# Simulating the Transition from the First Stars to the Second Stars

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## The First Stars: a well posed problem

#### • Initial conditions courtesy of The Big Bang





Thanks, WMAP.

## The First Stars: a well posed problem

- Gas Chemistry: Big Bang Nucleosynthesis
  - H, He, and a little D and Li (metal-free)
  - D and Li relatively unimportant
  - Radiative cooling
    - $T > 10^4$  K: atomic H, He
    - $T < 10^4 K H_2$

- Formation controlled by H<sub>2</sub> physics (Bromm et al. 2002; Abel et al. 2002)
  - H<sub>2</sub> only low-temperature coolant
  - lowest rot. trans. ~512 K  $\rightarrow$  T<sub>min</sub> ~ 200 K
  - states in LTE at n ~  $10^4$  cm<sup>-3</sup>
    - NLTE:  $\Lambda \propto n^2$  (efficient)
    - LTE:  $\Lambda \propto n$  (inefficient)









# Key Questions:

- How does the addition of metals alter the star-formation process?
- What chemical abundance is required to form the first low-mass stars?
- How rapid was the transition from Pop III to Pop II?
- What was the IMF of the first generation of Pop II stars? Is the IMF metallicity dependent?

## The First Metals

- Cooling
  - fine-structure
  - molecules
  - dust (also helps form H<sub>2</sub>)



## The Transition to Pop II



## The Transition to Pop II



#### Simulations with Enzo (Bryan & Norman 1997, O'Shea et al. 2004)

- Pre-enriched first star simulations
  - identical cosmological initial conditions (300h<sup>-1</sup> kpc box)
  - $Z = 0, 10^{-6}, 10^{-5}, 10^{-4}, 10^{-3.5}, 10^{-3} Z_{\odot}$
  - solar abundance patterns
  - non-eq H/He chemistry + tabulated metal cooling from Cloudy (all metals through Zn) (BDS, Sigurdsson, & Abel 2008)

#### Simulations with Enzo (Bryan & Norman 1997, O'Shea et al. 2004)

- refinement:
  - baryon/DM overdensity: 4/8
  - jeans length: 16 cells
  - time-step  $< t_{cool}$
- stop after 24 levels of refinement (10<sup>10</sup> dynamic range)

• z<sub>col</sub> ~ 15























#### $Z = 10^{-3} Z_{\odot}$ without CMB!











# Three Modes of Star Formation

- Z < Z<sub>cr</sub>: primordial (M<sub>char</sub> = few x 10<sup>3</sup> M<sub>☉</sub>) cooling cannot prevent loitering phase, collapse proceeds like metal-free case.
- $Z_{cr} \leq Z \leq Z_{CMB}$ : metallicity-regulated (M<sub>char</sub> = few M<sub>o</sub>) cools past loitering phase, does not reach T<sub>CMB</sub>.
- $Z \ge Z_{CMB}$ : CMB-regulated ( $M_{char} = \text{few x } 10^2 M_{\odot}$ ) cools rapidly to  $T_{CMB}$  where frag. stops.



### Caveats

- $M_{char} \neq M_{star}$ : stellar mass depends on accretion and feedback
- rotation: less rotation  $\rightarrow$  less fragmentation
- SN ejecta will not have solar pattern
- metal mixing important
- dust not included
- radiation field unknown



• the real ICs? - need Pop III supernovae



# Summary



- fragmentation suppressed by CMB
- 2 threshold metallicities: Z<sub>cr</sub> and Z<sub>CMB</sub>
- 3 modes of star formation: primordial, metallicity-regulated, and CMB regulated
- star formation modes time dependent top heavy IMF in distant past
- when was star formation 'normal?'
  - GMCs:T ~ 10 K (easy answer)
  - $T_{CMB} = 10 \text{ K}$  at  $z \sim 2.6 (2.6 \text{ Gyr after BB})$

# Metal Cooling: The Method

- Before the simulation:
  - assume ionization equilibrium
  - pre-compute cooling rates over n,T,Z, etc.
  - subtract H/He cooling: only metals left
- During the simulation:
  - follow H/He chemistry: H, H<sup>+</sup>, H<sup>-</sup>, H<sub>2</sub>, H<sub>2</sub><sup>+</sup>, He, He<sup>+</sup>, He<sup>++</sup>, e<sup>-</sup>
  - calculate H/He cooling directly (H<sub>2</sub>!)
  - interpolate metal cooling from data-grid