

Outline

- What this is all about?
- Coupling of generated energy with hydrodynamic
- Selection of the isotopes
- The CNO-Cycle as an example
- Definition of the network
- Nuclear statistical equilibrium as an approximation

What this is all about?

Why are we doing all this?

- Because we can do it!
- We want to explain the origin of the isotopes
- Nucleosynthesis calculations can give clues to explosion models of supernovae when they are compared with observations
- It gives answers to other questions like, evolution of stars or chemical evolution of galaxies.

Energy generation

- Each reaction generates or consumes energy.
→ coupling with hydrodynamic
- Depending on the reaction type, the generated energy may be kinetic or in form of photons or neutrinos. E.g. $A(x,y)B$, $A(x,\gamma)B$, $A(x,\nu)B$ reactions.
- The energy in form of photons and neutrinos may or may not escape. This depends on the external conditions (density, temperature) of the astrophysical site.
- If they escape, this leads to an energy loss.

Coupling to hydrodynamic

Due to the energy generation, there exists an additional source term $\rho\dot{\epsilon}$ in the conservation law for the total specific energy:

$$\dot{\epsilon} = h(d(\text{Amount of isotopes}), T)$$

plus network equations, which will be defined later.

$$[\dot{\epsilon}] = \text{erg g}^{-1}\text{s}^{-1}$$

Problems due to the Coupling

- Problems are mainly computational limitations, because solving the network equations **can** be very time consuming. (This depends on the problem).
- Another problem is due to the fact, that solving the set of hydrodynamical equations **is** usually very time consuming.

⇒ you are forced to make some simplifications

Selection of the Isotopes

The first question you have to ask yourself is:
„Do I want to model a hydro static object (e.g. the sun) or
do I want to model an explosion (e.g. supernovae)?“

- In the first case the hydrodynamical timescales are long compared to nuclear timescales. \Rightarrow it is a good approximation to neglect some β – unstable isotopes.
- In the second case the hydrodynamical timescales can be compared to nuclear timescales. \Rightarrow you have to consider β – unstable isotopes, because in these cases the cross sections for strong reactions are high enough to be important.

The second Case

- In principle one has to consider more isotopes.
- This is usually not done due to computational limitations

Simplification:

A reduced network is used to get an approximation for the Energy generation.

E.g. in type Ia supernova simulations done at the MPA the “network” consists of the isotopes ^{12}C , ^{16}O , ^4He , “ ^{24}Mg ”, “ ^{56}Ni ”

How do you get information of the nucleosynthesis in such cases?

- When simulating an explosive event, one puts „tracer particles“ in the simulated fluid.
- These particles record temperature and density
- These information are sufficient to calculate the nucleosynthesis by solving the network equations (will be defined later) in a post process.

The first Case (E.g. CNO-Cycle)

- May start, if sufficient heavy isotopes are available when a star has formed.
- The temperatures are usually of the order of 10^6 K

The basic reactions (CN-Cycle):



weak reactions,
 ν escapes

NO-Cycle:



Isotope	Lifetime
^{13}N	870 sec
^{15}O	178 sec
^{17}F	95 sec

Lifetime of Isotope X

Definition: Lifetime of Isotope X $\tau(X) := \frac{X}{dX/dt} = \frac{X}{\dot{X}}$

Now consider a reaction: $X(p,\gamma)Y$

You can calculate the lifetime of X against proton capture

$$\tau(X) = \frac{1}{p \times \lambda_{pX}}$$

Where p stands for the number of protons per unit volume

and λ_{pX} is the temperature dependent (!) cross section for the reaction.

⇒ if τ is small, then a reaction is very likely

if τ is large, then a reaction may be neglected

How to model the CNO-Cycle?

$$d^{12}\text{C}/dt = \lambda_{p15}\text{H}^{15}\text{N} - \lambda_{p12}\text{H}^{12}\text{C}$$

$$= {}^{15}\text{N}/\tau_{15} - {}^{12}\text{C}/\tau_{12} +$$

$$d^{13}\text{N}/dt = {}^{12}\text{C}/\tau_{12} - {}^{13}\text{N}/\tau_{\beta}(13) - \cancel{{}^{13}\text{N}/\tau_p(13\text{N})}$$

$$d^{13}\text{C}/dt = {}^{13}\text{N}/\tau_{\beta}(13) - {}^{13}\text{C}/\tau_{13}$$

$$d^{14}\text{N}/dt = {}^{13}\text{C}/\tau_{13} - {}^{14}\text{N}/\tau_{14} + {}^{17}\text{O}/\tau_{17}$$

$$d^{15}\text{O}/dt = {}^{14}\text{N}/\tau_{14} - {}^{15}\text{O}/\tau_{\beta}(15) - \cancel{{}^{15}\text{O}/\tau_p(15\text{O})}$$

$$d^{15}\text{N}/dt = {}^{15}\text{O}/\tau_{\beta}(15) - {}^{15}\text{N}/\tau_{15}$$

$$d^{16}\text{O}/dt = {}^{15}\text{N}/\tau_{15} - {}^{16}\text{O}/\tau_{16}$$

$$d^{17}\text{F}/dt = {}^{16}\text{O}/\tau_{16} - {}^{17}\text{F}/\tau_{\beta}(17) - \cancel{{}^{17}\text{F}/\tau_p(17\text{F})}$$

$$d^{17}\text{O}/dt = {}^{17}\text{F}/\tau_{\beta}(17) - {}^{17}\text{O}/\tau_{17}$$

Due to the fact that

$$\tau_p \gg \tau_{\beta}$$

Only for temperatures of the order of 10^6 K !

Further simplifications

Consider the equation
$$\frac{d^{13}\text{N}}{dt} = \frac{^{12}\text{C}}{\tau_{12}} - \frac{^{13}\text{N}}{\tau_{\beta}(13)}$$

Suppose ^{12}C and the τ 's are constant over sufficiently long period of time.

Then this equation is self regulating in the sense that the left side is >0 , if there is not enough ^{13}N and vice versa.

\Rightarrow Is there a state where the left side is $=0$?

YES! At least as a very good approximation.

Further Simplifications

A solution of $\frac{d^{13}\text{N}}{dt} = \frac{^{12}\text{C}}{\tau_{12}} - \frac{^{13}\text{N}}{\tau_{\beta}(13)}$ with the mentioned assumptions is given by

$$^{13}\text{N}(t) = \frac{\tau_{\beta}(13)}{\tau_{12}} ^{12}\text{C} (1 - e^{-t/\tau_{\beta}(13)})$$

⇒ This equation reaches an equilibrium value ($dN/dt=0$) of $(^{13}\text{N}/^{12}\text{C}) = \tau_{\beta}(13)/\tau_{12}$ in times of the order of τ_{β}

⇒ In times of the order of minutes it is possible to eliminate the Isotopes ^{13}N , ^{15}O and ^{17}F from the system that describes the CNO-cycle

Result of the Simplification

$$d^{12}\text{C}/dt = {}^{15}\text{N}/\tau_{15} - {}^{12}\text{C}/\tau_{12}$$

$$d^{13}\text{C}/dt = {}^{12}\text{C}/\tau_{12} - {}^{13}\text{C}/\tau_{13}$$

$$d^{14}\text{N}/dt = {}^{13}\text{C}/\tau_{13} - {}^{14}\text{N}/\tau_{14} + {}^{17}\text{O}/\tau_{17}$$

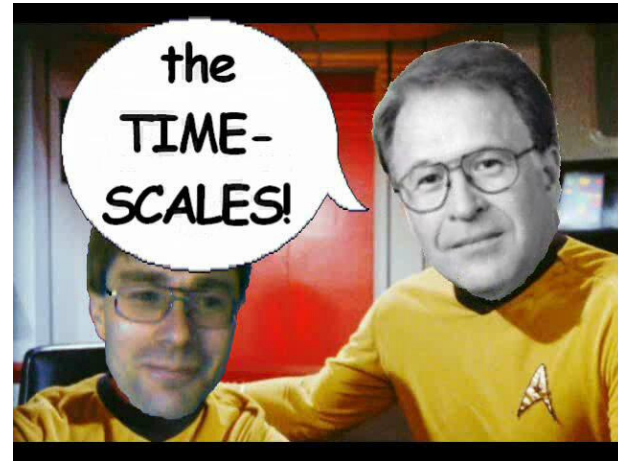
$$d^{15}\text{N}/dt = {}^{14}\text{N}/\tau_{14} - {}^{15}\text{N}/\tau_{15}$$

$$d^{16}\text{O}/dt = {}^{15}\text{N}/\tau_{15} - {}^{16}\text{O}/\tau_{16}$$

$$d^{17}\text{O}/dt = {}^{16}\text{O}/\tau_{16} - {}^{17}\text{O}/\tau_{17}$$

When you model non explosive astrophysical sites you can usually make some simplifications for your network

BUT



Definition of a Network

- The CNO-Cycle is just an example of a network.
- The isotopes in here are p and some C,N,O – isotopes; a relatively small number
- The relevant reactions are proton capture reactions and weak decays.

In general one has one deal with a lot more reactions like α captures, photo disintegrations, heavy ion reactions, “weak capture reactions”, decays and three body reactions

Definition of a Network

- If you want to model all these reactions, you have to write down a more general set of differential equations than it was done for the CNO-Cycle.

Suppose you want to consider m different isotopes. And let n_i be the number of isotope i per unit volume, then we obtain the following equation:

Definition of a Network

$$\dot{n}_i = \sum_j c_i^j \lambda_i^j n_j + \sum_{j,k} c_i^{j,k} \langle j, k \rangle n_j n_k + \sum_{j,k,l} c_i^{j,k,l} \langle j, k, l \rangle n_j n_k n_l$$

Statistical factors
to prevent
double counting

$$c_i^j := \pm N_i$$

$$c_i^{j,k} := \pm \frac{N_i}{N_j! N_k!}$$

$$c_i^{j,k,l} := \pm \frac{N_i}{N_j! N_k! N_l!}$$

$$d^{12}C/dt = {}^{15}N/\tau_{15} - {}^{12}C/\tau_{12}$$

$$d^{13}C/dt = {}^{12}C/\tau_{12} - {}^{13}C/\tau_{13}$$

$$d^{14}N/dt = {}^{13}C/\tau_{13} - {}^{14}N/\tau_{14} + {}^{17}O/\tau_{17}$$

$$d^{15}N/dt = {}^{14}N/\tau_{14} - {}^{15}N/\tau_{15}$$

$$d^{16}O/dt = {}^{15}N/\tau_{15} - {}^{16}O/\tau_{16}$$

$$d^{17}O/dt = {}^{16}O/\tau_{16} - {}^{17}O/\tau_{17}$$

Difficulties with this Form

$$\dot{Y}_i = \sum_j c_i^j \lambda_i^j Y_j + \sum_{j,k} c_i^{j,k} \rho N_A \langle j, k \rangle Y_j Y_k + \sum_{j,k,l} c_i^{j,k,l} (\rho N_A)^2 \langle j, k, l \rangle Y_j Y_k Y_l$$

- When you want to calculate a solution for an explosion, you encounter the problem, that total number of isotopes is changing.
→ you would run into unnecessary bookkeeping
- you have to introduce some new quantities
 - a) $\rho := (\sum_i n_i A_i) / N_A$
 - b) Mass fraction: $X_i := (n_i A_i) / (\rho N_A)$ with $\sum_i X_i = 1$
 - c) Specific abundances: $Y_i := X_i / A_i = n_i / (\rho N_A)$
- non linear & stiff \Rightarrow numerical solution is not trivial
- these equations define the nuclear network!

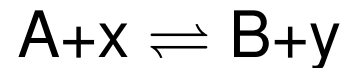
Summary 1

What do we know?

- We know that the last equation defines a nuclear network
- Depending on the situation (hydrostatic situation or explosion) the size of the network can vary.
- When you model the CNO-Cycle, 10 isotopes are sufficient
- But in case of (supernova) explosions you need about a factor of 30 to 50 more isotopes! (only in a post process)
- The energy generation of the network is coupled to hydrodynamic. → A reduced network is often used

The NSE

- The NSE is characterized by the fact that all strong reactions are in an equilibrium



where A,B,x,y can be every kind of particle!

- This can occur if the temperatures are so high that basically all Coulomb barriers can be overcome. (densities above 10^7 g/cm⁻³ assumed).
- This means corresponding reaction rates are identical (up to ± 1). → The network equations are replaced.

The NSE

- If you assume that ALL strong reactions are in an equilibrium, you may write for the chem. potential μ of a nucleus (A,Z) with Z protons and $A-Z$ neutrons

$$\mu(A,Z) = Z\mu_p + (A-Z)\mu_n + B(A,Z)$$

where $B(A,Z) = c^2(ZM_p + (A-Z)M_n - M(A,Z))$ is the binding energy of the nucleus (A,Z)

- Assuming a non relativistic Boltzmann gas you obtain for the number of nucleus (A,Z) per unit volume

$$n(A,Z) = G(A,Z) A^{3/2} h^{-3} (2\pi M_u kT)^{3/2} \exp(\beta\mu(A,Z))$$

The NSE

Using the last equation for the nucleus, protons and neutrons and the relation for the chemical you will get

$$n(A, Z) = G(A, Z) A^{3/2} \frac{n_p^Z n_n^{A-Z}}{2^A} \left(\frac{(2\pi M_u kT)^{3/2}}{h^3} \right)^{1-A} \exp\left(\frac{B(A, Z)}{kT}\right)$$

- The number densities are fixed by the neutron to proton ratio, the binding energy and the temperature.
- The neutron to proton ratio can only be changed by weak interactions.
- A “weak network” is solved first, then together with mass and charge conservation the new number densities are calculated.

Literature

- „Principles of Stellar Evolution and Nucleosynthesis“ – D.D. Clayton 1983 Univ. Chicago Press.
- „Nucleosynthesis and Supernovae“ – W.D. Arnett 1996 Princeton Univ. Press
- Thielemann et. al. Astro-ph/9802077 „Nucleosynthesis Basics and Applications to Supernovae
- Hillebrandt & Niemeyer „Type Ia Supernova Explosion Models“ Annu. Rev. As. Aphy. 2000. 38:191-230
- References in literature mentioned above