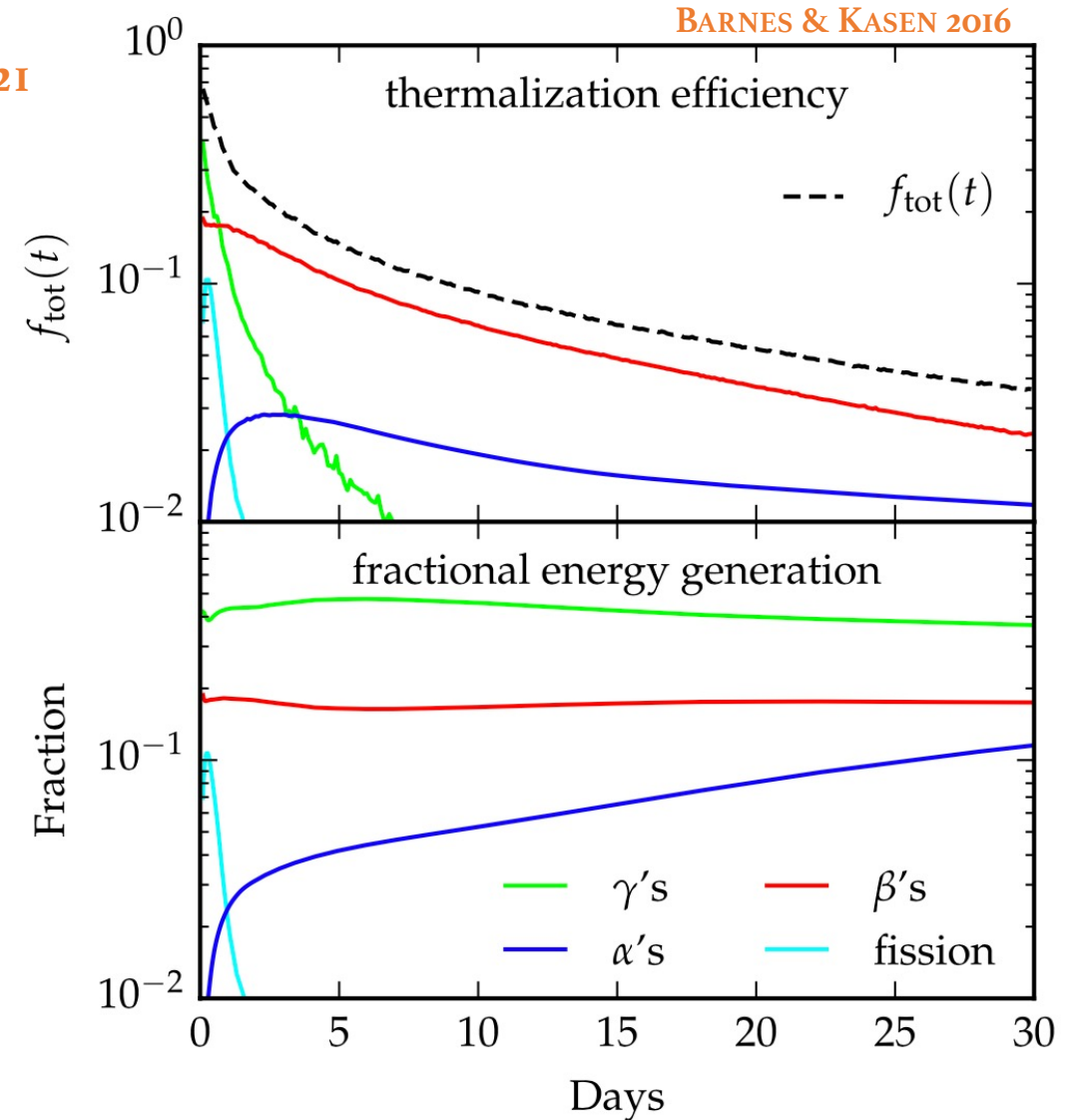
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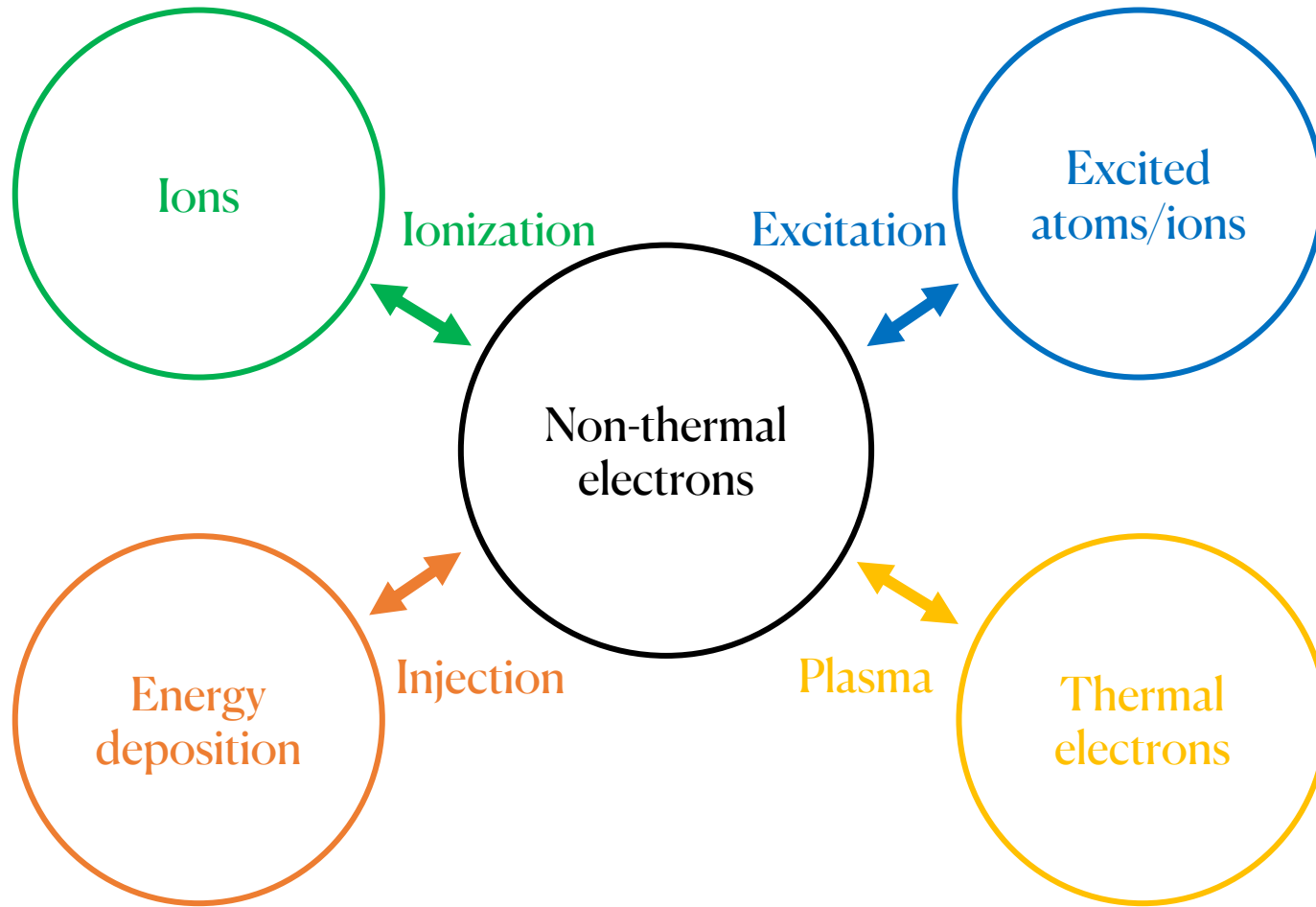
NLTE: Every energy deposition channel fraction:

- Heating
- Ionization
- Excitation



Non-thermal electron degradation

COMPETITION BETWEEN PROCESSES



Non-thermal electron degradation

COMPETITION BETWEEN PROCESSES

https://github.com/eliotayache/elec_degrad

Plasma losses: [Huba+2013](#)

Bremsstrahlung:

[Barnes&Kasen2016](#): Includes tabulated cross-sections from Seltzer&Berger for various compositions. Includes screening at low energies.

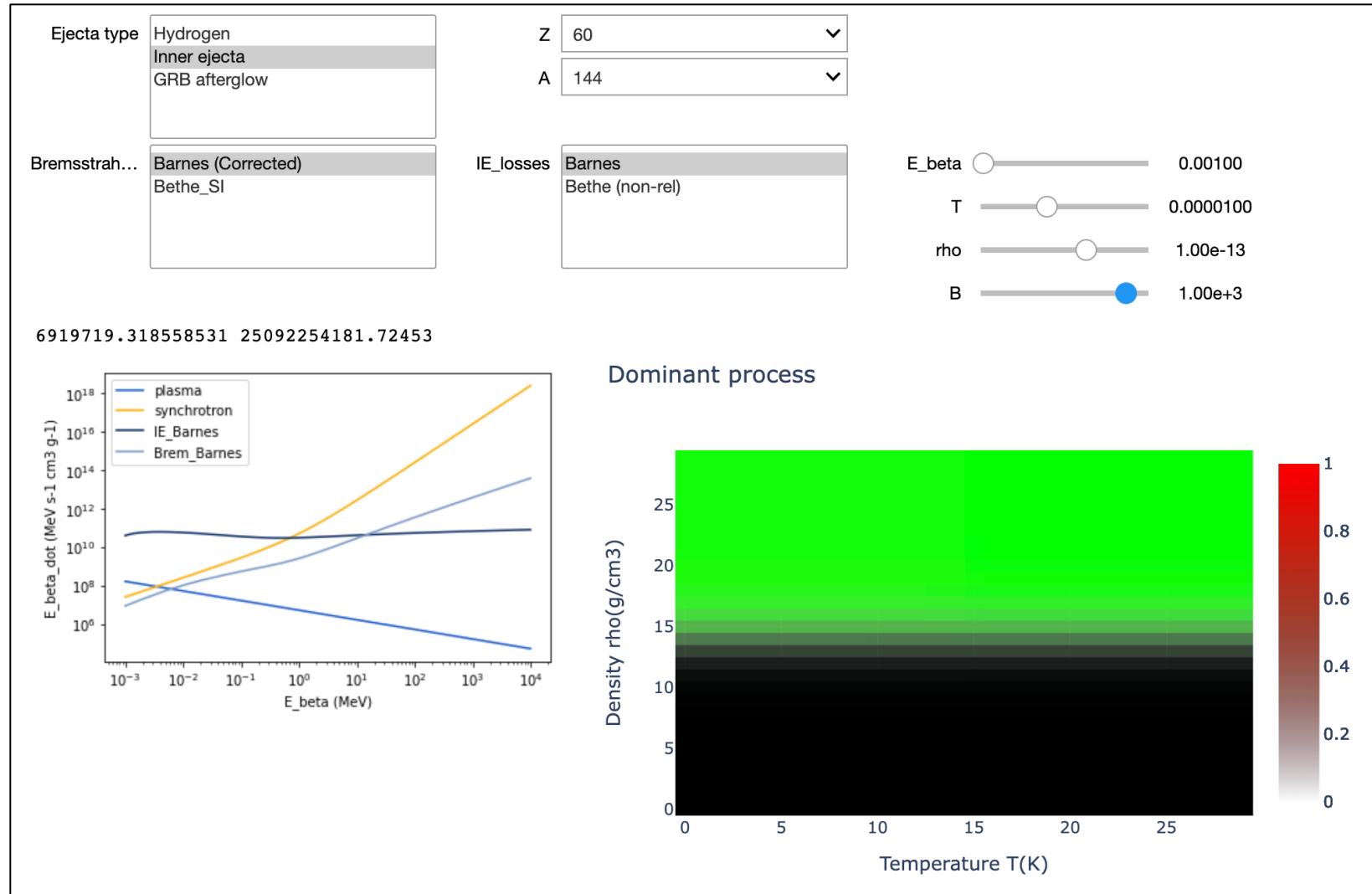
[Bethe1934](#): Does not include screening at low energies.

Ionization/Excitation:

[Barnes&Kasen2016](#): includes excitation and ionization for averaged composition
[Bethe formulation](#): non-relativistic. Only includes ionization.

Synchrotron

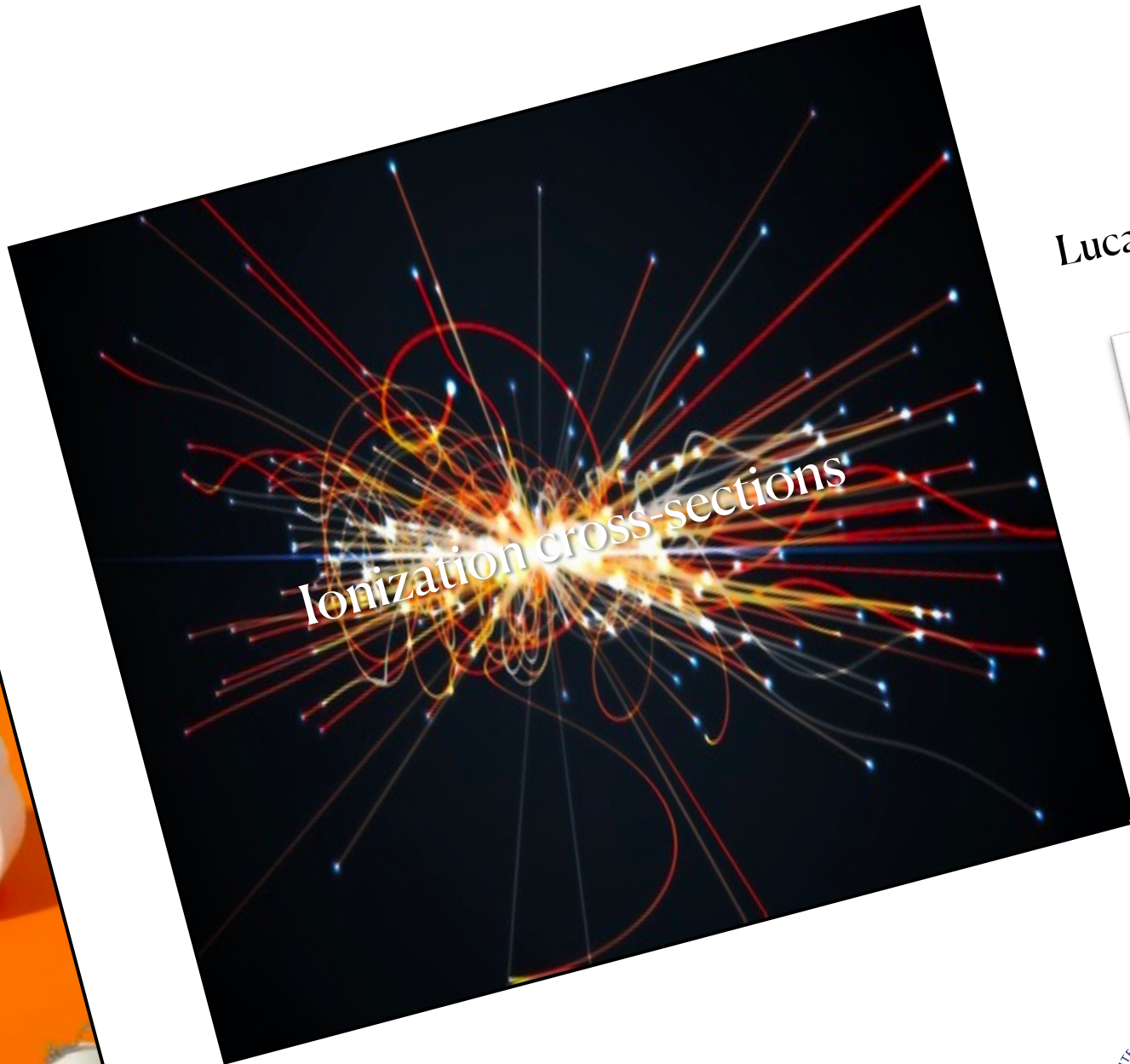
total synchrotron radiated energy for a single electron.



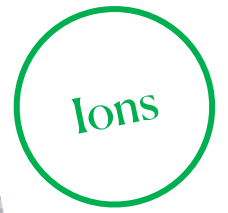
Non-thermal electron degradation

ASSUMPTIONS AND CHALLENGES

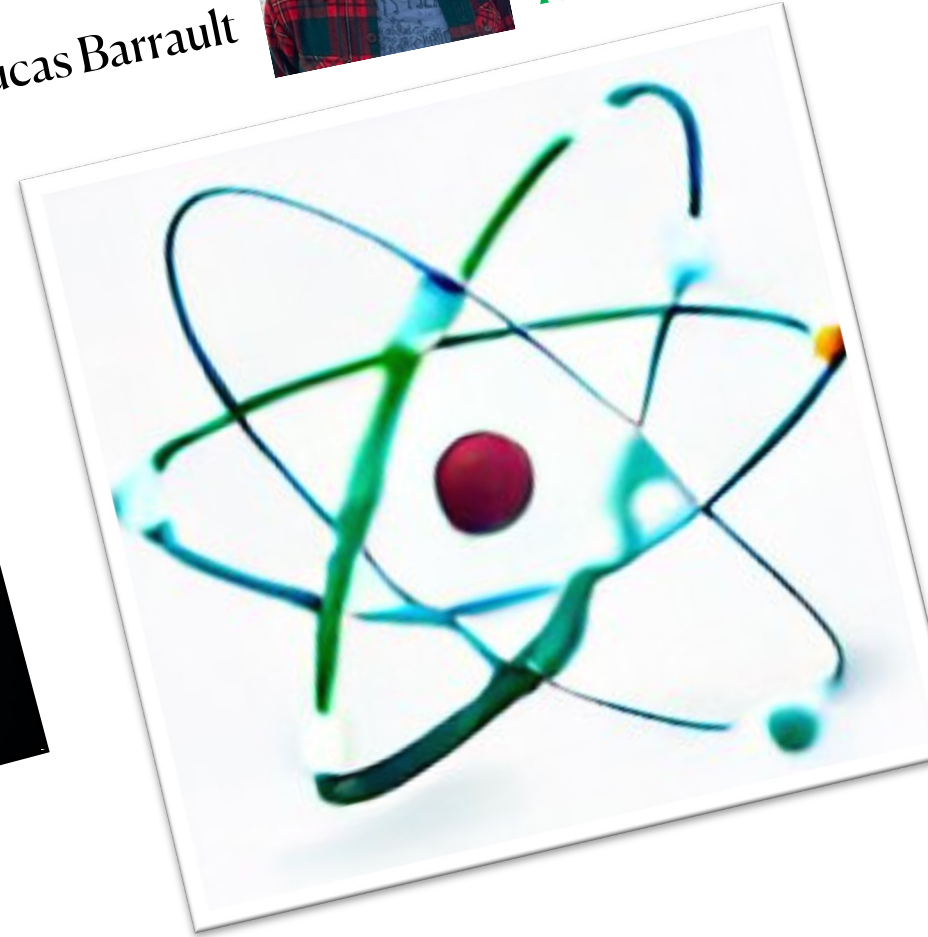




Lucas Barrault



Ionization



Ionization cross-sections

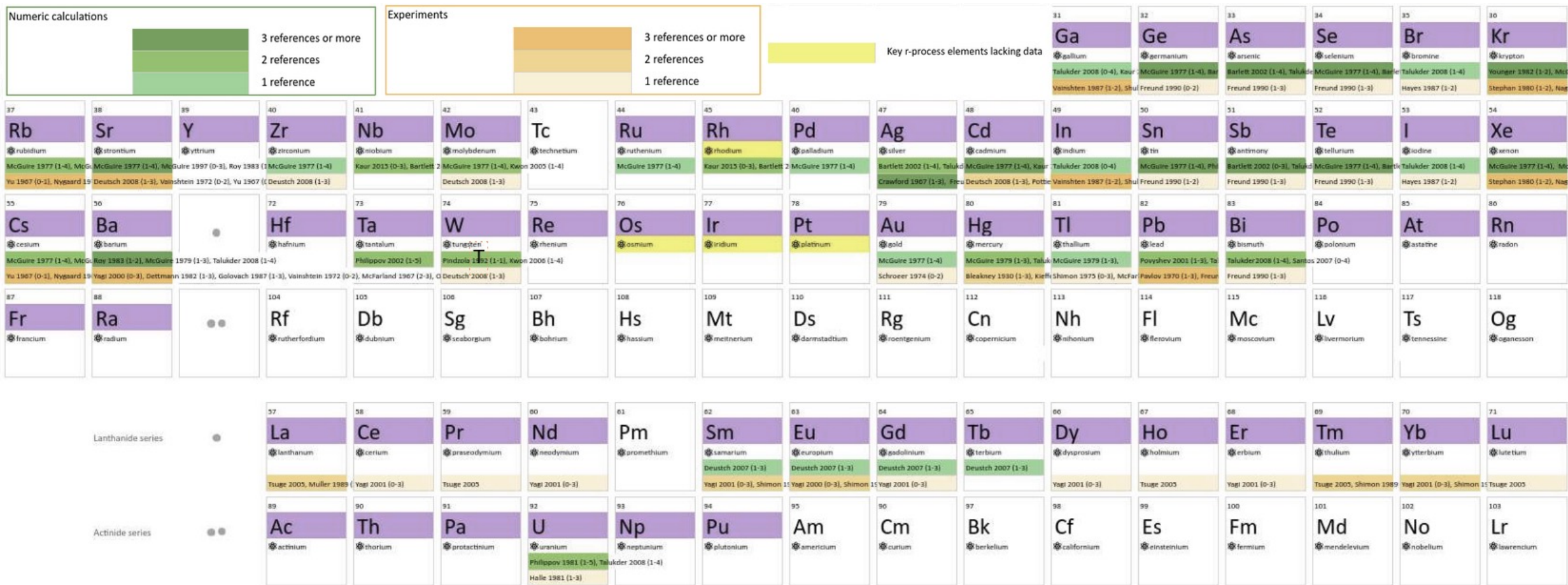
DATA DIMENSIONALITY

For **each element** the cross section value will depend on:

- **Ionization level** -> How many electrons does the initial ion have?
- **Subshell** -> Which subshell is loosing its electron? “partial” CS / “total” CS
- **Ion excitation state** -> What is the excitation structure of the resulting ion?
- **Incoming electron energy** -> How much energy is available? “total” CS
- **Secondary electron energy** -> “differential” CS

Ionization cross-sections

DATA AVAILABILITY



Ionization cross-sections

DATABASES

Experimental data

NIFS DATABASE.

<National Institute for Fusion Science>

Atomic and Molecular Numerical Databases

Cross Sections and Rate Coefficients for Ionization, Excitation, and Recombination by Electron Impact
Charge Transfer by Heavy Particle Collision, and Collision Processes of Molecules,
Sputtering Yields of Solids, and Back Scattering Coefficients from Solids

Bibliographic Databases

Fusion and Plasma Sciences, Atomic and Molecular Physics, and Atomic Collision Processes

Made by A&M Data Research Section

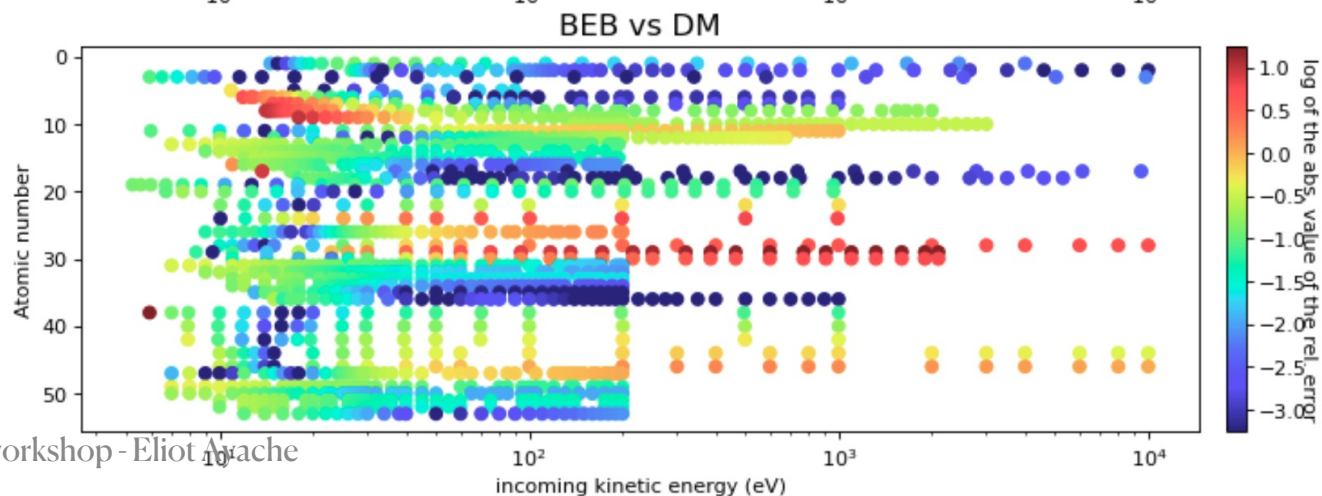
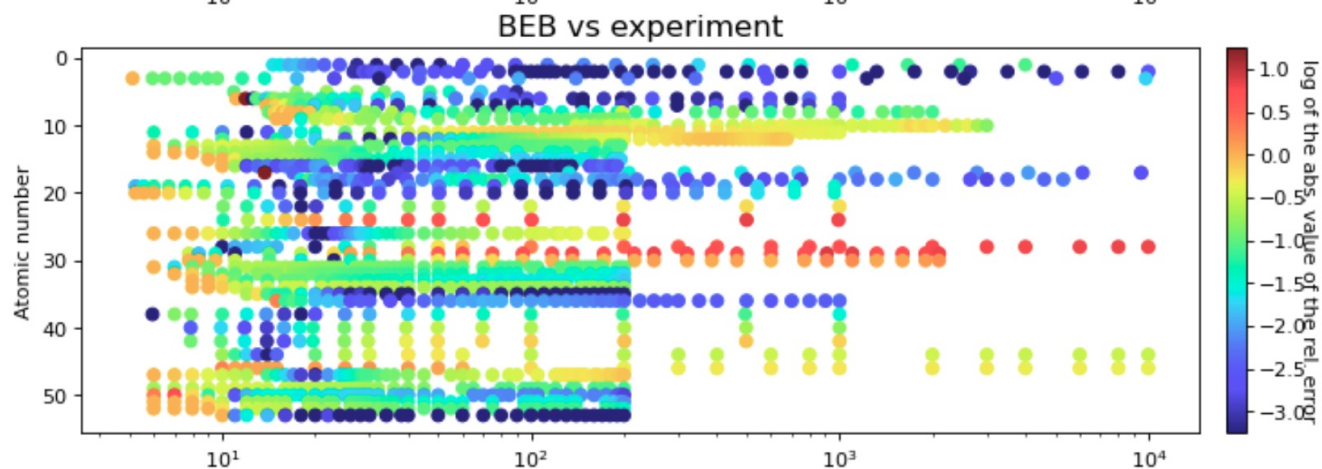
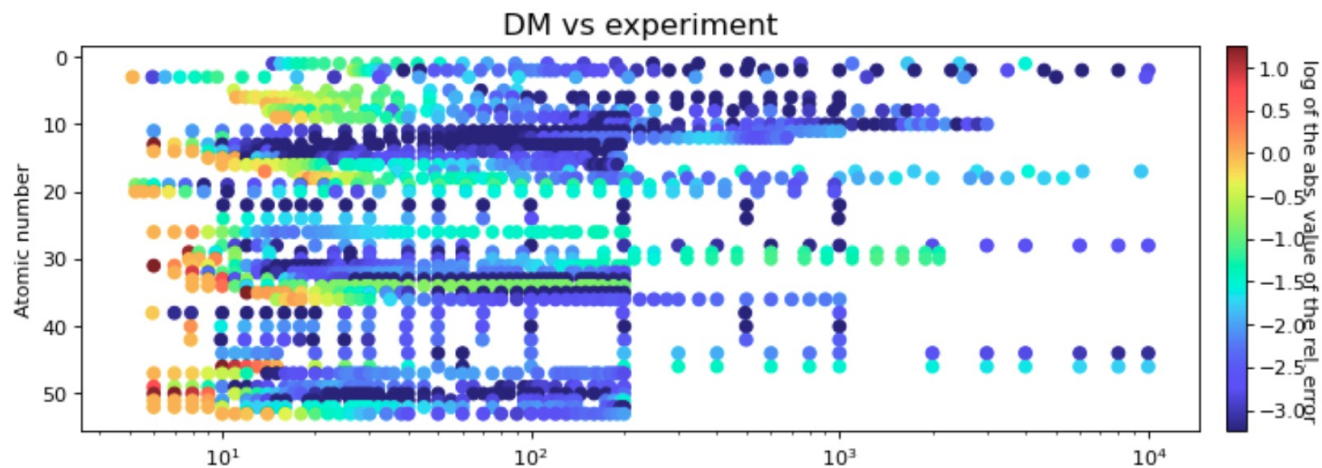
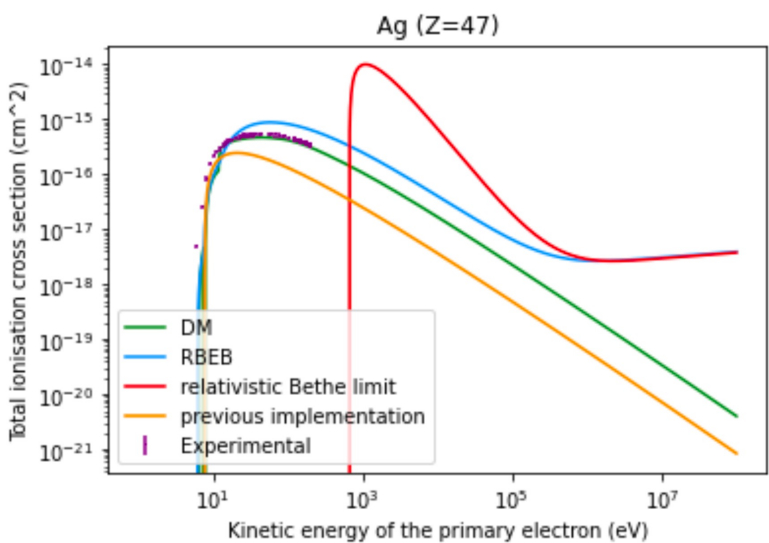
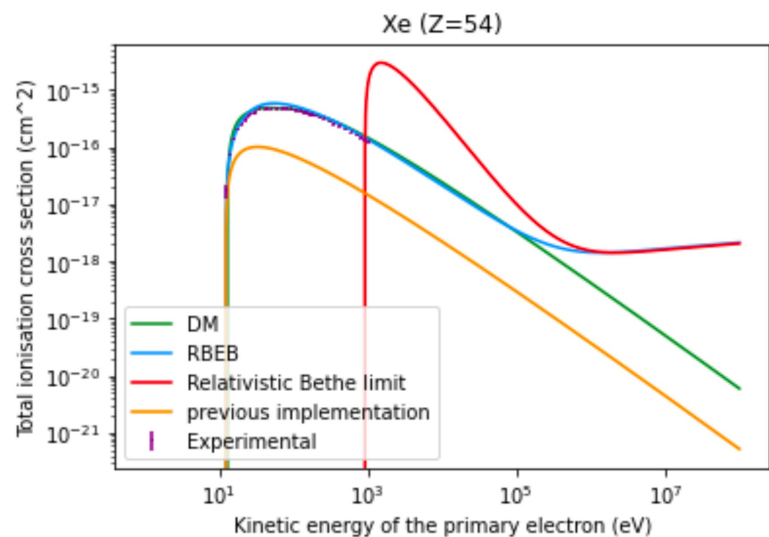
<https://dbshino.nifs.ac.jp/>

Model parameters

- Deutsch-Mark
- Binary Encounter

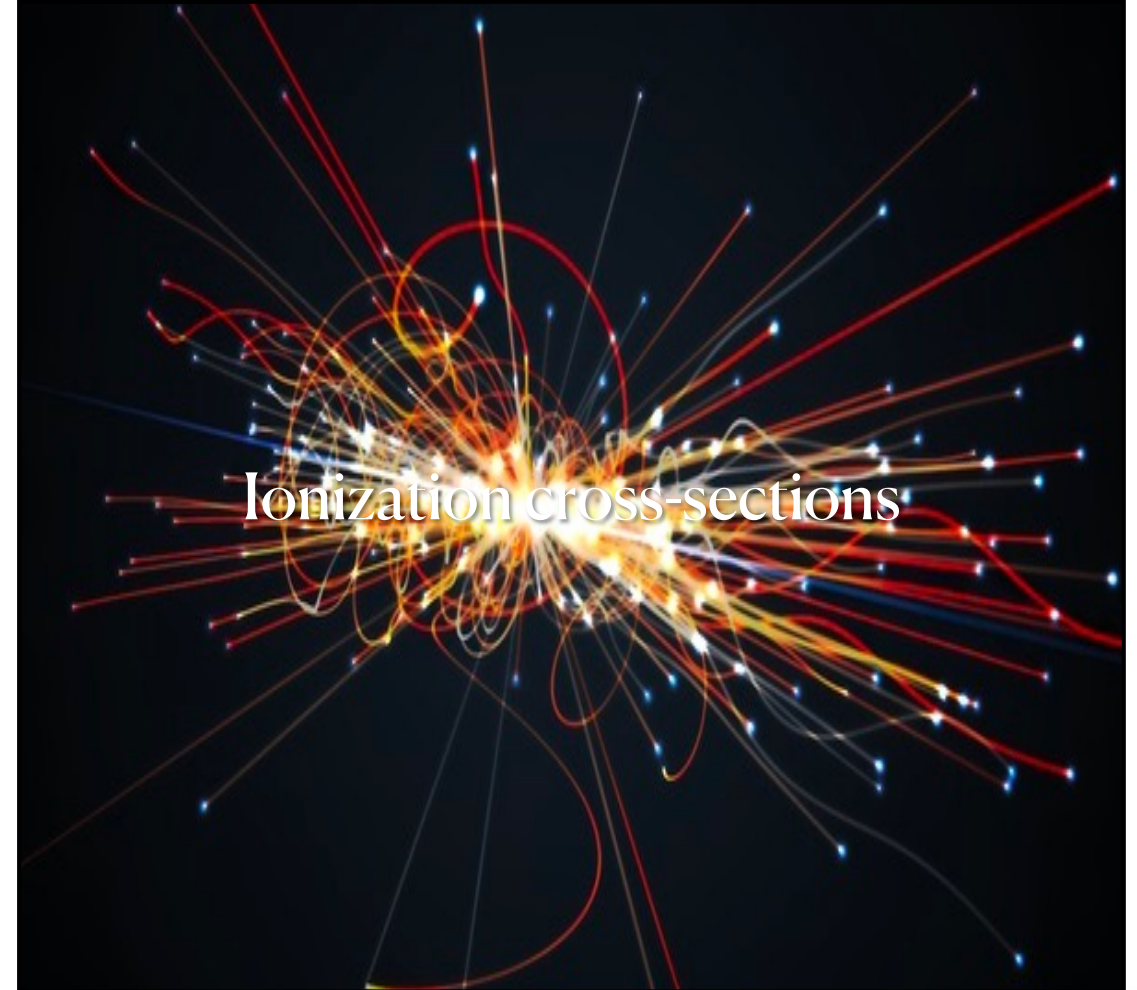
EADL -> Binding energies, kinetic energies for the subshells
NIST -> Ionization energy thresholds

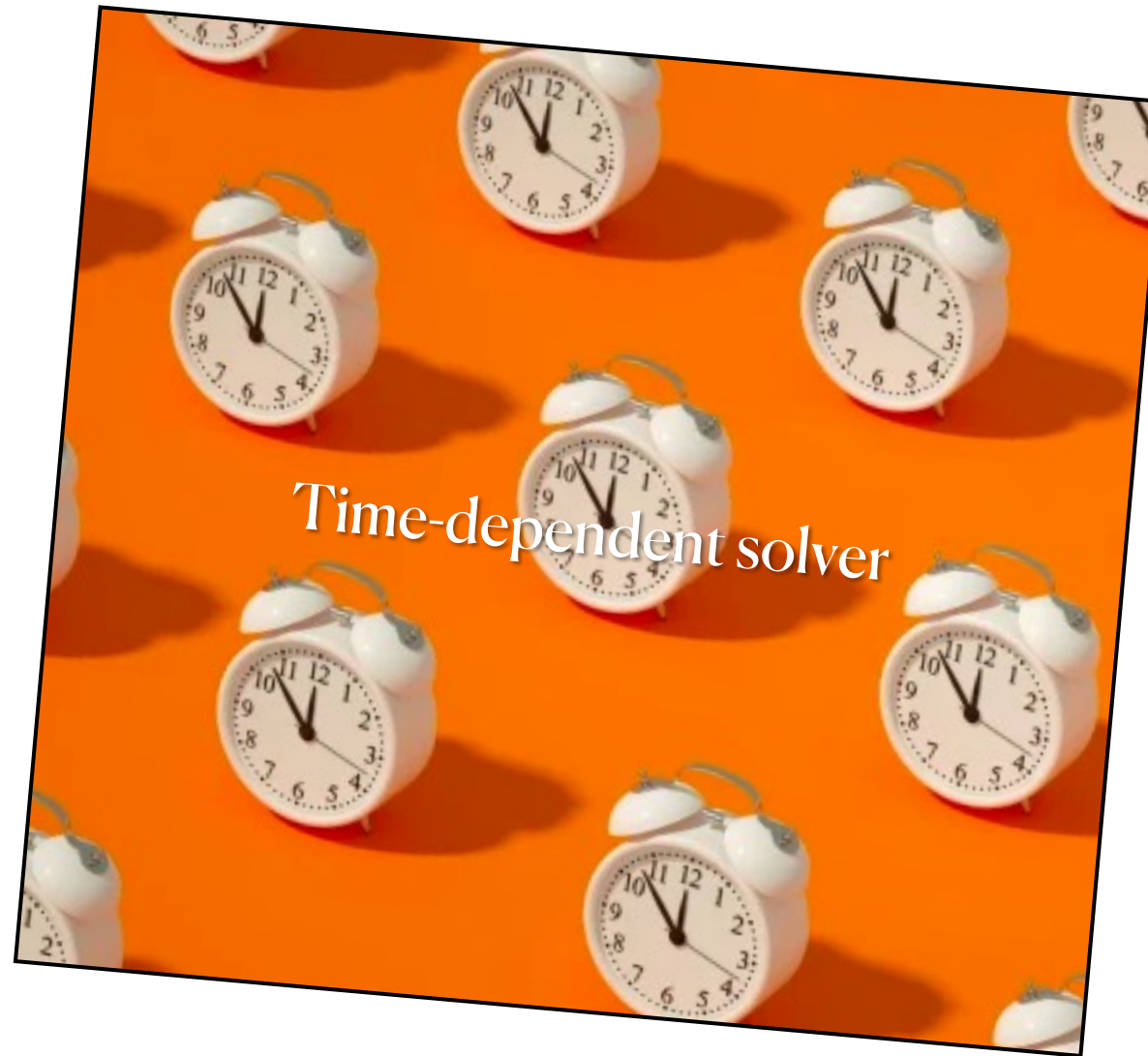
Model comparison



Non-thermal electron degradation

ASSUMPTIONS AND CHALLENGES





Time-dependent formulation

MATHEMATICAL FORMULATION

$$\frac{\partial F(E; t)}{\partial t} = n \left[\int F(E') R(E' \rightarrow E) dE' - \int F(E) R(E \rightarrow E'') dE'' \right] \quad \text{Vanilla Boltzmann equation}$$

$$\frac{1}{v_E} \frac{\partial z(E, t)}{\partial t} = \underbrace{nK z(E, t)}_{\text{Ionization}} + \underbrace{\frac{1}{v_E} \left[\frac{\partial z \mathcal{L}_e}{\partial E} - \frac{z \mathcal{L}_e}{v_E} \frac{\partial v_E}{\partial E} \right]}_{\text{Thermal losses}} + \underbrace{\frac{S(E, t)}{v_E}}_{\text{Energy deposition}} \quad \begin{array}{l} + \text{Excitation} \\ + \text{Adiabatic expansion} \\ + \dots \end{array}$$

Ionization

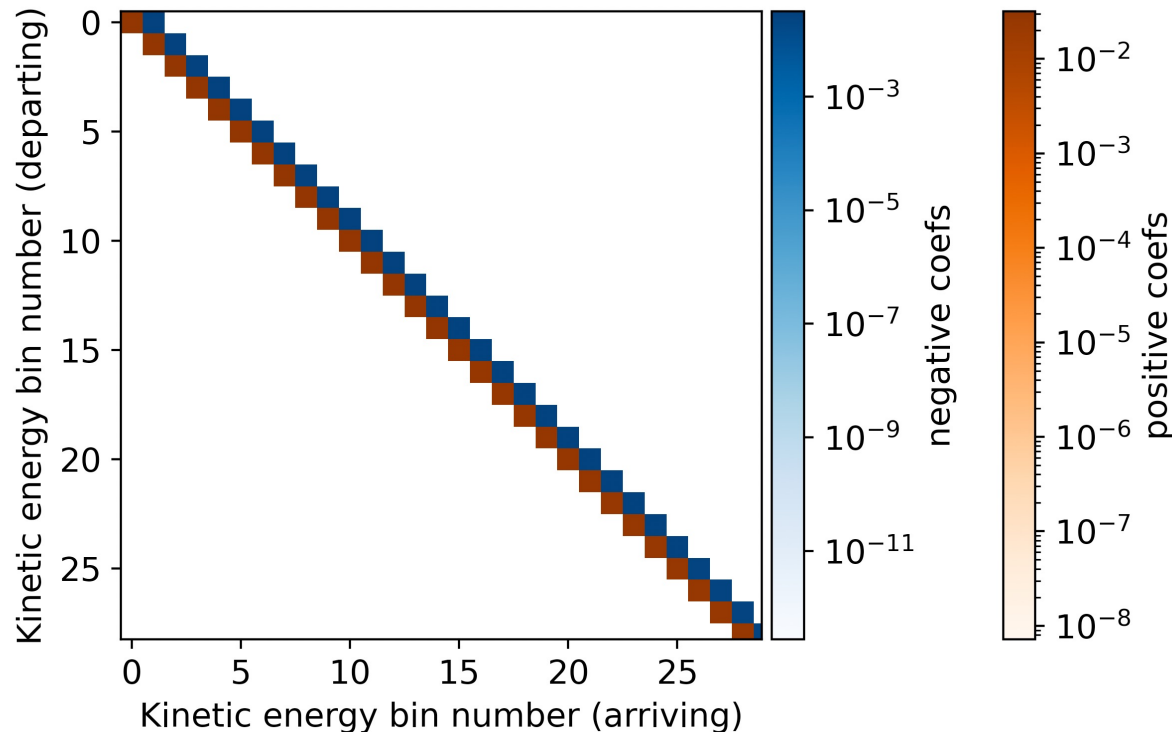
Thermal losses

Energy deposition

Time-dependent formulation

DISCRETIZATION

$$z_m^n \left[\frac{1}{v_m(t_n - t_{n-1})} - \frac{L_m}{v_m(E_m - E_{m-1})} + \frac{L_m}{v_m^2} \frac{dv}{dE} \Big|_m + n \sum_{m'} \sigma_{m \rightarrow m'} dE_{m'} \right] - n \sum_{m'} z_{m'}^n \sigma_{m' \rightarrow m} dE_{m'} + z_{m-1}^n \frac{L_{m-1}}{v_m(E_m - E_{m-1})} = \frac{z_m^{n-1}}{v_m(t_n - t_{n-1})}$$



“Upwinding” -> stability
Upper triangular -> energy resolution

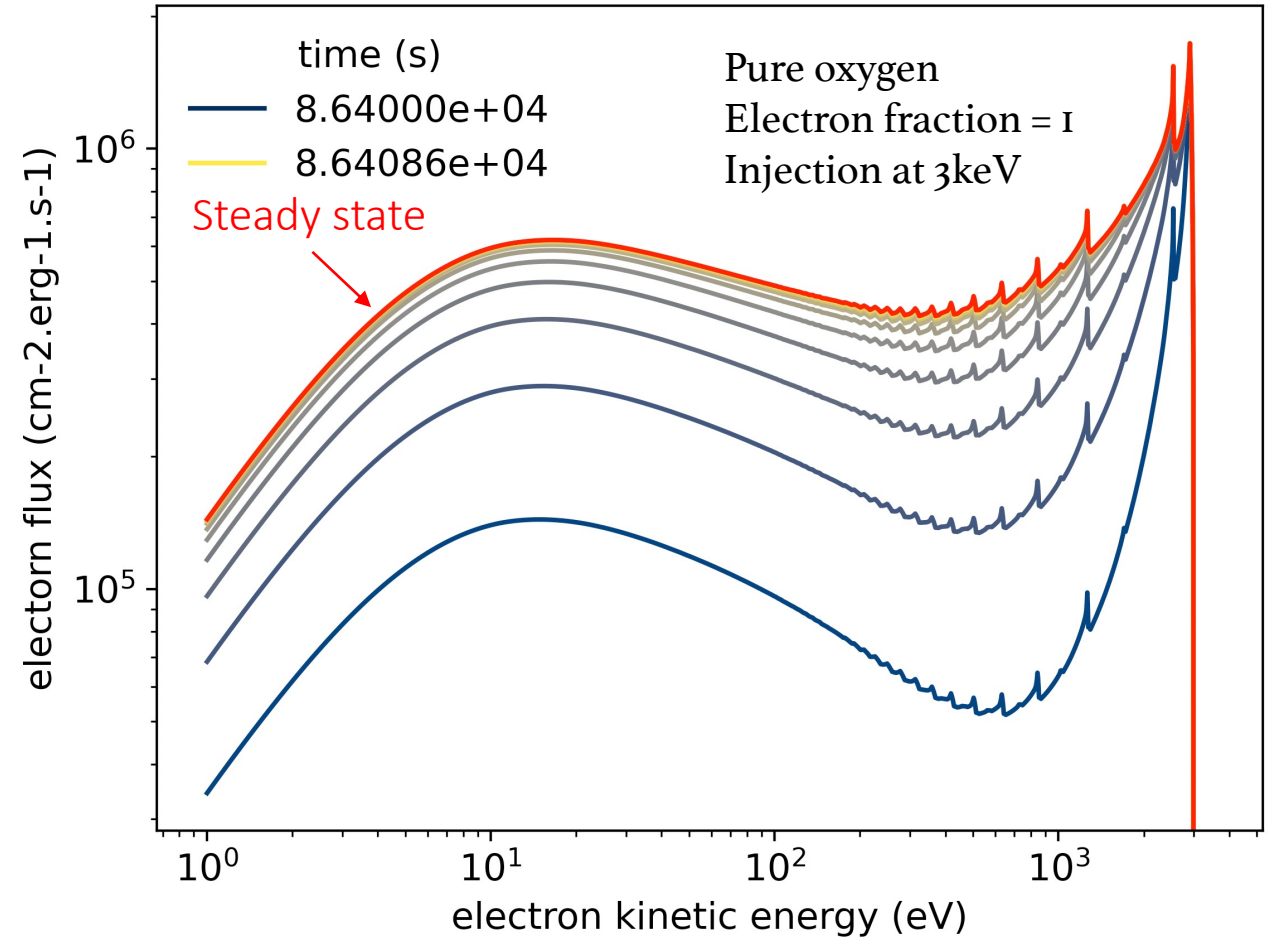
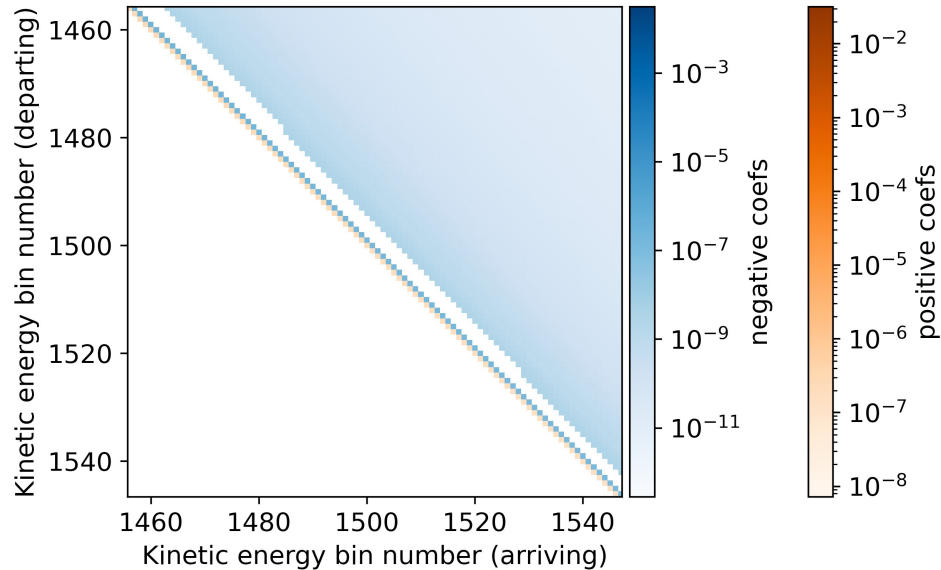
Thermal losses: **BARNES & KASEN 2016**

$$\dot{E}_\beta^{\text{pl}} = 7.7 \times 10^{-15} E_\beta^{-1/2} \times \frac{n_e}{1 \text{cm}^{-3}} \lambda_{ee} \left(1.0 - \frac{3.9 T}{7.7 E_\beta} \right) \text{MeVs}^{-1}$$

Preliminary results

TIME-DEPENDENT EVOLUTION

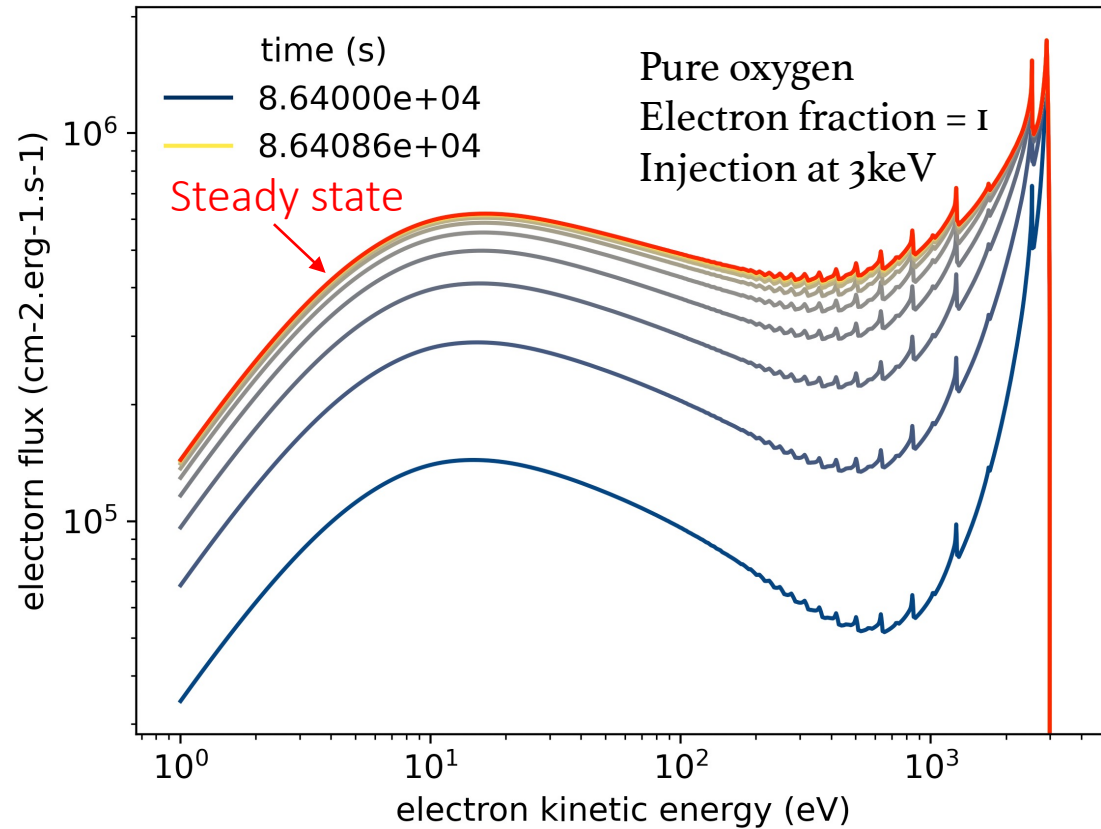
Converges onto steady state solution
Numerical artefacts: resolution dependent



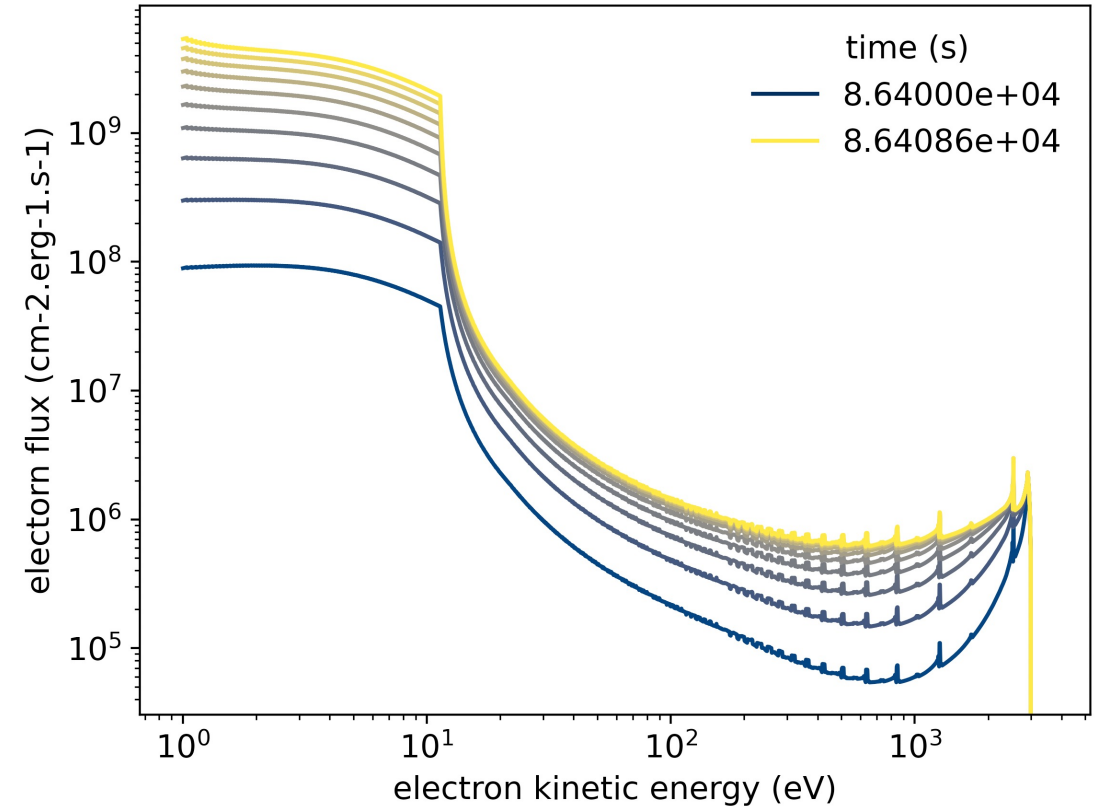
Preliminary results

TIME-DEPENDENT EVOLUTION

With thermal losses



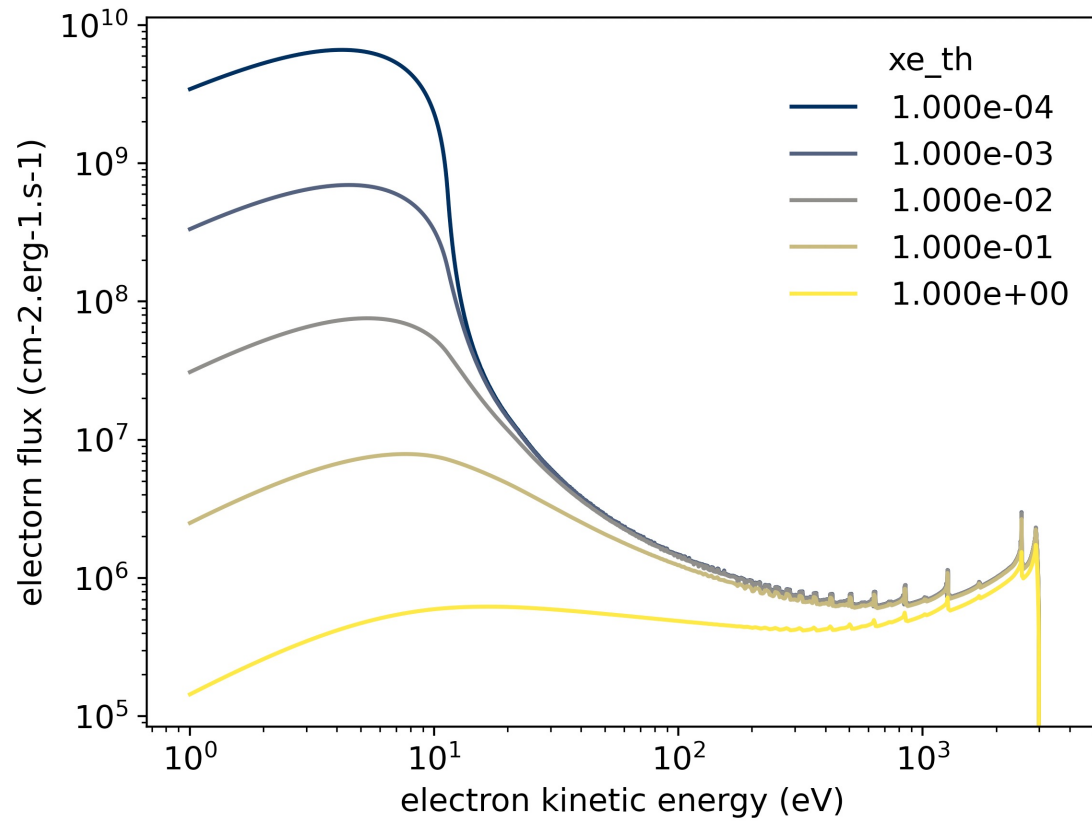
Without thermal losses



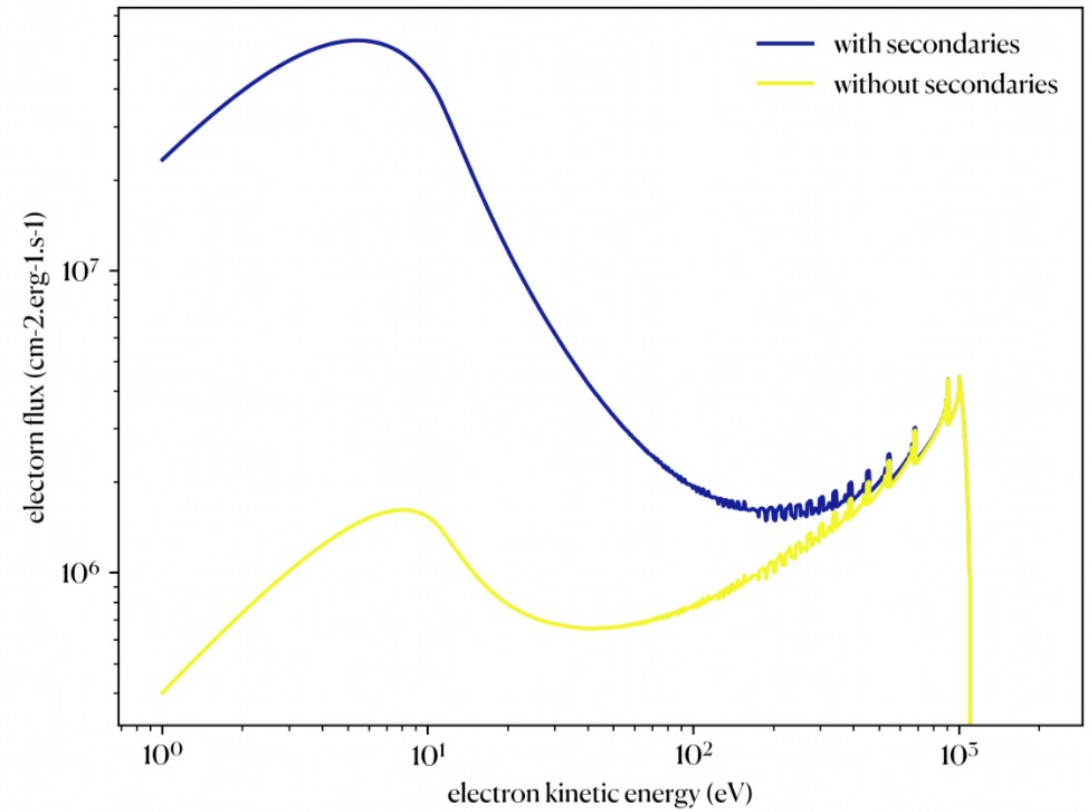
Preliminary results

PURE OXYGEN

Effect of electron fraction



Effect of secondaries



Conclusions

- Time-dependence needs implementing AT ALL STAGES of the NTLE processes
- More work is needed on measuring and modelling r-process cross-sections.
- Numerical non-thermal electron degradation is data heavy and computationally expensive, but can be carried out in a self-consistent way.
- Future work:
 - Adiabatic processes
 - Target excited states
 - Auger processes