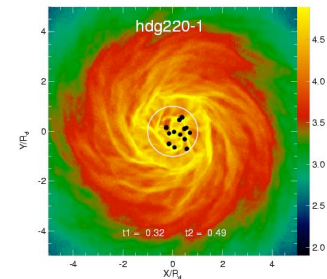
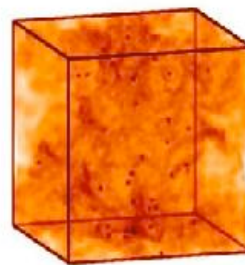
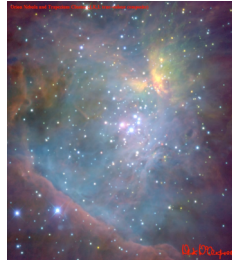
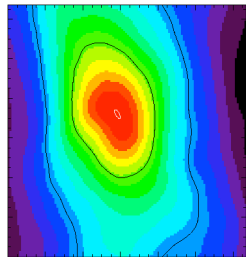
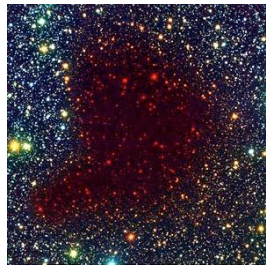


GRAVOTURBULENT STAR FORMATION with Smoothed Particle Hydrodynamics



Ralf Klessen

Emmy Noether Research Group (DFG)
Astrophysikalisches Institut Potsdam



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Collaborators

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Overview

1. Physics of star formation

- a) How do stars form?
- b) Where do stars form?
- c) Theory of *gravoturbulent star formation*

2. Numerical approach to star formation

- 1. Large-eddy simulations (*LES*) with smoothed particle hydrodynamics (*SPH*)
- 2. Transition to stellar-dynamics: introducing „*sink particles*“ to represent protostars (i.e. to describe *subgrid-scale physics*)

STAR FORMATION

Star formation in "typical" spiral:

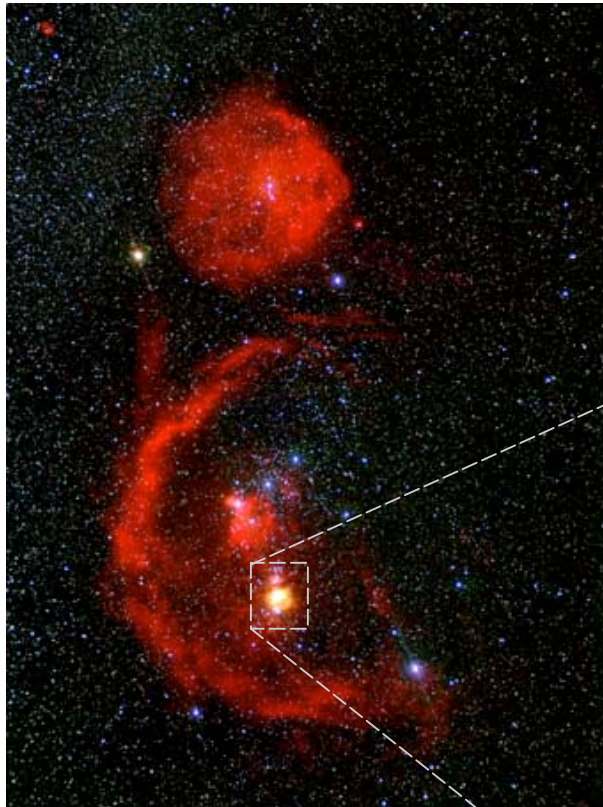


(from the Hubble Heritage Team)

NGC4622

- Star formation *always* is associated with *clouds of gas and dust*.
- Star formation is essentially a *local phenomenon* (on ~pc scale)
- **HOW** is star formation is *influenced* by *global* properties of the galaxy?

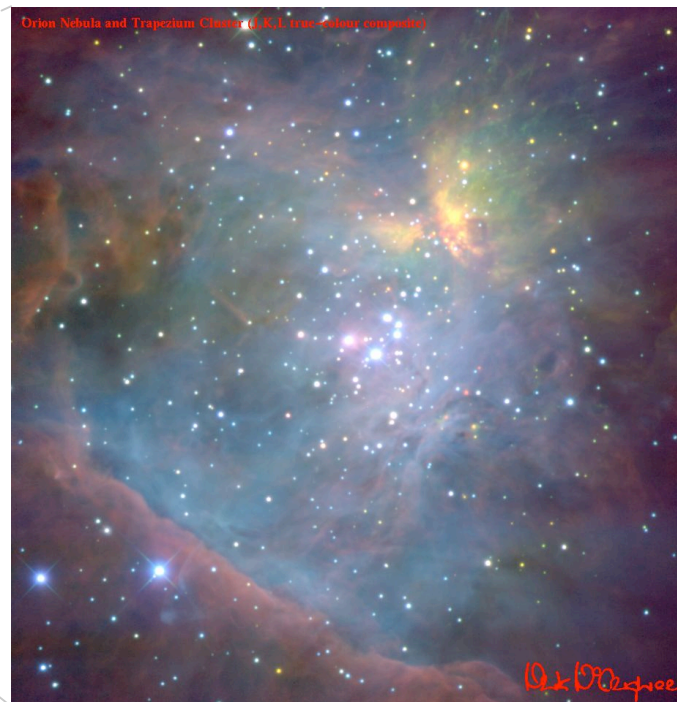
Local star forming region: The Trapezium Cluster in Orion



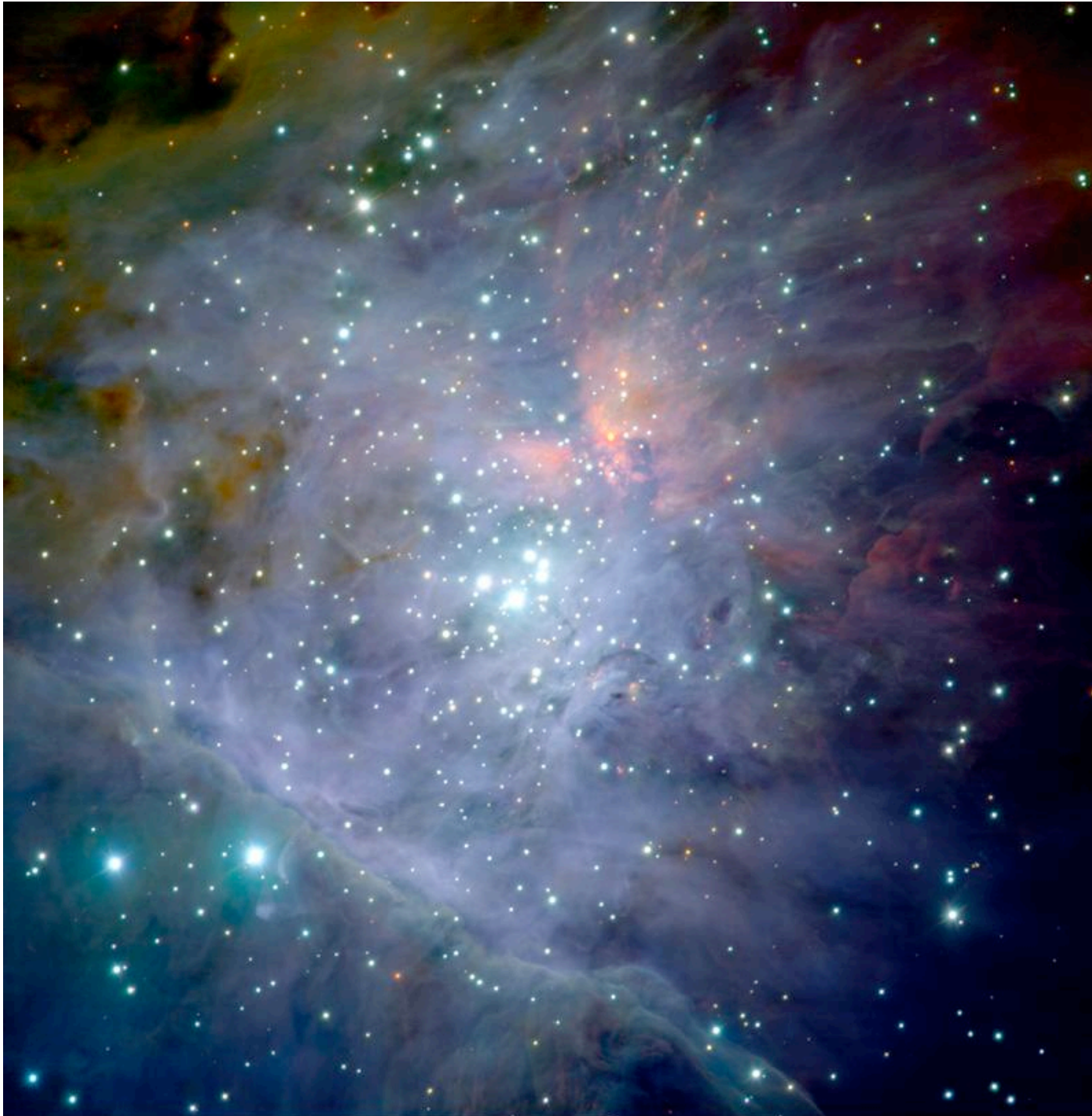
Orion molecular cloud

The Orion molecular cloud is the birth- place of several young embedded star clusters.

The Trapezium cluster is only visible in the IR and contains about 2000 newly born stars.



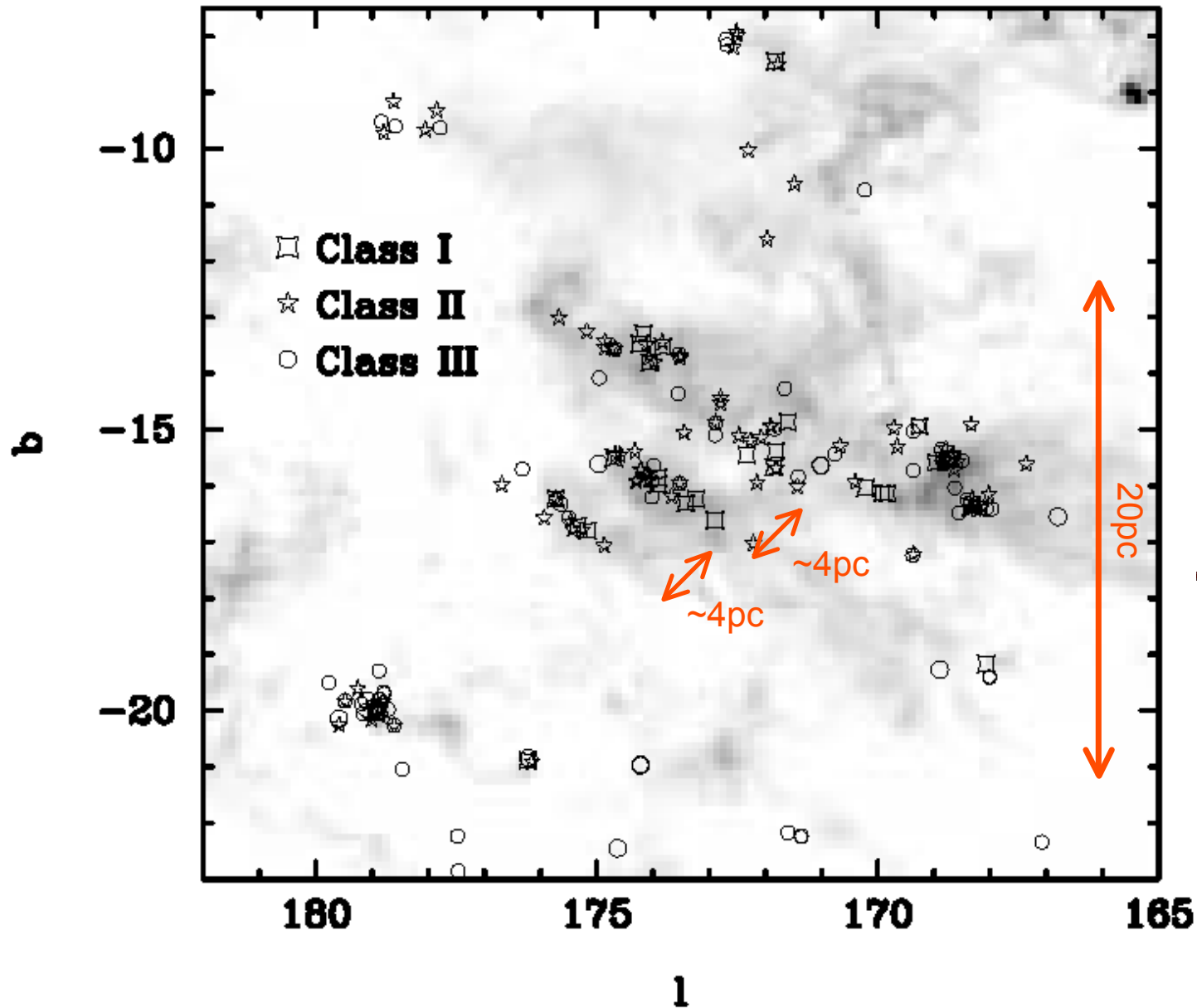
Trapezium cluster



Trapezium Cluster (detail)

- stars form in **clusters**
- stars form in **molecular clouds**
- (proto)stellar **feedback** is important

(color composite J,H,K
by M. McCaughrean,
VLT, Paranal, Chile)



Taurus molecular cloud

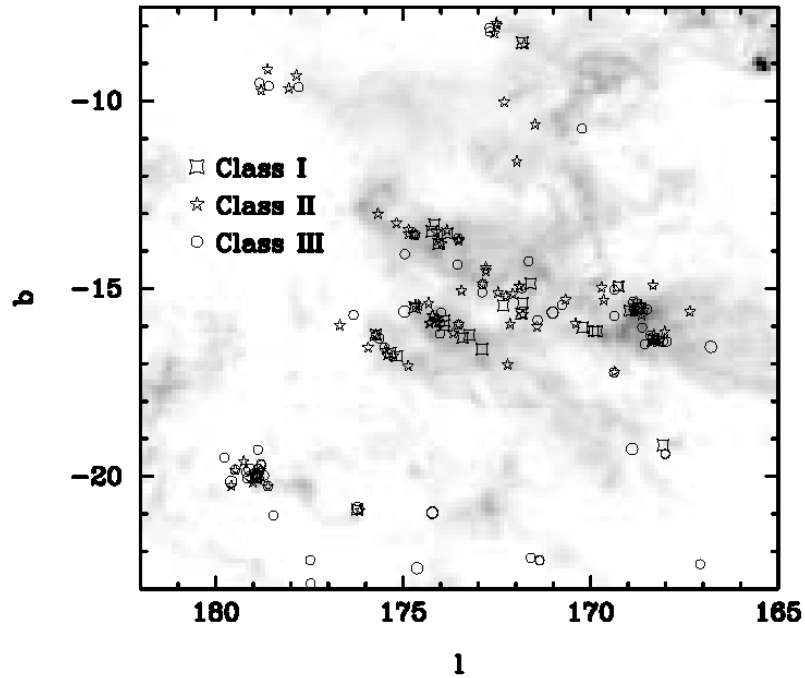
star-forming
filaments in the
Taurus cloud

- Structure and dynamics of young star clusters is coupled to *structure of mol. cloud*

Taurus molecular cloud

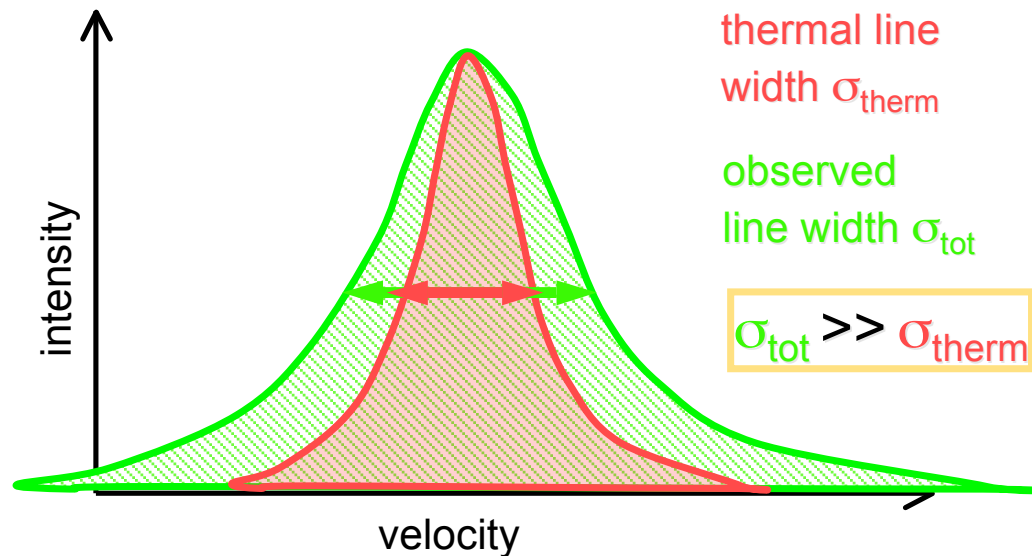
Star-forming filaments in *Taurus* cloud

(from Hartmann 2002)



- Structure and dynamics of young star clusters is coupled to *structure of molecular cloud*

- Structure and dynamics of *molecular cloud* is determined by *supersonic turbulence*



The star formation process

- *How* do stars form?
 - What determines *when* and *where* stars form?
 - What *regulates* the process and determines its *efficiency*?
 - How do *global* properties of the galaxy influence star formation (a *local* process)?
 - Are there different *modes* of SF?
(Starburst galaxies vs. *LSBs*, *isolated* SF vs. *clustered* SF)
- *What physical processes initiate and control the formation of stars?*

Gravoturbulent star formation

- New theory of star formation:

*Star formation is controlled
by interplay between
gravity and
supersonic turbulence!*

- Dual role of turbulence:
 - *stability on large scales*
 - *initiating collapse on small scales*

Gravoturbulent star formation

- New theory of star formation:

*Star formation is controlled
by interplay between
gravity and
supersonic turbulence!*

- Validity:

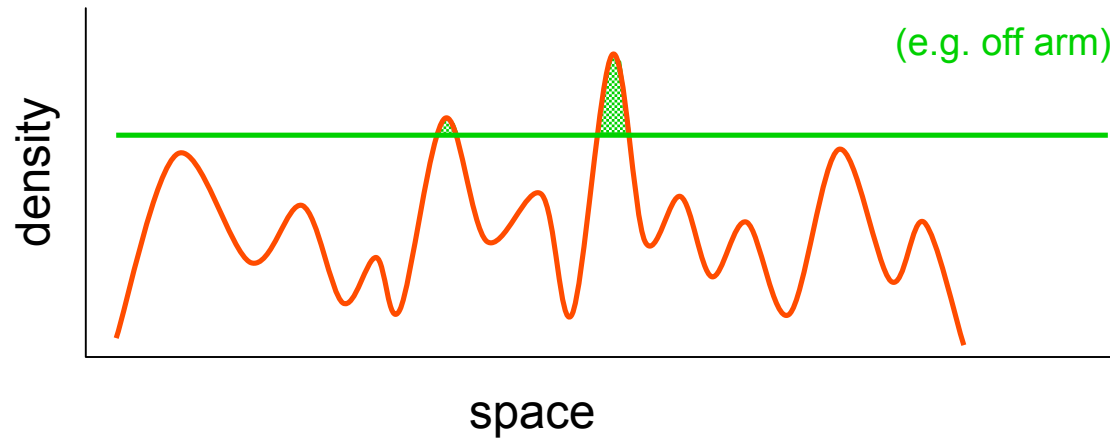
This hold on *all* scales and applies to build-up of stars and star clusters within molecular clouds as well as to the formation of molecular clouds in galactic disk.

Gravoturbulent Star Formation

- *Supersonic turbulence* in the galactic disk creates strong **density fluctuations** (in shocks: $\delta\rho/\rho \approx \mathcal{M}^2$)
 - chemical phase transition: atomic \rightarrow molecular
 - cooling instability
 - gravitational instability
- Cold *molecular clouds* form at the high-density peaks.
- *Turbulence* creates density structure, *gravity* selects for collapse
—————→ **GRAVOTUBULENT FRAGMENTATION**

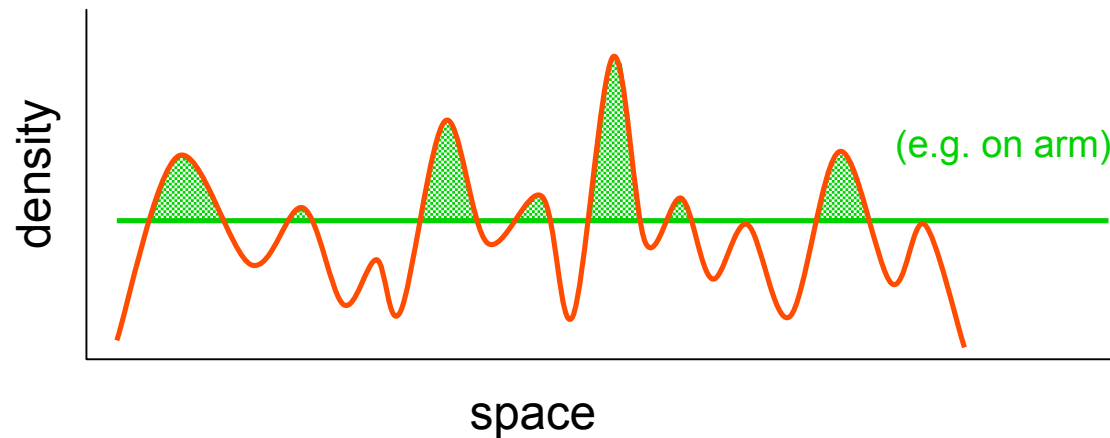
- *Turbulent cascade*: Local compression *within* a cloud provokes collapse \rightarrow individual *stars* and *star clusters*

Star formation on *global scales*



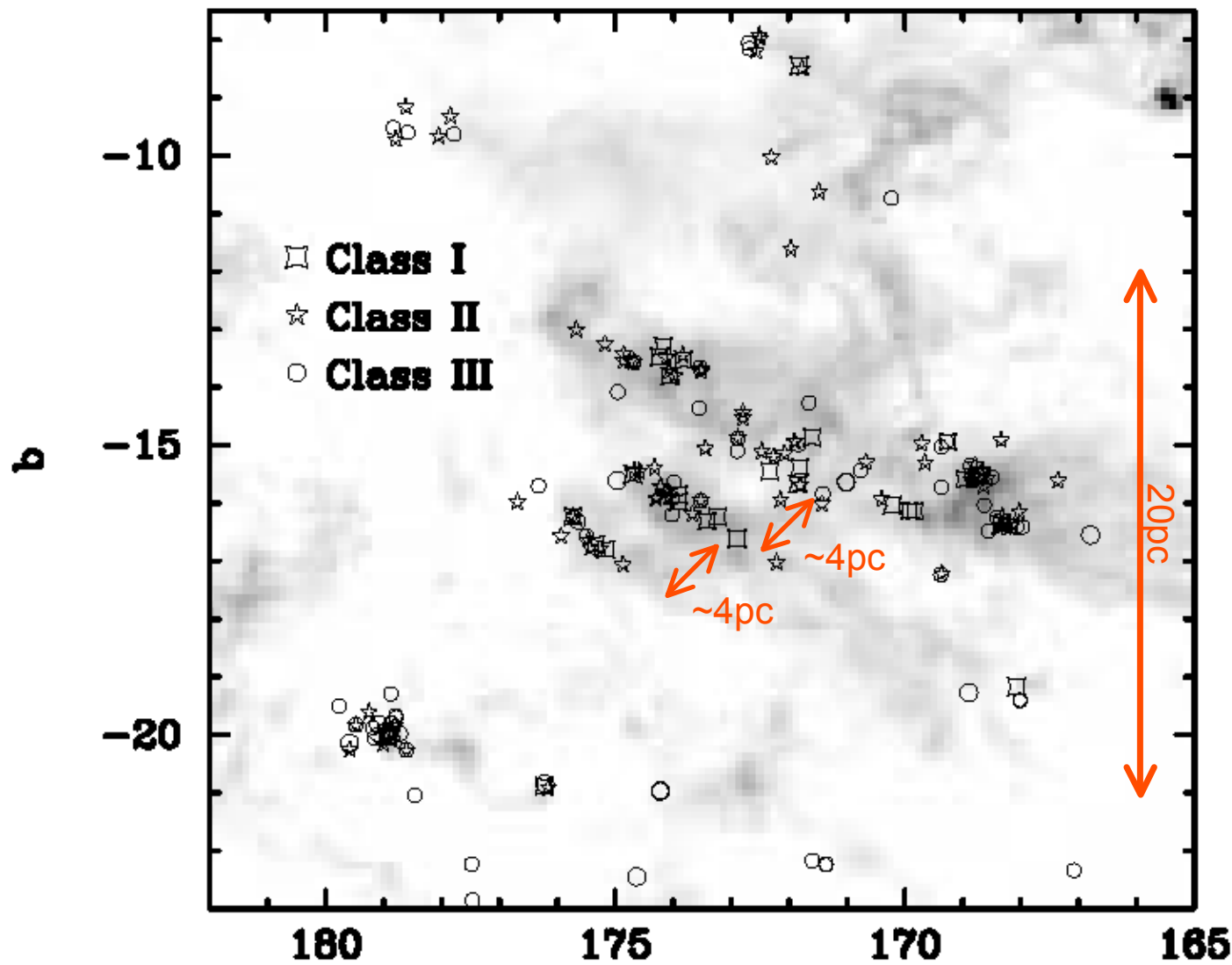
density fluctuations in warm atomic ISM caused by supersonic turbulence

some are dense enough to form H₂ within “reasonable timescale”
→ molecular clouds



external perturbations (i.e. potential changes) increase likelihood

Approaching
the problem

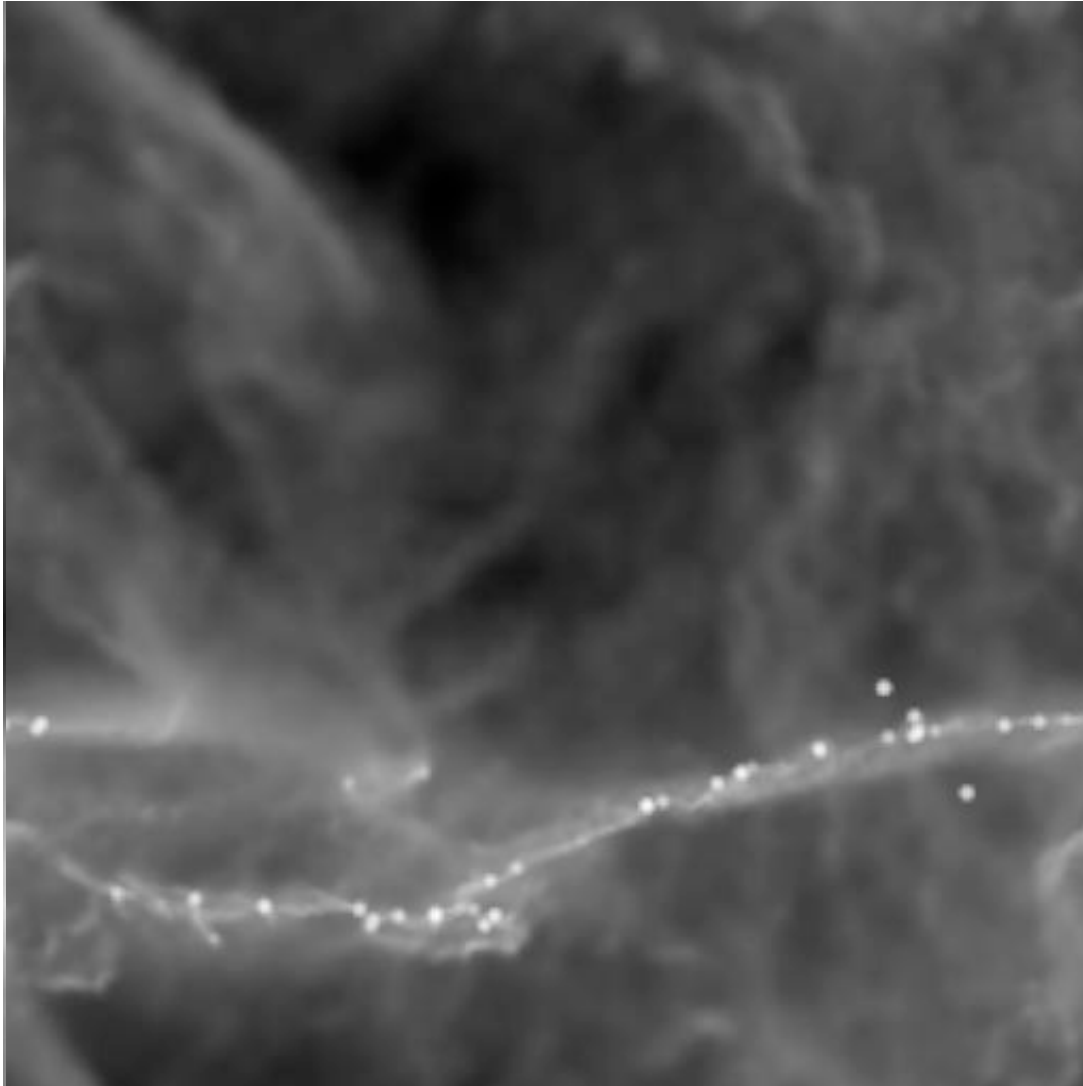


Taurus SF cloud

Star-
forming
filaments in
the *Taurus*
molecular
cloud

(from Hartmann 2002, ApJ)

Gravoturbulent fragmentation



Gravoturbulent fragmentation in molecular clouds:

- SPH model with 1.6×10^6 particles
- large-scale driven turbulence
- Mach number $\mathcal{M} = 6$
- periodic boundaries
- physical scaling:

“Taurus”:

- density $n(\text{H}_2) \approx 10^2 \text{ cm}^{-3}$:
- $L = 6 \text{ pc}$, $M = 5000 M_\odot$

What can we learn from that?

- *global properties* (statistical properties)
 - SF efficiency
 - SF time scale
 - IMF
 - description of self-gravitating turbulent systems (pdf's, Δ -var.)
 - **chemical mixing properties**
- *local properties* (properties of individual objects)
 - properties of individual clumps (e.g. shape, radial profile)
 - accretion history of individual protostars (dM/dt vs. t , j vs. t)
 - binary (proto)stars (eccentricity, mass ratio, etc.)
 - SED's of individual protostars
 - dynamic PMS tracks: $T_{\text{bol}}-L_{\text{bol}}$ evolution

Turbulent diffusion I

- Observations of young star clusters exhibit an enormous degree of chemical homogeneity (e.g. in the Pleiades: Wilden et al. 2002)
- Star-forming gas must be well mixed.
- How does this constrain models of interstellar turbulence?
- → Study mixing in supersonic compressible turbulence.....

Turbulent diffusion II

- Method:

- second central moment of displacement:

$$\xi_{\vec{r}}^2(t - t') = \left\langle [\vec{r}_i(t) - \vec{r}_i(t')]^2 \right\rangle_i$$

- classical diffusion equation:

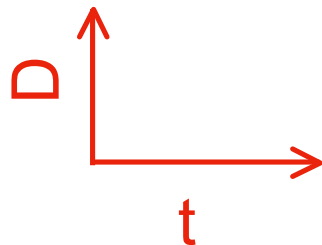
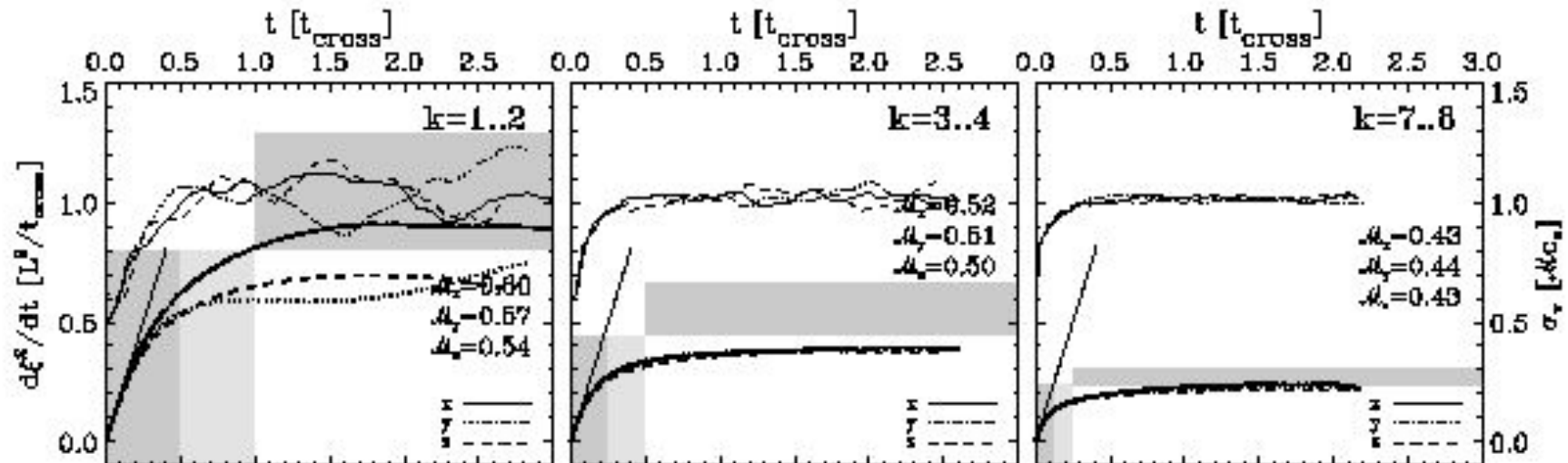
$$\frac{dn}{dt} = D \vec{\nabla}^2 n$$

- relation between D and ξ :

$$D(t - t') = \frac{d\xi_{\vec{r}}^2(t - t')}{dt} = 2 \left\langle \vec{v}_i(t - t') \cdot \vec{r}_i(t - t') \right\rangle_i$$

Turbulent diffusion III

- Time evolution of diffusion coefficient
(mean motion corrected).



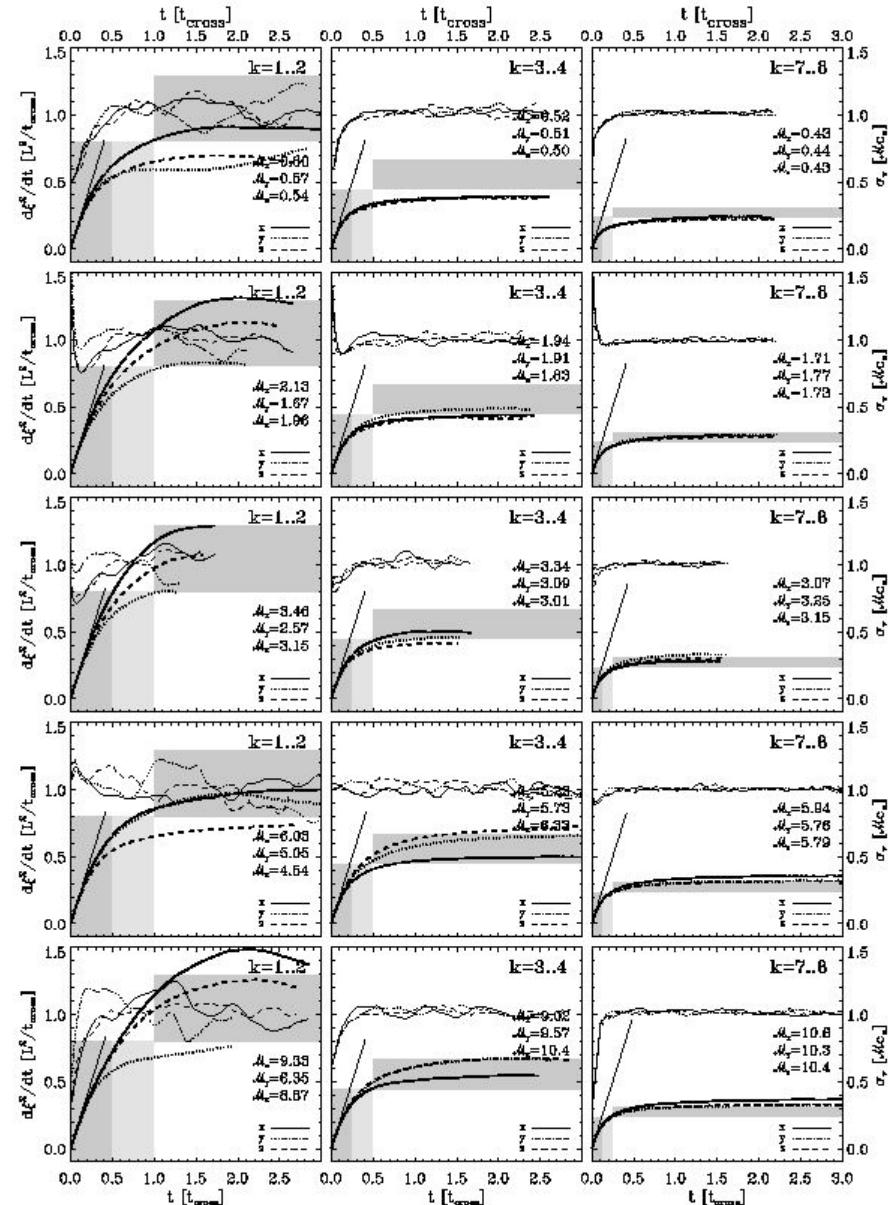
Turbulent diffusion IV

- Mean-motion corrected diffusion
- Simple mixing-length approach works!

- $D(t) \approx v_{\text{rms}}^2 t \quad t < \tau$

- $D(t) \approx v_{\text{rms}}^2 \tau$
 $= v_{\text{rms}} \ell \quad t > \tau$

- With v_{rms} = rms velocity and $\ell = L/k =$ shock sep.

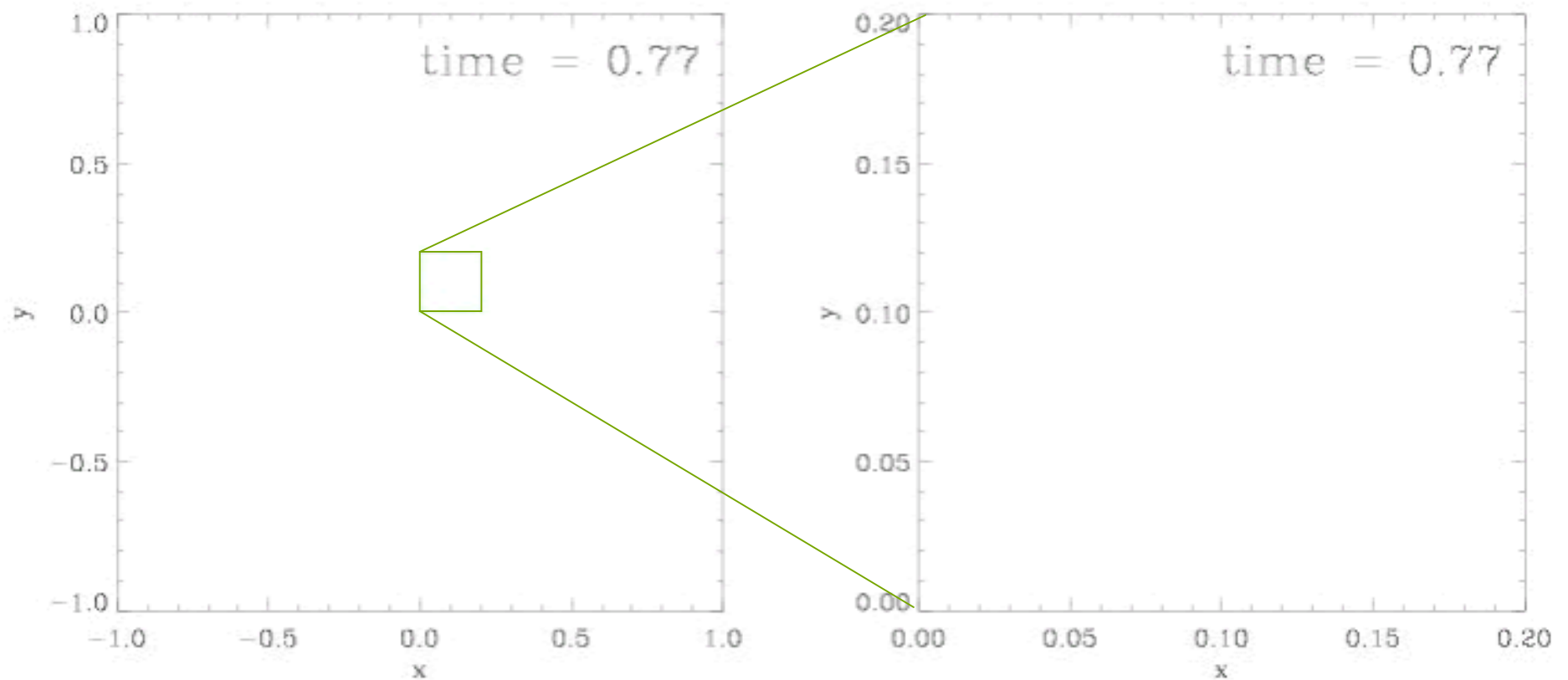


What can we learn from that?

- *global properties* (statistical properties)
 - SF efficiency and timescale
 - stellar mass function -- IMF
 - dynamics of young star clusters
 - description of self-gravitating turbulent systems (pdf's, Δ -var.)
 - chemical mixing properties
- *local properties* (properties of individual objects)
 - properties of individual clumps (e.g. shape, radial profile)
 - accretion history of individual protostars (dM/dt vs. t , j vs. t)
 - binary (proto)stars (eccentricity, mass ratio, etc.)
 - SED's of individual protostars
 - dynamic PMS tracks: $T_{\text{bol}}-L_{\text{bol}}$ evolution

Star cluster formation

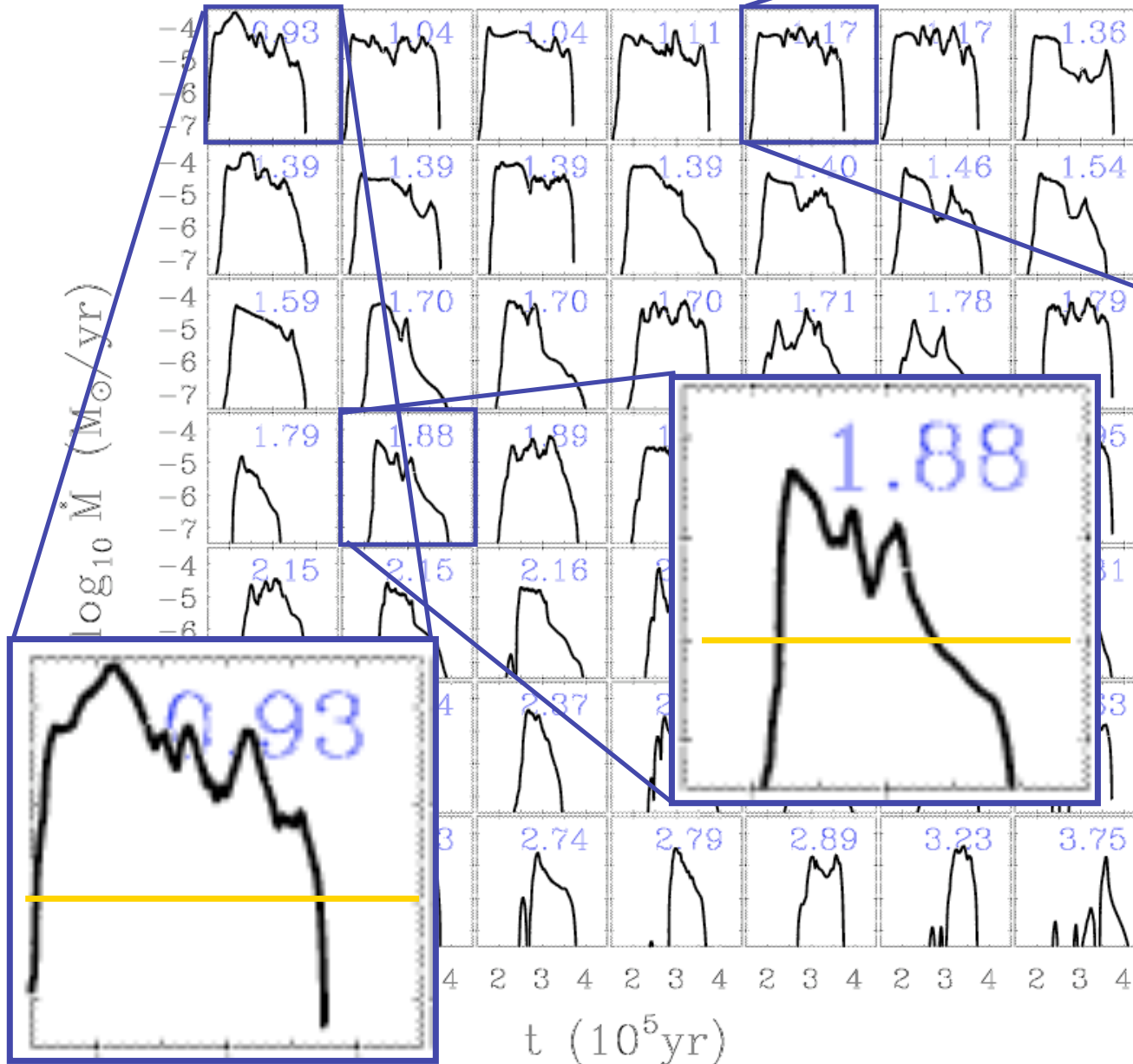
Most stars form in clusters → *star formation = cluster formation*



Trajectories of protostars in a nascent dense cluster created by gravoturbulent fragmentation

(from Klessen & Burkert 2000, ApJS, 128, 287)

Accretion rates in clu

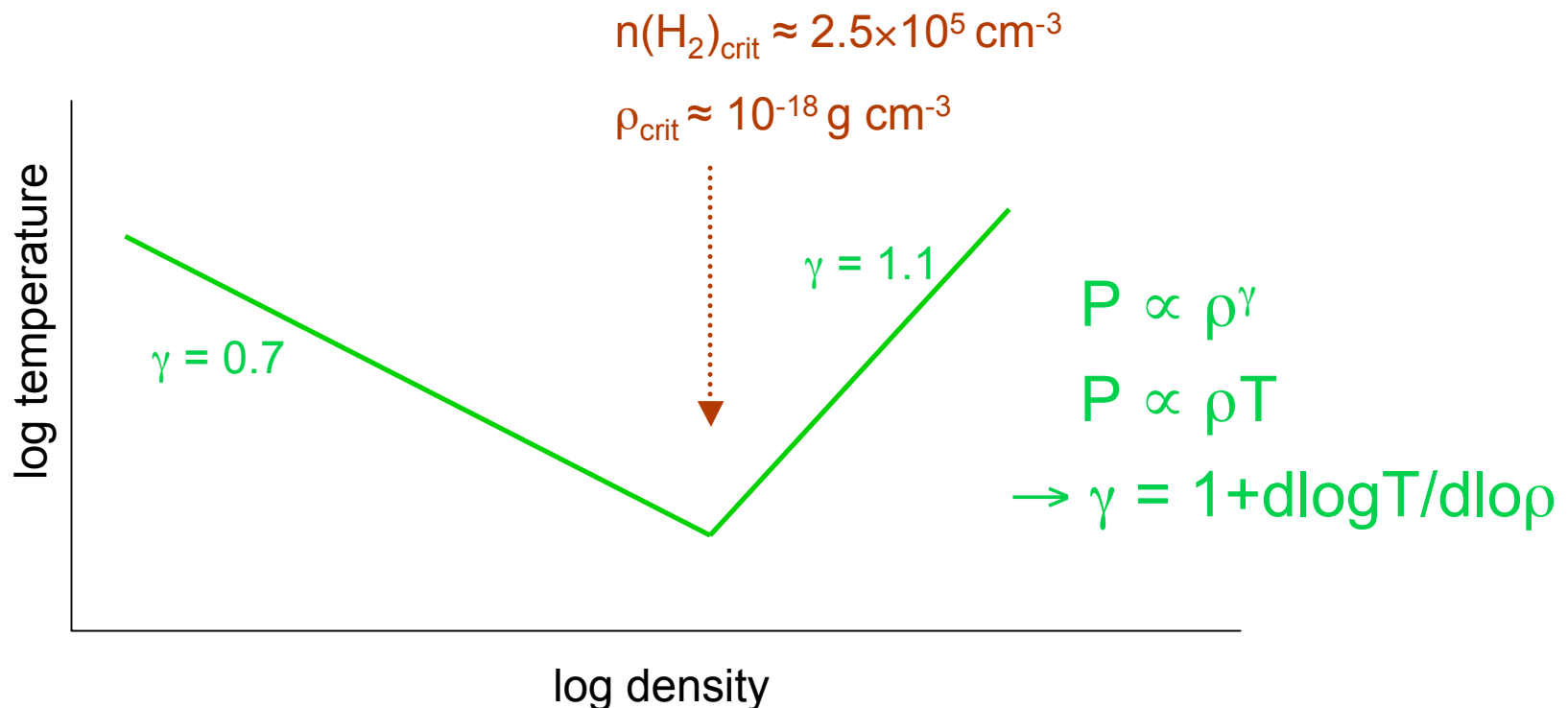


Mass accretion rates *vary with time* and are strongly *influenced* by the *cluster environment*.

(Klessen 2001, ApJ, 550, L77; also Schmeja & Klessen, 2004, A&A, 419, 405)

Influence of EOS

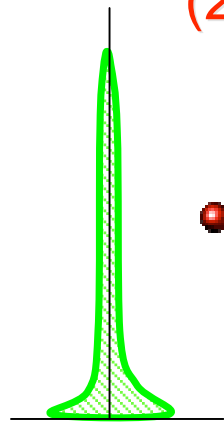
- But EOS depends on *chemical state*, on *balance* between *heating* and *cooling*



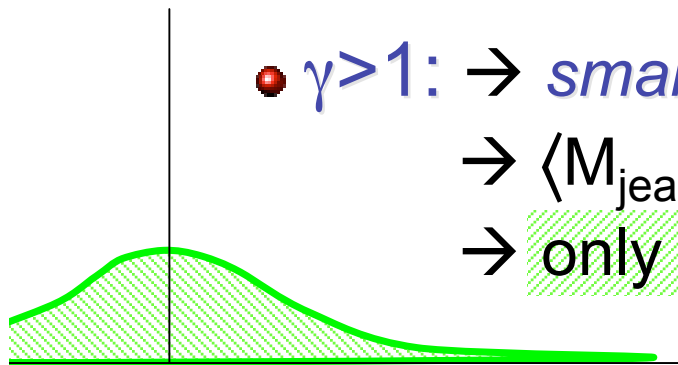
Influence of EOS

$$(1) \quad p \propto \rho^\gamma \quad \rightarrow \quad \rho \propto p^{1/\gamma}$$

$$(2) \quad M_{\text{jeans}} \propto \gamma^{3/2} \rho^{(3\gamma-4)/2}$$



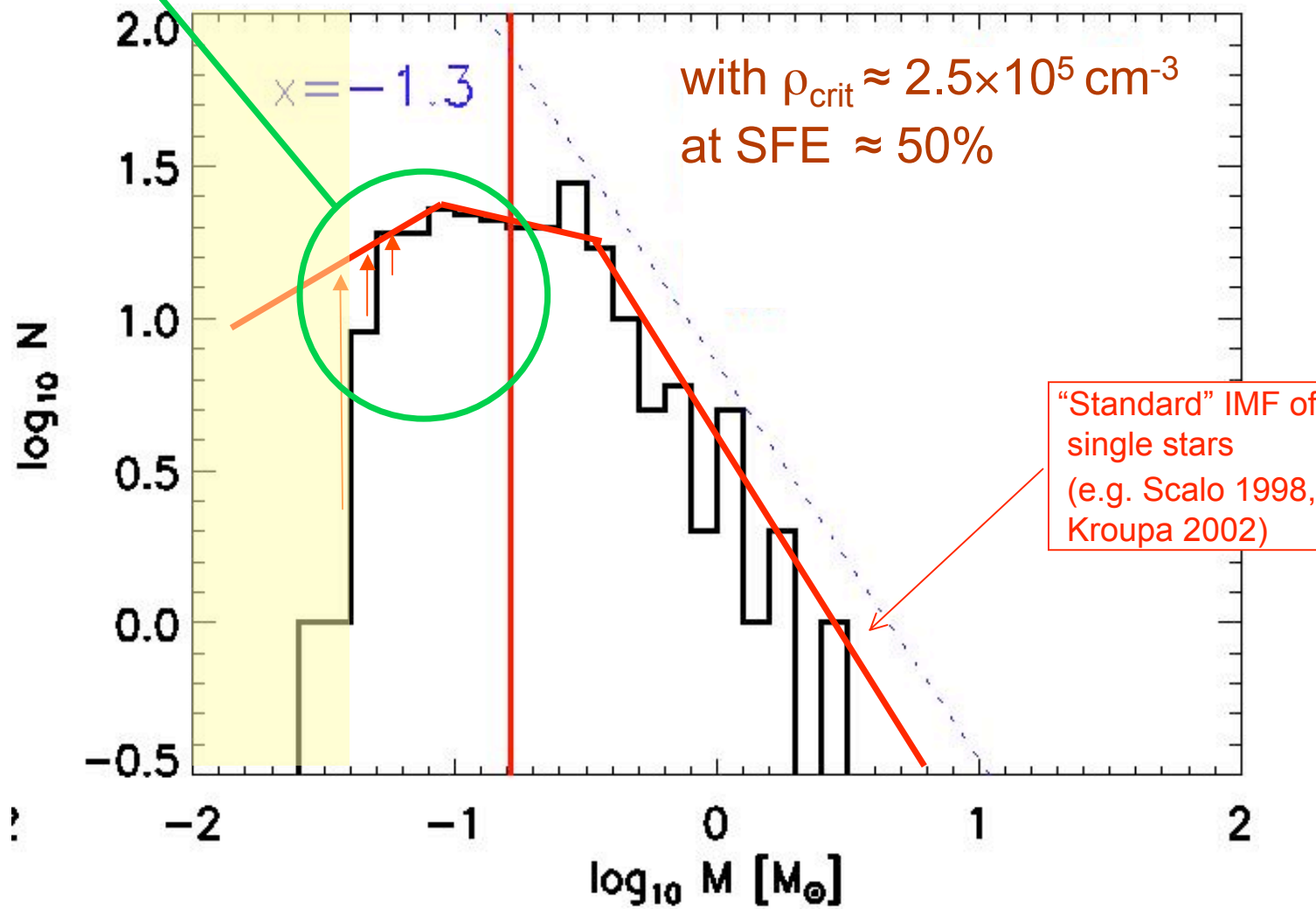
- $\gamma < 1$: \rightarrow *large* density excursion for given pressure
 \rightarrow $\langle M_{\text{jeans}} \rangle$ becomes small
 \rightarrow number of fluctuations with $M > M_{\text{jeans}}$ is large



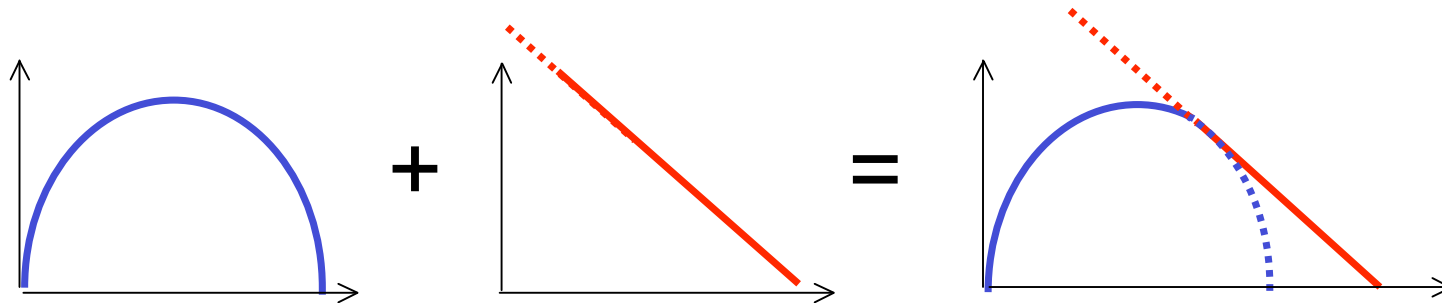
- $\gamma > 1$: \rightarrow *small* density excursion for given pressure
 \rightarrow $\langle M_{\text{jeans}} \rangle$ is large
 \rightarrow only few and massive clumps exceed M_{jeans}

Mass spectrum

sufficient # of brown dwarfs



Plausibility argument for shape



- Supersonic turbulence is scale free process

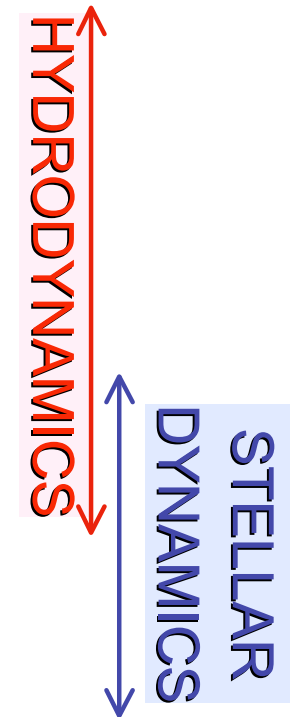
→ *POWER LAW BEHAVIOR*

- *But also:* turbulence and fragmentation are highly stochastic processes → central limit theorem

→ *GAUSSIAN DISTRIBUTION*

IMF: Summary

- To get the stellar mass function (IMF) we need to:
 - describe **supersonic turbulence** (LES)
 - include **self-gravity**
 - model **thermodynamic balance** of the gas (heating, cooling, time-dependent chemistry, EOS)
 - follow formation of **compact collapsed cores** (transition from hydro to stellar dynamics)
 - treat **stellar dynamical processes** (protostellar collisions, ejection by close encounters)



NUMERICS

Goal

- We want to understand the formation of star clusters in turbulent interstellar gas clouds.

--> We want to describe the transition from a hydrodynamic system (the self-gravitating gas cloud) to one that is dominated by (collisional) stellar dynamics (the final star cluster).

- How can we do that?

Numerical approach I

- Problem of star formation is very complex. It involves many scales (10^7 in length, and 10^{20} in density) and many physical processes → NO analytic solution
→ NUMERICAL APPROACH
- BUT, we need to...
 - solve the MHD equations in 3 dimensions
 - solve Poisson's equation (self-gravity)
 - follow the full turbulent cascade (in the ISM + in stellar interior)
 - include heating and cooling processes (EOS)
 - treat radiation transfer
 - describe energy production by nuclear burning processes

Numerical approach II

- Simplify!
Divide problem into little bits and pieces.....
- **GRAVOTURBULENT CLOUD FRAGMENTATION**
- We try to...
 - solve the HD equations in 3 dimensions
 - solve Poisson's equation (self-gravity)
 - include a (humble) approach to supersonic turbulence
 - describe perfect gas (with polytropic EOS)
 - follow collapse: include "sink particles"
(this will "handle" our subgrid-scale physics)

Intermezzo: HD & SPH

- derivation of equations of hydrodynamics
 - Boltzmann equation for 1D distribution function:

$$\frac{df}{dt} \equiv \frac{\partial f}{\partial t} + \dot{\vec{q}} \cdot \vec{\nabla}_{\vec{q}} f_{\vec{p}} + \dot{\vec{p}} \cdot \vec{\nabla}_{\vec{p}} f$$

- moments of distribution function:
density ρ , momentum \vec{p} , energy ε
- SPH: smoothed particle hydrodynamics
 - particle-based scheme to solve eqn.'s of hydrodynamics
 - thermodynamic behavior --> equation of state (EOS)

Hydrodynamics

- gases and fluids are *large* ensembles of interacting particles
- \longrightarrow state of system is described by location in $6N$ dimensional phase space $f^{(N)}(\vec{q}_1 \dots \vec{q}_N, \vec{p}_1 \dots \vec{p}_N) d\vec{q}_1 \dots d\vec{q}_N d\vec{p}_1 \dots d\vec{p}_N$
- time evolution governed by 'equation of motion' for $6N$ -dim probability distribution function $f^{(N)}$
- $f^{(N)} \rightarrow f^{(n)}$ by integrating over all but n coordinates \longrightarrow BBGKY hierarchy of equations of motion (after Born, Bogoliubov, Green, Kirkwood and Yvon)
- physical observables are typically associated with 1- or 2-body probability density $f^{(1)}$ or $f^{(2)}$
- at lowest level of hierarchy: 1-body distribution function describes the probability of finding a particle at time t in the volume element $d\vec{q}$ at \vec{q} with momenta in the range $d\vec{p}$ at \vec{p} .
- **Boltzmann equation** – equation of motion for $f^{(1)}$

$$\begin{aligned} \frac{df}{dt} &\equiv \frac{\partial f}{\partial t} + \dot{\vec{q}} \cdot \vec{\nabla}_{\vec{q}} f + \dot{\vec{p}} \cdot \vec{\nabla}_{\vec{p}} f \\ &= \frac{\partial f}{\partial t} + \vec{v} \cdot \vec{\nabla}_{\vec{q}} f + \vec{F} \cdot \vec{\nabla}_{\vec{p}} f = f_c \end{aligned}$$

- Boltzmann equation

$$\begin{aligned}\frac{df}{dt} &\equiv \frac{\partial f}{\partial t} + \dot{\vec{q}} \cdot \vec{\nabla}_{\vec{q}} f + \dot{\vec{p}} \cdot \vec{\nabla}_{\vec{p}} f \\ &= \frac{\partial f}{\partial t} + \vec{v} \cdot \vec{\nabla}_{\vec{q}} f + \vec{F} \cdot \vec{\nabla}_{\vec{p}} f = f_c\end{aligned}$$

→ first line: transformation from comoving to spatially fixed coordinate system.

→ second line: velocity $\vec{v} = \dot{\vec{q}}$ and force $\vec{F} = \dot{\vec{p}}$

→ all higher order terms are 'hidden' in the collision term f_c

- observable quantities are typically (velocity) moments of the Boltzmann equation, e.g.

→ density:

$$\rho = \int m f(\vec{q}, \vec{p}, t) d\vec{p}$$

→ momentum:

$$\rho \vec{v} = \int m \vec{v} f(\vec{q}, \vec{p}, t) d\vec{p}$$

→ kinetic energy density:

$$\rho \vec{v}^2 = \int m \vec{v}^2 f(\vec{q}, \vec{p}, t) d\vec{p}$$

- in general: the i -th velocity moment $\langle \xi_i \rangle$ (of $\xi_i = m\vec{v}^i$) is

$$\langle \xi_i \rangle = \frac{1}{n} \int \xi_i f(\vec{q}, \vec{p}, t) d\vec{p}$$

with the mean particle number density n defined as

$$n = \int f(\vec{q}, \vec{p}, t) d\vec{p}$$

- the equation of motion for $\langle \xi_i \rangle$ is

$$\int \xi_i \left\{ \frac{\partial f}{\partial t} + \vec{v} \cdot \vec{\nabla}_q f + \vec{F} \cdot \vec{\nabla}_p f \right\} d\vec{p} = \int \xi_i \{f_c\} d\vec{p},$$

which after some complicated rearrangement becomes

$$\frac{\partial}{\partial t} n \langle \xi_i \rangle + \vec{\nabla}_q (n \langle \xi_i \vec{v} \rangle) + n \vec{F} \langle \vec{\nabla}_p \xi_i \rangle = \int \xi_i f_c d\vec{p}$$

(Maxwell-Boltzmann transport equation for $\langle \xi_i \rangle$)

- if the RHS is zero, then ξ_i is a conserved quantity. This is only the case for first three moments, **mass** $\xi_0 = m$, **momentum** $\vec{\xi}_1 = m\vec{v}$, and **kinetic energy** $\xi_2 = m\vec{v}^2/2$.
- MB equations build a hierarically nested set of equations, as $\langle \xi_i \rangle$ depends on $\langle \xi_{i+1} \rangle$ via $\vec{\nabla}_q (n \langle \xi_i \vec{v} \rangle)$ and because the collision term cannot be reduced to depend on ξ_i only.
 - need for a closure equation
 - in hydrodynamics this is typically the equation of state.

assumptions

- **continuum limit:**

- distribution function f must be a 'smoothly' varying function on the scales of interest → local average possible
- stated differently: the averaging scale (i.e. scale of interest) must be larger than the mean free path of individual particles
- stated differently: microscopic behavior of particles can be neglected
- concept of fluid element must be meaningful

- **only 'short range forces':**

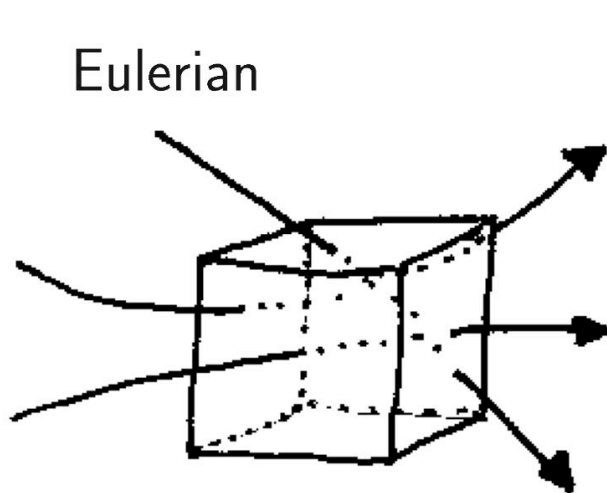
- forces between particles are short range or saturate → collective effects can be neglected
- stated differently: correlation length of particles in the system is finite (and smaller than the scales of interest)

limitations

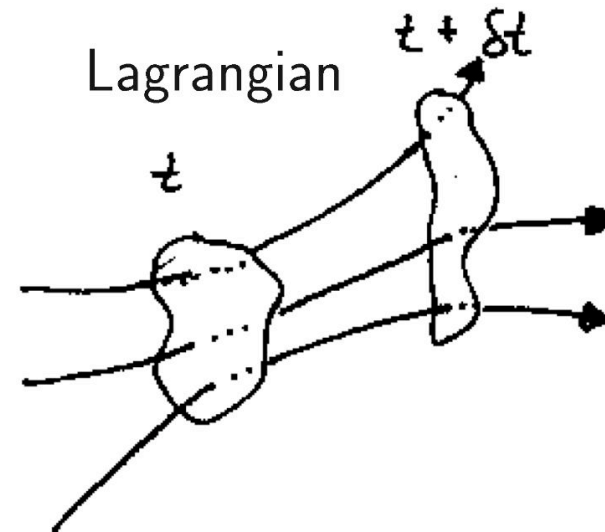
- shocks (scales of interest become smaller than mean free path)
- phase transitions (correlation length may become infinite)
- description of self-gravitating systems
- description of fully fractal systems

the equations of hydrodynamics

- hydrodynamics \equiv book keeping problem
One must keep track of the 'change' of a fluid element due to various physical processes acting on it. How do its 'properties' evolve under the influence of compression, heat sources, cooling, etc.?
- Eulerian vs. Lagrangian point of view



consider spatially fixed volume element



following motion of fluid element

- hydrodynamic equations = set of equations for the five conserved quantities ($\rho, \rho\vec{v}, \rho\vec{v}^2/2$) plus closure equation (plus transport equations for 'external' forces if present, e.g. gravity, magnetic field, heat sources, etc.)

- equations of hydrodynamics

$$\frac{d\rho}{dt} = \frac{\partial\rho}{\partial t} + \vec{v} \cdot \vec{\nabla}\rho = -\rho\vec{\nabla} \cdot \vec{v} \quad (\text{continuity equation})$$

$$\frac{d\vec{v}}{dt} = \frac{\partial\vec{v}}{\partial t} + (\vec{v} \cdot \vec{\nabla})\vec{v} = -\frac{1}{\rho}\vec{\nabla}p - \vec{\nabla}\phi + \eta\vec{\nabla}^2\vec{v} + \left(\zeta + \frac{\eta}{3}\right)\vec{\nabla}(\vec{\nabla} \cdot \vec{v})$$

(Navier-Stokes equation)

$$\frac{d\epsilon}{dt} = \frac{\partial\epsilon}{\partial t} + \vec{v} \cdot \vec{\nabla}\epsilon = T\frac{ds}{dt} - \frac{p}{\rho}\vec{\nabla} \cdot \vec{v} \quad (\text{energy equation})$$

$$\vec{\nabla}^2\phi = 4\pi G\rho \quad (\text{Poisson's equation})$$

$$p = \mathcal{R}\rho T \quad (\text{equation of state})$$

$$\vec{F}_B = -\vec{\nabla} \frac{\vec{B}^2}{8\pi} + \frac{1}{4\pi} (\vec{B} \cdot \vec{\nabla}) \vec{B} \quad (\text{magnetic force})$$

$$\frac{\partial \vec{B}}{\partial t} = \vec{\nabla} \times (\vec{v} \times \vec{B}) \quad (\text{Lorentz equation})$$

ρ = density, \vec{v} = velocity, p = pressure, ϕ = gravitational potential, ζ and η viscosity coefficients, $\epsilon = \rho \vec{v}^2 / 2$ = kinetic energy density, T = temperature, s = entropy, \mathcal{R} = gas constant, \vec{B} = magnetic field (cgs units)

- mass transport – continuity equation

$$\frac{d\rho}{dt} = \frac{\partial\rho}{\partial t} + \vec{v} \cdot \vec{\nabla}\rho = -\rho\vec{\nabla} \cdot \vec{v}$$

(conservation of mass)

- transport equation for momentum – Navier Stokes equation

$$\frac{d\vec{v}}{dt} = \frac{\partial \vec{v}}{\partial t} + (\vec{v} \cdot \vec{\nabla}) \vec{v} = -\frac{1}{\rho} \vec{\nabla} p - \vec{\nabla} \phi + \eta \vec{\nabla}^2 \vec{v} + \left(\zeta + \frac{\eta}{3} \right) \vec{\nabla} (\vec{\nabla} \cdot \vec{v})$$

momentum change due to

→ pressure gradients: $(-\rho^{-1} \vec{\nabla} p)$

→ (self) gravity: $-\vec{\nabla} \phi$

→ viscous forces (internal friction, contains $\text{div}(\partial v_i / \partial x_j)$ terms):
 $\eta \vec{\nabla}^2 \vec{v} + \left(\zeta + \frac{\eta}{3} \right) \vec{\nabla} (\vec{\nabla} \cdot \vec{v})$

(conservation of momentum, general form of momentum transport: $\partial_t(\rho v_i) = -\partial_j \Pi_{ij}$)

- transport equation for internal energy

$$\frac{d\epsilon}{dt} = \frac{\partial \epsilon}{\partial t} + \vec{v} \cdot \vec{\nabla} \epsilon = T \frac{ds}{dt} - \frac{p}{\rho} \vec{\nabla} \cdot \vec{v}$$

- follows from the thermodynamic relation $d\epsilon = T ds - p dV = T ds + p/\rho^2 d\rho$ which describes changes in ϵ due to entropy changes and to volume changes (compression, expansion)
- for adiabatic gas the first term vanishes ($s = \text{constant}$)
- heating sources, cooling processes can be incorporated in ds (conservation of energy)

- closure equation – equation of state
 - general form of equation of state $p = p(T, \rho, \dots)$
 - ideal gas: $p = \mathcal{R}\rho T$
 - special case – isothermal gas: $p = c_s^2 T$ (as $\mathcal{R}T = c_s^2$)

Note:

- in reality, computing the EOS is VERY complex!
- depends on detailed *balance* between *heating* and *cooling*
- these depend on *chemical composition* (which atomic and molecular species, dust)
- and on the ability to radiate away „cooling lines“ and black body radiation
 - > problem of *radiation transfer* (see, e.g., IPAM III)

In general:

- the „standard way“ of solving the equations of (magneto) hydrodynamics is using finite differences on a grid
- alternative use particle-based scheme: *SPH*
- see IPAM workshop I

SPH

concept of SPH

- 'invented' independently by Lucy (1977) and Gingold & Monaghan (1977)
- originally proposed as Monte Carlo approach to calculate the time evolution of gaseous systems
- more intuitively understood as interpolation scheme:

The fluid is represented by an ensemble of particles i , each carrying mass m_i , momentum $m_i\vec{v}_i$, and hydrodynamic properties (like pressure p_i , temperature T_i , internal energy ϵ_i , entropy s_i , etc.). The time evolution is governed by the equation of motion plus additional equations to modify the hydrodynamic properties of the particles. Hydrodynamic observables are obtained by a local averaging process.

properties of local averaging processes

- local averages $\langle f(\vec{r}) \rangle$ for any quantity $f(\vec{r})$ can be obtained by convolution with an appropriate smoothing function $W(\vec{r}, \vec{h})$:

$$\langle f(\vec{r}) \rangle \equiv \int f(\vec{r}') W(\vec{r} - \vec{r}', \vec{h}) d^3 r' .$$

the function $W(\vec{r}, \vec{h})$ is called smoothing kernel

- the kernel must satisfy the following two conditions:

$$\int W(\vec{r}, \vec{h}) d^3 r = 1 \quad \text{and} \quad \langle f(\vec{r}) \rangle \longrightarrow f(\vec{r}) \quad \text{for} \quad \vec{h} \rightarrow 0$$

the kernel W therefore follows the same definitions as Dirac's delta function $\delta(\vec{r})$: $\lim_{h \rightarrow 0} W(\vec{r}, h) = \delta(\vec{r})$.

- most SPH implementations use spherical kernel functions

$$W(\vec{r}, \vec{h}) \equiv W(r, h) \quad \text{with} \quad r = |\vec{r}| \quad \text{and} \quad h = |\vec{h}| .$$

(one could also use triaxial kernels, e.g. Martel et al. 1995)

properties of local averaging processes

- as the kernel function W can be seen as approximation to the δ -function for small but finite h we can expand the averaged function $\langle f(\vec{r}) \rangle$ into a Taylor series for h to obtain an estimate for $f(\vec{r})$; if W is an even function, the first order term vanishes and the errors are second order in h

$$\langle f(\vec{r}) \rangle = f(\vec{r}) + \mathcal{O}(h^2)$$

this holds for functions f that are smooth and do not exhibit steep gradients over the size of W (\rightarrow problems in shocks).

(more specifically the expansion is $\langle f(\vec{r}) \rangle = f(\vec{r}) + \kappa h^2 \vec{\nabla}^2 f(\vec{r}) + \mathcal{O}(h^3)$)

properties of local averaging processes

- within its intrinsic accuracy, the smoothing process therefore is a linear function with respect to summation and multiplication:

$$\langle f(\vec{r}) + g(\vec{r}) \rangle = \langle f(\vec{r}) \rangle + \langle g(\vec{r}) \rangle$$

$$\langle f(\vec{r}) \cdot g(\vec{r}) \rangle = \langle f(\vec{r}) \rangle \cdot \langle g(\vec{r}) \rangle$$

(one follows from the linearity of integration with respect to summation, and two is true to $\mathcal{O}(h^2)$)

- derivatives can be ‘drawn into’ the averaging process:

$$\frac{d}{dt} \langle f(\vec{r}) \rangle = \left\langle \frac{d}{dt} f(\vec{r}) \right\rangle$$

$$\vec{\nabla} \langle f(\vec{r}) \rangle = \langle \vec{\nabla} f(\vec{r}) \rangle$$

Furthermore, the spatial derivative of f can be transformed into a spatial derivative of W (no need for finite differences or grid):

$$\vec{\nabla} \langle f(\vec{r}) \rangle = \langle \vec{\nabla} f(\vec{r}) \rangle = \int f(\vec{r}') \vec{\nabla} W(|\vec{r} - \vec{r}'|, h) d^3 r' .$$

(shown by integrating by parts and assuming that the surface term vanishes; if the solution space is extended far enough, either the function f itself or the kernel approach zero)

properties of local averaging processes

- basic concept of SPH is a **particle representation** of the fluid
→ *integration* transforms into *summation* over discrete set of particles; example density ρ :

$$\langle \rho(\vec{r}_i) \rangle = \sum_j m_j W(|\vec{r}_i - \vec{r}_j|, h) .$$

in this picture, the mass of each particle is smeared out over its kernel region; the density at each location is obtained by summing over the contributions of the various particles → ***smoothed particle hydrodynamics!***

the kernel function

- different functions meet the requirement $\int W(|\vec{r}|, h) d^3r = 1$
and $\lim_{h \rightarrow 0} \int W(|\vec{r} - \vec{r}'|, h) f(\vec{r}') d^3r' = f(\vec{r})$:

→ Gaussian kernel:

$$W(r, h) = \frac{1}{\pi^{3/2} h^3} \exp\left(-\frac{r^2}{h^2}\right)$$

- *pro*: mathematically sound
- *pro*: derivatives exist to all orders and are smooth
- *contra*: all particles contribute to a location

→ spline functions with compact support

the kernel function

- different functions meet the requirement $\int W(|\vec{r}|, h) d^3r = 1$
and $\lim_{h \rightarrow 0} \int W(|\vec{r} - \vec{r}'|, h) f(\vec{r}') d^3r' = f(\vec{r})$:

→ the standard kernel: **cubic spline**

with $\xi = r/h$ it is defined as

$$W(r, h) \equiv \frac{1}{\pi h^3} \begin{cases} 1 - \frac{3}{2}\xi^2 + \frac{3}{4}\xi^3, & \text{for } 0 \leq \xi \leq 1; \\ \frac{1}{4}(2 - \xi)^3, & \text{for } 1 \leq \xi \leq 2; \\ 0, & \text{otherwise.} \end{cases}$$

- *pro*: compact support → all interactions are zero for $r > 2h$ → number of particles involved in the average remains small (typically between 30 and 80)
- *pro*: second derivative is continuous
- *pro*: dominant error term is second order in h

the fluid equations in SPH

- there is an infinite number of possible SPH implementations of the hydrodynamic equations!
- some notation: $h_{ij} = (h_i + h_j)/2$, $\vec{r}_{ij} = \vec{r}_i - \vec{r}_j$, $\vec{v}_{ij} = \vec{v}_i - \vec{v}_j$, and $\vec{\nabla}_i$ is the gradient with respect to the coordinates of particle i ; all measurements are taken at particle positions (e.g. $\rho_i = \rho(\vec{r}_i)$)
- *general form of SPH equations:*

$$\langle f_i \rangle = \sum_{j=1}^{N_i} \frac{m_j}{\rho_j} f_j W(r_{ij}, h_{ij})$$

the fluid equations in SPH

- *density* — continuity equation (conservation of mass)

$$\rho_i = \sum_{j=1}^{N_i} m_j W(r_{ij}, h_{ij})$$

or
$$\frac{d\rho_i}{dt} = \sum_{j=1}^{N_i} m_j \vec{v}_{ij} \cdot \vec{\nabla}_i W(r_{ij}, h_{ij})$$

(the second implementation is almost never used, see however Monaghan 1991 for an application to water waves)

important

density is needed for **ALL** particles **BEFORE** computing other averaged quantities → at each timestep, SPH computations consist of **TWO** loops, first the *density* is obtained for each particle, and then in a second round, all *other* particle properties are updated.

the fluid equations in SPH

- *pressure* is defined via the *equation of state* (for example for isothermal gas $p_i = c_s^2 \rho_i$)

the fluid equations in SPH

- *velocity* — Navier Stokes equation (conservation of momentum)

$$\frac{d\vec{v}}{dt} = \frac{\partial\vec{v}}{\partial t} + (\vec{v} \cdot \vec{\nabla})\vec{v} = \sum_i \vec{F}_i = \vec{F}_{\text{pressure}} + \vec{F}_{\text{viscosity}} + \vec{F}_{\text{gravity}}$$

rate of change of momentum of fluid element depends on sum of all forces acting on it.

the fluid equations in SPH

- *velocity* — **Navier Stokes equation** (conservation of momentum)

→ consider for now *only* pressure contributions: Euler's equation

$$\frac{d\vec{v}}{dt} = \frac{\partial\vec{v}}{\partial t} + (\vec{v} \cdot \vec{\nabla})\vec{v} = -\frac{1}{\rho} \vec{\nabla} p = -\vec{\nabla} \left(\frac{p}{\rho} \right) - \frac{p}{\rho^2} \vec{\nabla} \rho \quad (*)$$

here, the identity $\vec{\nabla}(p\rho^{-1}) = \rho^{-1}\vec{\nabla}p - p\rho^{-2}\vec{\nabla}\rho$ is used

→ in the SPH formalism this reads as

$$\frac{d\vec{v}_i}{dt} = - \sum_{j=1}^{N_i} m_j \left(\frac{p_i}{\rho_i^2} + \frac{p_j}{\rho_j^2} \right) \vec{\nabla}_i W(r_{ij}, h_{ij})$$

where the first term in (*) is neglected because it leads to surface terms in the averaging procedure; it is assumed that either the pressure or the kernel becomes zero at the integration border; if this is not the case *correction terms* need to be added above.

the fluid equations in SPH

- *velocity* — **Navier Stokes equation** (conservation of momentum)
→ the SPH implementation of the standard artificial viscosity is

$$\vec{F}_i^{\text{visc}} = - \sum_{j=1}^{N_i} m_j \Pi_{ij} \vec{\nabla}_i W(r_{ij}, h_{ij}),$$

where the viscosity tensor Π_{ij} is defined by

$$\Pi_{ij} = \begin{cases} (-\alpha c_{ij} \mu_{ij} + \beta \mu_{ij}^2) / \rho_{ij} & \text{for } \vec{v}_{ij} \cdot \vec{r}_{ij} \leq 0, \\ 0 & \text{for } \vec{v}_{ij} \cdot \vec{r}_{ij} > 0, \end{cases}$$

where

$$\mu_{ij} = \frac{h \vec{v}_{ij} \cdot \vec{r}_{ij}}{\vec{r}_{ij}^2 + 0.01 h^2}.$$

with $\vec{r}_{ij} = \vec{r}_i - \vec{r}_j$, $\vec{v}_{ij} = \vec{v}_i - \vec{v}_j$, mean density $\rho_{ij} = (\rho_i + \rho_j)/2$, and mean sound speed $c_{ij} = (c_i + c_j)/2$.

the fluid equations in SPH

- *velocity* — **Navier Stokes equation** (conservation of momentum)
 - if self-gravity is taken into account, the gravitational force needs to be added on the RHS

$$\vec{F}_G = -\vec{\nabla}\phi_i = -G \sum_{j=1}^N \frac{m_j}{r_{ij}^2} \frac{r_{ij}}{r_{ij}}$$

note that the sum needs to be taken over *ALL* particles ←
computationally expensive

- set together, the momentum equation is

$$\frac{d\vec{v}_i}{dt} = - \sum_{j=1}^{N_i} m_j \left(\frac{p_i}{\rho_i^2} + \frac{p_j}{\rho_j^2} + \Pi_{ij} \right) \vec{\nabla}_i W(r_{ij}, h_{ij}) - \nabla \phi_i$$

the fluid equations in SPH

- *energy equation* (conservation of momentum)

→ recall the hydrodynamic energy equation:

$$\frac{d\epsilon}{dt} = \frac{\partial\epsilon}{\partial t} + \vec{v} \cdot \vec{\nabla}\epsilon = \frac{ds}{dt} - \frac{p}{\rho} \vec{\nabla} \cdot \vec{v}$$

→ for *adiabatic* systems ($c = \text{const}$) the SPH form follows as

$$\frac{d\epsilon_i}{dt} = \frac{p_i}{\rho_i^2} \sum_{j=1}^{N_i} m_j \vec{v}_{ij} \cdot \vec{\nabla}_i W(r_{ij}, h_{ij}),$$

(note that the alternative form

$$\frac{d\epsilon_i}{dt} = \frac{1}{2} \sum_{j=1}^{N_i} m_j \left(\frac{p_i}{\rho_i^2} + \frac{p_j}{\rho_j^2} \right) \vec{v}_{ij} \cdot \vec{\nabla}_i W(r_{ij}, h_{ij})$$

can lead to unphysical solutions, like negative internal energy)

the fluid equations in SPH

- *energy equation* (conservation of momentum)

→ *dissipation* due to (artificial) viscosity leads to a term

$$\frac{d\epsilon_i}{dt} = \frac{1}{2} \sum_{j=1}^{N_i} m_j \Pi_{ij} \vec{v}_{ij} \cdot \vec{\nabla}_i W(r_{ij}, h_{ij})$$

→ the presence of *heating* sources or *cooling* processes can be incorporated into a function Γ_i .

→ altogether:

$$\frac{d\epsilon_i}{dt} = \frac{p_i}{\rho_i^2} \sum_{j=1}^{N_i} m_j \vec{v}_{ij} \cdot \vec{\nabla}_i W_{ij} + \frac{1}{2} \sum_{j=1}^{N_i} m_j \Pi_{ij} \vec{v}_{ij} \cdot \vec{\nabla}_i W_{ij} + \Gamma_i$$

can lead to unphysical solutions, like negative internal energy)

fully conservative formulation using Lagrange multipliers

- the Lagrangian for compressible flows which are generated by the thermal energy $\epsilon(\rho, s)$ acts as effective potential is

$$\mathcal{L} = \int \rho \left\{ \frac{1}{2} v^2 - u(\rho, s) \right\} d^3r.$$

equations of motion follow with $s = \text{const}$ from

$$\frac{d}{dt} \frac{\partial \mathcal{L}}{\partial \vec{v}} - \frac{\partial \mathcal{L}}{\partial \vec{r}} = 0$$

- after some SPH arithmetics, one can derive the following acceleration equation for particle i

$$\frac{d\vec{v}_i}{dt} = - \sum_{j=1}^{N_i} m_j \left\{ \frac{1}{f_i \rho_i^2} p_i \vec{\nabla}_i W(r_{ij}, h_i) + \frac{1}{f_j \rho_j^2} p_j \vec{\nabla}_i W(r_{ij}, h_j) \right\}$$

where

$$f_i = \left[1 + \frac{h_i}{3\rho_i} \frac{\partial \rho_i}{\partial h_i} \right]$$

Large-eddy simulations

- We use **LES** to model the large-scale dynamics
- Principal problem: only large scale flow properties
 - Reynolds number: $Re = LV/\nu$ ($Re_{nature} \gg Re_{model}$)
 - dynamic range much smaller than true physical one
 - need **subgrid model** (in our case simple: only dissipation) more complex when processes (chemical reactions, nuclear burning, etc) on subgrid scale determine large-scale dynamics
- Also: stochasticity of the flow \Rightarrow unpredictable when and where “interesting things” happen
 - occurrence of localized collapse
 - location and strength of shock fronts
 - etc.

LES with SPH

- For self-gravitating gases **SPH** is probably okay ...
 - fully Lagrangian (particles are free to move where needed)
 - good resolution in high-density regions (in collapse)
 - particle based --> good for transition from hydrodynamics to stellar dynamics
- BUT:
 - low resolution in low-density region
 - difficult to reach very high levels of refinement (however, particle splitting may be promising path)
 - dissipative and need for artificial viscosity
 - how to handle subgrid scales?

Gravoturbulent SF with SPH

- Comparison between particle-based and grid-based methods: **SPH** vs. **ZEUS**

Klessen, Heitsch, Mac Low (2000)

Heitsch, Mac Low, Klessen (2001)

Ossenkopf, Klessen, Heitsch (2001)

- Both methods are complementary...

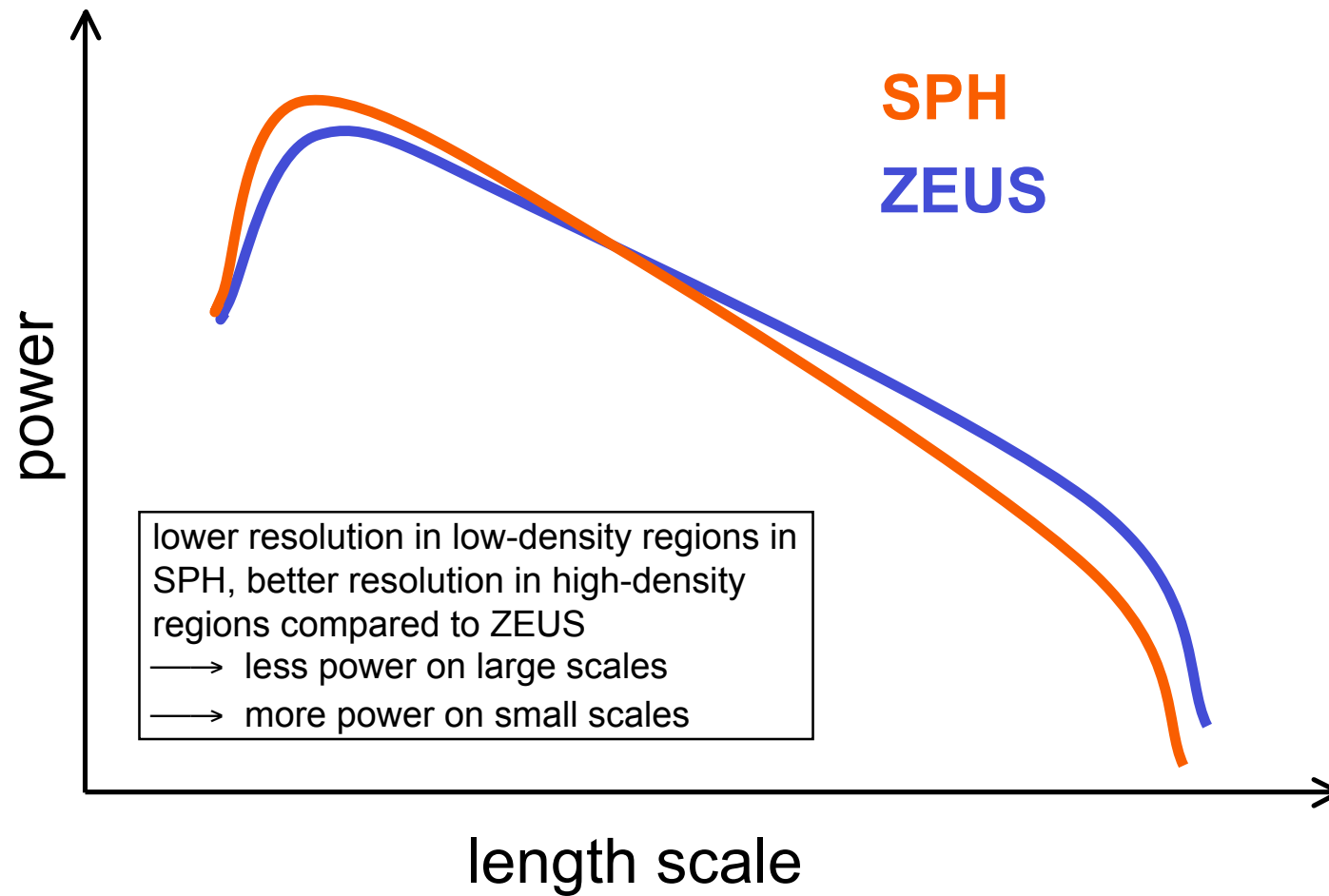
→ **Bracketing reality!**

- As a crude estimation:

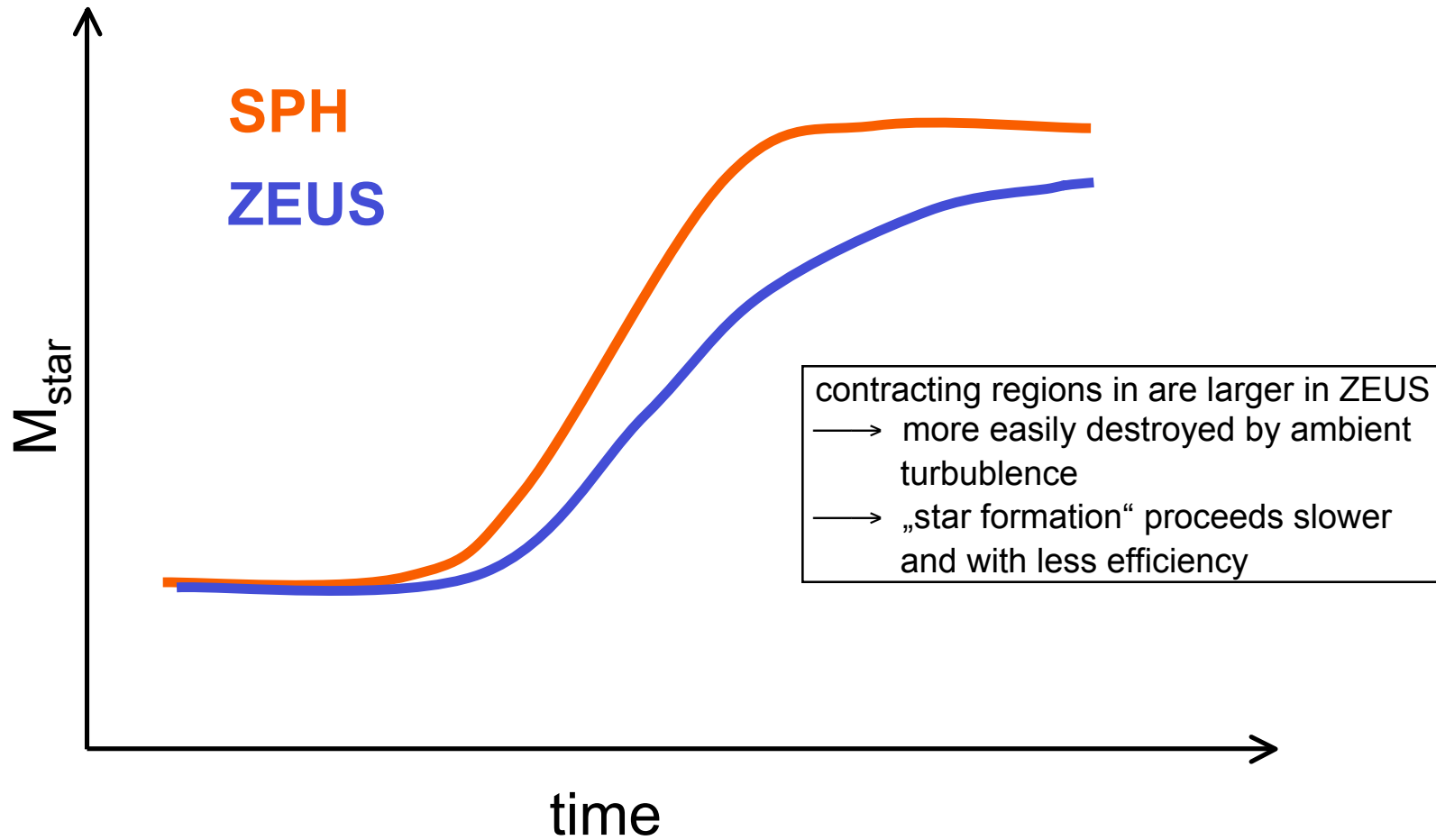
SPH is better in high-density regions

ZEUS is better in low-density regions

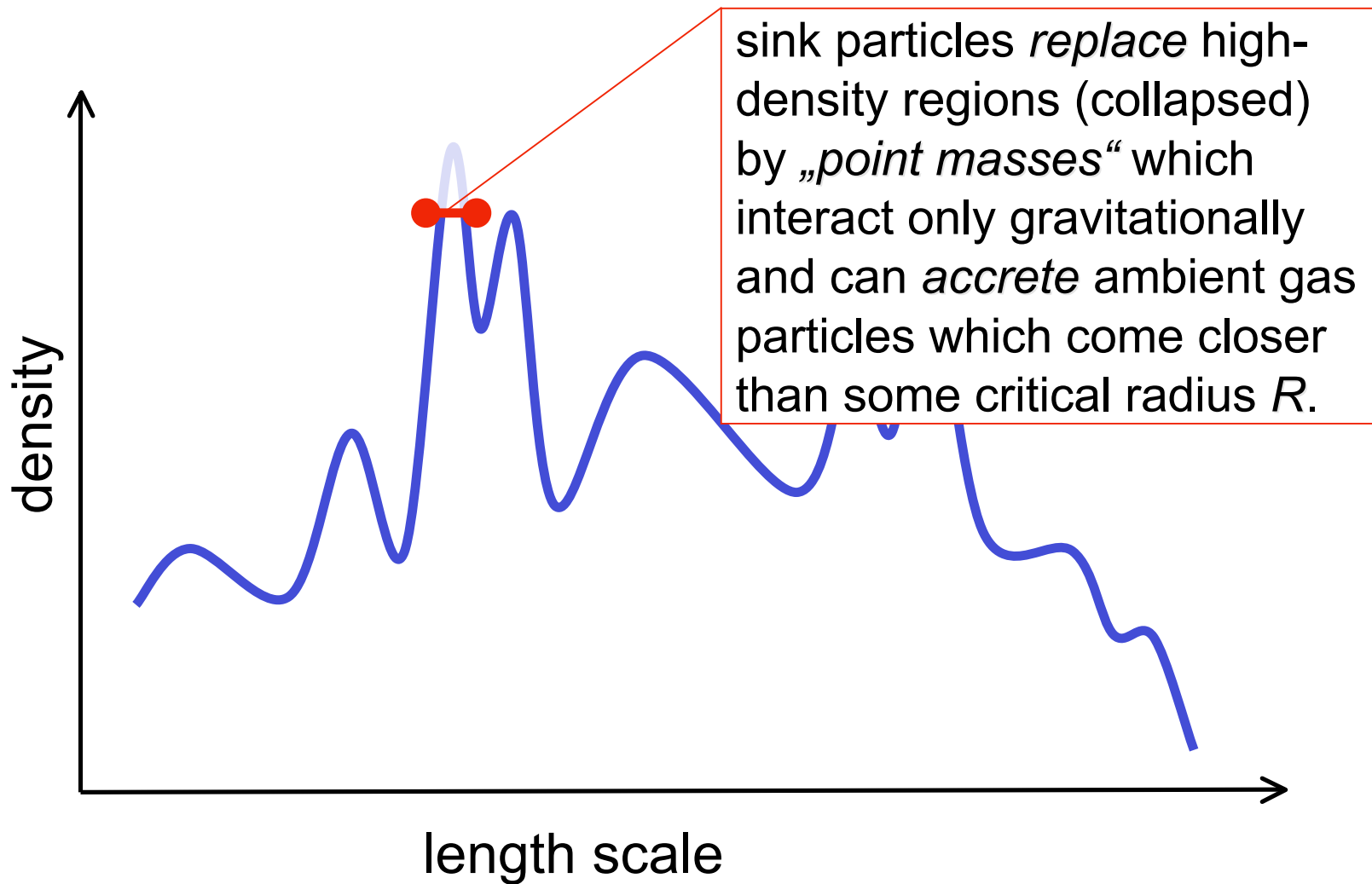
SPH vs. ZEUS



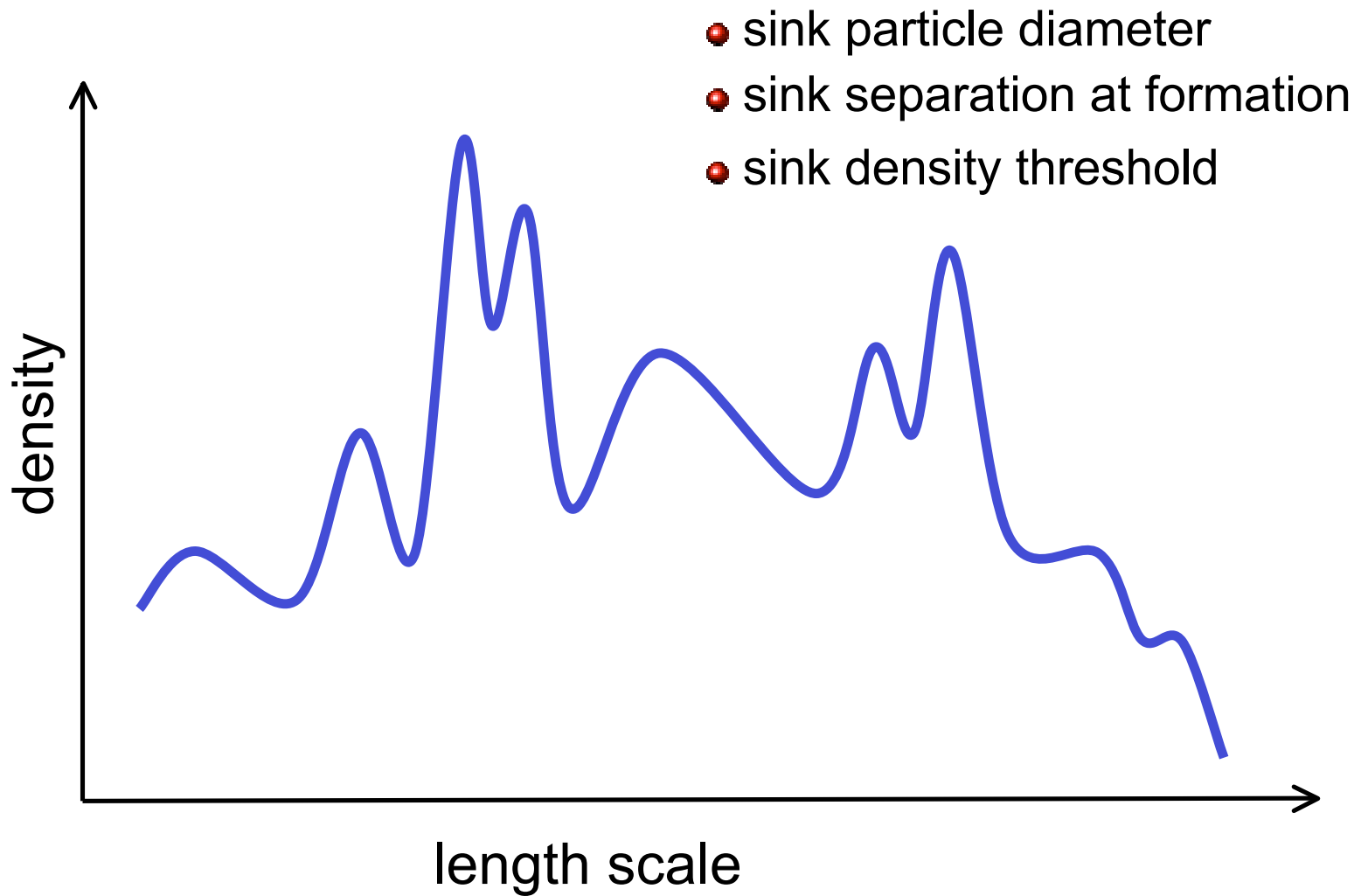
SPH vs. ZEUS



SPH with sink particles I

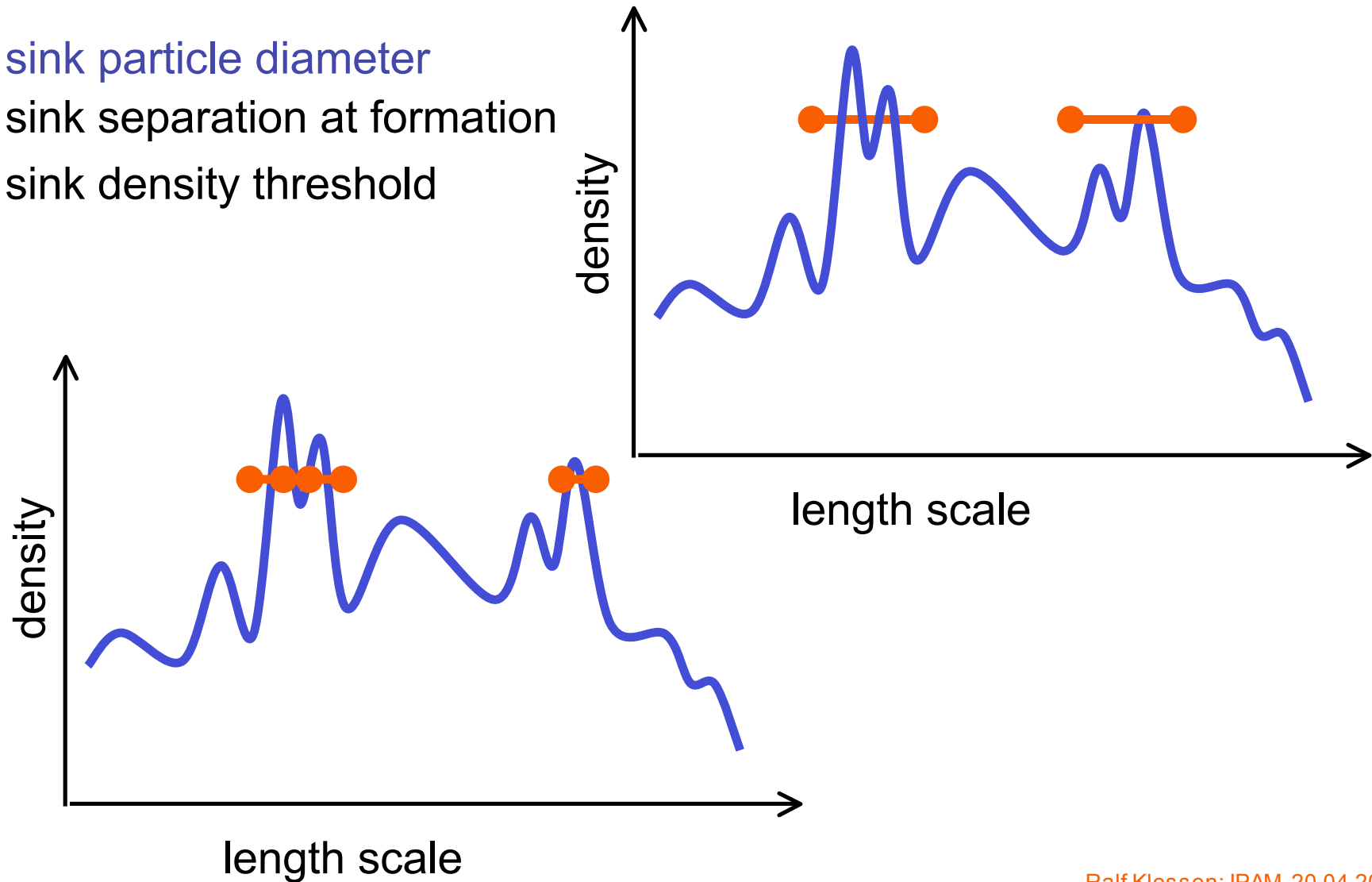


SPH with sink particles I



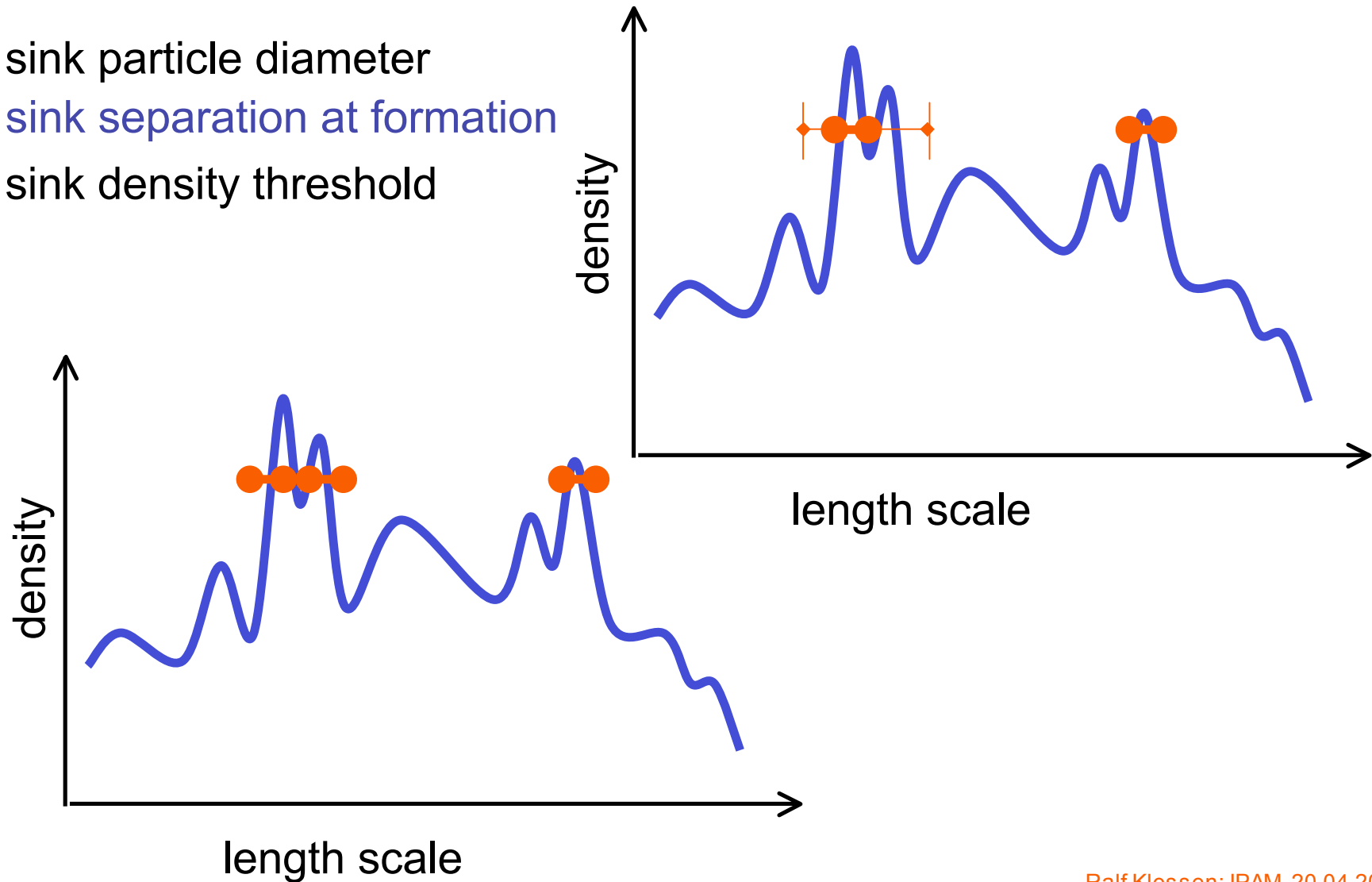
SPH with sink particles II

- sink particle diameter
- sink separation at formation
- sink density threshold



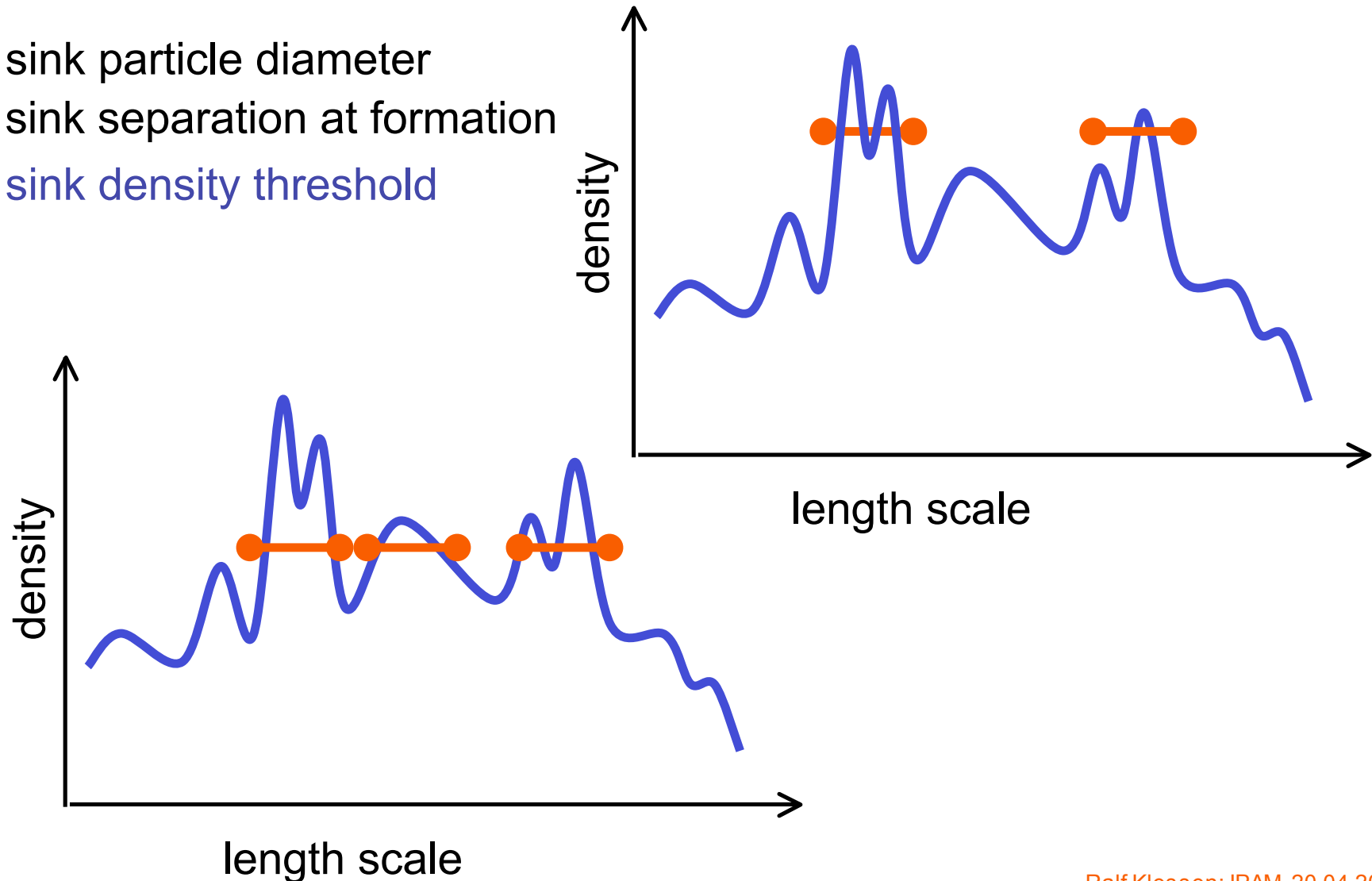
SPH with sink particles III

- sink particle diameter
- sink separation at formation
- sink density threshold



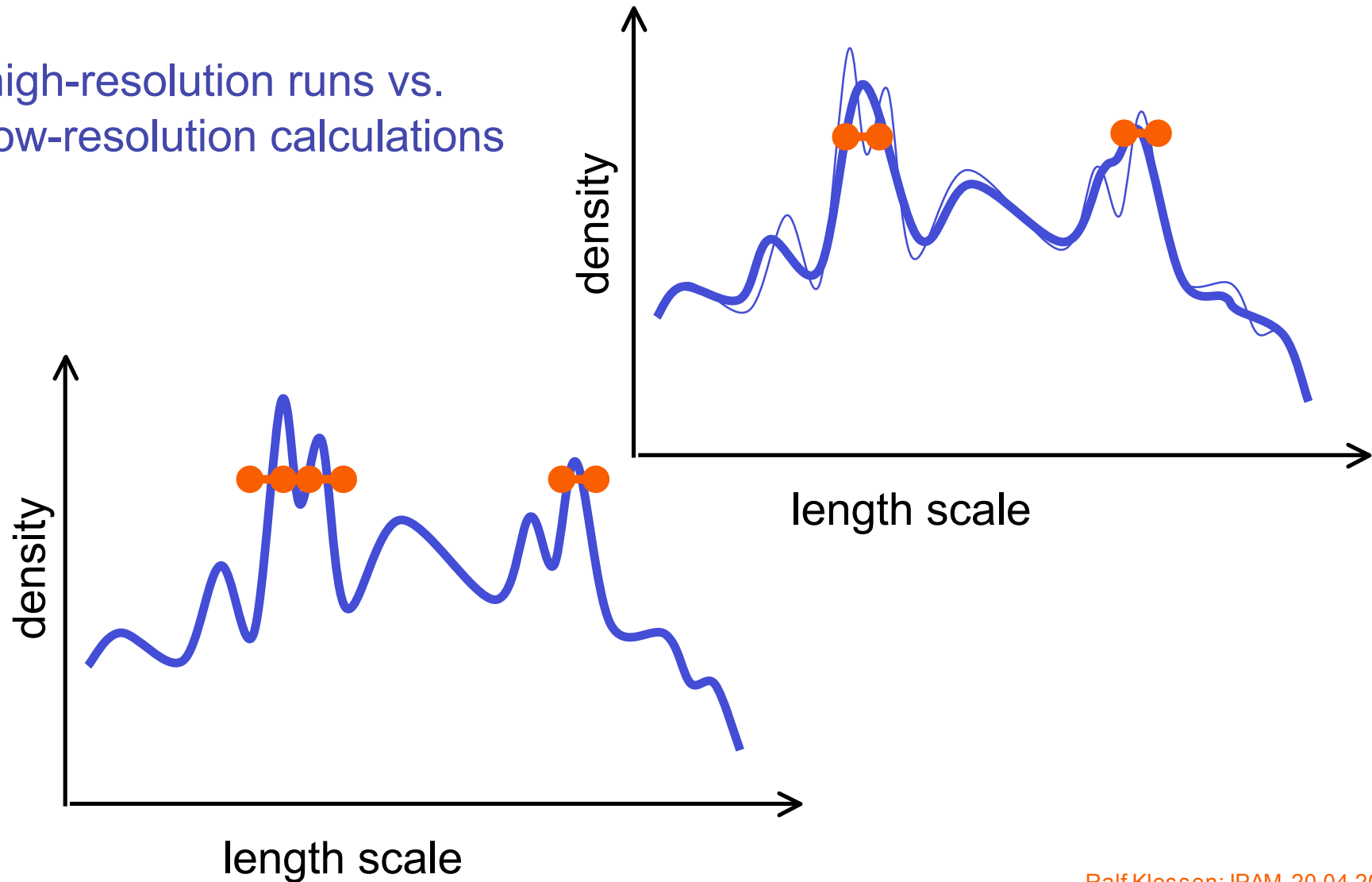
SPH with sink particles IV

- sink particle diameter
- sink separation at formation
- sink density threshold



SPH with sink particles V

high-resolution runs vs.
low-resolution calculations



Some final remarks...

- *GRAVOTURBULENT STAR FORMATION:*

This dynamic theory can explain and reproduce many features of star-forming regions on small as well as on large galactic scales.

- Some open questions:

- role of magnetic fields?
- role of thermodynamic state of the gas?
- what drives turbulence?
- how are small scales (local molecular clouds) connected to large-scale dynamics?
- what terminates star formation locally?

Some final remarks...

● *NUMERICS:*

SPH appears able to describe gravoturbulent fragmentation and star formation in molecular clouds.

- Pro:
 - *Lagrangian* character of method.
 - can resolve *large density contrasts*.
 - good for transition from hydro- to stellar dynamics
--> accreting sink particles describe protostars
- Con:
 - low resolution in low-density regions.
 - difficulties with shock-capturing and treating B-fields.
- Next steps:
 - particle-splitting to locally increase resolution,
 - GPM, XSPH with “physical” viscosity

Outlook & First Examples

- **WHAT WE REALLY NEED: more physics!!**
We need good *subgrid-scale models* for unresolved scales in our calculations.
- 2 Examples:
 - LOCAL SCALES: so far, we use dumb sink particles to protostellar collapse --> we do not know how the star “inside” forms and how it backreacts onto the ambient environment
--> combine 3D hydro with 1D/2D PMS models
 - GALACTIC SCALES: can gravoturbulent models give us some handle on star-formation efficiency?
--> some thoughts...

Example 1

Towards a complete picture...

COMBINE:

- **3D** hydrodynamic simulations of the *turbulent fragmentation* of *entire molecular cloud regions*.

(using SPH with GRAPE: Klessen & Burkert 2000, 2001, Klessen, Heitsch & Mac Low 2000, Klessen 2001b)

WITH:

- detailed **2D** hydrodynamic modelling of *protostellar accretion disks*. (e.g. Yorke & Bodenheimer 1999)

→ with rad. transfer → **SED, T_{bol} , L_{bol} etc.**

AND/OR:

- implicit **1D** radiation-hydrodynamic scheme with time-dependent convection and D network following the *collapse of individual cores towards the MS* (Wuchterl & Tscharnuter 2002)
→ **PMS tracks, absolute stellar ages for cluster stars**

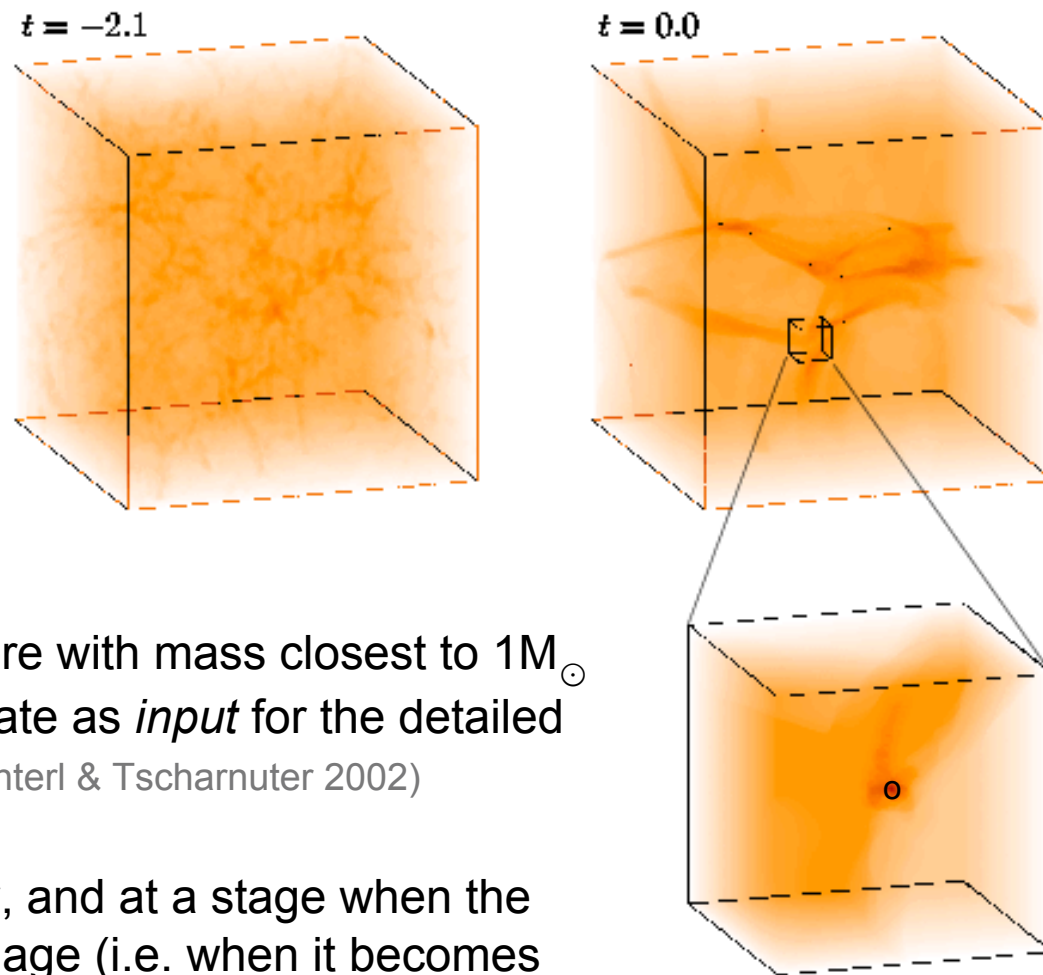
Formation of a $1M_{\odot}$ star

Dynamical evolution of a molecular cloud region of size $(0.32\text{pc})^3$ containing $200 M_{\odot}$ of gas
(from Klessen & Burkert 2000)

Within two free-fall times the system builds up a *cluster* of deeply embedded accreting *protostellar cores*.

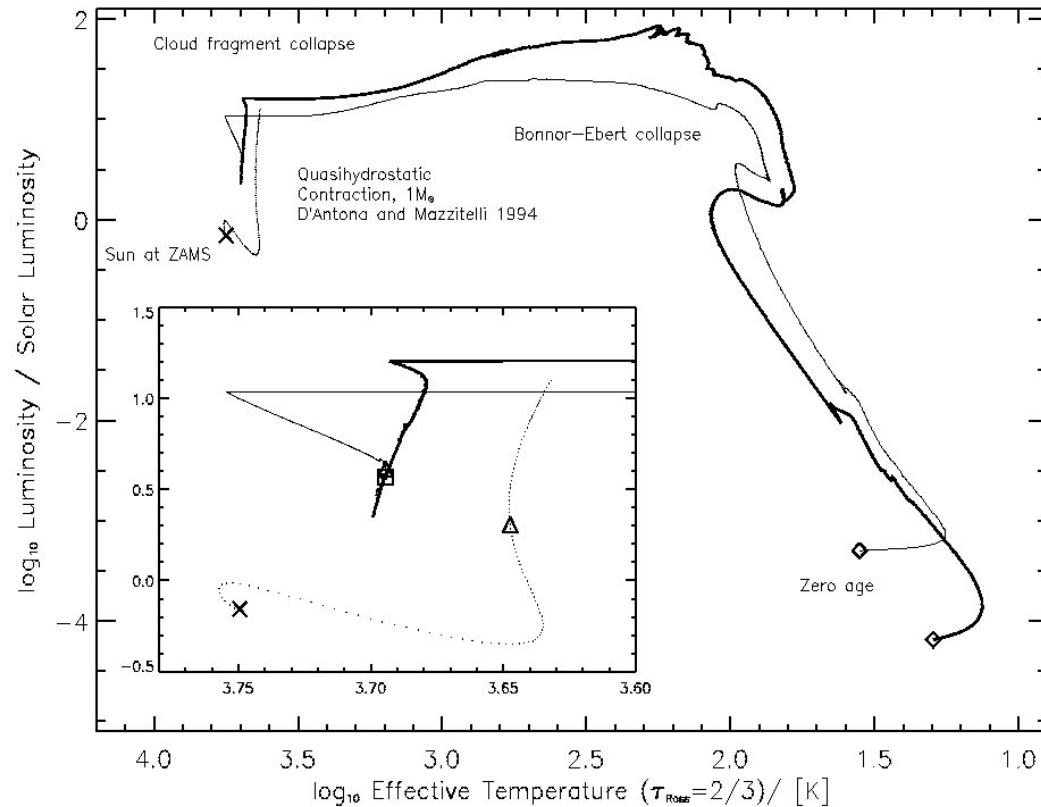
We select the protostellar core with mass closest to $1M_{\odot}$ and use its mass accretion rate as *input* for the detailed 1D-RHD calculation (see Wuchterl & Tscharnuter 2002)

The system is shown initially, and at a stage when the $1M_{\odot}$ -fragment reaches zero age (i.e. when it becomes optically thick for the first time).



(from Wuchterl & Klessen 2001)

PMS tracks



- core in cluster environment
- isolated Bonnor-Ebert core
- D'Antona & Mazzitelli track

For ages less than 10^6 years, different collapse conditions lead to different evolutionary tracks. Later the dynamical tracks converge.

There are large differences to the hydrostatic track, due to different stellar structure.

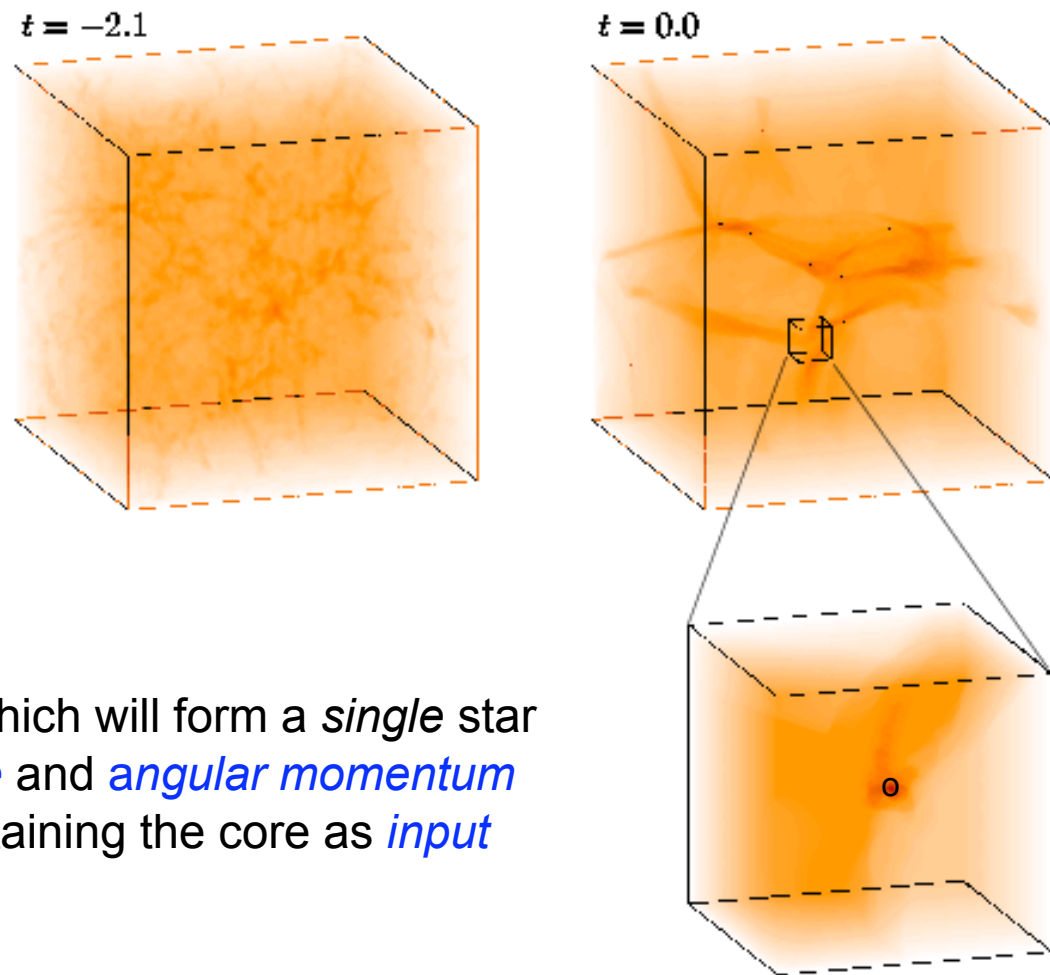
(from Wuchterl & Klessen 2001)

SED's of a $1 M_{\odot}$ star

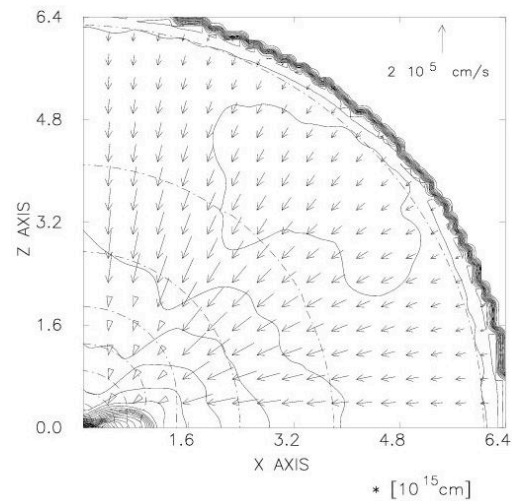
Dynamical evolution of a molecular cloud region of size $(0.32\text{pc})^3$ containing $200 M_{\odot}$ of gas
(from Klessen & Burkert 2000)

Within one to two free-fall times the system builds up a *cluster* of deeply embedded accreting *protostellar cores*.

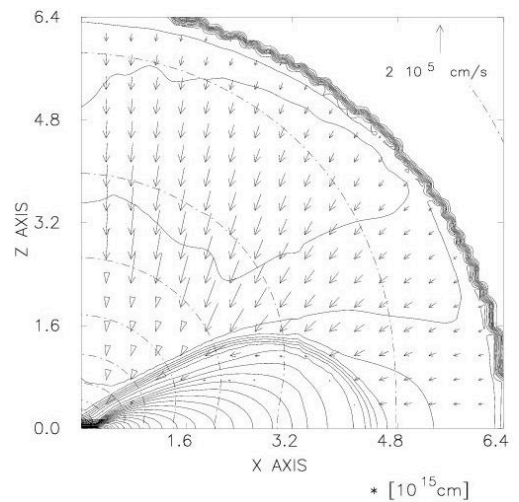
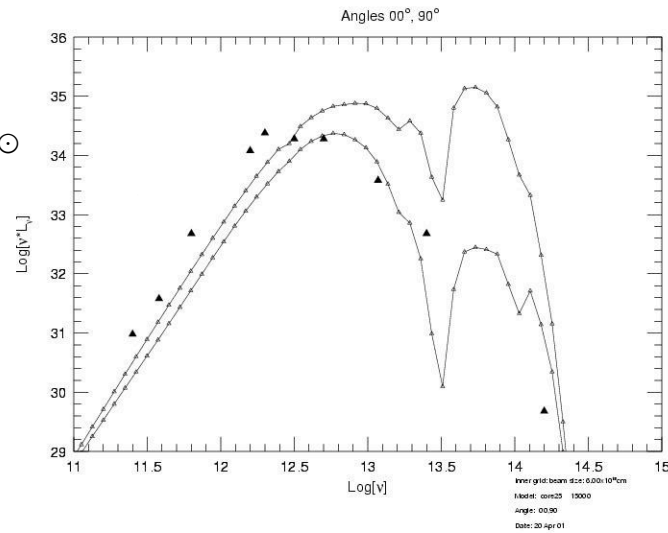
We select a protostellar core which will form a *single* star and use its *mass accretion rate* and *angular momentum gain* into a control volume containing the core as *input* for a detailed 2D calculation
(Bodenheimer & Klessen, in preparation)



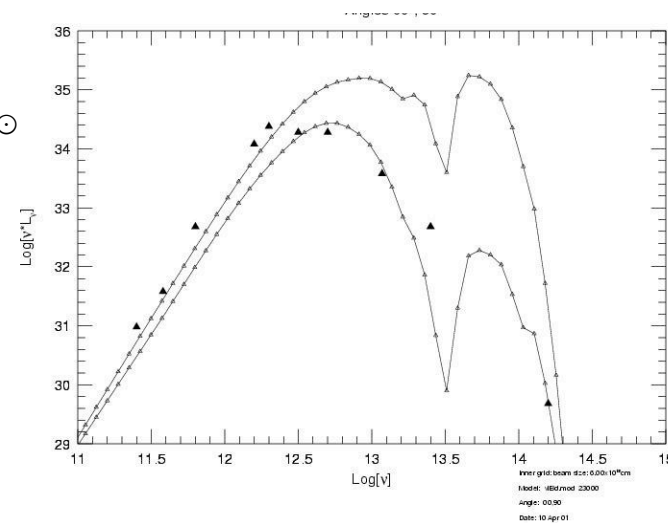
SED's of a $1 M_{\odot}$ star



$t \approx 15000\text{yr}$
 $M_{\text{tot}} \approx 0.75 M_{\odot}$



$t \approx 30000\text{yr}$
 $M_{\text{tot}} \approx 0.89 M_{\odot}$



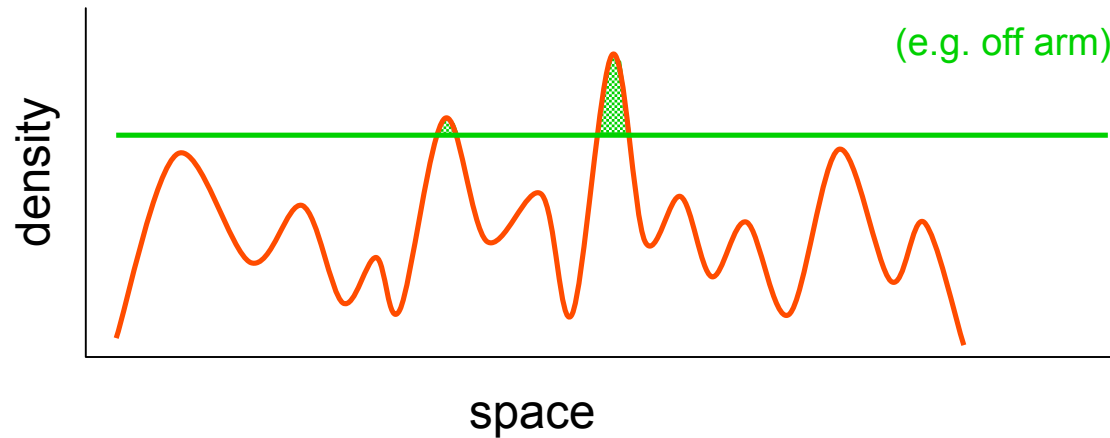
(Bodenheimer & Klessen, in preparation)

Example 2

Star formation on *global scales*

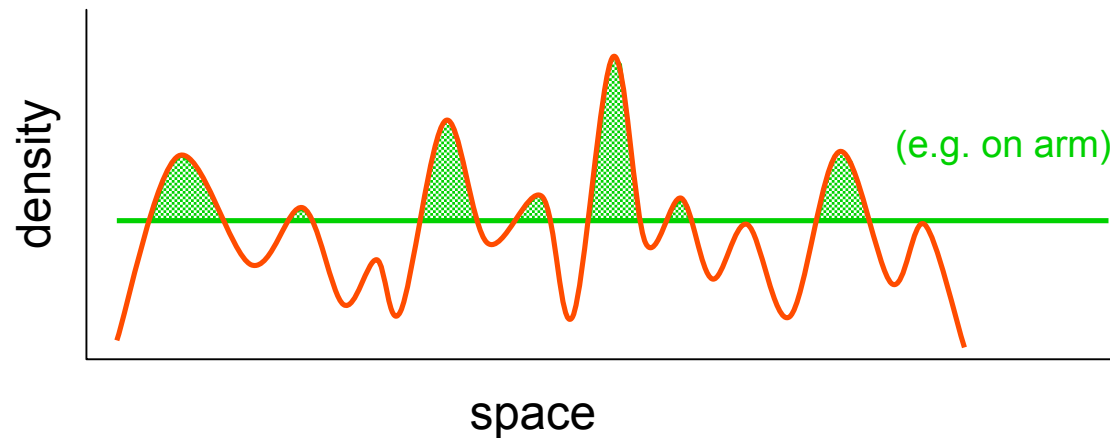
- *SF on global scales* = *formation of molecular clouds*
- MC's form at *stagnation points* of *convergent large-scale flows* (need $\sim 0.5 \text{kpc}^3$ of gas) \rightarrow high density \rightarrow enhanced cooling \rightarrow fast H_2 formation & gravitational instability \rightarrow local collapse and star formation
- External perturbations *increase* the local likelihood of MC formation (e.g. in spiral density waves, galaxy interactions, etc.)

Star formation on *global scales*



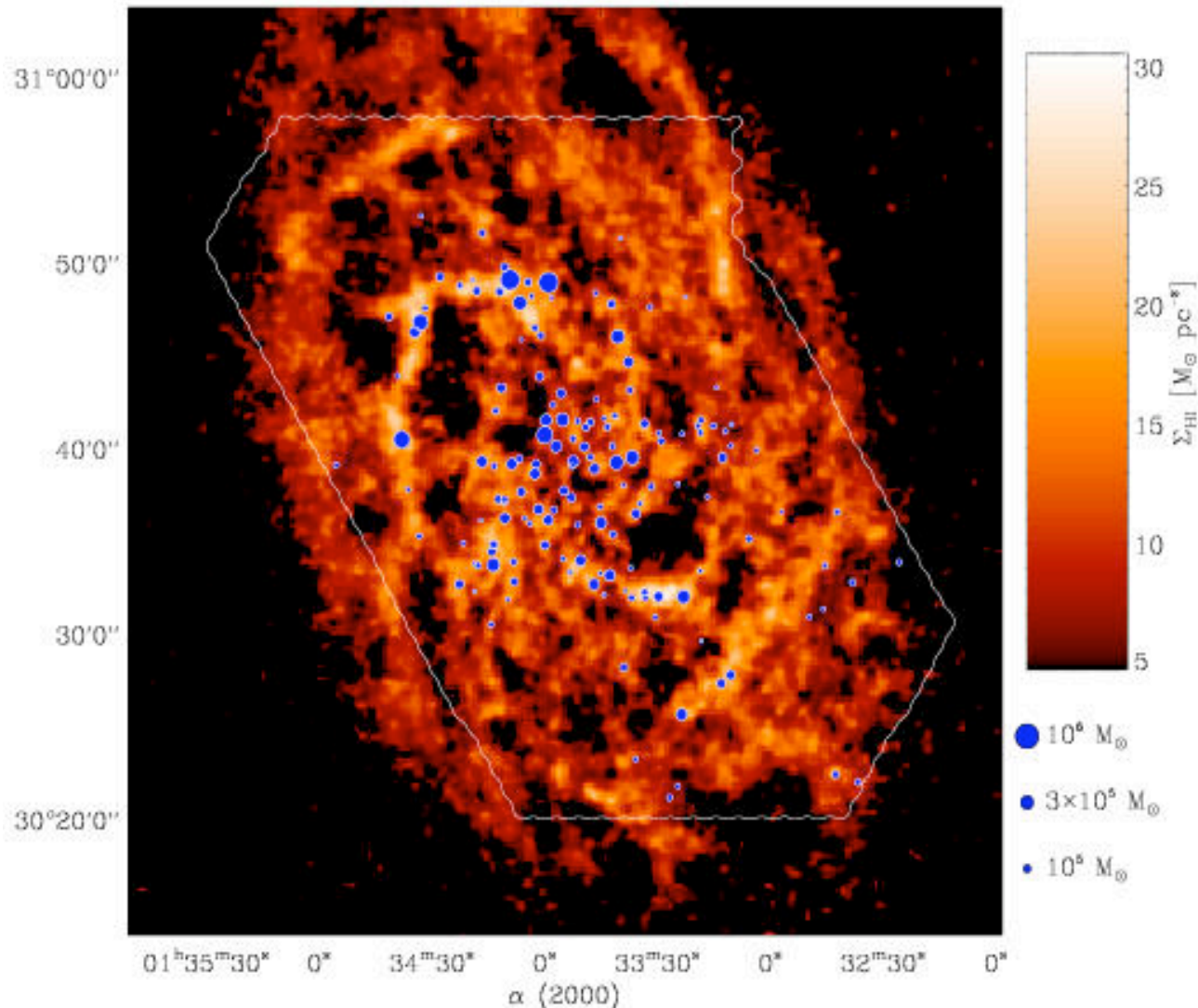
density fluctuations in warm atomic ISM caused by supersonic turbulence

some are dense enough to form H₂ within “reasonable timescale”
→ molecular clouds



external perturbations (i.e. potential changes) increase likelihood

Correlation between H₂ and HI



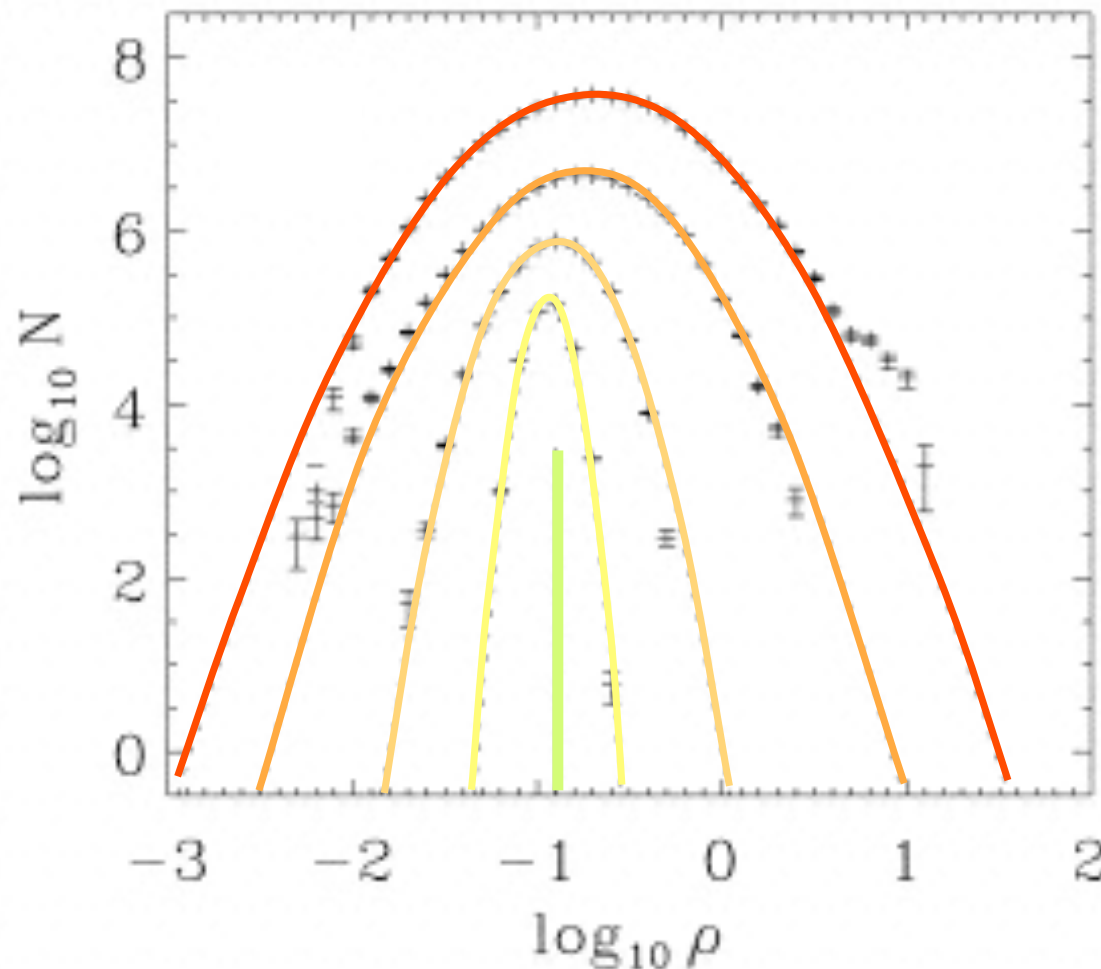
Compare H₂ - HI
in M33:

- H₂: BIMA-SONG Survey, see Blitz et al.
- HI: Observations with Westerbork Radio T.

H₂ clouds are seen
in regions of high
HI density
(in spiral arms and
filaments)

(Deul & van der Hulst 1987, Blitz et al. 2004)

Star formation on *global scales*



probability
distribution
function of
density
(ρ -pdf) for
decaying
supersonic
turbulence

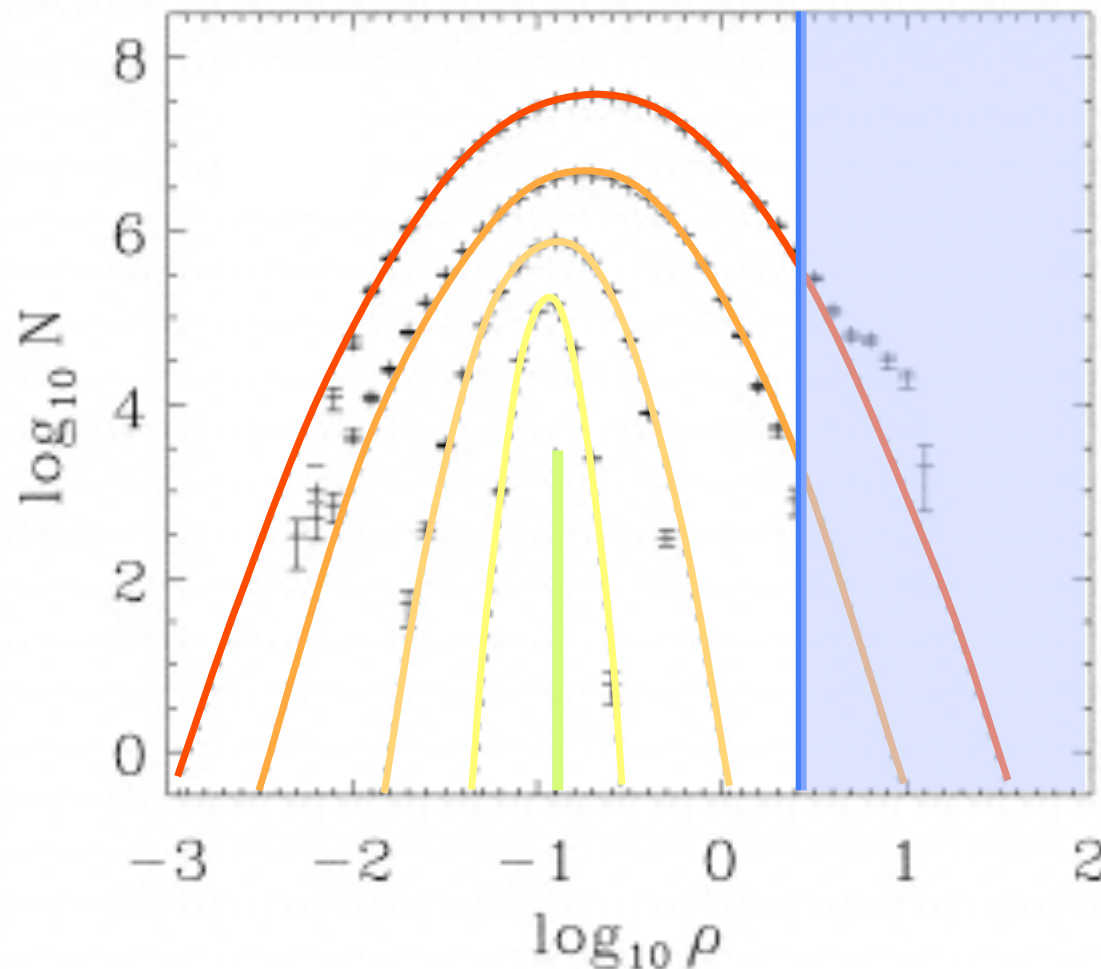
*varying rms Mach
numbers:*

M1 > **M2** >
M3 > **M4** > **0**

mass weighted ρ -pdf, each shifted by $\Delta \log N = 1$

(from Klessen, 2001)

Star formation on *global scales*



mass weighted ρ -pdf, each shifted by $\Delta \log N = 1$

(from Klessen, 2001; rate from Hollenback, Werner, & Salpeter 1971)

H₂ formation rate:

$$\tau_{\text{H}_2} \approx \frac{1.5 \text{ Gyr}}{n_{\text{H}} / 1 \text{ cm}^{-3}}$$

For $n_{\text{H}} \geq 100 \text{ cm}^{-3}$,
H₂ forms within
10 Myr, this is
about the lifetime
of typical MC's.

*What fraction of
the galactic ISM
reaches such
densities?*

THANKS