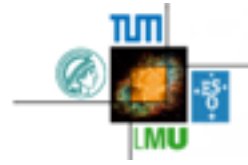




Excellence Cluster Universe



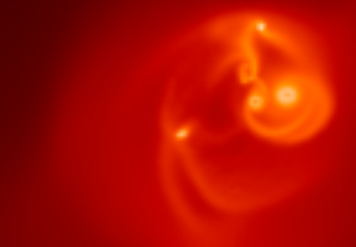
# N-body algorithms in GANDALF

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David Hubber

USM, LMU, München  
Excellence Cluster Universe,  
Garching bei München, Germany

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# Collisional vs. Collisionless N-body dynamics

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- N-body algorithms are usually divided up into two main classes :
  - **Collisional** : N-body particles are central point masses which can have strong 2-body interactions (e.g. stellar encounters)
    - NBODY6, Starlab/kira
  - **Collisionless** : N-body particles have a smoothed potential so only feel long-range potential forces (e.g. cold-dark matter fluid)
    - GADGET 2/3, GASOLINE
- Both ‘versions’ of N-body simulations can be realised in GANDALF
- However, the collisional N-body dynamics is **only realised designed for relatively small N-body systems** and not for large-N systems (e.g. the million body problem)

# Simple collisionless N-body integrators

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- **Collisionless** N-body integrators in GANDALF use the same algorithms as the SPH particles, i.e.

- Leapfrog kick-drift-kick (i.e. lfkdk)

$$\mathbf{r}_i(t + \Delta t) = \mathbf{r}_i(t) + \mathbf{v}_i(t) \Delta t + \frac{1}{2} \mathbf{a}_i(t) \Delta t^2$$

$$\mathbf{v}_i(t + \Delta t) = \mathbf{v}_i(t) + \frac{1}{2} (\mathbf{a}_i(t) + \mathbf{a}_i(t + \Delta t)) \Delta t$$

- Leapfrog drift-kick-drift (i.e. lfdkd)
- These integrators are **symplectic**, i.e. have very good conservation properties, particularly angular momentum

# Simple collisionless N-body integrators

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- **Simplest way** to simulate collisionless N-body is to use SPH particles with self gravity but hydro\_forces switched off!

```
hydro_forces = 0  
self_gravity = 1
```

- Developing multi-species in GANDALF in order to have cdm particles, i.e. self-gravitating but no hydro forces, as well as hydro particles

```
template <int ndim>  
struct Particle  
{  
    bool active;           ///< Flag if active (i.e. recompute step)  
    bool potmin;          ///< Is particle at a potential minima?  
    int iorig;            ///< Original particle i.d.  
    int itype;            ///< SPH particle type  
    etc..  
};
```

```
part.itype = gas;  
part.itype = cdm;
```

# Simple collisional N-body integrators

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- **Collisional** N-body integrators are more demanding because
  - Stars may have rather violent 2-body (or 3-body) interactions
  - Requires much higher accuracy with the integrations
- Simplest integrators are the same as the collisionless code
  - Leapfrog kick-drift-kick (i.e. Ifkdk)
  - Leapfrog drift-kick-drift (i.e. Ifdkd)

# More sophisticated N-body integrators

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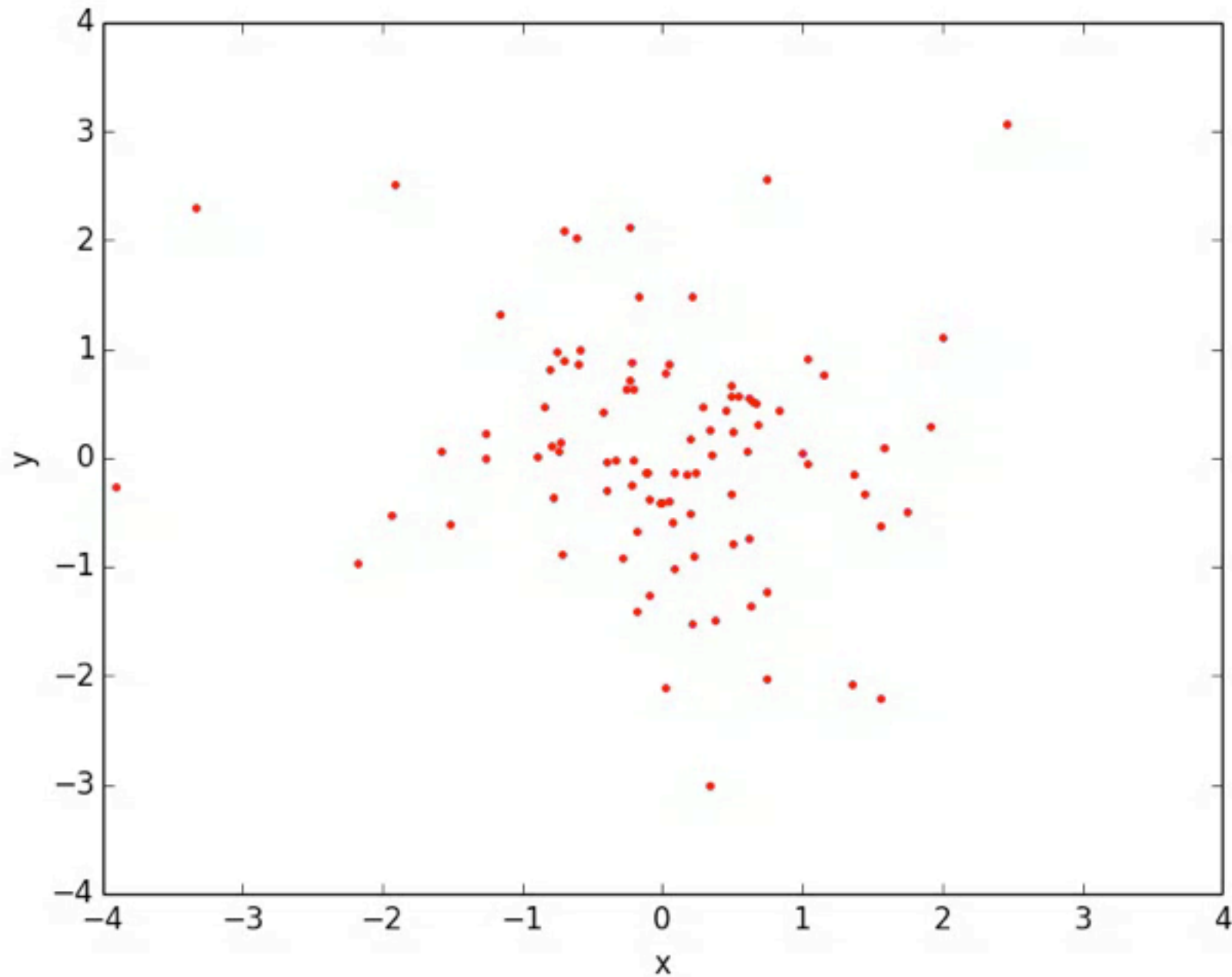
- For more accuracy, we can use :
  - 4th, 6th and 8th-order Hermite scheme (Makino & Aarseth 1992)
  - KS-regularisation
- Hermite schemes compute both the force AND the force derivative

$$\mathbf{a}_s = -G \sum_{t=1}^N m_t \phi'(\mathbf{r}_{st}, \bar{h}_{st}) \hat{\mathbf{r}}_{st} - G \sum_{i=1}^N m_i \phi'(\mathbf{r}_{si}, \bar{h}_{si}) \hat{\mathbf{r}}_{si}$$

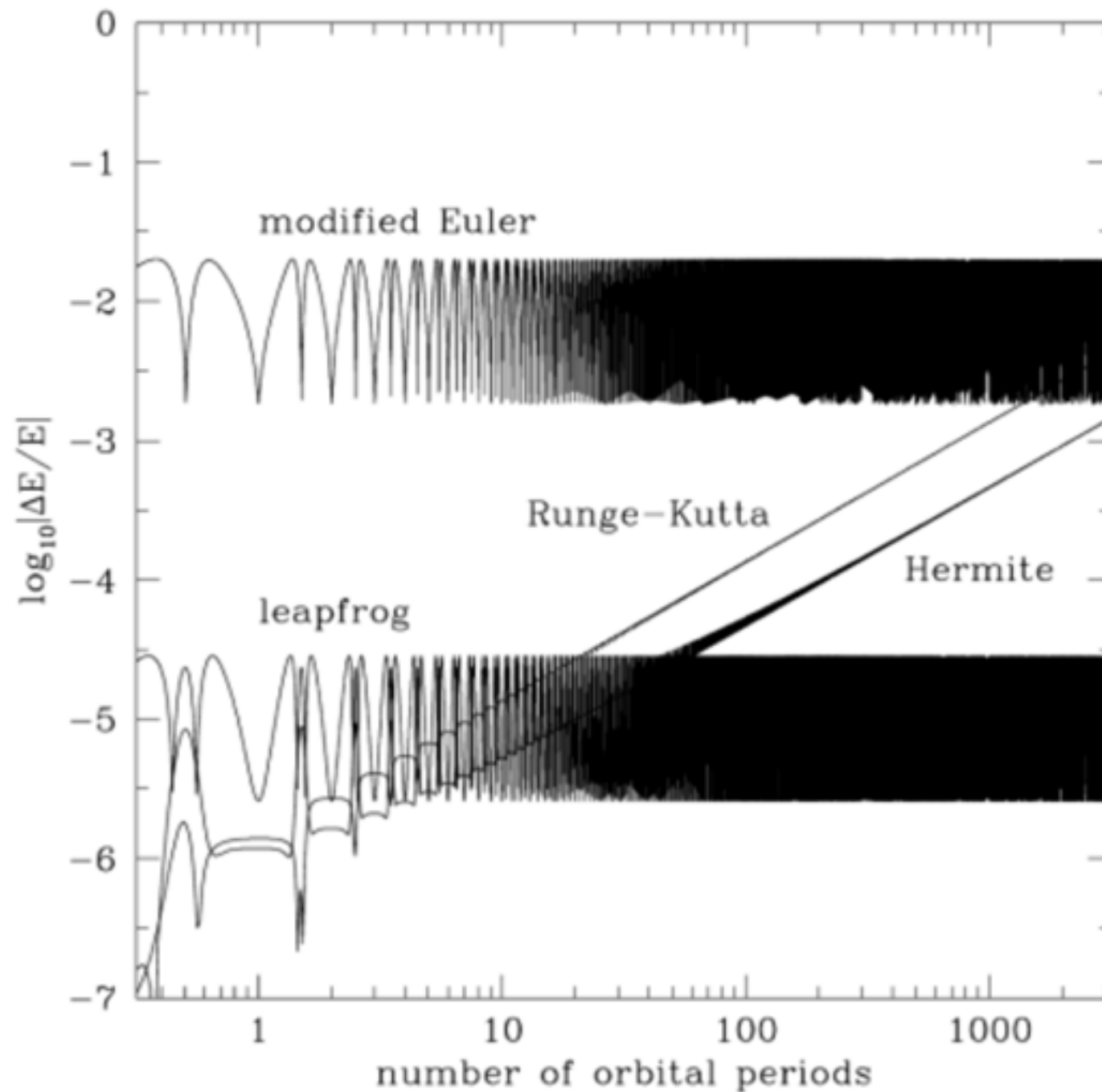
$$\begin{aligned} \dot{\mathbf{a}}_s = & -G \sum_{t=1}^N \frac{m_t \phi'(\mathbf{r}_{st}, \bar{h}_{st})}{|\mathbf{r}_{st}|} \mathbf{v}_{st} + 3G \sum_{t=1}^N \frac{m_t (\mathbf{r}_{st} \cdot \mathbf{v}_{st}) \phi'(\mathbf{r}_{st}, \bar{h}_{st})}{|\mathbf{r}_{st}|^3} \mathbf{r}_{st} \\ & - 4\pi G \sum_{t=1}^N \frac{m_t (\mathbf{r}_{st} \cdot \mathbf{v}_{st}) W(\mathbf{r}_{st}, \bar{h}_{st})}{|\mathbf{r}_{st}|^2} \mathbf{r}_{st} . \end{aligned}$$

# A simple example : A plummer sphere ( $N = 100$ )

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# Energy errors in N-body codes

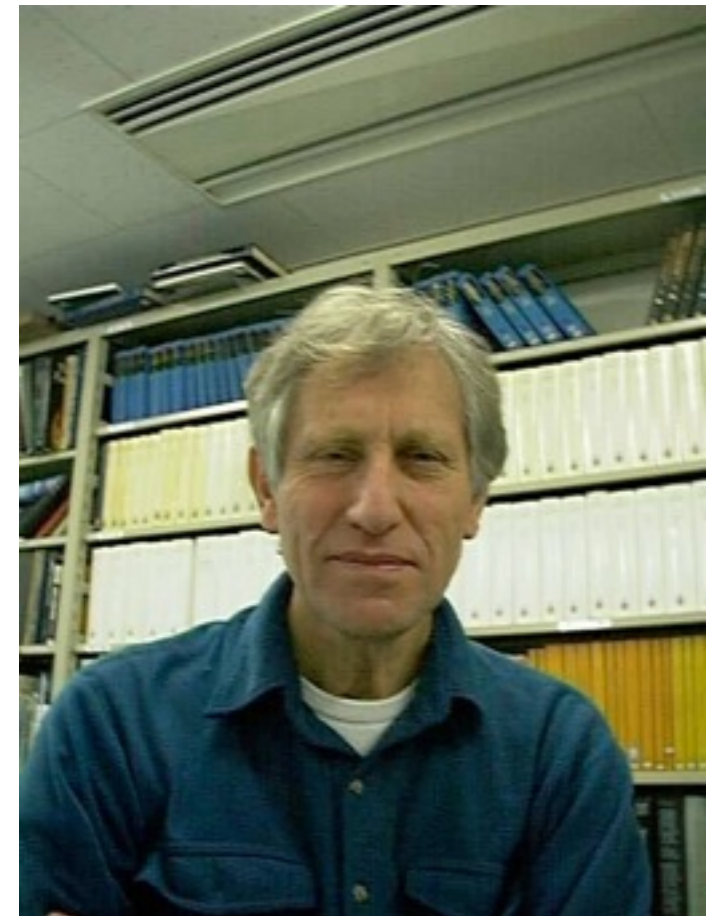




# What about 'Regularisation'?

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- **KS-Regularisation** is a powerful technique used in some N-body codes to :
  - (i) allow very accurate integration of very close 2-body encounters
  - (ii) therefore eliminate the need for softening/smoothing of grav. forces
- Some reasons not to use it
  - **Extremely** complicated
  - Hard to combine other physics (e.g. gas forces)
  - There are alternatives these days, not quite as accurate but much easier to implement
- Will I get hunted down by Sverre Aarseth if I don't use it??
  - Hopefully not



# Sub-systems

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- If binary or higher-order multiple systems form, then the simulation may progress slower and slower
  - **Spends a lot of CPU effort integrating the binary system** with short timesteps as the **rest of the simulation proceeds very slowly**
  - Most of the time, **the binary motion can be isolated and simulated as a separate system** (with or without external perturbations)
- If a binary is identified (as in the previous slide), then
  - Binary motion is integrated separately
  - Rest of simulation interacts with centre-of-mass of binary

# Hydrodynamics + N-body

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- GANDALF employs a hybrid scheme for modelling the evolution of a gaseous stellar cluster
  - Gas is modelled with SPH particles using 2nd order Leapfrog scheme
  - N-body particles are modelled with 4th-order Hermite scheme
  - Derived **coupling terms that maintains energy conservation**

# Possible challenges to hybrid scheme

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- **N-body codes usually require high accuracy** (e.g. total energy conserved to less than 0.001% accuracy), but hydro-codes usually operate with much higher error tolerances.

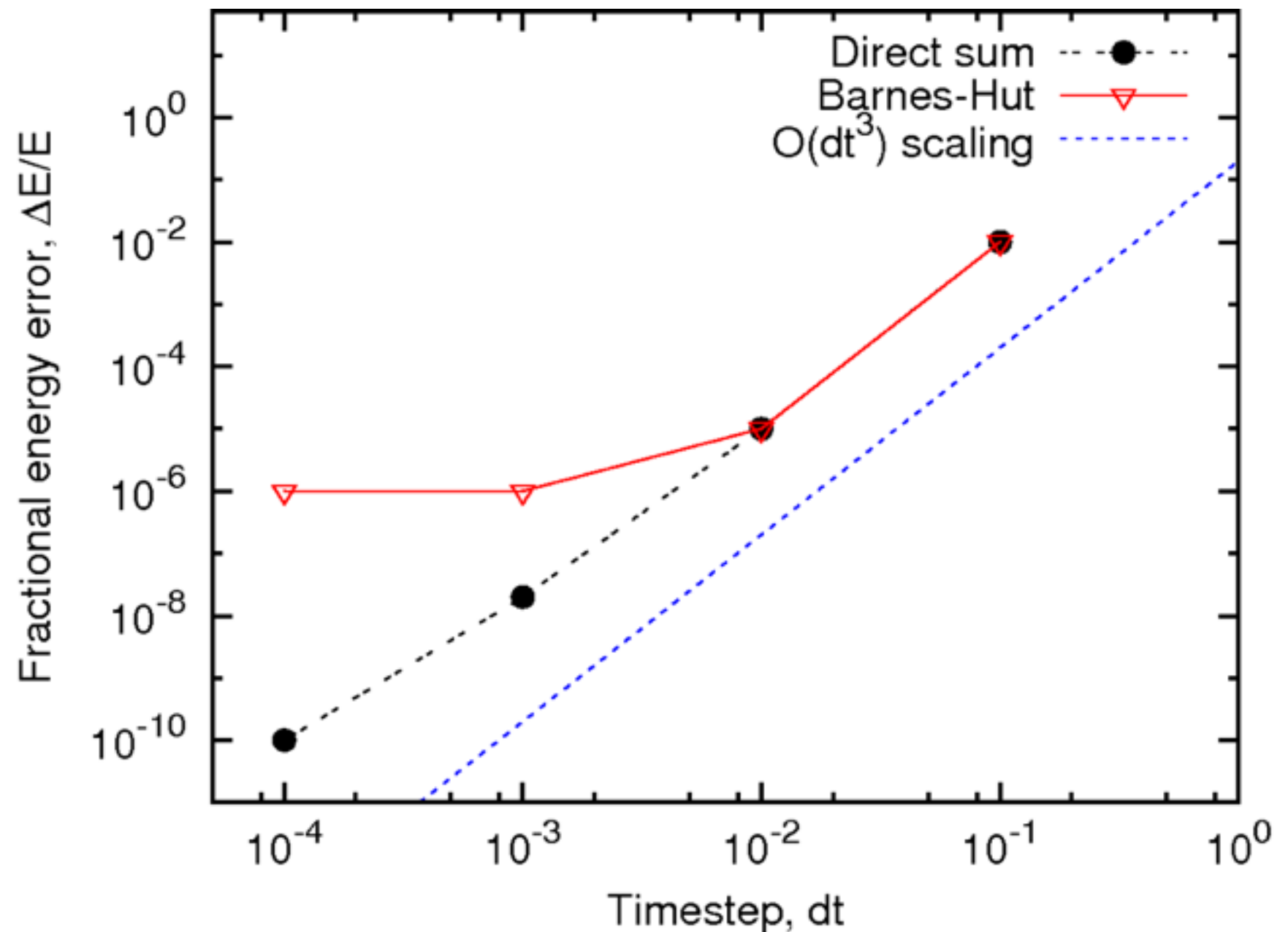
challenges. A simple workaround has been proposed by [200]. For each and every simulation the conservation of energy, momentum and angular momentum should be monitored. Reducing the time step size and increasing the force accuracy, say, if a tree is used for gravity, should improve the conservation properties. A correct code should ensure conservation to better than 1% over several thousand time steps.

Rosswog (2009)

- However, modern SPH schemes derived via Lagrangian mechanics can, in principle, conserve momentum, angular momentum and energy to rounding error given a robust integration scheme.

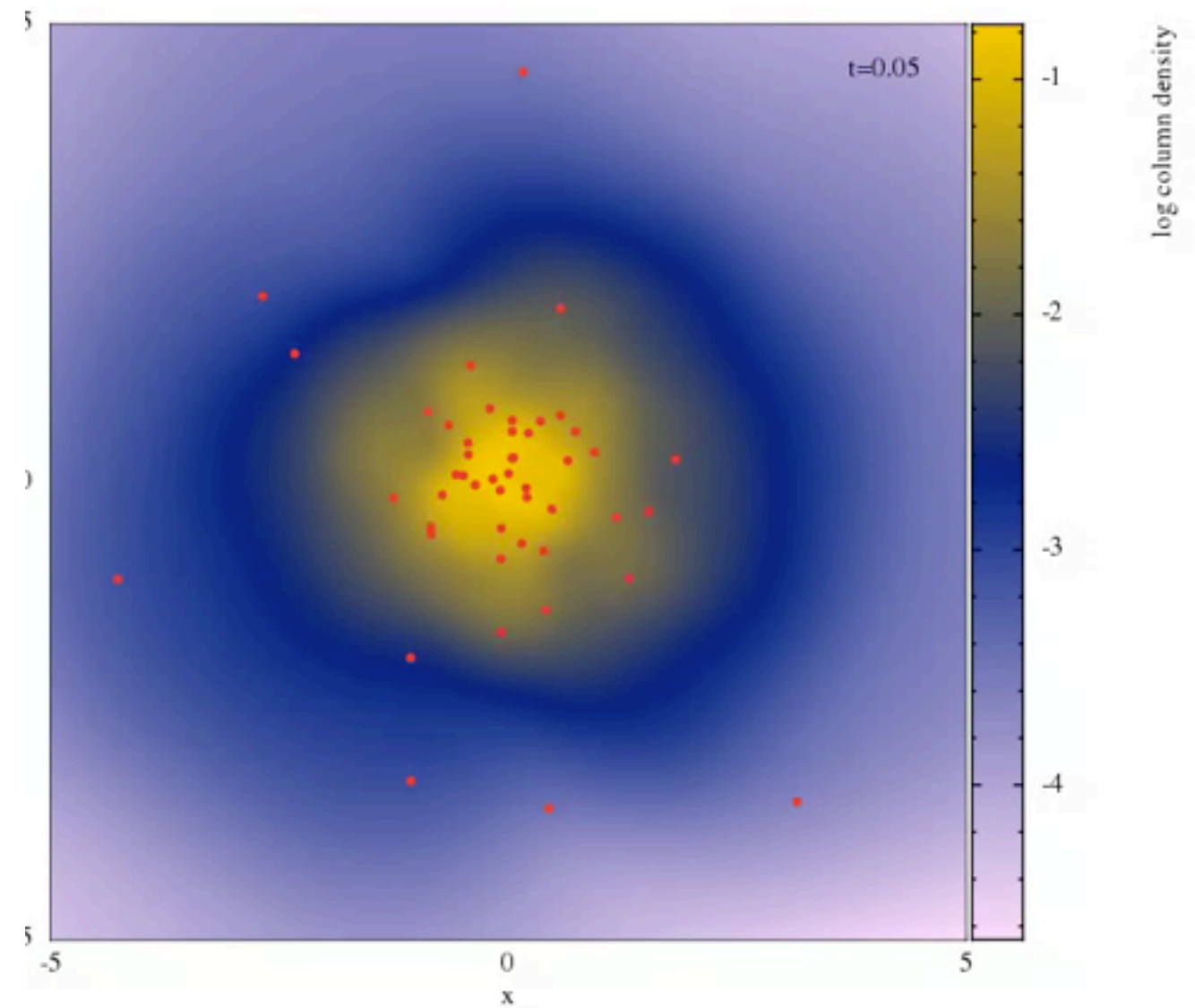
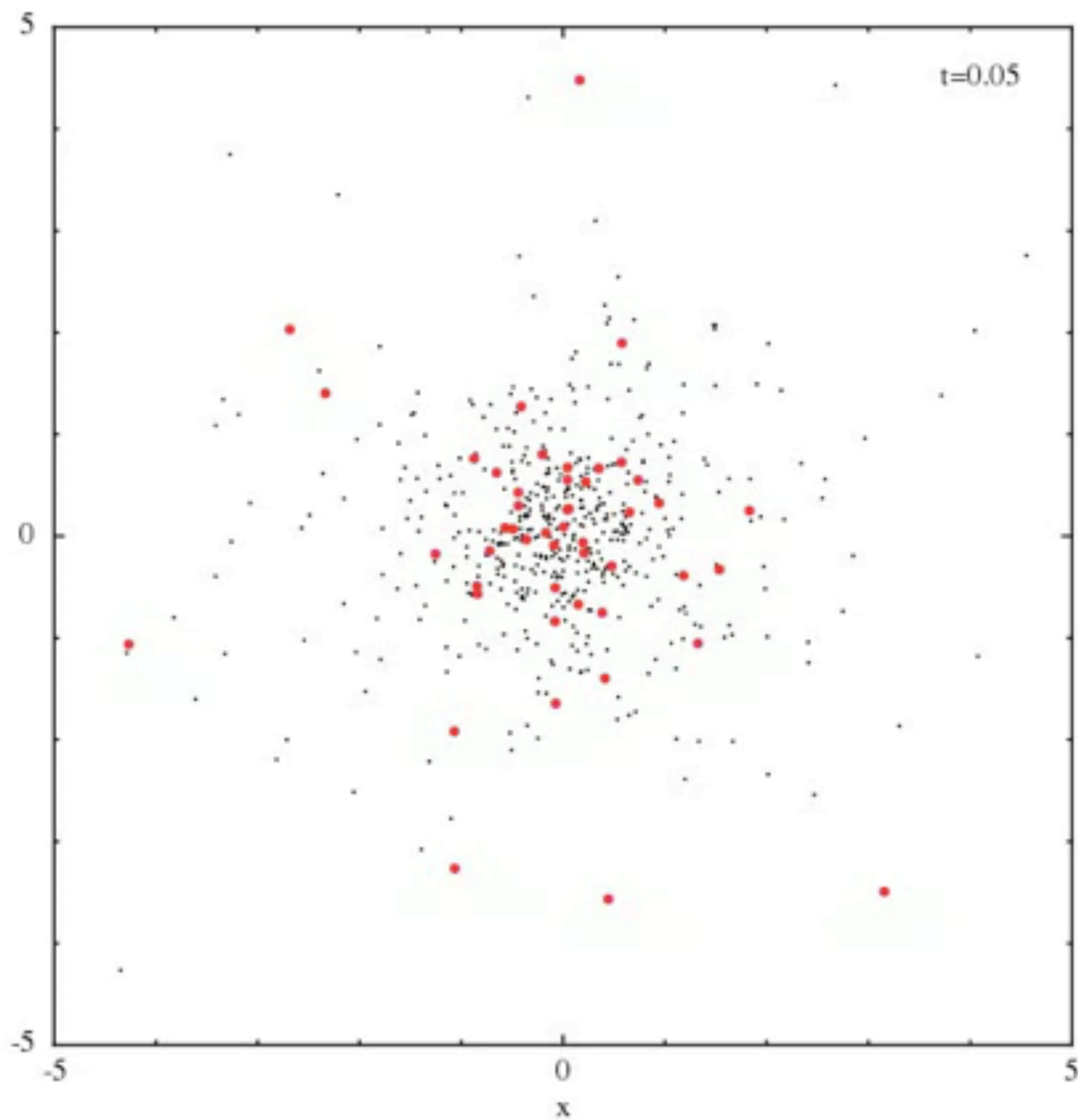
# Errors in SPH/N-body codes

- Integration (truncation) error
  - SPH - 2nd-order Leapfrog
  - N-body - 4th-order Hermite
- Block timesteps
- Gravity tree errors



# Gaseous Plummer spheres

- A Plummer sphere can be combined with a  $n=5$  polytrope to produce a stable ‘gaseous cluster’.



# Modelling star formation : Sink particles

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- Modelling how low-density gas collapses into stars is a very expensive process
  - Can perhaps investigate a single star in detail
  - **Almost impossible with current capabilities to model a cluster of fully formed stars**
- Bate, Bonnel & Price (1995) introduced **dynamical sink particles**, to mimic the formation of a star and to capture the effects of any subsequent accretion
  - Sinks are created like little black holes / vacuum cleaners that sweep up any gas that enters it
  - Allows simulations to run fast enough to follow large-scale cluster formation





# Sink particles : formation criteria

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- The choice of formation criteria is crucial for obtaining converged simulations
- We use the following criteria

- Exceeds a density threshold

$$\rho_i > \rho_{\text{SINK}}$$

- Gravitational potential minima

$$\phi_i < \text{MIN} \{ \phi_j \}$$

- Doesn't overlap with existing sink

$$|\mathbf{r}_i - \mathbf{r}_{s'}| > X_{\text{SINK}} h_i + R_{s'}$$

- There's an additional criterion which should be implemented soon

- Hills sphere criteria

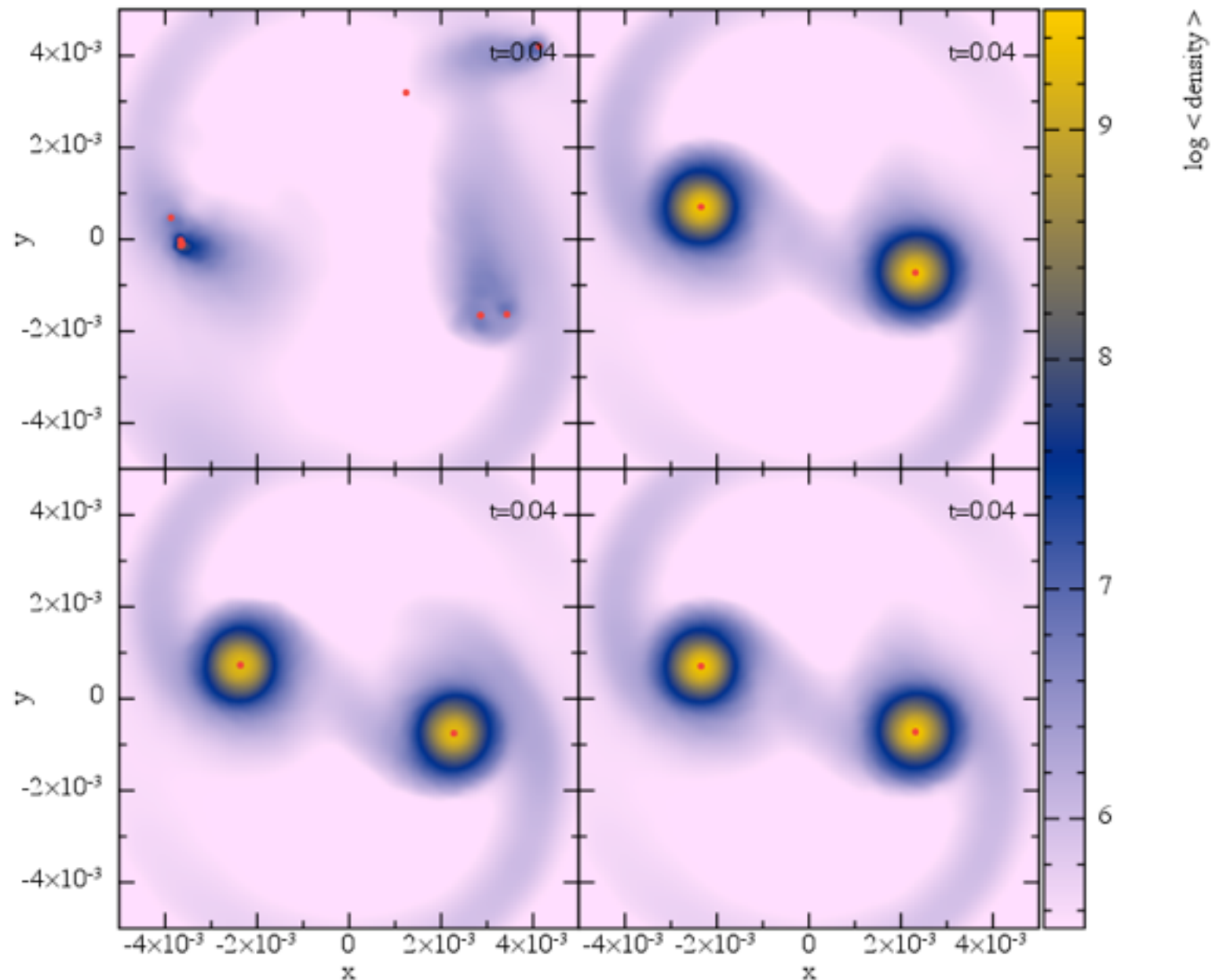
$$\rho_i > \rho_{\text{HILL}} \equiv \frac{3 X_{\text{HILL}} (-\Delta \mathbf{r}_{is'} \cdot \Delta \mathbf{a}_{is'})}{4 \pi G |\Delta \mathbf{r}_{is'}|^2}$$

# Sink particles : formation criteria

Density criterion

Low sink density

High sink density

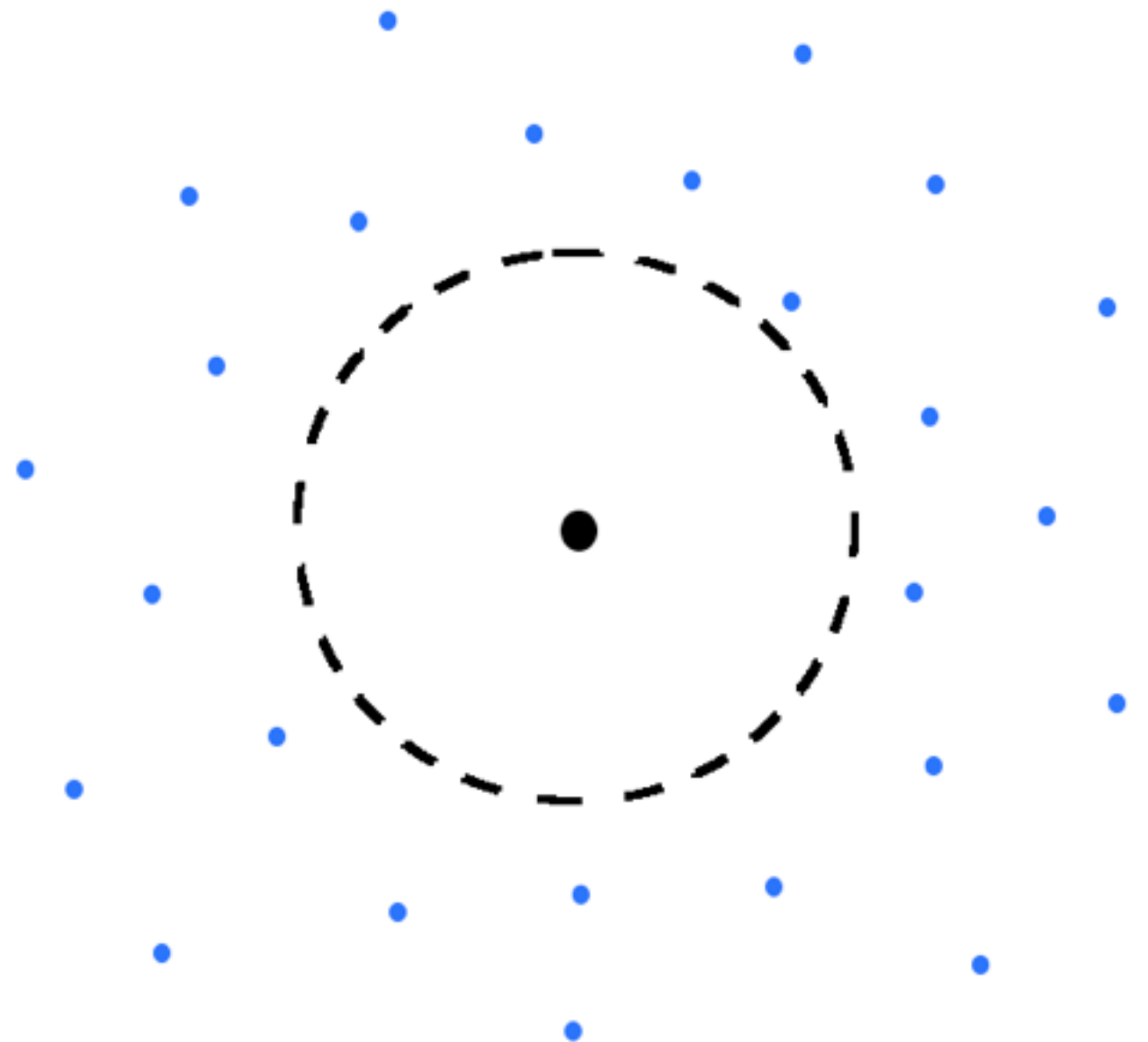


Density & potential  
minimum criteria

# Modelling accretion

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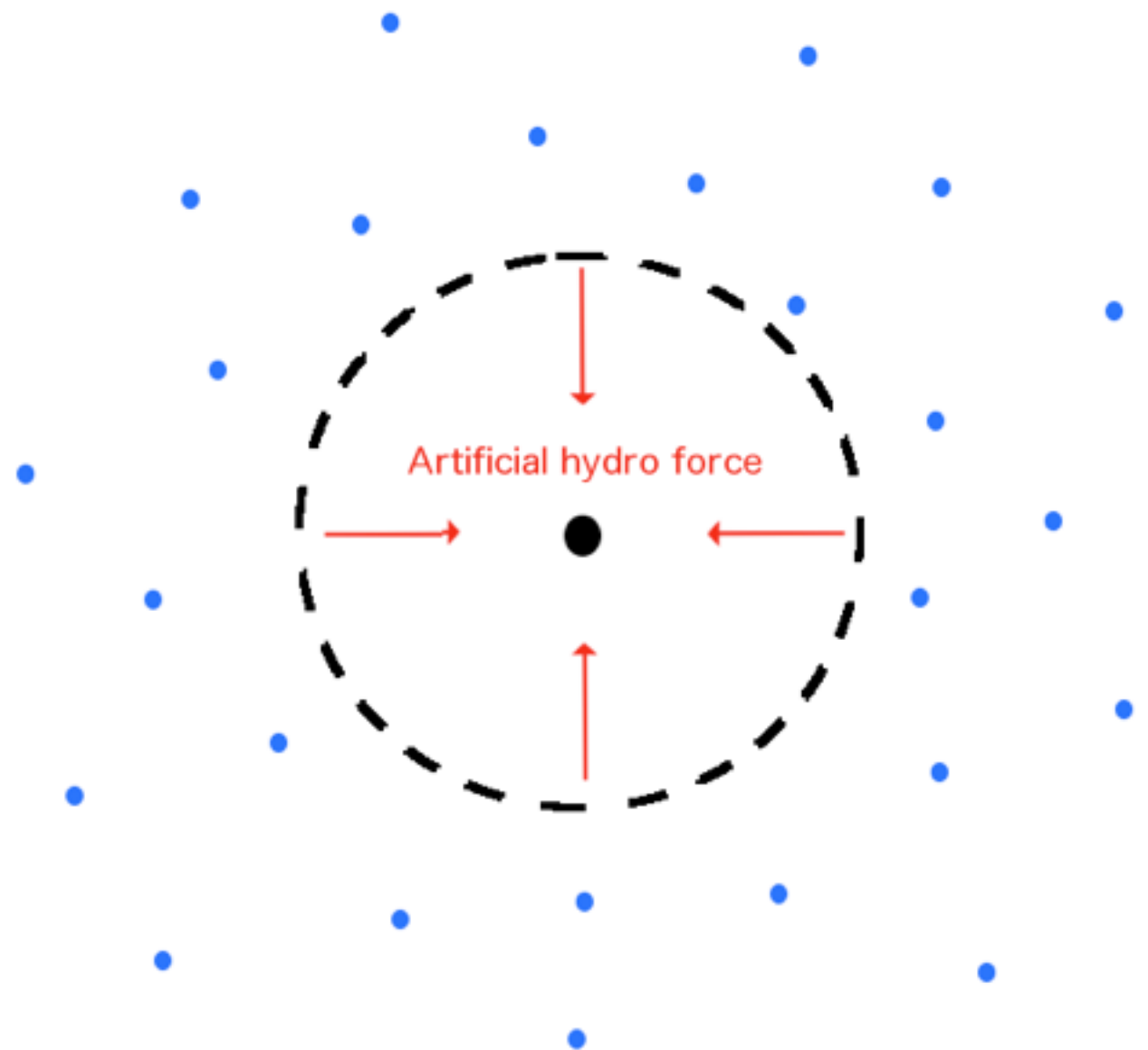
- Accretion is modelled by removing particles from the simulation that
  - Enter the sink accretion radius
  - Are gravitationally bound to the sink
- Generally leads to an empty 'exclusion' zone inside the sink that is devoid of any SPH particles



# Artificial boundary forces

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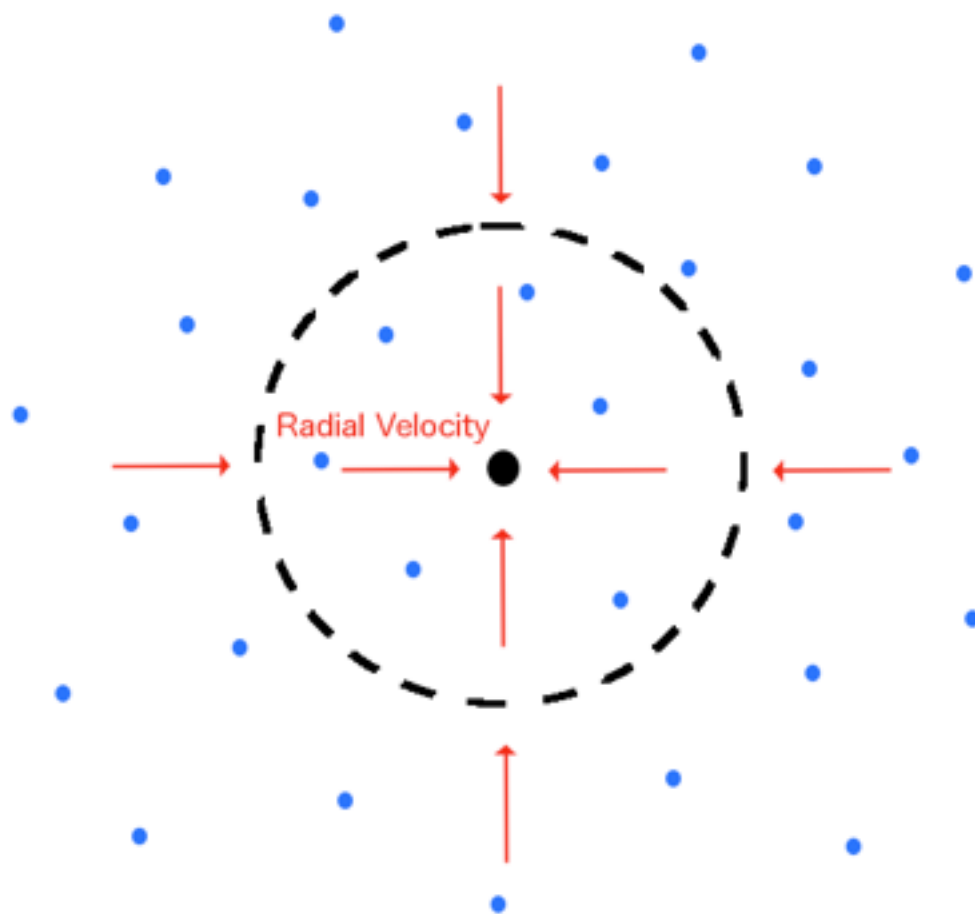
- Particles just outside the accretion radius see no neighbours inside the sink
  - Discontinuous sampling of density field
  - All SPH properties are incorrect, in particular the hydro forces
  - Leads to artificial outward pressure gradient, and therefore artificial inward hydro force
- BBP95 originally suggested using some correction terms to account for missing neighbours
  - Does not work so well and is not used any more (as far as I know)



# Spherical and disc accretion

- For sub-grid accretion model, we consider two limiting cases

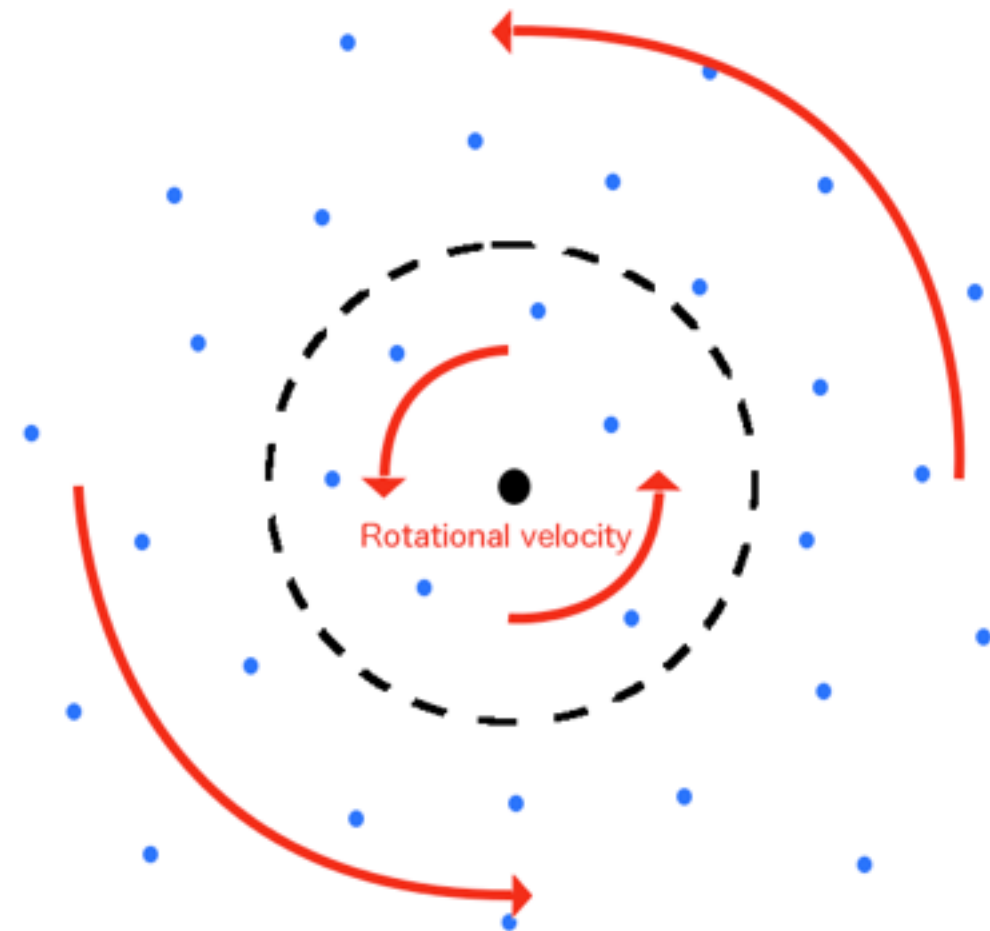
Spherical accretion



$$\dot{M}(r) = -4\pi r^2 \rho(r) v_{\text{RAD}}(r)$$

$$\langle t_{\text{RAD}} \rangle_s = \frac{\sum_j \{m_j\} W}{4\pi \sum_j \{|\Delta \mathbf{r}_{js}| \Delta \mathbf{r}_{js} \cdot \Delta \mathbf{v}_{js} m_j W(|\Delta \mathbf{r}_{js}|, H_s)\}}$$

Disc accretion



$$t_{\text{SS}} = \frac{(G M_* R_d)^{1/2}}{\alpha_{\text{SS}} a^2}$$

$$\langle t_{\text{DISC}} \rangle = \frac{(GM_s)^{1/2}}{\alpha_{\text{SS}} W} \sum_j \left\{ \frac{|\Delta \mathbf{r}_{js}|^{1/2} m_j W(|\Delta \mathbf{r}_{js}|, H_s)}{\rho_j a_j^2} \right\}$$

# Computing the accretion timescale

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- We compute the ratio of rotational energy to gravitational energy of particles inside the sink as an indicator of which limiting case is applicable

$$f = \text{MIN} \left\{ \frac{2E_{\text{ROT}}}{|E_{\text{GRAV}}|}, 1 \right\}$$

- If  $f = 1$ , particles are in **rotational equilibrium** :  $t_{\text{ACC}} \rightarrow \langle t_{\text{DISC}} \rangle_s$
- If  $f = 0$ , particle motion is **purely radial** :  $t_{\text{ACC}} \rightarrow \langle t_{\text{RAD}} \rangle_s$
- To deal with intermediate cases that also give the correct limiting behaviour, we use

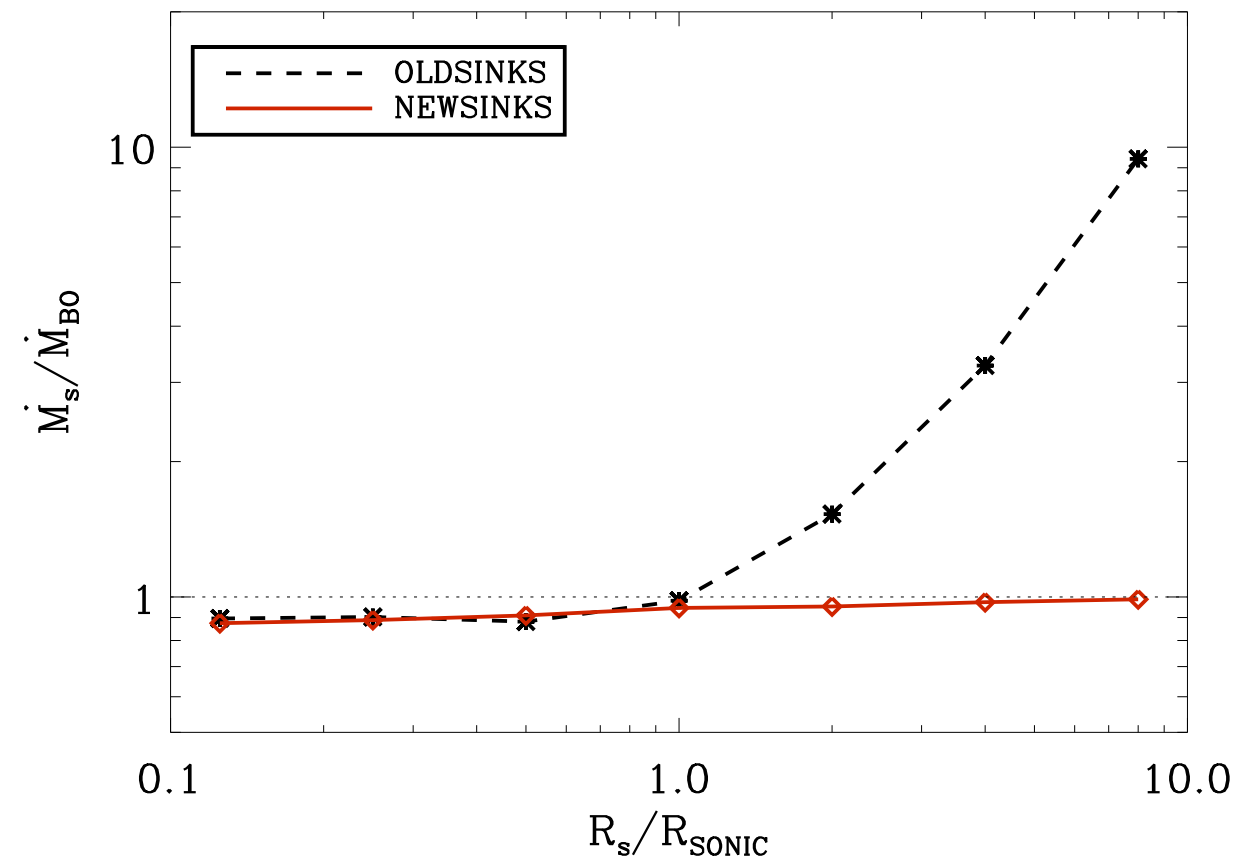
$$t_{\text{ACC}} = \langle t_{\text{RAD}} \rangle_s^{(1-f)} \langle t_{\text{DISC}} \rangle_s^f$$

- The mass accreted in the current timestep is then

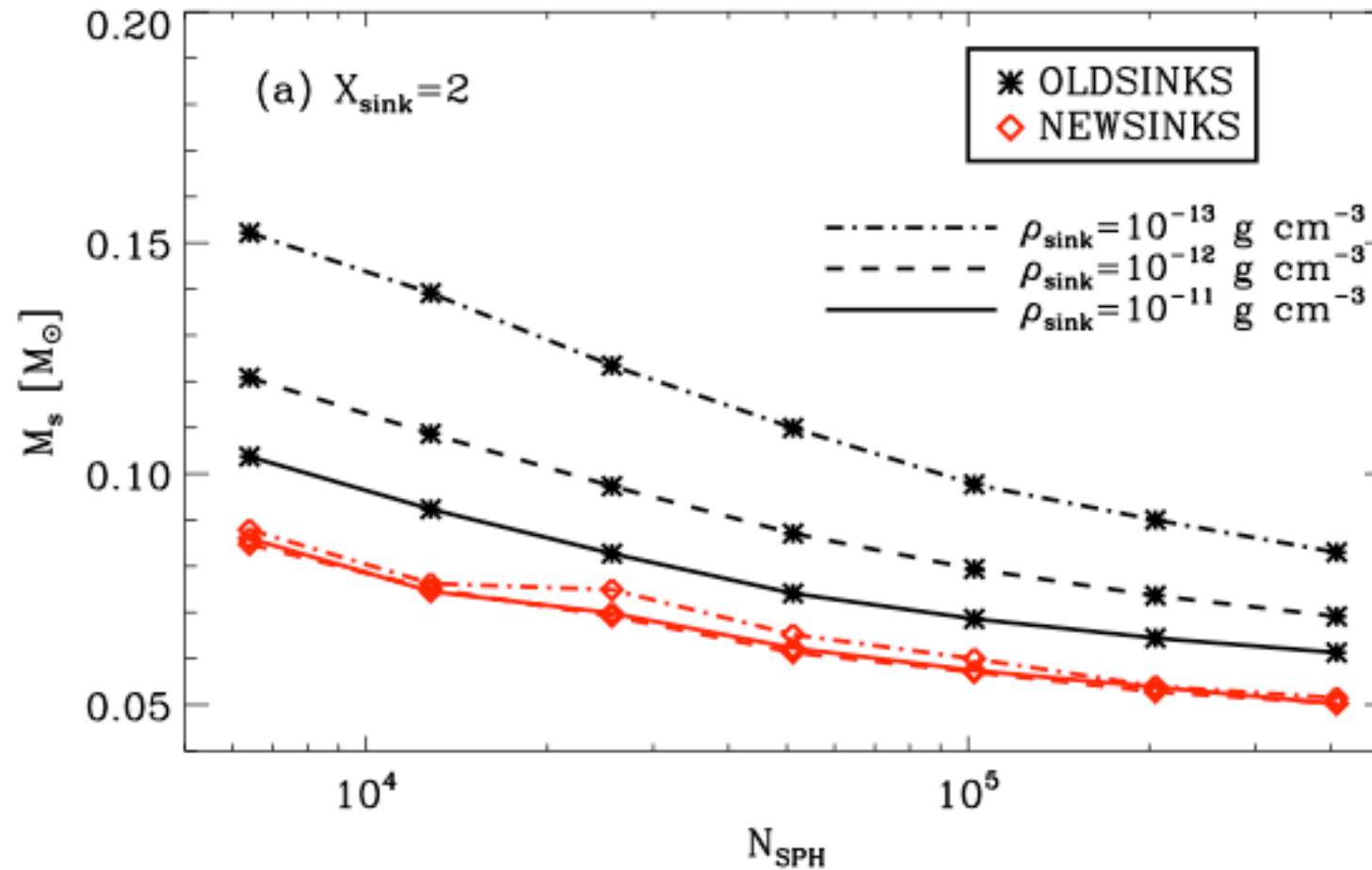
$$\delta M_{\text{ACC}} = M_{\text{INT}} \left[ 1 - \exp \left( -\frac{\delta t_s}{t_{\text{ACC}}} \right) \right]$$

# Bondi accretion (Spherical accretion)

- In Bondi accretion, the sonic point defines the radius where the inflow velocity is equal to the local sound speed
- For **large radii, the inflow is subsonic** (both hydro and gravity forces important)
- For small radii, the inflow is supersonic (only gravity important)
- Old sinks are correct for small radii since the lack of hydro forces is unimportant. For large radii, the lack of hydro forces leads to incorrect accretion rates
- New sinks give correct accretion rates for all sink radii
- Note : For monatomic gases, the sonic radius is zero. Therefore, for old sinks the accretion rate is always wrong.



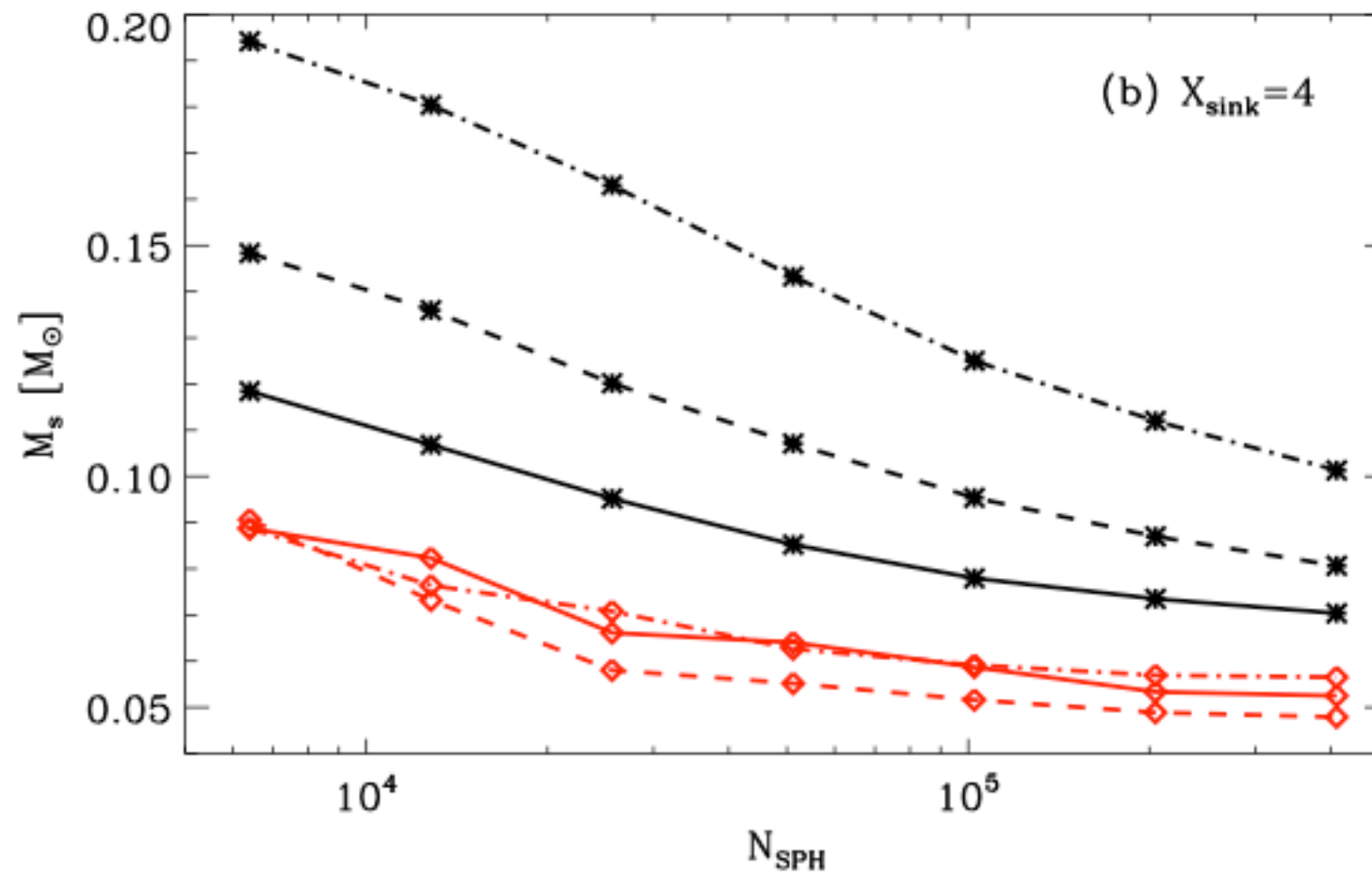
# Boss-Bodenheimer test : Convergence of sink properties



- For old sinks, the total mass contained in a sink varies greatly depending on the formation density, and hence the sink radius.
- For new sinks, although the results vary with resolution (external hydrodynamics), they are essentially independent of sink density/radius.



# Boss-Bodenheimer test : Convergence of sink properties



- For larger sinks (same formation density), old sinks have even larger masses, but new sinks are still converged at the same masses

# Future development

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- GANDALF will allow both collisional and collisionless N-body simulations (but far more optimised for collisionless)
- Collisional N-body will be optimised in the future, particularly with the sub-systems and binary integrators
- Sink particles currently only implemented in SPH schemes
  - Will be added to Meshless scheme soon