# Types of Equilibria

- Steady Flow of Reactions
- Chemical Equilibrium of Reactions
- Complete Chemical Equilibrium (NSE)
- Clusters of Chemical Equilbrium (QSE)
- QSE Clusters linked by Steady Flow

# CNO(I)-Cycle in Steady Flow

The CNO-Cycles in Hydrogen Burning

cycle	reaction sequence
CNOI	$^{12}C(p,\gamma)^{13}N(e^+\nu)^{13}C(p,\gamma)^{14}N(p,\gamma)^{15}O(e^+\nu)^{15}N(p,\alpha)^{12}C$
CNOII	$^{15}N(p,\gamma)^{16}O(p,\gamma)^{17}F(e^{+}\nu)^{17}O(p,\alpha)^{14}N$
CNOIII	$17O(p,\gamma)^{18}F(e^+\nu)^{18}O(p,\alpha)^{15}N$
CNOIV	$^{18}\mathrm{O}(\mathrm{p},\gamma)^{19}\mathrm{F}(\mathrm{p},lpha)^{16}\mathrm{O}^{-10}$

$$\begin{split} Y_{1} = & \bar{\rho} N_{A} \langle 12, 1 \rangle Y_{12} Y_{1} - \rho N_{A} \langle 13, 1 \rangle Y_{13} Y_{1} - \rho N_{A} \langle 14, 1 \rangle Y_{14} Y_{1} \\ & - \rho N_{A} \langle 15, 1 \rangle Y_{15} Y_{1} \\ & = -4 C_{CNO} = -4 \rho N_{A} \langle 14, 1 \rangle Y_{14} Y_{1} = -\frac{1}{\tau_{1,14}} Y_{1} \end{split}$$

 $\dot{Y}_4 = \rho N_A \left< 15, 1 \right> Y_{15} Y_1 = C_{CNO}$ 

the network entry for nuclei with mass numbers A=12, 13, 14, 15 is governed in each case by a production reaction (proton reaction on A-1) and a distruction reaction (proton reaction on A). In case of a steady flow they cancel and lead to Y=0 for all A, linking all of these terms and identical to (A=14 is useful as this encounters the slowest reaction and essentially all mass assembles in <sup>14</sup>N)

$$C_{CNO} = 
ho N_A \langle 14, 1 \rangle Y_{14} Y_1$$
 $Y_{14} \approx rac{1.4 imes 10^{-2}}{14}$ 

summing all mass fractions of CNO nuclei for solar metallicity

### s-process and steady flow



shown are s-, r-, and p-only nuclei!

### s-process and steady flow



## The sigma\*N-curve



a complete steady flow is not given, but in between magic numbers (where the neutron capture cross sections are small) almost attained!

#### Käppeler et al. (2007): needed cross sections with uncertainties between 1 and 5% for complete set of isotopes from <sup>12</sup>C to <sup>210</sup>Po, including unstable samples (branching points)

![](_page_5_Figure_1.jpeg)

### s- and r-decomposition

![](_page_6_Figure_1.jpeg)

the almost constant sigma\*N-curve leads to a large odd-even staggering in the abundances (due to the odd-even staggering in n-capture cross sections!

# Steady flows and chem. equilibrium in stellar burning

pp-cycles and CNO-cycle lead to steady flows in H-burning

- 1. Hydrogen Burning $T = (1-4)x10^7 K$ pp-cycles-> $^{1}H(p,e^+\nu)^2 H$ CNO-cycle-> slowest reaction $^{14}N(p,\gamma)^{15}O$
- 2. Helium Burning  $T=(1-2)x10^{8}K$ <sup>4</sup>He+<sup>4</sup>He  $\Leftrightarrow$  <sup>8</sup>Be <sup>8</sup>Be( $\alpha,\gamma$ )<sup>12</sup>C[( $\alpha,\gamma$ )<sup>16</sup>O] <sup>14</sup>N( $\alpha,\gamma$ )<sup>18</sup>F( $\beta$ <sup>+</sup>)<sup>18</sup>O( $\alpha,\gamma$ )<sup>22</sup>Ne( $\alpha,n$ )<sup>25</sup>Mg

<sup>4</sup>He+<sup>4</sup>He → <sup>8</sup>Be is in chemical equilibrium released neutrons lead to steady flow in neutron capture

# Complete chem. equilibrium (NSE)

![](_page_8_Figure_1.jpeg)

Si-burning in stellar evolution and expl. Si-burning at high densities lead to NSE!

# QSE in explosive Si-burning

![](_page_9_Figure_1.jpeg)

# **QSE** Formalism

light group

$$Y_{NSE}(^{A}Z) = C(^{A}Z)Y_{n}^{N}Y_{p}^{Z}$$

$$Si-group$$

$$Y_{QSE,Si}(^{A}Z) = \frac{C(^{A}Z)}{C(^{28}Si)}Y(^{28}Si)Y_{p}^{Z-14}Y_{n}^{N-14}$$

$$Y_{RG} = \sum_{i\in Lt \ group} N_{i}Y_{i} + \sum_{i\in Si \ group} (N_{i} - 14)Y_{i} + \sum_{i\in Fe \ group} (N_{i} - 28)Y_{i}$$

$$Y_{ZG} = \sum_{i\in Lt \ group} Z_{i}Y_{i} + \sum_{i\in Si \ group} (Z_{i} - 14)Y_{i} + \sum_{i\in Fe \ group} (Z_{i} - 28)Y_{i}$$

$$Y_{SiG} = \sum_{i\in Si \ group} Y_{i},$$

$$Y_{SiG} = \sum_{i\in Si \ group} Y_{i},$$

$$Y_{FeG} = \sum_{i\in Fe \ group} Y_{i}.$$

 $\theta = \left(\frac{m_u k_B T}{2\pi\hbar^2}\right)^{3/2}$ 

binding energy differences, i.e. masses enter directly time evolution for those quantities which are in equilibrium and the individual abundances of nuclei with slow reactions which link equilibrium groups (Hix, Parete-Koon, Freiburghaus, Thielemann 2007)

# Obtaining equilibrium at high T

![](_page_11_Figure_1.jpeg)

at T=4 GK the equilibrium description is correct after about 10<sup>-3</sup> s!

## Incomplete Si-burning with freeze-out

![](_page_12_Figure_1.jpeg)

### Normal and alpha-rich freeze-out

![](_page_13_Figure_1.jpeg)

# Interim conclusions

- steady flows are approached in many hydrostatic burning stages during stellar evolution, including the s-process. They are determined by rates (often the smallest ones), which are/can be related to small Q-values.
- NSE/QSE equilibria are obtained in hydrostatic Si-burning and in explosive burning. Abundance distribution depends directly on mass differences, but for these applications mostly close to stability.
- How about QSE-equilibria linked by steady flows (and far from stability)?

### The classical r-process

- Assume conditions where after a charged-particle freeze-out the heavy QSE-group splits into QSE-subgroups containing each one isotopic chain Z, and a high neutron density is left over
- these QSE-groups are connected by beta-decays from Z to Z+1
- neutrons are consumed to form heavier nuclei
- is a steady flow of beta-decays conceivable?

### s- and r-decoposition

![](_page_16_Figure_1.jpeg)

$$\begin{split} Y(Z,A) &= -\lambda_{\beta^-}(Z,A)Y(Z,A) - \rho N_A < \sigma v >_{n,\gamma} Y_n Y(Z,A) \\ &= -\lambda_{\beta^-}(Z,A)Y(Z,A) - < \sigma v >_{n,\gamma} n_n Y(Z,A) \\ &= -\frac{1}{\tau_{\beta}}Y(Z,A) - \frac{1}{\tau_{n,\gamma}}Y(Z,A). \end{split}$$

which timescale is shorter? neutron capture inverse proportional to  $n_n$  !

Heavy Elements are made by slow and rapid neutron capture events

# High neutron densities lead to nuclei far from stability

![](_page_17_Figure_1.jpeg)

Nuclear Reactions to be considered:  $(n, \gamma)$ ,  $(\gamma, n)$  $(\beta, xn)$ ,  $(\beta, f)$ , (n, f), inelastic  $\nu$ -scattering,  $(\nu_e, e^-)$  ....

### The classical r-process

How to predict abundance changes?

- $\dot{Y}(Z,A) = \sum \lambda_{Z',A'} Y_{Z',A'} + \sum \rho N_A < \sigma v >_{Z',A'} Y_{Z',A'} Y_n$ with  $n_n = \rho N_A Y_n$
- $\dot{Y}(Z,A) \approx \lambda_{\gamma}(Z,A+1)Y(Z,A+1) \langle \sigma v \rangle_{Z,A} Y_{Z,A}n_n$  in case  $(n,\gamma)$ ,  $(\gamma,n)$  rates dominate
- $\dot{Y}(Z,A) = 0$  in chemical equilibrium, •  $Y(Z,A+1)/Y(Z,A) = f(n_n,T,S_n)$  due to detailed balance relation between  $\lambda_{\gamma}(Z,A+1)$  and  $\langle \sigma v \rangle_{Z,A}$

• abundance maxima for all Z's at same  $S_n$ 

$$\begin{split} \frac{Y(Z,A+1)}{Y(Z,A)} &\neq \frac{\langle \sigma v \rangle_{n,\gamma} (A)}{\lambda_{\gamma,n}(A+1)} n_n \qquad \frac{2G(Z,A)}{G(Z,A+1)} [\frac{A}{A+1}]^{3/2} [\frac{m_u kT}{2\pi\hbar^2}]^{3/2} \langle \sigma v \rangle_{n,\gamma} (A) \exp(-S_n(A+1)/kT) \\ &\qquad \frac{Y(Z,A+1)}{Y(Z,A)} = n_n \frac{G(Z,A+1)}{2G(Z,A)} \left[\frac{A+1}{A}\right]^{3/2} \left[\frac{2\pi\hbar^2}{m_u kT}\right]^{3/2} \exp(S_n(A+1)/kT) \end{split}$$

### classical calculation with $n_n$ =const and T=const

![](_page_19_Figure_1.jpeg)

## A=80 and 195 peaks

![](_page_20_Figure_1.jpeg)

three components produce the A=80, 130, and 195 peaks during "comparable" timescales (for the first time experimental half-lives and masses are known in the r-process path at A=80 and 130)!

### Following three $S_n$ 's for timescales $t_1, t_2, t_3$ Kratz, Bitouzet, Thielemann, Möller, Pfeiffer and permutations 1993-1999

![](_page_21_Figure_1.jpeg)

![](_page_22_Figure_0.jpeg)

# Multi-components and steady beta-flow

![](_page_23_Figure_1.jpeg)

(where long half-lives are encountered).

# Explosive H/He-burning and the onset of the rp-process

![](_page_24_Figure_1.jpeg)

detailed modeling is more complex and identifies reaction sensitivities and drip-line dependence (Fisker, FKT, Wiescher 2005, Fisker, Schatz, FKT 2008)

## Test calculations as functions of T

![](_page_25_Figure_1.jpeg)

early evaluations: van Wormer et al. (1994) T=1.5 10<sup>8</sup> K hot CNO cycle no break-out yet

for T=2-3 10<sup>8</sup> K more CNO-type cycles develop up to Ca

Uncertainties in:  ${}^{14}O(\alpha, p){}^{17}F, {}^{17}F(p, \gamma){}^{18}Ne,$   ${}^{18}Ne(\alpha, p){}^{21}Na$  ${}^{15}O(\alpha, \gamma){}^{19}Ne, {}^{19}Ne(p, \gamma){}^{20}Na$ 

### $T = 4 \times 10^8 K$

![](_page_26_Figure_1.jpeg)

### $T = 6 \ge 10^8 K$

![](_page_27_Figure_1.jpeg)

### $T = 8 \ge 10^8 K$

![](_page_28_Figure_1.jpeg)

### $T = 1.5 \ 10^9 K$

![](_page_29_Figure_1.jpeg)

# Break-out from hot CNO at 4-5 10<sup>8</sup> K

Wiescher, Görres, Thielemann, van Wormer, Schatz, Rembges ..

![](_page_30_Figure_2.jpeg)

at Z above 20 a sequence of rapid proton captures and beta-decays (rp-process), for smaller target charges (Coulomb barriers) (a,p) reactions possible

#### rp-process (encounters p-drip line, endpoint Sb-Te cycle)

![](_page_31_Figure_1.jpeg)

Si

9

10 11 12 13

Al Mg 15 16

14

develops features of **QSE-groups along isotonic lines** with links via beta-decays (waiting points) and alpha-induced reactions (tested for energy generation by Rembges et al. 1997)