Moving mesh cosmology: the hydrodynamics of galaxy formation

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Accepted 2012 June 4. Received 2012 April 10; in original form 2011 September 16

ABSTRACT
We present a detailed comparison between the well-known smoothed particle hydrodynamics (SPH) code GADGET and the new moving-mesh code AREPO on a number of hydrodynamical test problems. Through a variety of numerical experiments with increasing complexity we establish a clear link between simple test problems with known analytic solutions and systematic numerical effects seen in cosmological simulations of galaxy formation. Our tests demonstrate deficiencies of the SPH method in several sectors. These accuracy problems not only manifest themselves in idealized hydrodynamical tests, but also propagate to more realistic simulation set-ups of galaxy formation, ultimately affecting local and global gas properties in the full cosmological framework, as highlighted in companion papers by Vogelsberger et al. and Keres et al. We find that an inadequate treatment of fluid instabilities in GADGET suppresses entropy generation by mixing, underestimates vorticity generation in curved shocks and prevents efficient gas stripping from infalling substructures. Moreover, in idealized tests of inside-out disc formation, the convergence rate of gas disc sizes is much slower in GADGET due to spurious angular momentum transport. In simulations where we follow the interaction between a forming central disc and orbiting substructures in a massive halo, the final disc morphology is strikingly different in the two codes. In AREPO, gas from infalling substructures is readily depleted and incorporated into the host halo atmosphere, facilitating the formation of an extended central disc. Conversely, gaseous sub-clumps are more coherent in GADGET simulations, morphologically transforming the central disc as they impact it. The numerical artefacts of the SPH solver are particularly severe for poorly resolved flows, and thus inevitably affect cosmological simulations due to their inherently hierarchical nature. Taken together, our numerical experiments clearly demonstrate that AREPO delivers a physically more reliable solution.

Key words: methods: numerical – galaxies: formation – cosmology: theory.

1 INTRODUCTION
Numerical simulations have become an indispensable tool for studying astrophysical phenomena. Over the last decade, the numerical accuracy and fidelity of simulation methods have undergone drastic improvements, both in terms of resolvable dynamic range and of the complexity of the physical processes that are now routinely incorporated into the codes. This remarkable progress in numerical techniques coupled with the ever increasing power of high performance computing platforms has led to major advances in a number of topics, such as star formation (Klessen, Krumholz & Heitsch 2011), accretion disc dynamics and associated jet phenomena (De Villiers & Hawley 2003; Gammie, McKinney & Tóth 2003), supernova explosions (Janka et al. 2007), black hole coalescence (Pretorius 2007) and cosmic structure formation (Springel et al. 2005).

Clearly, an in-depth understanding of many astrophysical processes, particularly those that are inherently non-linear, relies crucially on the accuracy and realism of the numerical approach. While the latter is largely determined by the degree of adequateness, comprehensiveness and consistency of the physical model assumed, the former depends sensitively on the discretization scheme adopted for the equations, which with increasing resolution should converge to

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the correct solution. None the less, even for some elementary physical problems, with known analytic solutions, simulation methods can sometimes produce poorly converged results, or even converge to a solution which is, however, different from the expected result (Springel 2010b).

In the context of hydrodynamical cosmological simulations, an important example is given by the seminal work of Frenk et al. (1999) (the Santa Barbara Comparison Project. This study performed a detailed comparison between 12 different simulation codes that tracked the non-radiative evolution of a forming galaxy cluster in a cold dark matter (CDM) cosmology. The initial conditions were generated independently by each group from a provided linear theory density or displacement field. Very different numerical methods, which can be broadly classified into smoothed particle hydrodynamics (SPH) and mesh-based techniques, employing also different gravity solvers and different effective resolutions, were then used by the groups to follow the formation and evolution of the target object. Reassuringly, the Santa Barbara Comparison Project has shown that the global properties of the simulated object, both in terms of dark matter and gas distribution, were in reasonable agreement among the codes. However, detailed gas properties of the Santa Barbara cluster, notably in the central region, exhibited a much poorer level of consistency. In particular, Frenk et al. (1999) pointed out that there is a systematic discrepancy in the central entropy profiles predicted by SPH and mesh-based codes, with the former producing power-law entropy profiles all the way to the centre and the latter yielding some form of entropy core.

Even though progress in numerical techniques has led to substantial improvements in current simulation codes compared to those considered in Frenk et al. (1999), the systematic difference in the predicted central cluster entropy still persists today, and its origin has not been fully understood thus far. Agertz et al. (2007) have performed a set of numerical experiments with SPH and mesh codes, where they have simulated the evolution of a cold, dense blob in a hot wind tunnel. By comparing the outcomes from different methods (and at various resolutions) against the characteristic blob disruption time-scale expected analytically, they conclude that in SPH codes the development of fluid instabilities can be numerically hampered. The suppression of dynamical fluid instabilities, such as Kelvin–Helmholtz, Rayleigh–Taylor and Richtmyer–Meshkov instabilities, leads to less efficient mixing of fluid elements with different specific entropy, and hence also inhibits entropy generation through mixing in the simulated system. In simulations of colliding isolated galaxy clusters, Mitchell et al. (2009) have shown that the central entropy profiles obtained with the SPH code GADGET and the mesh code FLASH show a similar level of discrepancy as found by Frenk et al. (1999), which they attribute to the different levels of mixing.

There have been a number of attempts to improve the description of mixing in SPH by modifying the standard set of discretized equations (see e.g. Price 2008; Wadsley, Veeravalli & Couchman 2008; Heß & Springel 2010) or by incorporating a Riemann solver in place of the artificial viscosity (e.g. Inutsuka 2002; Cha & Whitworth 2003; Murante et al. 2011). For example, Price (2008) has shown that the introduction of an artificial thermal conductivity term can improve the behaviour of fluid elements at contact discontinuities, which in turn leads to better developed Kelvin–Helmholtz instabilities. More generally, Wadsley et al. (2008) suggested that a physical modelling of heat diffusion is necessary when simulating high Reynolds number flows (both for SPH and mesh-based methods), and that by applying such an approach a flat entropy core is likely a more correct solution for non-radiative galaxy cluster simulations.

Differences in the hydrodynamical solver between SPH and mesh-based codes are possibly not the only reason for the systematically different central entropy profiles in the Santa Barbara Comparison Project. As mentioned above, the groups involved in this study did not use identical initial conditions, and did not perform their simulations at equal numerical resolutions and with identical gravity solvers, which opens up the possibility of additional sources for discrepancies. A more recent code comparison study by Heitmann et al. (2008) evolved uniform, dark matter only cosmological boxes in an ΛCDM universe with 10 different codes, starting from the same initial conditions. They showed that for large systems there is a reassuring agreement in the halo mass functions between the codes, as well as in the internal structures of haloes in the outer regions. However, the study by Heitmann et al. (2008) also revealed significant discrepancies for small haloes, demonstrating that the typical root grid resolution commonly adopted in adaptive mesh refinement (AMR) codes for simulations of cosmic structure formation is overly coarse and leads to a suppression of low-mass halo formation (see also O’Shea et al. 2005). This emphasizes the need to simultaneously strive for high accuracy both in the gravity solver and in the hydrodynamics. In addition to the high accuracy of the gravity and hydro solver, it is also of prime importance to adopt sufficiently high mass and spatial resolution for the studied problem at hand. For example, if besides pure hydrodynamics other physical processes, such as star formation and associated feedback, are modelled in cosmological simulations, it is necessary to resolve all star-forming haloes with a sufficiently large number of resolution elements to reach convergent results (Springel & Hernquist 2003b).

In the present work we perform a detailed comparison study between the widely used SPH code GADGET and the novel moving-mesh code AREPO on a number of hydrodynamical test problems with increasing levels of complexity. We have adopted a combination of existing test problems and newly devised numerical experiments1 in order to provide a clear link between the results of our test problems and those of full cosmological simulations, which are discussed in detail in our companion papers (Vogelsberger et al. 2012, hereafter Paper I) and (Keres et al. 2012, hereafter Paper II). Our work thus contributes to the understanding of the significant differences in baryon properties found in cosmological simulations between mesh-based and SPH techniques, both at a global level (see Paper I) and at the scale of individual galaxies (see Paper II).

A unique advantage of our comparison study lies in the fact that simulations with GADGET and AREPO can be started from identical initial conditions and that both codes employ the same gravity solver. Note that in case of AREPO gravitational softenings for the gas can be kept fixed as is the case of GADGET or can be computed adaptively determined by the cell size. In the numerical experiments with gas self-gravity we explore both fixed and adaptive gas softening, where in the case of adaptive softenings a floor equal to the GADGET gas softenings is set. We find that these different choices of gravitational softenings in AREPO do not affect our findings. This allows us to isolate cleanly how the different hydro solvers used by GADGET and AREPO impact gas properties. We first consider elementary hydrodynamical numerical tests such as a strong 1D Sod shock tube test, a 2D implosion test (which is widely used for benchmarking mesh codes, but has rarely been considered in SPH), and the ‘blob’ test (Agertz et al. 2007). We then perform a number of isolated or

1 High-resolution images and movies of various numerical experiments are available for download at website http://www.cfa.harvard.edu/itc/research/movingmeshcosmology

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merging halo simulations in the non-radiative regime, aimed at understanding how shocks and fluid instabilities affect their gaseous atmospheres. Finally, we study the differences between GADGET and AREPO in radiative simulations, where we follow inside-out disc formation and interactions between the central discs and orbiting substructures.

This paper is organized as follows. In Section 2, we provide a brief overview of the numerical codes used and the types of simulations performed. Section 3 represents the core of the paper where all of our numerical experiments are discussed. Finally, we summarize our findings in Section 4.

2 METHODOLOGY

2.1 Numerical codes

2.1.1 GADGET

GADGET is a massively parallel TREEPM-SPH code widely used in numerical astrophysics. In this study we adopt the latest GADGET-3 version. A detailed description of an earlier version can be found in Springel (2005). GADGET is fully adaptive in time and space, and in its entropy formulation for SPH (Springel & Hernquist 2002) manifestly conserves both energy and entropy in the absence of artificial viscosity. Gravitational forces are computed with an octree method (Barnes & Hut 1986; Hernquist 1987). To speed up the computation, long-range forces can be optionally evaluated with a PM method, with the tree being restricted to short-range forces only.

In all our tests, we adopt as standard value for the artificial viscosity strength $\alpha = 1.0$, and we use 64 neighbours for kernel interpolation in 3D simulation, unless we specifically vary these parameters to assess their effect.

2.1.2 Other SPH implementations

As mentioned in the Introduction, there have been a number of recent proposals to improve the standard SPH implementation in various ways, for example by invoking a time-dependent artificial viscosity (Morris & Monaghan 1997; Dolag et al. 2005), a modified density estimate (e.g. Ritchie & Thomas 2001; Heß & Springel 2010; Saitoh & Makino 2012), a decoupling of the hot and cold neighbours in multiphase flows (Marri & White 2003; Okamoto et al. 2003), an artificial thermal conductivity (Price 2008), a modified equation of motion (Heß & Springel 2010; Abel 2011), an explicit modelling of mixing (Wadsley et al. 2008), an enlarged neighbour number combined with a different kernel shape (Read, Hayfield & Agertz 2010), or a replacement of the artificial viscosity by a Riemann solver (e.g. Inutsuka & Makino 2012). While some of these modifications of the standard SPH formalism deliver more accurate results in targeted numerical experiments, no consensus has yet emerged whether any of these approaches (or a combination thereof) is sufficiently robust for cosmological applications and leads to universally more accurate results in galaxy formation simulations. Therefore, in this work we focus on the traditional SPH implementation rather than on the various possible modifications proposed recently. We note that this standard formulation of SPH has also been typically employed in state-of-the-art cosmological SPH calculations (e.g. Crain et al. 2009; Di Matteo et al. 2012). Furthermore, as discussed in Paper I, there are other, generic issues with SPH that have not been resolved by any of the above modifications. These issues ultimately mean that SPH does not currently have a formal convergence condition, which also complicates rigorous evaluations of variants of the standard SPH algorithm.

2.1.3 AREPO

AREPO (Springel 2010a) is a new massively parallel simulation code, which uses the same gravity solver as GADGET (augmented with the possibility of adaptive gravitational softenings for the gas), but employs a completely different method for the evolution of the fluid. It adopts a second-order accurate finite-volume technique, where the solution of the Euler equations is computed by an un-split Godunov method equipped with an exact Riemann solver. Throughout we use the default choice of the slope limiter in AREPO which prevents the linear reconstruction to over- or undershoot the maximum/minimum values of neighbouring cells, as described in detail in Springel (2010a). Unlike standard finite-volume codes used in numerical astrophysics, AREPO solves the equations on an unstructured Voronoi mesh, which is allowed to freely move with the fluid. The resulting quasi-Lagrangian nature of AREPO automatically guarantees spatial adaptivity and greatly reduces numerical diffusivity even in the presence of large bulk flows. Compared to standard Eulerian mesh codes, AREPO has the advantage of being fully Galilean invariant (as is GADGET as well), it is less prone to advection errors and over-mixing, and preserves contact discontinuities better.

In this study, we employ a mesh regularization method in AREPO by default, based on a Lloyd algorithm (for details see Springel 2010a), which guarantees that the geometric centre of each Voronoi cell is sufficiently close to the cell’s mesh-generating point to ensure good accuracy of the spatial reconstruction. In cosmological simulations (see Paper I), an alternative regularization criterion has proved to be advantageous, based on the maximum opening angle under which a cell face is seen from the mesh-generating point. We have checked for a number of test runs presented in this study that this alternative regularization method does not affect any of the results described here.

For most of the simulations presented here we do not use the possibility of mesh refinement and de-refinement operations (see Springel 2010a), except in our numerical experiments with star formation presented in Section 3.4.3, where we employ it for verification of our findings. Mesh de-/refinements are used to constrain the mass of cells to a small range around a target value (equal to the gas particle mass in the matching GADGET run). This restricts the spectrum of star particle masses which are generated from gaseous cells, and thus ensures that N-body heating effects are minimized. Also, as our default choice we use the energy formulation of AREPO. We verified for each numerical experiment that there is no significant spurious transfer of kinetic into thermal energy.

2.2 Types of simulations

2.2.1 Physical processes

We perform both radiative and non-radiative hydrodynamical simulations, where in the former case gas is represented by a primordial mixture of hydrogen and helium in an optically thin limit (Katz, Weinberg & Hernquist 1996). In simulations with radiative cooling, we employ a subresolution multiphase model for star formation.
and associated supernova feedback (Springel & Hernquist 2003a). We slightly modify the behaviour of this model for gas elements which are hot but already above the density threshold for star formation, by allowing them to settle quickly on to the effective equation of state: if their newly estimated temperature from radiative cooling would fall below the temperature of the multiphase medium, we set it equal to the multiphase temperature. This change has been introduced to make the subgrid star formation module consistent with its current implementation in AREPO. We also perform some simulations with the subresolution star formation model where spawning of star particles is intentionally prevented, but the cold, dense gas above the density threshold for star formation is still governed by the effective equation of state. These simulations are particularly useful for understanding the development of dynamical instabilities between cold and hot media, and have the advantage of not being prone to fragmentation which might affect pure cooling runs.

Note that we deliberately consider only very simplified baryonic physics implementations in all presented numerical experiments, as detailed above. This is for two reasons. First, we need to make sure that gas cooling and star formation processes are treated on an equal footing as possible in order to allow a meaningful comparison of the two codes. Thus, we adopt a simple multiphase model for star formation which is largely insensitive to the detailed structure of the gas on very small scales where gas joins the interstellar medium and is described by a relatively stiff effective equation of state. Secondly, our aim is to isolate in an as clean manner as possible the differences between simulated systems with GADGET and AREPO stemming from the discretization of the fluid equations alone. This goal is given precedence over trying to reproduce observational findings through a more sophisticated modelling of additional physics. While it is likely that such additional physical processes will change the properties of some of our simulated systems by possibly different degrees in GADGET and AREPO, it is of significant interest in its own right to disentangle numerical uncertainties from uncertainties in the physical modelling of star formation and associated feedback processes.

2.2.2 Gravitational softenings

For simplicity, many of the numerical experiments presented in this study have been evolved without gas self-gravity. For simulations where self-gravity of the gas is nevertheless included, we note that GADGET employs constant gravitational softening for gas particles in the manner of Hernquist & Katz (1989), while AREPO uses either the same constant gravitational softening or an adaptive softening determined by the cell size with a minimum softening value set to the fixed softening of the matching GADGET run. We have verified that this does not lead to substantial differences for any of the tests presented in this study.

2.2.3 Effective hydro resolution

Even though GADGET and AREPO calculations use the same gravity solver and can be initiated from identical initial conditions, due to the very different nature of the hydro solver it is not straightforward to define unambiguously a comparison strategy at the same or equivalent hydro resolution. In this study we choose to keep the number of resolution elements the same in the both codes, i.e. to have the same number of SPH particles as Voronoi cells, corresponding to a comparable mass resolution in the gas. This allows us to adopt indeed the same initial conditions and to also have similar mass resolution in the stellar component in those simulations where star formation is included. Furthermore, the CPU costs of the two codes are then roughly comparable, as discussed in Paper I.

We note, however, that this choice implies that the ‘effective’ spatial resolution of GADGET is lower than that of the matching moving-mesh calculation, given that fluid properties in SPH are evaluated by kernel averaging over a somewhat larger number of neighbours than needed in AREPO for its stencil of neighbouring cells, for example for gradient estimates. For this reason, we perform all of our numerical experiments at a number of different resolutions, which also help us to gain some insight into the convergence properties of the two codes. None the less, it is important to stress that the convergence properties of the SPH method are still not well understood. For example, one would ultimately require that the number of neighbours be increased with increasing total particle number (see e.g. Rasio 2000; Read et al. 2010; Robinson & Monaghan 2011, and discussion in Paper I), but the appropriate scaling of the neighbour number with increasing resolution is currently unknown. It is common practice, which we adopt as well, to simply always keep the number of neighbours fixed when the total particle number is varied, even though the discretized representation of the density field is not guaranteed to converge to its underlying smooth distribution with this choice.

2.3 Initial conditions generation

For all the tests presented in this study (except for the “blob” experiment of Agertz et al. 2007, for which we take publicly available initial conditions), we generate the initial conditions in terms of possibly different degrees in GADGET and AREPO, it is of significant interest in its own right to disentangle numerical uncertainties from uncertainties in the physical modelling of star formation and associated feedback processes.

3 RESULTS

3.1 Strong shocks and interacting curved shocks in multidimensions

3.1.1 Strong shock in 1D

As an introductory problem, we consider a strong shock with a Mach number of $M = 6.3$ in one dimension. The initial conditions have been constructed as follows: in a computational domain of length $L_x = 10.0$, $N = 200$ particles (cells) have been placed on a regular grid such that for $x < 5.0$ the pressure and density are $P = 30.0$ and $\rho = 1.0$, while for $x \geq 5.0$ they are $P = 0.14$ and $\rho = 0.125$, respectively. We adopt an adiabatic index of $\gamma = 1.4$, and
assume that the fluid is initially at rest. In the test run with GADGET, we adopt a standard value of the artificial viscosity equal to $\alpha = 1.0$, and we vary the neighbour number $N_{\text{ngb}}$ by setting it to 5, 7, 11 or 15, appropriate for the 1D nature of the test.

In Fig. 1, we show the gas density, velocity, entropy (i.e. $P/\rho^\gamma$) and pressure at time $t = 0.25$ for GADGET with $N_{\text{ngb}} = 5$ (left-hand panels), and for AREPO (middle panels). Blue symbols give the values of individual particles/cells, dashed red lines represent the initial conditions, while dotted red lines are the analytic solution to this Riemann problem. It can be seen that in both GADGET and AREPO the post-shock properties of the fluid are captured well, but the shock and the contact discontinuity are significantly broader in GADGET which also shows a characteristic ‘pressure blip’ at the contact discontinuity (see also Springel 2010b). Contrary to what one may perhaps suppose, the post-shock oscillations present in GADGET are not caused by inadequate artificial viscosity, but are instead induced by an inaccurate treatment of the sharp contact discontinuity of the initial conditions. To demonstrate this point, we have performed exactly the same shock tube test but this time smoothing the initial contact discontinuity over five particles in density and internal energy with a Hann window function, so as to reduce its sharpness. Green cross symbols in the third row of Fig. 1 (see also the right-hand panels where we zoom into the region around the shock) illustrate how the gas entropy is affected by this choice of smoothed initial conditions. Post-shock oscillations and the ‘spike’ in entropy at the contact discontinuity are greatly reduced.

It is interesting to note that also in the case of AREPO the ‘spike’ in the entropy at the contact discontinuity is cured by our smoothed initial conditions. This feature is absent in the results of standard grid codes, as we have verified by running this test with a static mesh option in AREPO (and by running it with ENZO), which tends to broaden contact discontinuities (depending on the advection speed) and thus largely wash out this feature. This can be seen in the lower right-hand panel of Fig. 1 where we show the results from the AREPO run with a static mesh with magenta triangle symbols. The much lower numerical diffusivity of AREPO for contact discontinuities preserves the initial start-up feature to much higher accuracy, but at the same time this can lead to a larger ‘wall heating’ effect (Rider 2000) than in static grid codes, which tend to smooth out at some level the initial start-up errors at the contact discontinuity (see also the description of the Noh problem in Springel 2010a).

When we increase the number of neighbours from $N_{\text{ngb}} = 5$ (as shown in Fig. 1) to 15 in the shock tube tests performed with GADGET (but keeping the total number of particles constant), the spatial region over which the contact discontinuity and the shock are broadened increases progressively, but the pressure jump between the post-shock and pre-shock gas outside of the broadened region remains the same, yielding consistent Mach numbers. Based on these tests, we later discuss in more detail the consequences of...
shock broadening in \textit{gadget} when simulating the radial infall of gas into static dark matter haloes in Section 3.3.2.

3.1.2 Interacting shocks in 2D

While 1D shock tube tests are an essential basic benchmark for hydrodynamical code performance, they are far less demanding than multidimensional flow problems where complex interactions of non-planar shocks may occur, as is the case in realistic structure formation simulations. Such tests have however not been examined widely in the SPH literature thus far. We hence perform the so-called ‘implosion test’ in two dimensions (Hui, Li & Li 1999).\footnote{See also ‘Comparison of several difference schemes on 1D and 2D test problems for the Euler equations’ by R. Liska and B. Wendroff on http://www-troja.ifi.cuni.cz/~fiska/CompareEuler/compareE and http://www.astro.princeton.edu/~jstone/Athena/tests/} To set up this test problem, we select a computational domain \(L_x = L_y = 0.3\) with \(N_x = N_y = 200\) particles or cells, respectively (we have also run higher resolution versions with \(N_x = N_y = 400\) and 800). The initial pressure and density are \(P = 1.0\) and \(\rho = 1.0\) for \(x + y > 0.15\), while \(P = 0.14\) and \(\rho = 0.125\) otherwise. The adiabatic index is \(\gamma = 1.4\) and the fluid is initially at rest. We performed this test with \textit{gadget}, adopting an artificial viscosity of \(\alpha = 1.0\) and \(N_{\text{high}} = 22\) smoothing neighbours (suitable for this 2D test). While previously this test has been considered using reflective boundary conditions, we here adopt periodic boundary conditions due to their more straightforward implementation in SPH codes.

In Fig. 2, we show density maps for simulations performed with \textit{gadget} (left-hand panels) and with \textit{arepo} (right-hand panels), at times \(t = 0.1, 0.3, 0.5\) and 0.7 for \(N_x = N_y = 200\). The complex, evolving gas density structure is caused by the continuous interaction of shocks throughout the computational domain. Initially, due to the discontinuity in density and pressure along \(x + y > 0.15\), a planar mild shock front develops perpendicular to the \(x = y\) diagonal travelling towards the origin. Given that we have adopted periodic boundary conditions, the gas will interact on all four sides of the simulated box, and in particular, interacting shock waves in the lower left-hand corner result in the formation of a narrow jet along the \(x = y\) diagonal (which is clearly visible only in the \textit{arepo} run). As the travelling shocks accelerate the fluid in the regions of density discontinuity a Richtmyer–Meshkov instability develops, as manifested by ‘mushroom-like’ features in the gas distribution.

While the global gas density distribution in Fig. 2 agrees reasonably well in \textit{gadget} and \textit{arepo} runs (i.e. with respect to the shape of the regions with different densities and the magnitude of the density differences), there are several significant differences: (i) shocks and contact discontinuities are much sharper in \textit{arepo}, in agreement with our findings in Section 3.1.1; (ii) the density distribution in \textit{gadget} appears not only smoothed out but it is also noisier, as evidenced by the graininess of the density maps, which is caused by intrinsic noise in multidimensional flows in SPH (Springel 2010b); (iii) most striking, perhaps, is the quite different appearance of the low density gas in the bottom left corner – a narrow, extended, dense jet along the diagonal is largely absent in \textit{gadget} (due to the broadening of the contact discontinuity) and Richtmyer–Meshkov instabilities along discontinuities are suppressed. We have verified that higher resolution simulations with \(N_x = N_y = 400\) and 800) result in sharper shocks and contact discontinuities (with a somewhat feeble jet forming), but they cannot satisfactorily cure the absence of well-developed Richtmyer–Meshkov instabilities. This is in agreement with previous studies (e.g. Agertz et al. 2007; Springel 2010b) indicating fundamental limitations of the SPH method in its widely used form. We also note that the moving mesh in \textit{arepo} does an excellent job of maintaining symmetry across the \(x = y\) diagonal. Furthermore, the level of numerical diffusion of discontinuities is very low, as indicated by the narrowness and length of the jet.

In addition to the density fields it is also instructive to analyse the vorticity distribution in the implosion test, as shown in Fig. 3. To construct vorticity maps we either (i) first compute spatially adaptive velocity field maps for each component on a uniform grid (in our case only \(x\) and \(y\) components) and then finite difference them to obtain the curl or (ii) we compute spatially adaptive vorticity maps, where the curl has been evaluated directly in the code. For \textit{gadget} runs, vorticity maps computed with method (i) are shown in the upper panels of Fig. 3 for \(N_x = N_y = 200\) and \(N_x = N_y = 800\). Using method (ii) vorticity maps look essentially the same. In the case of \textit{arepo} (shown in the lower panels for the same run with \(N_x = N_y = 200\)) the vorticity maps are also very similar when computed with method (i) or (ii). However, regardless of the exact details of the vorticity map generation there are marked differences in the vorticity field between \textit{gadget} and \textit{arepo}. For the \(N_x = N_y = 200\) run, the vorticity map in \textit{gadget} is largely featureless and noisy, with artificial suppression of vorticity generation at locations where surfaces of constant density and constant pressure are not aligned (i.e. baroclinic source term \(\propto \nabla \rho \times \nabla p\)). Strikingly, even in the case of the high-resolution \textit{gadget} run with \(N_x = N_y = 800\) (upper right-hand panel) vorticity generation in the regions with large baroclinic term is largely suppressed compared to the \textit{arepo} simulation, even though the latter has a much smaller number of resolution elements.

It is also worth noting that the overall magnitude of the vorticity is relatively high in \textit{gadget} with respect to \textit{arepo} in the regions which should be characterized by having low vorticity. This is due to the noise alluded to above, visible both in the density and vorticity maps. This noise is inherently present in multidimensional SPH simulations. It is caused by inaccurate pressure gradient estimates and is particularly evident in the subsonic regime (for further details see e.g. Springel 2010b; Bauer & Springel 2012). The jitter in gas velocities caused by the noisy gradient estimates introduces an artificial vorticity ‘floor’ throughout the simulated volume of the box. The poor treatment of vorticity in \textit{gadget} has also direct consequences for the effectiveness of the widely used Balsara switch (Balsara 1995) for artificial viscosity which relies on evaluation of velocity divergence and curl.

Hence, from the results of the ‘implosion test’ we conclude that while \textit{gadget} can accurately capture the main fluid properties in the case of dynamically interacting shocks with complicated geometries, there are systematic biases in detailed aspects, related to the development of fluid instabilities and to the level of fluid mixing at different entropies. These unwanted features lead to the suppression of angular momentum transport by vortices and to the damping of turbulence in the wake of curved shocks (see also Bauer & Springel 2012).

3.2 Dynamical fluid instabilities

3.2.1 Blob experiment

Our findings from the implosion test (Section 3.1.2) indicate that there are fundamental differences in the treatment of fluid instabilities between \textit{gadget} and \textit{arepo}. To investigate this issue in detail, we first consider the so-called ‘blob’ test as proposed by Agertz...
Figure 2. Implosion test in two dimensions at times $t = 0.1, 0.3, 0.5$ and 0.7. Left-hand panels: GADGET ($\alpha = 1.0, N_{\text{ngb}} = 22$). Right-hand panels: AREPO moving mesh. For each row, the density scale is the same in the left-hand and right-hand panels, covering the density range $\rho_{\text{gas}} = 1.2 - 0.4$.

e et al. (2007), which has been analysed with many different codes by now (Agertz et al. 2007; Heß & Springel 2010; Murante et al. 2011). The idea behind this test is to simulate the evolution of a dense cold blob immersed in a hot windtunnel, mimicking in a simplified way the motion of a satellite galaxy through the intracluster medium (ICM; see also Heß & Springel 2012). If the relative velocity between the blob and the surrounding medium is supersonic, a bow shock will develop in front of the blob. Additionally, dynamical
Figure 3. Vorticity maps of the implosion test in two dimensions at time \( t = 0.3 \). Upper panels: GADGET runs at two different resolutions with \( N_x = N_y = 200 \) (left-hand panel) and \( N_x = N_y = 800 \) (right-hand panel). For both resolutions, vorticity maps are computed by finite differencing the velocity field. Lower panels: vorticity maps of the AREPO run with \( N_x = N_y = 200 \) computed by finite differencing the velocity field (left-hand panel) and by calculating vorticity in the code based on a discretized curl operator directly applied to the Voronoi cells (right-hand panel).

instabilities (mostly of Kelvin–Helmholtz and Rayleigh–Taylor types) will grow in the subsonic part of the flow between the bow shock and the surface of the blob (see Agertz et al. 2007, for a detailed description). These instabilities will greatly influence the evolution of the blob, leading to its eventual disintegration on the characteristic Kelvin–Helmholtz time-scale set by the initial conditions.

To simulate this problem, we adopt the initial conditions used in the original Agertz et al. (2007) paper, which are publicly available.\(^4\) The simulated domain consists of a periodic box with extensions \( L_x = 2000 \text{ kpc} \), \( L_y = 2000 \text{ kpc} \) and \( L_z = 8000 \text{ kpc} \), and the blob is initially placed at \( x_c = y_c = z_c = 1000 \text{ kpc} \). The blob has a radius of \( R_{\text{blob}} = 197 \text{ kpc} \), and it is 10 times colder and denser than the surrounding medium, such that pressure equilibrium is ensured. The density and temperature of the external medium are \( \rho_{\text{medium}} = 3.13 \times 10^{-8} \text{ M}_\odot \text{ kpc}^{-3} \) and \( T_{\text{medium}} = 10^7 \text{ K} \), respectively, and it is moving with a constant velocity \( v_{\text{medium}} = 1000 \text{ km s}^{-1} \) along the \( z \)-axis. The adiabatic index is set to \( \gamma = 5/3 \). Following Agertz et al. (2007), we define the characteristic Kelvin–Helmholtz time-scale as \( t_{\text{KH}} = 1.6 t_{\text{cr}} \), where \( t_{\text{cr}} \) is the blob crushing time defined as \( t_{\text{cr}} = 2R_{\text{blob}}(\rho_{\text{blob}}/\rho_{\text{medium}})^{1/2}/v_{\text{medium}} \). For these initial conditions, the characteristic Kelvin–Helmholtz time-scale is \( t_{\text{KH}} \sim 1.98 \text{ Gyr} \).

We have performed the blob test at three different resolutions with GADGET and AREPO. The resolutions used are \( 32 \times 32 \times 128 \), \( 64 \times 64 \times 256 \) and \( 128 \times 128 \times 512 \). The two higher resolutions correspond exactly to the resolution of the simulations used in Agertz et al. (2007), while we constructed the initial conditions for the lowest resolution run by subsampling.

In Fig. 4 we show projected surface density maps of a thin slice \((\Delta y = 100 \text{ kpc})\) centred around the blob position. Left-hand panels illustrate GADGET runs at times \( t = t_{\text{KH}} \), \( t = 2 t_{\text{KH}} \) and \( t = 3 t_{\text{KH}} \), while the right-hand panels are for the simulations with AREPO in the moving-mesh mode. The position and the shape of the bow shock are rather similar between GADGET and AREPO runs, especially at early times, but the shock is broader and less crisp in GADGET. In the run with GADGET, the blob acquires a cap-like appearance caused by the internal shock which compresses it, and it undergoes continuous ablation due to the low-pressure region which forms in the wake of the blob (Agertz et al. 2007). While initially the blob evolution is similar in AREPO, after \( t = t_{\text{KH}} \) well-developed Kelvin–Helmholtz and Rayleigh–Taylor instabilities lead to an efficient shredding of the blob, mixing it with the external medium. In Fig. 5 we also show projected surface density maps centred around the

\(^4\) http://www.astrosim.net/
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3.3 Shocks and fluid instabilities in isolated halo models

3.3.1 Isolated haloes in hydrostatic equilibrium

To explore more directly the impact of shocks and fluid instabilities in the context of structure formation, we have devised a number of idealized test problems involving isolated halo models. Here we briefly describe how we set up and verify hydrostatic equilibrium configuration of the initial conditions, which form the backbone for a series of numerical experiments discussed in Sections 3.3 and 3.4.

The initial conditions are constructed by populating static or live dark matter potentials with gas particles, whose positions are drawn randomly. For the dark matter distribution we assume a Hernquist profile (Hernquist 1990) so that we can easily generate self-consistent models when we use live haloes.

blob position for the $32 \times 32 \times 128$ and $64 \times 64 \times 256$ runs at $t = 3t_{\text{KH}}$.

We quantify the time evolution of the blob in Fig. 6. Here we show the remaining blob mass fraction as a function of time, where the material associated with the blob is selected such as to satisfy $T < 0.9T_{\text{medium}}$ and $\rho > 0.64\rho_{\text{blob}}$, as has been done in previous studies. We compare GADGET and moving-mesh AREPO simulations, performed at three different resolutions (as indicated on the legend). While for our highest resolution runs there is a broad agreement between GADGET and AREPO for $t < t_{\text{KH}}$, blob mass fractions are systematically different afterwards. This is exactly when the transition in the mass-loss rate from the ablation-dominated to the fluid instability-dominated regime occurs. In the latter regime, AREPO clearly delivers a physically more trustworthy solution. The moving-mesh AREPO simulation agrees qualitatively very well with the outcome of Eulerian grid codes studied in Agertz et al. (2007). However, as noted by Springel (2011), there is a small but systematic difference with the moving-mesh code delivering slightly higher remaining blob mass for $t > 1.5 \times t_{\text{KH}}$. As expected, the GADGET simulation results agree well with the findings of Agertz et al. (2007) for the other SPH codes. This indicates that the inaccuracies we find for GADGET are inherent to the standard SPH method, and have prompted a number of works suggesting possible improvements to the SPH method (e.g. Price 2008; Wadsley et al. 2008; Heß & Springel 2010; Saitoh & Makino 2012, see also Section 2.1.2).
Figure 5. Projected surface density maps in units of $M_\odot$ kpc$^{-2}$ at $t = 3 t_{\text{KH}}$ for the low (top row) and intermediate-resolution simulation (bottom row) with GADGET and AREPO. The thickness of the slices is $\Delta y = 100$ kpc and they are centred on $y_c = 1000$ kpc. Even at our lowest resolution, AREPO captures the dynamical evolution of the cold blob much more accurately.

Figure 6. The remaining blob mass fraction as a function of time in units of $t_{\text{KH}}$. GADGET and AREPO results at three different resolutions are shown, as indicated in the legend.

profile traces the dark matter at large radii but is slightly softened in the centre, i.e.,

$$\rho_{\text{gas}}(r) = \frac{M_{\text{vir}}}{(2\pi a^3)(x + x_0)(x + 1)^5},$$

where $M_{\text{vir}}$ is the virial mass of the system, $a$ is the Hernquist scale length parameter, $x = r/a$, and $x_0 = 0.01$ is the softening scale length parameter for the gaseous halo. For the simulations without any net rotation, the initial gas velocities are set to zero, while for the simulation with non-vanishing angular momentum we assign gas velocities such that the halo is characterized by the dimensionless spin parameter

$$\lambda = \frac{|J|}{\sqrt{2} GM_{\text{vir}}^5/2},$$

where $J$ represents the angular momentum, $E$ is the total energy of the halo, and we assume solid body rotation.

For validation purposes, we evolve isolated haloes with GADGET and AREPO for 2.45 Gyr with gas self-gravity and no radiative losses. These test runs confirm that the gas is in very good hydrostatic equilibrium within the dark matter potential. The differences in gas density, temperature and entropy between GADGET and AREPO are within a few per cent throughout the whole halo at the final time in the case of static dark matter haloes (see Fig. A1 in Appendix A; for live haloes see Section 3.5).

We also observe that even though the initial gas velocities are zero (in the case with $\lambda = 0$) some small gas velocities develop over time ($\sigma_{\text{gas,3D}} \sim 30$ km s$^{-1}$). This is primarily caused by the initial Poisson sampling of gas positions, which implies an initial state that is not perfectly relaxed. While this numerical artefact could be avoided by explicitly relaxing the initial conditions, we note that these residual gas velocities do not have any significant bearing on our results: the total gas kinetic energy $E_{\text{kin}}$ is less than 0.005 of either the total potential or the total internal gas energy at the final time, as shown in Fig. A2 in Appendix A. It is, however, interesting to note that while the total $\sigma_{\text{gas}}$ is very similar between GADGET and AREPO, there are some systematic differences in the radial profiles of $\sigma_{\text{gas}}$, which are caused by dissipation of gas motions on different spatial scales (see the bottom panel of Fig. A1).

3.3.2 Radial collapse of cold gas in a static dark matter halo

We now analyse how differences in shock capturing between GADGET and AREPO affect the radial infall of gas in dark matter haloes,
a problem of direct cosmological interest. For this purpose we intentionally adopt a set-up as simple as possible in order to isolate differences between the codes driven by the shock treatment only. We initially set up gas in hydrostatic equilibrium and at rest within a static Hernquist potential with mass $M_{\text{halo}} = 10^{14} M_\odot$, scale length $a = 176 \text{kpc}$ and a gas fraction of $f_{\text{gas}} = 0.17$. We consider both an analytic gravitational potential and a potential with a centrally softened core. We introduce a small modification in the codes such that gas self-gravity is switched off – the gas only feels the static dark matter potential and hydrodynamical forces in a purely non-radiative regime. We then artificially reduce the internal energies of gas particles/cells so as to displace the gas from the equilibrium solution. The newly assigned gas temperature is $4.7 \times 10^3 \text{K}$, equal for all resolution elements. This test is hence analogous to the well-studied Evrard collapse (Evrard 1988), but it is even simpler in nature because we intentionally neglect gas self-gravity.

As a consequence of the dramatic reduction in its temperature, the gas will suddenly lose pressure support and radially free-fall towards the centre. As the gas collapses, a radial shock develops in the centre, steepening while it propagates outwards and ploughing through the remainder of the outer material which is still falling in. The Mach number of the shock varies over the range $\sim 3-8$, well matched to the 1D shock tube test problem described in Section 3.1.1. Finally, as the shock propagates beyond the virial radius of the halo, the gas will reach a new hydrostatic equilibrium solution within the static dark matter potential. In Fig. 7, we show radial profiles of gas density, temperature and entropy computed at time $t = 0.1 \text{Gyr}$ after the start of the gas collapse for dark matter haloes with softened potentials. For each code ($\text{GADGET}$, $N_{\text{gas}} = 64$, $\alpha = 1.0$: blue lines; $\text{AREPO}$: red lines) we perform three runs at different resolutions: $N_{\text{gas}} = 10^4$ and $r_{\text{soft}} = 14.0 \text{kpc}$ (dotted lines), $N_{\text{gas}} = 10^5$ and $r_{\text{soft}} = 6.5 \text{kpc}$ (dashed lines) and $N_{\text{gas}} = 10^6$ and $r_{\text{soft}} = 3.0 \text{kpc}$ (continuous lines). Note that the $r_{\text{soft}}$ values indicate the spatial scale over which we smooth the analytic dark matter potential. They do not necessarily represent the minimum spatial resolution of these simulations, given that the gas is not self-gravitating and that the dark matter halo is rigid. The smoothing lengths of gas particles in the central region are of the order of $r_{\text{smhl}} \sim 1 \text{kpc}$ in $\text{GADGET}$ for the lowest resolution run with $N_{\text{gas}} = 10^4$, while the typical central cell sizes in $\text{AREPO}$ are a few hundred pc.

From Fig. 7 it can be seen that there are systematic differences in the gas properties predicted by the simulations with $\text{GADGET}$ and $\text{AREPO}$. In particular, the gas entropy distribution is broader in $\text{GADGET}$ both in pre- and post-shock gas in all three simulations, while for the two lower resolution runs there is a slight mismatch in the exact position of the shock front and in its strength (similar to the findings for the Evrard collapse in Springel 2010a), which is minimized in the case of $N_{\text{gas}} = 10^6$ particles. Note that the differences in shock front position between different resolution simulations for a given code are not driven by resolution effects but by different spatial softening of the central potential, which affects the gas collapse in the innermost regions. We have checked this explicitly by performing runs with $N_{\text{gas}} = 10^4$, $10^5$ and $10^6$ for both codes, but this time simulating cold gas collapse within an analytic Hernquist potential, as shown in Fig. 8. In this case the shock properties for different resolution runs are almost identical for a given code, indicating that the simulation with $N_{\text{gas}} = 10^4$ resolution elements is in principle sufficient for capturing the shock position accurately. None the less,  

$^5$ We modify the analytic Hernquist potential by convolving it with the spline softened potential in the centre, as we describe in detail in Section 3.5.

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**Figure 7.** Radial profiles of gas density, temperature and entropy at time $t = 0.1 \text{Gyr}$ for $\text{GADGET}$ (blue lines) and for $\text{AREPO}$ (red lines). For each code we show three different resolution runs: $N_{\text{gas}} = 10^4$ and $r_{\text{soft}} = 14.0 \text{kpc}$ (dotted lines), $N_{\text{gas}} = 10^5$ and $r_{\text{soft}} = 6.5 \text{kpc}$ (dashed lines) and $N_{\text{gas}} = 10^6$ and $r_{\text{soft}} = 3.0 \text{kpc}$ (continuous lines). Vertical black lines with the same style indicate the softening scales of the dark matter potential (note that this does not correspond strictly to the spatial resolution limit because the gas is not self-gravitating). The black vertical dot–dashed line denotes the virial radius of the system.
regardless of the resolution, differences between GADGET and AREPO in the pre- and post-shock gas persist also in the case of analytic Hernquist potentials.

The differences in entropy content of pre- and post-shock gas in GADGET and AREPO are in part due to the large shock broadening in SPH, as discussed in Sections 3.1.1 and 3.1.2. In fact, if we adopt $N_{\text{ngb}} = 32$ (which is considered the minimum number still permissible in 3D simulations) instead of $N_{\text{ngb}} = 64$, the gas entropy profile in GADGET becomes less broad, but is still not as sharp as in AREPO, indicating that simulations with a larger number of particles are needed in GADGET than in AREPO to recover shock features with the same accuracy.

Moreover, there is another numerical effect leading to spatially different entropy generation in GADGET: in the converging subsonic part of the flow, artificial viscosity (as implemented in GADGET) leads to artificial dissipation which increases the entropy in the pre-shock gas. While this feature is clearly visible in fig. 40 of Springel (2010a) for the case of the Evrard collapse, here we see that it also enlarges the central entropy in GADGET. The reason for this is the following: as soon as the gas is brought out of equilibrium and starts free-falling towards the centre, the gas entropy will be boosted in the central region due to an active artificial viscosity in the converging flow, creating an entropy bump that extends up to several kpc (or even several tens of kpc in the case of potentials with large cores) away from the centre, even though the shock has not fully formed yet at this point. This artificial entropy generation is much smaller in AREPO. As the shock forms and propagates outwards, it will lead to additional physical dissipation of much higher magnitude, bringing the entropy profiles of GADGET and AREPO into better agreement. Interestingly, in the case of the Evrard collapse, the initial difference in central entropy profiles is minimized with time due to the gas self-gravity (as we checked explicitly by running a simulation with exactly the same gas configuration but with gas self-gravity and without static dark matter potential), while it persists in our test runs even when the new equilibrium solution of the system is reached. The central entropy is higher in GADGET by a factor of $\sim 1.2$ and $\sim 1.5$ for analytic and softened potentials, respectively. Thus, it follows that differences in the gas properties due to different spatial dissipation of kinetic energy in GADGET and AREPO can be aggravated in the case of non-self-gravitating gas.

Note, however, that the difference in the central entropy profiles between the two codes goes in the opposite direction to what is found in non-radiative simulations of hierarchically forming galaxy clusters, where the central entropy is higher in mesh-based calculations. This indicates that accretion shocks during cosmological structure formation do not seem to be the likely cause of this central entropy discrepancy.

### 3.3.3 Infall of two gaseous spheres in a static dark matter halo

We now further increase the complexity of the problem by considering two gaseous spheres instead of one, collapsing into one common static dark matter halo placed in-between the two spheres. Each sphere has an initial spatial displacement from the centre of the halo. This test is similar in spirit to a number of previous works which analysed collisions of two galaxy clusters in isolation (see e.g. Ricker & Sarazin 2001; Ritchie & Thomas 2002; McCarthy et al. 2007; Springel & Farrar 2007; Mitchell et al. 2009; ZuHone 2011), but here we devise a cleaner set-up in order to minimize additional possible numerical effects (e.g. gravitational $N$-body heating and differences due to gas self-gravity). As in Section 3.3.2, we simulate a static analytic Hernquist dark matter halo (with exactly the same parameters) but instead of one cold gaseous sphere we generate two identical cold spheres separated by 1.2 Mpc along the $x$-axis, and we again neglect any radiative losses and gas self-gravity. The gaseous spheres are constructed in the same way as in Section 3.3.2, i.e. gas is first set up to be in hydrostatic equilibrium within a static Hernquist dark matter potential, and then its internal energy is reduced to $4.7 \times 10^4$ K. For each code we perform runs with different gas particle numbers, i.e. $N_{\text{gas}} = 10^2$, $N_{\text{gas}} = 10^3$ and $N_{\text{gas}} = 10^4$ per sphere.

Under the gravitational pull from the central dark matter potential the two cold spheres collapse towards its centre and violently collide. The interesting aspect of this problem is that radial symmetry is broken and the gas interaction results in much more complicated shock geometries. This is illustrated in Fig. 9, where we show a time sequence of projected gas density maps (first two columns) and mass-weighted entropy maps (last two columns) for runs with $N_{\text{gas}} = 2 \times 10^4$ performed with the two codes. While the GADGET simulation shows qualitatively similar gas structures, the detailed properties substantially differ.

Initially, as the spheres start to fall in, the gas is compressed in the centre of the halo, generating a spherical overdensity, which is somewhat broader and less peaked in GADGET, largely due to a poorer effective spatial resolution and non-negligible artificial viscosity (see Section 3.3.2). Also, during this initial stage, more entropy is produced in the central region in our SPH calculation. As more gas falls in, a shock develops which rapidly assumes a cocoon-like geometry elongated perpendicular to the direction of the collapse (see top panels of Fig. 9). From the gas density maps it can be seen that the shock front is narrower and sharper in AREPO, whereas in GADGET it has a more splotchy-like appearance, caused by kernel averaging. High entropy plumes propagating outwards along the $y$-axis are clearly visible in the right-hand panels, while the central entropy in GADGET within $\sim 250$ kpc is still higher than in the moving-mesh simulation. As the cocoon propagates against the infalling material, gas in the very centre is pushed perpendicular to the $x$-axis, generating a dense sheet-like region (see the second row of Fig. 9). Dynamical fluid instabilities at the boundary of this
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Figure 9. Projected surface density maps (first two columns; in units of $M_\odot \text{ kpc}^{-2}$) and mass-weighted entropy maps (last two columns; in internal units) for GADGET and AREPO simulations, showing the collision of two gaseous spheres with $N_{\text{gas}} = 2 \times 10^6$ at times $t = 0.54 \text{ Gyr}$ (first row), $t = 1.2 \text{ Gyr}$ (second row), $t = 2.9 \text{ Gyr}$ (third row) and $t = 15 \text{ Gyr}$ (fourth row). The plotted spatial domain is $0.6 \times 0.6 \times 1 \text{ Mpc}$.

dense region induce typical mushroom-like morphologies (cap-like in projection). However, even for our highest resolution simulations with $N_{\text{gas}} = 2 \times 10^6$ particles, there are some marked differences between the two codes. Mushroom-like features (corresponding to the red–orange colours in the density maps) originating at the very boundary of the dense central region are more coherent in GADGET, while in AREPO they break up and mix more efficiently with the surrounding medium. This also leads to the more efficient mixing of different entropy gas in the very core in the moving-mesh simulation, as can be seen from the entropy maps.

The differences in the fluid properties at the early stages of the simulated system as described above are, however, not the only reason why the thermodynamic properties of the gas are systematically discrepant between the two codes when a new equilibrium state is reached. With time, dense shells of gas completely disperse by mixing with the surrounding material in AREPO, while in the case of GADGET filaments and blobs of dense gas survive and gradually sink back to the centre (see the third row of Fig. 9). This buoyantly driven deposition of low entropy material in GADGET causes even larger differences between the final entropy distributions, which are illustrated in Fig. 10 (see also the bottom row of Fig. 9). The radial entropy profiles are computed once the system has reached hydrostatic equilibrium at time $t \sim 15 \text{ Gyr}$ from the start of the simulation. The blue lines denote GADGET results at three different resolutions (dotted lines: $N_{\text{gas}} = 2 \times 10^4$, dashed lines: $N_{\text{gas}} = 2 \times 10^5$, solid lines: $N_{\text{gas}} = 2 \times 10^6$), while the red lines are for the runs with AREPO. For $N_{\text{gas}} \geq 2 \times 10^5$, both codes seem to produce converged entropy profiles, but they converge to very different results. While in GADGET the entropy profiles steadily decrease towards the centre, in the moving-mesh code a large entropy core is produced. This systematic difference in the central entropy profiles is in good agreement with the previous study by Mitchell et al. (2009) of idealized major merger simulations of haloes in the non-radiative regime.
three components of the velocity vector for each blob, i.e. the reminder of the intracluster gas is initially at rest. We set all cluster-centric distance of 700 kpc and thickness of 150 kpc, and to $10^3$ (low-resolution simulation), and $10^4$ (dotted lines), and $N_{\text{gas}} = 2 \times 10^5$ (dashed lines) and $N_{\text{gas}} = 2 \times 10^6$ (continuous lines). The vertical dotted line indicates the virial radius of the underlying dark matter halo. While for $N_{\text{gas}} \geq 2 \times 10^5$ the entropy profiles seem converged for each code, they converge to a very different result.

3.3.4 Generalized blob test: non-radiative case

We now devise a numerical experiment to capture the evolution of cold, dense blobs in a more realistic setting, rather than a uniform density wind tunnel (see Section 3.2.1). For this purpose we consider our default isolated halo with a static, analytic Hernquist profile for the dark matter component, gas in hydrostatic equilibrium, and we include gas self-gravity, but neglect any radiative losses. We additionally populate this halo with 10 blobs, with the following properties: the gas pressure within blobs is set to be 0.01 of the maximum intracluster pressure within 800 kpc radius from the centre; the radius of each blob is $R_{\text{blob}} = 20$ kpc; the total mass of all blobs is 20 per cent of the total intracluster gas mass, i.e. $M_{\text{blobs}} = 0.2 M_{\text{gas}} M_{\text{halo}}$; the adiabatic index is the same for the blobs and for the surrounding gas, $\gamma = 5/3$.

We place the blobs randomly within a spherical shell with a cluster-centric distance of 700 kpc and thickness of 150 kpc, and particles belonging to the blobs are drawn randomly as well. Apart from positions, we also assign bulk velocities to the blobs, while the reminder of the intracluster gas is initially at rest. We set all three components of the velocity vector for each blob, i.e. $v_r$ (only inward radial velocity), $v_\theta$, and $v_\phi$, assuming a random distribution for each velocity component starting from a characteristic velocity value of 200 km s$^{-1}$. Blob velocity values range then from $\sim 230$ to $\sim 510$ km s$^{-1}$, with the average velocity of all blobs being $v_{\text{mean}} \sim 400$ km s$^{-1}$. In this way, individual blobs will not reach the cluster centre all at the same time, giving them more realistic orbits than purely radial ones. Initially, the blobs are roughly in pressure equilibrium with the surrounding gaseous halo. We perform this numerical experiment at two different resolutions: $N_{\text{gas}} = 10^4$, $N_{\text{blob}} = 10^3$ (low-resolution simulation), and $N_{\text{gas}} = 10^5$, $N_{\text{blob}} = 10^4$ (higher resolution simulation).

In Fig. 11, we show the time evolution of the dense blobs moving through the isolated halo, for the higher resolution run. In the top panels, the projected surface density map is plotted at the initial time, where both the intracluster gas structure and the blob properties are exactly the same in the two codes. Initially, for $t \leq 1$ Gyr, the blobs are moving on almost identical orbits in GADGET and AREPO, and they also have very similar morphologies. As the blobs start approaching the inner cluster region, well-defined bow shocks develop ahead of each blob and ram-pressure stripping ablates the blobs. Additionally, the Kelvin–Helmholtz and Rayleigh–Taylor instabilities arise which tend to disrupt the blobs on a characteristic time-scale of several Gyr, as discussed in Section 3.2.1 (but note that here gas is self-gravitating). The lower rate of ram-pressure stripping and suppression of fluid instabilities in GADGET has a significant effect not only on blob morphologies but also on their orbits. At $t = 1.37$ Gyr, as illustrated in the middle panels, it is still possible to cross-identify each blob in GADGET with the respective blob in the AREPO run, but the blobs in GADGET have lost less material and thus appear denser and some of them are closer to the centre.

This different loss of blob material leads to systematically diverging orbits due to higher buoyancy and dynamical friction forces acting on each blob in GADGET. This can be clearly seen in the bottom panels where the blobs in GADGET are markedly more concentrated and have essentially all fallen to the innermost cluster region, while in AREPO they are at much larger cluster-centric distances being gradually eroded. Moreover, given that the blobs in the moving-mesh calculation are ablated more efficiently and that therefore the dynamical friction exerted on the blobs is lower, they have higher velocities when passing at the pericentre and thus can reach larger distances after the first passage, as visible in the bottom panels of Fig. 11. As the gas is self-gravitating, the blobs are also subject to tidal stripping when passing close to the innermost regions which contributes to the ablation of the blob material.

To illustrate more clearly the differences in mass loss of the blobs in GADGET and AREPO, in Fig. 12 we show projected maps of gas material at $t = 1.37$ Gyr (corresponding to the middle panels of Fig. 11), which initially belonged to the blobs. To compute these maps in the case of GADGET, we show integrated mass along the line of sight ($\Delta z = 2$ Mpc) for all particles initially contained in the blobs. In the case of AREPO, we use a tracer field to follow spreading of fluid elements initially within the blobs: each cell is characterized by an additional scalar field $\text{Tracer}$ which at $t = 0$ is equal to 1 for all blob cells and 0 elsewhere. The tracer field essentially evolves as a dye cast on the moving fluid, and at some time $t > 0$ the $\text{Tracer}$ value indicates the mass fraction of the material which initially was in the blobs for each cell. In the right-hand panel of Fig. 12, we plot the projected density-weighted tracer field. The dynamical range is the same in both panels, ranging from the maximum of the projected quantity to $10^{-7}$ of this maximum value, using a logarithmic colour mapping. As anticipated, cold dense blobs in GADGET lose much less material while in the run with AREPO the lost blob material is significantly more diffuse. In the wake of the infalling blobs prominent tails develop, extending up to several 100 kpc. The mass deposition of blob material is clearly much more spatially extended in the moving-mesh code and occurs over larger cluster-centric distances, allowing fluids with different entropies to intermingle and to affect host halo properties on larger scales.

The formation of bow shocks in front of the moving blobs also implies that vorticity will be generated due to the baroclinic term. To explore this issue in more detail, in Fig. 13 we show vorticity maps for the run with GADGET (left-hand panel) and for the simulation with AREPO (right-hand panel) at $t = 1.37$ Gyr (corresponding to the middle panels of Fig. 11). The vorticity maps are constructed by first evaluating projected mass-weighted velocity maps for the
Figure 11. Projected surface density maps in units of $M_\odot \text{kpc}^{-2}$ at the initial time (top panels), at $t = 1.37 \text{Gyr}$ (middle panels), and at $t = 2.65 \text{Gyr}$ (bottom panels) for GADGET ($\alpha = 1.0, N_{\text{ngb}} = 64$) and AREPO. The plotted spatial domain is $2 \times 2 \times 2 \text{Mpc}$, so as to encompass initially the whole halo whose virial radius is $R_{200} = 755 \text{kpc}$. While in the beginning the orbits and the morphologies of the blobs are very similar in the two codes, for $t \geq 1 \text{Gyr}$ they begin to diverge, with blobs sinking to the centre faster in GADGET.
Figure 12. Spatial distribution of gas at $t = 1.37$ Gyr that was contained in the blobs at time $t = 0$. The plotted spatial domain is $2 \times 2 \times 2$ Mpc (for details on map-making see the main text). Most of the material stays confined within blobs in the case of GADGET, with a relatively small fraction lost to their wake. The spatial distribution of blob material is markedly different in AREPO, showing much more stripped gas that creates prominent tails which extend up to several 100 kpc.

Figure 13. Projected mass-weighted vorticity maps in units of km s$^{-1}$ kpc$^{-1}$ (absolute value of the $z$-component only) for a run with GADGET (left-hand panel) and with AREPO (right-hand panel). The maps are computed at time $t = 1.37$ Gyr from the start of the simulation. The thickness of the projected region is $\Delta z = 1$ Mpc. In the wake of the bow shocks produced by the moving blobs, turbulence is generated. The spatial extent of the turbulent wakes is significantly larger in AREPO.

$x$ and $y$-components, and then by taking the absolute value of the $z$-component of the curl. Note that while the positions and the structure of the bow shocks are rather similar in the two codes, as expected, there is a striking difference between the spatial extent of the high-vorticity regions produced in the wake of the bow shocks (denoted with green colours), where turbulent motions should be generated. For example, focusing on the rightmost blob centred on $x \sim 250$ kpc and $y \sim -100$ kpc in projection, the mean projected vorticity value in its wake is a factor of $\sim 2$ higher in the moving-mesh calculation. This clearly suggests that the suppression of vorticity generation in GADGET will have an impact on the level of turbulence injected by curved shocks that are associated with galaxy or sub-halo motions through the ICM, producing a bias in the amount of non-thermal pressure support in galaxy clusters (see also e.g. Dolag et al. 2005; Iapichino et al. 2011; Vazza et al. 2011a).

As the blobs orbit at larger cluster-centric distances in AREPO for a longer time and mix with the surrounding medium more efficiently than is the case for GADGET, they lead to higher entropy generation over a wider range of radii. This is shown in Fig. 14, where we compute radial entropy profiles for all intracluster gas including the blobs, at the final time $t \sim 10$ Gyr when the system has reached hydrostatic equilibrium. Both in the low- and high-resolution runs with AREPO, the gas entropy profile is significantly higher in the inner regions up to $r \sim 100$ kpc. Instead, in GADGET, the entropy content of the dense blobs is increased less as they sink towards the centre, such that they settle on a lower adiabat, corresponding to smaller radii.
We also note that these systematic differences in entropy profiles are entirely due to the different hydro solvers employed by GADGET and AREPO, and not due to the different choice of gravitational softenings for gas (fixed in GADGET and adaptive in AREPO, but with a floor set equal to the GADGET value). In fact, if in Fig. 14, the green line shows the entropy profile obtained from an identical low-resolution AREPO simulation where instead of an adaptive a constant gravitational softening is used for the gas, in the same way as in GADGET.

It is also instructive to analyse how the entropy profiles of the ICM evolve with time. In Fig. 14 we show radial entropy profiles at the initial time (grey lines) for both resolutions. The initial entropy profile of the run with a higher particle number extends further inwards due to the better spatial resolution, while some differences in the outer regions are due to different positions of the blobs which are drawn randomly. The kink in the initial entropy profile for \( r \geq 500 \) kpc is caused by the blobs which populate this region and have low entropy content with respect to the surrounding ICM. During the first 0.5 Gyr, the entropy profiles in GADGET and AREPO do not change noticeably, remaining nearly identical to each other and to the values prescribed by the initial conditions. At 1 Gyr, the gas entropy starts rising in the very centre in the run with AREPO (it is higher by a factor of \( \leq 3 \) with respect to the initial value), while it takes almost twice as much time in the run with GADGET before the central entropy rises by the same amount. At that time, however, the central entropy profile in AREPO is already about a factor of 6 higher than it was initially, and this systematic difference in the central entropy values persists until the final simulated time.

At time \( t \sim 2 \) Gyr, another interesting process occurs: as several blobs reach the cluster core region for the first time, dissipation of their kinetic energy leads to a gradual expansion of the whole gaseous halo (compare middle to bottom panels of Fig. 11) which needs to readjust itself to find a new equilibrium solution within the static dark matter potential.

Interestingly, in a recent paper by Vazza et al. (2011b) non-radiative cosmological simulations performed with the GADGET and ENZO codes show that infalling satellites in GADGET sink to smaller radii and are characterized by lower entropy content than is the case for ENZO, confirming the importance of the physical processes discussed here in the full cosmological setting. This indicates that at least part of the systematic difference in central entropy values found between SPH and grid codes in the Santa Barbara cluster comparison project (Frenk et al. 1999) is due to the different treatment of stripping and mixing of the cold material from the infalling satellites.

### 3.4 Radiative gas cooling in isolated haloes

#### 3.4.1 Radiative gas cooling in non-rotating haloes

We now investigate the time evolution of an isolated \( 10^{14} \) M\(_{\odot} \) mass halo subject to radiative cooling and star formation. We first simulate a halo without any net rotation and compare gas and stellar properties of this system between the two codes. The results illustrated here are for haloes with static dark matter potentials, but we find very similar results if we consider live haloes instead.

In Fig. 15 we show star formation rates as a function of time for our simulated object consisting of \( N_{\text{gas}} = 10^4 \) or \( N_{\text{gas}} = 10^6 \) particles/cells, calculated with either GADGET or AREPO. Over the whole simulated time-span of 10 Gyr the star formation rates are very similar in the two codes. This indicates that not only do the implemented gas cooling rates match very well, but also that the subgrid model for star formation and supernova feedback is consistent between the codes, even at a relatively low numerical resolution (a similar conclusion is reached for mergers of isolated galaxies by Hayward et al., in preparation).

![Figure 14](image1.png)

**Figure 14.** Radial entropy profiles at time \( t \sim 10 \) Gyr after the start of the simulation, at which point all the blobs have been disrupted and the system has reached a new equilibrium. For each code (GADGET: blue lines; AREPO: red lines) two resolution simulations are shown: \( N_{\text{gas}} = 10^4 \) and \( N_{\text{blob}} = 10^3 \) (dashed lines), and \( N_{\text{gas}} = 10^5 \) and \( N_{\text{blob}} = 10^4 \) (continuous lines). Grey curves with the same line styles represent the initial entropy profiles of these simulations. An additional green line shows the AREPO simulation with \( N_{\text{gas}} = 10^4 \) and \( N_{\text{blob}} = 10^3 \), but with the fixed gravitational softening as in GADGET. AREPO runs with adaptive and fixed softenings for the gas produce very similar entropy profiles. Vertical lines with the same line styles denote the gravitational softenings and vertical dotted line the virial radius of the system. The interaction of the moving dense blobs with the ICM leaves systematically different imprints on the gas entropy profiles when simulated with AREPO or GADGET.

![Figure 15](image2.png)

**Figure 15.** Time evolution of the star formation rate of a \( M_{\text{vir}} = 10^{14} \) M\(_{\odot} \) isolated halo which radiatively cools and has negligible spin. Illustrated are low \( (N_{\text{gas}} = 10^4; \) dotted lines) and high \( (N_{\text{gas}} = 10^6; \) continuous lines) resolution simulations with GADGET (blue) and AREPO (red). The agreement in star formation rates is very good over the whole simulated time-span.
with GADGET (blue) and AREPO (red). In the inset, the ratio of \( M_{\text{cold}} \) resolution runs. The distribution of gas entropies is almost identical from the start of the simulation for the low-resolution and high-resolution runs. The gaseous halo is allowed to radiatively cool but there is no net rotation. The vertical dotted lines indicate the gravitational softening used in the GADGET run, which correspond to the floor values of the adaptive softenings in the moving-mesh calculations.

To demonstrate more clearly that gas radiative cooling and star formation proceed in a very similar manner in our isolated haloes in the absence of any net rotation, in Fig. 16 we plot the total mass of cold baryons (stars and gas with entropy \( < 10^5 \) in internal units) as a function of time for GADGET (blue lines) and AREPO (red lines), at two different resolutions, i.e. \( N_{\text{part}} = 10^4 \) (dotted lines) and \( N_{\text{part}} = 10^5 \). For the same number of SPH particles as cells used in AREPO the amount of cold baryons matches to within a few per cent between the two codes, as is evident from the inset in the plot where we show the ratio of \( M_{\text{cold}} \) values found with our moving-mesh code and with GADGET. In the case of live haloes we find that the total mass of cold baryons exhibits the same level of agreement.

Furthermore, in Fig. 17 we show a 2D histogram of intracluster gas entropies as a function of cluster-centric distance, after 5 Gyr from the start of the simulation for the low-resolution and high-resolution runs. The distribution of gas entropies is almost identical in the two codes for \( r > r_{\text{soft}} \) (note that below \( r_{\text{soft}} \), the simulation results are not trustworthy due to the limited gravitational resolution on these scales). It can be seen that there is some difference in the gas entropy close to the virial radius of the system, which is caused by different boundary conditions (vacuum for GADGET and a uniform low-resolution grid for AREPO).

These results are in line with the expectation that gas cooling and condensation in this simulated system are determined entirely by the gas properties at the cooling radius (Bertschinger 1989; White & Frenk 1991; Hernquist & Springel 2003) which corresponds very closely between the simulation codes.

### 3.4.2 Radiative gas cooling in rotating haloes

Even though gas cooling and star formation proceed in a remarkably similar way in GADGET and the moving-mesh code for haloes with vanishing spins, this is not guaranteed to remain the case once some degree of rotation is included. We therefore simulate exactly the same isolated haloes as in the previous section, but imposing a certain level of gas rotation within the static dark matter potential. To highlight the effect, we use a large spin parameter equal to \( \lambda = 0.4 \).

Fig. 18 shows the time evolution of the total mass in cold baryons (stars and gas with entropy \( < 10^5 \) in internal units) for the rotating haloes simulated with GADGET (blue curves) and AREPO (red curves) at different resolutions. Comparing \( M_{\text{cold}} \) with our findings from Fig. 16 for non-rotating haloes indicates that overall less gas cools from the hot phase if the gas spins. This effect is not surprising given our initial conditions. Even though the gas density and temperature distribution are initially identical, in the simulations where there is considerable spin the gas will be subject to centrifugal accelerations, preventing it from collapsing radially. The (partial) centrifugal support will tend to reduce the gas densities and hence the cooling rates.

More importantly, from Fig. 18 it can be seen that there is poorer agreement in the amount of cold baryons between GADGET and AREPO for a rotating gaseous halo. The discrepancy is larger for the low-resolution run with \( N_{\text{gas}} = 10^4 \), where the final \( M_{\text{cold}} \) value after 10 Gyr is about 30 per cent higher in the moving-mesh code. The reason for this discrepancy is twofold: at low resolution, GADGET underestimates the cooling rate somewhat, while AREPO overestimates it. The SPH result turns out to be more stable at poor resolution than the mesh-based calculation. Here it is advantageous for SPH that even for a few particles a clearly defined phase boundary between cold and hot gas is maintained (in fact, a ‘pressure blip’ in SPH leads to a sampling gap at this boundary) whereas in AREPO this boundary is blurred at low resolution, causing slightly elevated cooling. This trend is also confirmed by the gas radial velocities which are least negative in the low-resolution GADGET run and most negative in the low-resolution AREPO simulation. The radial velocities systematically differ within a few 100 kpc and throughout most
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Figure 18. Time evolution of the cold baryonic mass (stars and gas with entropy < 10^5 in internal units) for a M_{vir} = 10^{14} M_⊙ isolated halo which radiatively cools and has a spin of λ = 0.4. Illustrated are low (N_{gas} = 10^4; dotted lines) and high (N_{gas} = 10^6; continuous lines) resolution simulations with GADGET (blue) and AREPO (red). In the inset, the ratio of M_{cold} values found in AREPO and GADGET is shown for both numerical resolutions. The agreement between the amount of cold baryons formed is poorer for N_{gas} = 10^4, but improves for the high-resolution run, with M_{cold} in AREPO being ∼10–15 per cent higher.

of the simulated time-span. For N_{gas} = 10^6, the radial velocities obtained with the two codes are much closer, which also translates into a better match of the amount of cold baryons, with M_{cold} higher by ∼10–15 per cent in AREPO.

At least part of the remaining difference in M_{cold} between GADGET and AREPO is driven by the interaction between the gas which is cooling towards the centre and the cold gas that is already in the disc. This is shown in Fig. 19, where we plot 2D histograms of the gas entropy as a function of cluster-centric distance at t = 5 Gyr for the low- (left-hand panel) and high-resolution simulations (right-hand panel). In the moving-mesh code the gas surrounding the central disc has a range of entropy values, with some fluid elements exhibiting very large entropies due to the shock in the immediate vicinity of the disc. In contrast, in the runs with GADGET and especially at low resolution, there is a clear gap between cold material in the disc and hotter gas in the halo (see e.g. Agertz et al. 2007; Springel 2010b). Similarly, as described in Section 3.3.2, SPH particles are affected by artificial viscosity in the converging part of the flow which can slightly offset cooling losses. Note that for N_{gas} = 10^6 the difference in gas entropy structure around the disc is lower between the codes, as expected, given that unwanted artificial viscosity effects are reduced and that the gap between the cold and hot gas phases due to repulsive pressure forces is smaller.

From Fig. 19 it is also evident that the extent of the cold disc is different between the codes, especially at low resolutions. To quantify this important effect, in Table 1 we summarize the main properties of the forming disc. For N_{gas} = 10^6 the half-mass radius of the gaseous disc in GADGET is almost a factor of 2 smaller than in the higher resolution run, while in the moving-mesh code R_{gas,HM} is 70 per cent of the value we obtain with N_{gas} = 10^4. This indicates that the convergence rate of the gas disc size is slower in the case of GADGET, due to spurious transfer of angular momentum from the cold to the hot phase (Okamoto et al. 2003) and due to the artificial viscosity in the case of poorly resolved discs (note, however, that the total angular momentum is manifestly conserved in GADGET). In the case of the stellar discs, they are essentially not resolved in our low-resolution runs (R_{stars,HM} ∼ r_{soft} = 14 kpc), while the half-mass radius is ∼40 per cent higher in AREPO at higher resolution. Contrary to GADGET, total angular momentum conservation is not automatically guaranteed in the moving-mesh code, particularly for discs resolved with only a small number of cells. None the less, it is reassuring that R_{gas,HM} increases with higher resolution in AREPO (attesting that spurious angular momentum transport inwards is probably small) and that the value of R_{gas,HM} obtained with both codes for N_{gas} = 10^6 is relatively close. This indicates that the
Table 1. Gaseous and stellar properties of the disc at \( t = 5 \) Gyr which forms in an isolated \( 10^{14} \, M_{\odot} \) halo which radiatively cools and rotates. For simulations with two different resolutions with \textsc{gadget} and \textsc{arepo}, we list the maximum radial extent of the gaseous disc, its half-mass radius and the total mass within the half-mass radius, in the third, fourth and fifth columns, respectively. Stellar half-mass radius and the total stellar mass enclosed within it are shown in the sixth and seventh columns. In the eighth column, the total mass of cold baryons is given, being equal to \( 2 \times (M_{\text{gas, HI}} + M_{\text{stars, HI}}) \). All gas particles/cells with entropy less than \( 10^{5} \) (in internal units) are assumed to be part of the disc, while for the stellar disc we consider all star particles that form in the simulated volume.

<table>
<thead>
<tr>
<th>Code</th>
<th>( N_{\text{gas}} )</th>
<th>( R_{\text{gas, max}} )</th>
<th>( R_{\text{gas, HM}} )</th>
<th>( M_{\text{gas, HM}} )</th>
<th>( R_{\text{stars, HM}} )</th>
<th>( M_{\text{stars, HM}} )</th>
<th>( M_{\text{cold}} )</th>
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</thead>
<tbody>
<tr>
<td></td>
<td>( \times 10^{10} , M_{\odot} )</td>
<td>( \text{kpc} )</td>
<td>( \text{kpc} )</td>
<td>( \times 10^{10} , M_{\odot} )</td>
<td>( \text{kpc} )</td>
<td>( \times 10^{10} , M_{\odot} )</td>
<td>( \times 10^{10} , M_{\odot} )</td>
</tr>
<tr>
<td>\text{GADGET}</td>
<td>( 10^{4} )</td>
<td>40.0</td>
<td>30.3</td>
<td>27.0</td>
<td>13.5</td>
<td>215.3</td>
<td>484.6</td>
</tr>
<tr>
<td>\text{GADGET}</td>
<td>( 10^{6} )</td>
<td>79.0</td>
<td>54.7</td>
<td>9.8</td>
<td>6.8</td>
<td>238.3</td>
<td>496.2</td>
</tr>
<tr>
<td>\text{AREPO}</td>
<td>( 10^{4} )</td>
<td>79.0</td>
<td>47.4</td>
<td>27.7</td>
<td>18.1</td>
<td>267.9</td>
<td>591.2</td>
</tr>
<tr>
<td>\text{AREPO}</td>
<td>( 10^{6} )</td>
<td>85.0</td>
<td>67.1</td>
<td>12.8</td>
<td>9.7</td>
<td>263.1</td>
<td>551.8</td>
</tr>
</tbody>
</table>

3.4.3 Generalized blob test: radiative case

Here we consider the evolution of cold, dense blobs embedded in a galaxy cluster as in Section 3.3.4, but now the whole system is allowed to radiatively cool. To add an additional layer of realism the blobs are constructed to be similar to cosmological substructures: they are equipped with their own dark matter halo, and stars may form out of their gas during the simulated time-span. More specifically, each blob is represented by a live Hernquist dark matter halo and gas in hydrostatic equilibrium. The virial mass of each blob is \( 2 \times 10^{13} \, M_{\odot} \), the scale length parameter is \( a = 41.4 \, \text{kpc} \), and the gas fraction amounts to \( f_{\text{gas}} = 0.17 \). As before, we populate our default \( 10^{13} \, M_{\odot} \) halo (that has a static dark matter potential) with 10 identical substructures. The procedure for assigning blob positions and velocities is the same as in Section 3.3.4, but here we use a different random number seed which leads to the different initial positions and velocities with respect to Section 3.3.4. The positions of blob centres are in the range of \( \sim 650 \) to \( \sim 750 \, \text{kpc} \), while the characteristic blob velocities range from 200 to 500 km s\(^{-1}\). Also, in Section 3.3.4 the gas in the halo is initially at rest while here the gas has considerable angular momentum, which contributes to the relative velocity between the blobs and the gas. We simulate the evolution of this system at three different resolutions: \( N_{\text{gas}} = 10^{4} \), \( N_{\text{blob}} = 10^{2} \) (low resolution), \( N_{\text{gas}} = 10^{5} \), \( N_{\text{blob}} = 10^{4} \) (intermediate resolution), and \( N_{\text{gas}} = 10^{6} \), \( N_{\text{blob}} = 10^{4} \) (high resolution), with gas self-gravity, and with cooling and the subgrid model for star formation.

Additionally, we perform two simulations at intermediate resolution where the standard subgrid model for star formation is included as well, but where we simply prevent any star particles from being spawned out of the star-forming phase, denoted by EOS. This simulation set-up is particularly useful for following the thermodynamical evolution of cold and hot gas for many Gyr without the dense cold gas being subject to fragmentation that would likely occur in pure cooling runs. For numerical experiments with \textsc{arepo}, as a default choice, we have not used mesh refinement and de-refinement. However, we have performed extra runs at intermediate resolution where we adopt a de-/refinement strategy to limit the mass range of gas cells (within a factor of 2 of the gas particle mass in the matching \textsc{gadget} run) which automatically imposes a narrow range of stellar masses as well. For these runs we have increased the number of gas cells/particles in each blob to \( 2.5 \times 10^{3} \), such that the gas particle/cell mass in blobs is exactly identical to the one in the parent halo (which is optimal for our de-/refinement method). These test runs recover very closely all of our results with the default set-up, confirming that \( N \)-body heating effects (e.g. due to the spectrum of star particle masses) are not very important here.

In Fig. 20, we show projected surface density maps at \( t = 0.9 \), 2.3 and 6 Gyr for simulations where we use the subgrid model for star formation, but prevent the spawning of star particles. Initially, the evolution of the blobs proceeds in a very similar fashion in simulations with \textsc{gadget} and \textsc{arepo}. However, already after less than a Gyr (see top panels of Fig. 20) blobs in the moving-mesh code are much more affected by ram pressure and dynamical fluid instabilities, which cause efficient gas stripping. As the blobs reach the very inner regions, they interact with the forming disc. In \textsc{gadget} simulations, the blobs have a significantly more damaging effect on the disc, simply because they are less gas depleted. This is clearly visible in the central panels of Fig. 20, where in \textsc{gadget} feeble spiral features are present, while in \textsc{arepo} the disc is more extended and spiral arms are well developed. Additionally, in the moving-mesh simulation a large fraction of the stripped material is deposited in the main halo, and is gradually accreted on to the central disc, promoting its growth. Blobs in \textsc{gadget} are more coherent and eventually lose their angular momentum due to hydro-dynamical friction. As they merge with the central disc, an ellipsoidal structure is formed (see bottom panels of Fig. 20). Note that due to the static dark matter halo the dynamical friction force arises only due to the gas in the main halo. Thus, we expect the blobs will lose their angular momentum on an even faster time-scale in simulations with live dark matter haloes, as is the case for cosmological runs. After 6 Gyr, the difference between \textsc{gadget} and \textsc{arepo} simulations is significant: while in the latter case an extended disc forms, with gaseous spiral arms reaching up to 60 kpc away from the centre, in \textsc{gadget} we are left with a flattened, amorphous blob.

In numerical experiments where we do not suppress star particles from forming, we find very similar results. In Fig. 21, we show projected stellar density maps at \( t = 6 \) Gyr using our intermediate-resolution simulations (which both in resolution and in time match...
Figure 20. Projected surface density maps in units of $M_\odot \text{kpc}^{-2}$ at times $t = 0.9 \text{ Gyr}$ (top panels), $t = 2.3 \text{ Gyr}$ (central panels) and $t = 6 \text{ Gyr}$ (bottom panels) for a $M_{\text{vir}} = 10^{14} M_\odot$ isolated halo which rotates, radiatively cools, and has 10 orbiting substructures. The thickness of the projection is $\Delta z = 2 \text{ Mpc}$. Although we have used our subgrid model for star formation in these simulations, spawning of star particles has been intentionally prevented here. Gas stripping from the orbiting blobs is found to be very different in the two numerical techniques. In GADGET, blobs largely survive, and when they interact with the central disc they tend to disrupt it, whereas in AREPO, the blobs are more easily shredded and have a less damaging effect on the forming disc. In fact, some of the stripped blob material ends up contributing to the extended gaseous disc.

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the bottom panels of Fig. 20). Clearly, the stellar distribution in GADGET is more centrally concentrated, forming a small disc. On the contrary, a well-defined extended stellar disc assembles in the moving-mesh calculation, endowed with a central bar. We quantify the properties of the gaseous and stellar disc in Table 2. The convergence rate of the gas disc size in GADGET is even slower than in the simulations without blobs, because in addition to the spurious transfer of the angular momentum the damaging effect of the
The interaction of cold blobs with the surrounding medium influences the global star formation rates as well. In Fig. 22, we compare the star formation histories in simulations without blobs (left-hand panel; see Section 3.4.2) with the numerical tests where haloes are populated with 10 substructures (right-hand panel). In the case without blobs, the star formation rates are somewhat higher in AREPO especially for the low-resolution run due to the more efficient gas cooling and thus larger amounts of cold gas available for star formation. For the high-resolution run the differences in star formation rates between the codes are smaller, given that the difference in the amount of cold gas between the codes amounts to ∼ 30 per cent. The difference in the star formation rates between the codes becomes more pronounced once we include the blobs. Initially, as can be seen from the inset plot where we show the star formation rate from the blobs and from the disc separately, the star formation rate within blobs contributes more than 50 per cent to the total star formation rate, and it is very similar in both codes. This induces a much higher peak in the total star formation rate in the AREPO blobs dramatically drops, and at t > 1 Gyr it is truncated altogether. This is due to ram-pressure stripping and dynamical fluid instabilities that efficiently remove the gas from the blobs. In fact, we can crudely estimate a characteristic Kelvin–Helmholtz time-scale at the initial time using the equation from Section 3.2.1 (but note that here gas is self-gravitating), where for the blob density we take the average density within the blob radius R_{blob} (which we varied from the scale length parameter to the virial radius) and for the surrounding medium density we take the typical halo density at the position of the blobs. With these assumptions, we obtain typical values for f_{KH} in the range of 1–3 Gyr, depending on the blob positions, relative velocities and the choice of R_{blob}. These
gravitational softening, the same as in the matching GADGET run. It
is a factor within the blobs is suppressed in AREPO. In GADGET simulations, star
formation proceeds in the central disc which contains a greater
amount of cold, star-forming gas which fills a larger area. In the inset plot, we compute the total star formation rate occurring within the blobs (dot–dashed and continuous lines) and the one coming from the disc (double dot–dashed and continuous lines) for our intermediate- and high-resolution simulations with GADGET (blue curves) and AREPO (red curves). In the right-hand panel we also show intermediate-resolution AREPO run with the fixed gravitational softening for the gas (green lines) which produces very similar star formation rates (both in the disc and the blobs) as the simulation with the adaptive softening.

$t_{KH}$ values are comparable to the time-scale on which star formation within the blobs is suppressed in AREPO. In GADGET simulations, star formation proceeds in the blobs even until 9 Gyr, albeit at a progressively reduced rate. For $t > 1$ Gyr the total star formation rate comes mainly from the central regions. The central discs that form in the AREPO simulations have higher star formation rates over many Gyr, which are typically larger by a factor of 2 than the GADGET results. Even in our highest resolution moving-mesh simulation the star-forming disc has roughly twice as large amount of cold gas, a difference which originates from the interaction of hotter infalling gas with the gas in the disc, as described in Section 3.4.2. However, also the star-forming gas is distributed over a larger area. In fact, while within the half-mass radius of the GADGET disc the star formation rate is $\sim$40 per cent higher in the moving-mesh run, outside of it is a factor $\sim$2.3 higher, contributing about half to the total AREPO star formation rate. The reason why cold, star-forming gas in GADGET is filling a smaller area in the disc and is more confined to the dense arm segments and blobs (the so-called ‘string-of-pears’, which can be also seen in the middle panel of Fig. 20), is due to the SPH surface tension originating at the interface between cold and hot media in relative motion. In the right-hand panel of Fig. 22, we also show an identical intermediate-resolution AREPO run, but with fixed gas gravitational softening, the same as in the matching GADGET run. It can be seen that the choice of gas gravitational softening in AREPO does not affect our results in any significant way: the evolution of the star formation rate both in the central disc and in the orbiting blobs remains very similar, indicating that the differences that we see between the two codes are indeed entirely driven by hydrodynamical effects.

These findings have immediate consequences for more realistic astrophysical situations. For example, Puchwein et al. (2010) have simulated a high-resolution sample of galaxy clusters with GADGET finding that up to 30 per cent of intracluster stars form in dense cold blobs – remnants of infalling satellites. Suppressed dynamical instabilities in GADGET enhance the probability of survival for these dense blobs, which can then serve as sites of star formation, thus possibly overpredicting the number of intracluster stars. In fact, if we compute the total mass of cold baryons $M_{\text{cold}}$ (stars plus gas above the density threshold for star formation) in two cosmological simulations (for further details see Paper I), where in one we prevent the spawning of stars and in the other we allow it (the simulations are otherwise identical), $M_{\text{cold}}$ is found to be very similar in GADGET regardless of whether stars are formed or not. Instead, in the simulations with AREPO, we find that $M_{\text{cold}}$ is systematically reduced at lower redshifts if the spawning of stars is prevented. This demonstrates explicitly that very low entropy material formed in GADGET cannot easily be shredded and mixed with higher entropy gas (at least in the absence of additional feedback processes such as galactic winds or black hole heating), so that $M_{\text{cold}}$ is preserved (a similar conclusion has been reached independently by Heß & Springel 2012). In contrast, in the simulations with AREPO if star formation is switched-off, some of the cold material can be stripped out of galaxies due to dynamical instabilities, returning it to diffuse form and lowering $M_{\text{cold}}$. Importantly, this also implies that the number of stars formed in AREPO will be much more sensitive to the characteristic time-scale for star formation and to the numerical resolution (needed to resolve fluid instabilities) than is the case for GADGET. In a recent paper by Agertz, Teyssier & Moore (2011), it has been shown that lower star formation efficiency leads to the production of larger discs in cosmological simulations. While this result is in line with our findings, it is important to recognize that the actual cause is quite different: in the study by Agertz et al. (2011), differences in the physical modelling of star formation in cosmological simulations affect the disc sizes,

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while here the cause lies in different accuracies of the hydro solvers involved.

3.5 Gravitational $N$-body heating

We now investigate potential artificial gas heating due to the Poisson noise induced by the finite number of particles in dark matter haloes. Steinmetz & White (1997) outlined the analytic theory of this effect, and confirmed it with non-radiative and radiative numerical experiments that quantified gravitational $N$-body noise present in structure formation simulations. For an equilibrium system they defined a characteristic $N$-body heating time-scale which is proportional to the cube power of the dark matter velocity dispersion and inversely proportional to the dark matter particle mass and dark matter density. Due to this inverse proportionality to the dark matter density, $N$-body heating is expected to be strongest in the innermost regions of haloes. From their analysis it follows that the dark matter particle mass adopted in numerical simulations should be lower than a critical mass which is of the order of a few times $10^9 M_\odot$ for galaxy clusters (see their equation 10), otherwise the radiative cooling losses may be overwhelmed by spurious heating.

Since nowadays the typical mass resolution of hydrodynamical cosmological simulations has improved dramatically, with simulated galaxy clusters containing several times $10^6$ up to $10^7$ particles in zoom-in runs (Dolag et al. 2009; Sijacki, Springel & Haehnelt 2009; Vazza et al. 2010), possible numerical artefacts due to a grainy dark matter distribution are rarely addressed in the literature (but see e.g. Kay et al. 2000; Borgani et al. 2006; Vazza 2011). Recently, Springel (2010a) has pointed out that mesh-based codes are potentially more severely affected by gravitational $N$-body heating than SPH codes, due to their better ability to detect weak shocks, which in this case is an unwanted feature.

Here we perform a number of idealized test problems aimed at explicitly addressing the $N$-body heating problem, and in particular to understand whether there are systematic differences between GADGET and AREPO in this respect. Note that intentionally all previous tests (except for Section 3.4.3) have been performed either without any dark matter component or with a static dark matter potential to avoid such possible Poisson-noise imprints.

We first consider our standard isolated halo, where we replace the static dark matter potential with a live dark matter halo in which dark matter particles have a characteristic velocity dispersion for the given mass of the system (but there is no net rotation). As in the case of the rigid potential, we populate the halo with gas particles in hydrostatic equilibrium which are initially at rest. We include self-gravity of the gas component and evolve the system non-radiatively for 10 Gyr.

In Fig. 23, we show radial profiles of gas density, temperature and entropy at the final time for both codes at two resolutions: $N_{\text{gas}} = N_{\text{DM}} = 10^4$ particles (dotted lines) and $N_{\text{gas}} = N_{\text{DM}} = 10^6$ particles (continuous lines). The centre of the system is defined by the position of the most bound particle. For cluster-centric distances greater than the gravitational softening value, $r_{\text{soft}}$, the entropy profiles of GADGET and AREPO agree to within $\sim 20$ per cent for the low-resolution runs, and to better than 10 per cent for the high-resolution simulation. For $r < r_{\text{soft}}$, the differences between the codes in the entropy profile amount up to a factor of 2, with the gas entropy in AREPO being systematically higher. An extra test simulation performed with AREPO in its dual entropy mode (where entropy instead of energy is explicitly conserved for cells whose Riemann problems with adjacent cells all have Mach number less than 1.1)
resulted in identical radial profiles as for the AREPO simulation in the standard energy mode. This clearly indicates that the small entropy core which develops in AREPO is not caused by weak shocks generated by the grainy dark matter potential. Rather, central gas particles/cells acquire small-scale velocities when simulated with live haloes. In the case of the moving-mesh code, these velocities lead to fluid mixing and temperature equilibration in the innermost regions, while in GADGET such mixing does not occur. Note, however, that in the gravitating systems, \( r_{\text{soft}} \) represents the minimum spatial scale below which the simulation results are not trustworthy due to discreteness effects. Therefore, provided that gravitational softening lengths are chosen conservatively, the gas properties of galaxy clusters in an equilibrium configuration simulated with GADGET and AREPO match very well.

We now consider a more challenging problem, where we simulate the infall of cold gas into dark matter haloes. This is the same problem as discussed in Section 3.3.2, but here we compare outcomes from live versus static dark matter haloes. To match numerical experiments with rigid and live haloes as closely as possible, we adopt the following procedure.

For live haloes, the motion of dark matter particles is prevented in the code to avoid local deformations of the gravitational potential which will not be present in the static case. Effectively, in this way the distribution of dark matter is ‘frozen’, and live haloes are coarse-grained representations of static haloes.

For static haloes, we convolve the analytic Hernquist potential with a spline softened potential in the centre, such that gas particles feel exactly the same gravitational acceleration as in the case of live haloes, where the central potential needs to minimize effects from two-body encounters. For the Hernquist potential, gravitational accelerations are given by

\[
g(r) = -\frac{GM_{\text{vir}}}{(r + a)^2}. \tag{3}\]

We modify the gravitational acceleration felt by gas particles in the code, viz.

\[
g_g(r) = \frac{GM_{\text{vir}}}{(\frac{32\pi}{15} \frac{a}{M_{\text{vir}}})^{-\frac{1}{2}} (r + a)^2 \exp(-h_1/r)}, \tag{4}\]

where \( h_0 \) and \( h_1 \) are free coefficients (to first approximation \( h_0 \sim h_1 \sim r_{\text{soft}} \)). We determine the value of \( h_0 \) and \( h_1 \) for three different resolution runs by fitting equation (4) to the gravitational acceleration of a live ‘frozen’ halo with 10^6 dark matter particles simulated three times assuming \( r_{\text{soft}} \) values appropriate for the numerical resolutions we want to investigate.

In Fig. 24 we illustrate the outcome of this method. In the upper panel, the green lines denote the radial gravitational acceleration of the live ‘frozen’ halo with 10^6 dark matter particles, adopting \( r_{\text{soft}} = 3, 6.5 \) and 14 kpc (indicated by numbers 1, 2 and 3), respectively. Blue dashed lines are our best-fitting functions to these simulations using equation (4). Dotted red lines indicate the radial acceleration in the case of the analytic Hernquist profile and of the spline-softened potential, which are the limiting cases of equation (4) for large and small radii, respectively. The bottom panel of Fig. 24 shows how accurate our modified gravitational acceleration is as a function of cluster-centric distance.

With the procedure described above, we perform simulations of the radial infall of cold gas into live (‘frozen’) haloes. We keep the number of gas resolution elements the same and equal to \( N_{\text{gas}} = 10^5 \) (corresponding to \( r_{\text{soft,gas}} = 3 \) kpc for GADGET), while gravitational softenings are adaptive in AREPO but with a floor of 3 kpc), while we increase the number of dark matter particles from \( N_{\text{DM}} = 10^4 \) (\( r_{\text{soft,DM}} = 14 \) kpc), to \( 10^5 \) (\( r_{\text{soft,DM}} = 6.5 \) kpc) and \( 10^6 \) (\( r_{\text{soft,DM}} = 3 \) kpc). For each live halo we run a matching static halo simulation where the number of gas particles/cells is kept the same, i.e. \( N_{\text{gas}} \) = 10^5. In all runs, gas self-gravity is neglected and there are no radiative losses. In close encounters of gas and dark matter, the effective gravitational softening is the maximum between \( r_{\text{soft,gas}} \) and \( r_{\text{soft,DM}} \), which is relevant length scale of the problem.

In Fig. 25 we show radial profiles of gas density, temperature and entropy at time \( t = 0.3 \) Gyr (analogous to Fig. 7). The left-hand panels are for simulations of live and static haloes with GADGET, while the right-hand panels show the results with AREPO. Due to gravitational N-body heating there are systematic differences in the central gas properties. The central gas density is lower, while the central gas temperature and entropy are higher for the live haloes, due to spurious transfer of energy from dark matter to gas, which heats the gas and makes it expand. In general, N-body heating effects are largest for \( N_{\text{DM}} = 10^4 \) and smallest for \( N_{\text{DM}} = 10^6 \), and they are confined to spatial regions within \( r_{\text{soft,DM}} \), which is reassuring.

We can see by comparing the left-hand to the right-hand panels that AREPO is indeed much more sensitive to spurious gravitational heating, as discussed in Springel (2010a), with the central entropy boosted by two orders of magnitude. In the right-hand panel of Fig. 25, we also show radial entropy profile obtained with AREPO in the dual entropy mode (where entropy instead of energy is explicitly conserved for cells with all Mach numbers less than 1.1) with \( N_{\text{DM}} = 10^6 \) particles (orange continuous line). This indicates that at least part of the central entropy core is generated by the weak shocks which gas cells experience when moving through the grainy dark matter potential.

At the final time \( t = 2.45 \) Gyr, when the system has reached an equilibrium state, the difference between the matching live and static halo runs for \( r > r_{\text{soft,DM}} \) is \( \sim 10, \sim 5 \) and \( \sim 5 \) per cent, respectively, for low, intermediate and high resolution in GADGET,
Figure 25. Radial profiles of gas density, temperature and entropy at time $t = 0.3$ Gyr for GADGET (left-hand panels; blue lines: static dark matter haloes; green lines: live dark matter haloes) and AREPO (right-hand panels; red lines: static dark matter haloes; green lines: live dark matter haloes). For each code we show three different resolution runs: $N_{DM} = 10^4$ and $r_{soft} = 14.0$ kpc (dotted lines), $N_{DM} = 10^5$ and $r_{soft} = 6.5$ kpc (dashed lines), and $N_{DM} = 10^6$ and $r_{soft} = 3.0$ kpc (continuous lines), while $N_{gas} = 10^6$ is kept fixed. In the right-hand panel we also show the entropy profile obtained with the moving-mesh code in the entropy mode for our highest resolution simulation (continuous orange line). The vertical black lines with the same style indicate the softening scales of the dark matter potential. The black vertical dot–dashed lines denote the virial radius of the system.

while it is somewhat higher in AREPO, i.e. $\sim 15–20$, 10 and 5 per cent, respectively. For $r < r_{soft}$, the ratios of central entropy values are $\leq 10$, $\leq 4$ and $\leq 2$ in GADGET and $\leq 300$, $\leq 150$ and $\leq 25$ in AREPO, for low-, intermediate- and high-resolution runs. Even though the amount of artificial heating is significantly larger in AREPO than it is for GADGET, the departures between static and live halo radial profiles always occur within $r_{soft,DM}$ for both codes. Provided that gravitational softening values for the dark matter component are chosen cautiously, our numerical experiments hence indicate that the systematic discrepancy in central entropy values found between SPH and mesh-based codes for the Santa Barbara cluster comparison project (Frenk et al. 1999; Springel 2010a) is unlikely to be due to effects from a grainy dark matter potential during cosmological structure formation.
4 DISCUSSION AND CONCLUSIONS

In this study we have carried out a detailed comparison between the SPH code GADGET and the new moving-mesh code AREPO on a number of hydrodynamical test problems, which are crucial for understanding cosmological simulations of galaxy formation. In a purely hydrodynamical regime without gas self-gravity or an external gravitational potential, we have first carried out a set of numerical experiments previously considered in the literature, some of which have rarely been shown for SPH codes, such as the 2D implosion test. We have then focused on idealized non-radiative galaxy cluster simulations, specifically aimed towards benchmarking differences in hydro solvers for problems with shocks and fluid instabilities. In simulations where radiative losses were included we have analysed the amount of baryons which cool from the hot halo atmospheres in GADGET and AREPO, both for rotating and non-rotating haloes. In the former case, we have also studied how the central baryonic discs form, and how orbiting substructures affect the disc morphology and the star formation rate. Finally, we have constructed special test problems designed to gauge the effect of gravitational $N$-body heating on the gaseous properties of the haloes. Our main conclusions are as follows.

(i) While post-shock fluid properties are captured well both in GADGET and AREPO in the case of 1D shock tube tests with high Mach numbers, the shocks are significantly broader in GADGET, and substantial post-shock oscillations develop, largely because of an inadequate treatment of the initial contact discontinuity. AREPO in the moving-mesh mode preserves the contact discontinuity much more accurately, but it is broadened if we employ a static mesh, due to larger numerical diffusivity in this case. Note that the more accurate treatment of the contact discontinuity in the moving-mesh code can lead to a larger ‘wall heating’ effect (Rider 2000) than in static grid codes, which tend to wash-out at some level the initial code can lead to a larger ‘wall heating’ effect (Rider 2000) than in accurate treatment of the contact discontinuity in the moving-mesh due to larger numerical diffusivity in this case. Note that the more accurately, but it is broadened if we employ a static mesh, due to larger numerical diffusivity in this case. Note that the more accurate treatment of the contact discontinuity in the moving-mesh code can lead to a larger ‘wall heating’ effect (Rider 2000) than in static grid codes, which tend to wash-out at some level the initial start-up errors at the contact discontinuity (see also description of the Noh problem in Springel 2010a).

(ii) For shocks with complicated geometries in multidimensions, differences between GADGET and AREPO are more striking. Even though the global fluid properties remain similar, the sampling of the fluid properties is much noisier in GADGET, and the development of dynamical fluid instabilities is inhibited. These problems are not specific to GADGET, but are inherent to the standard SPH method as a whole (see also Agertz et al. 2007; Springel 2010b, and discussion in Section 2.1.2). The suppression of fluid instabilities reduces the amount of entropy generation by mixing and artificially prolongs the lifetime of gaseous structures which are moving through a medium with a different density. Additionally, vorticity generation in the wake of curved shocks due to the baroclinic source term is largely suppressed in GADGET, which directly impacts angular momentum transfer by vortices and the level of generated turbulence.

(iii) These fundamental differences between GADGET and AREPO identified in simple hydrodynamical test problems affect the properties of gas in more realistic, cosmologically motivated simulations as well. Specifically, in non-radiative idealized simulations of merging galaxy clusters and in simulations of isolated haloes with orbiting gaseous substructures, we find that in AREPO: (i) an entropy core is produced in the centre due to more efficient fluid mixing, (ii) the gas stripping rate from the orbiting substructures is larger and (iii) more vorticity is produced in the wake of curved shocks. These findings are in line with results of previous works which simulated similar problems with AMR codes (e.g. Mitchell et al. 2009; Vazza 2011; Vazza et al. 2011b). Moreover, the unphysical dissipation of shocks and subsonic turbulence in GADGET, as shown by Bauer & Springel (2012) (see also section 4.2 of Paper I), leads to the heating of the halo outskirts rather than of the central regions.

(iv) In radiative simulations of isolated haloes without any net rotation, gas cooling and condensation into stars proceeds in a very similar fashion in GADGET and AREPO, given that the gas properties at the cooling radius match closely. However, for spinning haloes there is a net difference in the total amount of cold baryons, which is higher in AREPO, and this difference persists at a level of about 10–15 per cent in our highest resolution runs. Baryonic discs which form due to dissipative collapse of rotating gas are systematically larger in AREPO at low resolution, and the convergence rate of the gas disc sizes is higher.

(v) In numerical experiments where we follow the interaction between a forming central disc and 10 orbiting substructures in an isolated halo which radiatively cools, the final disc morphology is significantly different. While in AREPO an extended disc is produced with well-developed spiral arms and a central bar, in GADGET the disc is more centrally concentrated and amorphous. Orbiting substructures are more much more efficiently stripped of their gas content in the moving-mesh calculation and the material is incorporated into the host halo atmosphere. Instead, in GADGET gaseous substructures are more coherent, thus they lose more angular momentum from hydro-/dynamical friction, and when passing through the central disc they induce morphological transformations.

(vi) While star formation is more readily truncated in infalling substructures due to gas stripping, extended gaseous discs in AREPO have significantly larger star formation rates for many Gyr than is the case for GADGET.

(vii) Due to its better ability to detect weak shocks, AREPO is more sensitive to gravitational $N$-body heating (Springel 2010a). This is confirmed by our specifically designed numerical experiments, which allow us to quantify the magnitude of spurious gas heating for simulations of haloes with a finite number of dark matter particles with respect to the analytic dark matter potentials. None the less, the spatial extent of this artificial heating is reassuringly constrained to lie within the gravitational softening length for typical set-ups, which is a scale below which simulation results are not trustworthy due to resolution effects anyway.

The numerical experiments presented in this study clearly demonstrate that several important shortcomings of the SPH solver not only affect idealized test problems but are equally detrimental in more realistic set-ups relevant for structure formation, and ultimately in full cosmological simulations. This is especially the case because of: (i) the complicated flows involving multiphase media which are the norm in cosmological simulations and (ii) the hierarchical nature of structure formation where low-mass systems are always poorly resolved. In both of these regimes, the hydrodynamical solver of the standard SPH method exhibits the largest inaccuracies. On the other hand, our numerical tests confirm and significantly extend the findings of Springel (2010a) that the moving-mesh code AREPO delivers a physically more accurate representation of the evolution of inviscid gases. Note that in the current work we deliberately kept the modelling of the baryon physics at a very simple level, so as to isolate the differences between the hydro solvers in as clean a manner as possible. While inclusion of more realistic feedback mechanisms, such as supernovae winds and AGN heating, will likely modify the properties of the simulated galaxies significantly, it is of prime importance to disentangle numerical inaccuracies of the hydro solver from the uncertainties of the feedback physics modelling. It is hence clear that cosmological simulations with AREPO have the potential to provide a much more realistic description of structure formation in
the Universe (see also Papers I and II), something we will explore in more depth in forthcoming studies.

ACKNOWLEDGMENTS

We would like to thank Andrew MacFadyen and Daniel Eisenstein for very useful discussions and suggestions on the topic, and Manfred Kitzbichler and Diego Muñoz for carefully reading the manuscript. We would like to thank the anonymous referee for many constructive suggestions, which helped to improve the presentation of the results. DS acknowledges NASA Hubble Fellowship through grant HST-HF-51282.01-A. DK acknowledges NASA Hubble Fellowship through grant HST-HF-51276.01-A. The computations in this paper were performed on the Odyssey cluster supported by the FAS Science Division Research Computing Group at Harvard University.

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APPENDIX A: ISOLATED HALOES IN HYDROSTATIC EQUILIBRIUM

Here we show how gaseous spheres placed in hydrostatic equilibrium within a static Hernquist dark matter halo evolve with time to determine the level of accuracy of our initial conditions. The general set-up is described in detail in Section 3.3.1. Specifically for the figures presented here we assume that the gas is self-gravitating and that it has initially zero velocity.

In Fig. A1, we show gas density, temperature and 1D velocity dispersion radial profiles at $t = 0.05$ Gyr (dotted lines) and $t = 2.45$ Gyr (continuous lines) for GADGET (blue lines) and AREPO (red lines) with $N_{\text{gas}} = 10^6$ resolution elements. As discussed in Section 3.3.1, due to the Poisson sampling of the gas positions in the initial conditions, they are not perfectly relaxed, which leads to the development of small-scale random motions, as evidenced by a non-zero gas velocity dispersion. Over time, these residual gas velocities are dissipated, as can be seen from the bottom panel of Fig. A1. This leads to a slight readjustment of the temperature distribution, which is very similar in both codes. Note, however, that the kinetic energy of the random gas motions is at all times a very small fraction of both the potential and the internal energy, as illustrated in Fig. A2, amounting to only about 0.2 per cent of the total gas energy after 2.45 Gyr.

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Figure A1. Radial profiles of gas density, temperature and 1D velocity dispersion at time $t = 0.05$ Gyr (dotted lines) and $t = 2.45$ Gyr (continuous lines) for GADGET (blue lines) and AREPO (red lines) with $N_{\text{gas}} = 10^6$ in a static dark matter halo with a Hernquist profile. Vertical black lines indicate the softening scales of the gas equal to $r_{\text{soft}} = 3.0$ kpc. The black vertical dot-dashed line denotes the virial radius of the system.

Figure A2. Time evolution of the ratio of gas kinetic (dotted lines), potential (dashed lines) and internal (continuous lines) to the total gas energy in GADGET (blue) and AREPO (red) simulations of isolated haloes in hydrostatic equilibrium. Gas is self-gravitating and represented by $N_{\text{gas}} = 10^6$ resolution elements and is evolved within a static Hernquist dark matter halo.

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