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A HIGHER-ORDER UPWIND METHOD FOR VISCOELASTIC FLOW

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We present a conservative finite difference method designed to capture elastic wave propagation in viscoelastic fluids in two dimensions. We model the incompressible Navier–Stokes equations with an extra viscoelastic stress described by the Oldroyd-B constitutive equations. The equations are cast into a hybrid conservation form which is amenable to the use of a second-order Godunov method for the hyperbolic part of the equations, including a new exact Riemann solver. A numerical stress splitting technique provides a well-posed discretization for the entire range of Newtonian and elastic fluids. Incompressibility is enforced through a projection method and a partitioning of variables that suppresses compressive waves. Irregular geometry is treated with an embedded boundary/volume-of-fluid approach. The method is stable for time steps governed by the advective Courant–Friedrichs–Lewy (CFL) condition. We present second-order convergence results in $L^1$ for a range of Oldroyd-B fluids.

1. Introduction

The governing equations for viscoelastic flow of an Oldroyd-B fluid are the incompressible Navier–Stokes equations plus an extra viscoelastic stress described by the Oldroyd-B constitutive equations:

$$\frac{\partial \mathbf{u}}{\partial t} + (\mathbf{u} \cdot \nabla) \mathbf{u} - \frac{1}{\rho} \nabla \cdot \mathbf{\tau} = -\frac{1}{\rho} \nabla p + \frac{\mu_s}{\rho} \Delta \mathbf{u},$$  \hspace{1cm} (1)

$$\nabla \cdot \mathbf{u} = 0,$$  \hspace{1cm} (2)

$$\frac{\partial \mathbf{\tau}}{\partial t} + (\mathbf{u} \cdot \nabla) \mathbf{\tau} - (\nabla \mathbf{u}) \mathbf{\tau} - \mathbf{\tau} (\nabla \mathbf{u})^T = \frac{\mu_p}{\lambda} 2 \mathbf{D} - \frac{1}{\lambda} \mathbf{\tau},$$  \hspace{1cm} (3)

where $\mathbf{u}$ is the fluid velocity, $\mathbf{\tau}$ is the polymeric stress tensor, $p$ is the isotropic pressure, and $\mathbf{D} = [\nabla \mathbf{u} + (\nabla \mathbf{u})^T]/2$ is the rate-of-strain tensor. The parameters that describe the fluid are the density, $\rho$, relaxation time, $\lambda$, and the solvent and polymeric

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contributions to the total viscosity, $\mu = \mu_s + \mu_p$. The dimensionless parameters that characterize these types of flows are the Reynolds number, $Re = \frac{\rho U L}{\mu}$, and the Weissenberg number, $We = \frac{\lambda U}{L}$, where $U$ and $L$ are the characteristic velocity and length.

Though the Reynolds number and the Weissenberg number independently characterize viscoelastic flows, it is the elastic Mach number, $Ma = \sqrt{Re \cdot We}$, that is the critical parameter in determining well-posedness of the system. In particular, the system of equations exhibits a change in type from parabolic to hyperbolic when the elastic Mach number becomes supercritical ($Ma > 1$), admitting propagation of discontinuities. This mathematical behavior was alluded to in the experimental results of Ultman and Denn [33] and formally noted in [7, 18]. Joseph suggested that a method suitable for transonic flows might be needed to capture the transition to supercritical flows in viscoelasticity [17]. The analysis described in [30] capitalized on this concept in the design of a numerical algorithm that resolves unsteady elastic wave behavior in viscoelastic fluids.

In this paper, we extend the previous numerical algorithm [30] by leveraging the conservative hyperbolic formulation described therein to design a suitable higher resolution upstream method for the hyperbolic. In the original algorithm the Oldroyd-B equation is recast into a well-posed hyperbolic form with source terms using a stress-splitting technique; a Lax–Wendroff method is used to discretize the quasilinear form of the hyperbolic part in the context of a predictor-corrector projection method. (Projection methods are an approach to enforcing the constraint in incompressible flows [3, 2] and have proven to be successful in treating unsteady viscoelastic flows [19, 30].) Our new method uses a second-order Godunov method [5, 6], instead of Lax–Wendroff as in [30], to discretize the hyperbolic part of the equations, resulting in two immediate advantages. First, the maximum time step is increased by a factor of four to allow an advective CFL number restriction of $0 < CFL < 1$. Second, we can apply second-order conservative finite volume techniques which have been developed for hyperbolic conservation laws [6], elliptic equations [16], and parabolic equations [22] in an embedded boundary (EB) framework for irregular geometry. Our results are consistent with the modified equation analysis in these methods, and we obtain second-order solution error convergence in $L^1$ for a range of Oldroyd-B fluids.

## 2. Hyperbolic analysis

Through the introduction of the inverse deformation tensor, $g$, which links material (Lagrangian) coordinates, $X$, and spatial (Eulerian) coordinates, $x$, as in

$$g_{\alpha\beta} = \frac{\partial X_\alpha}{\partial x_\beta},$$

(4)
the advective part of the PDE for viscoelastic stress (3) may be put in conservation form. The quantity \( M \) is conserved:

\[
\frac{\partial M}{\partial t} + \nabla \cdot (u \otimes M) = g \left( -\frac{1}{\lambda} \tau + \left( \frac{\mu_p}{\lambda} - \rho a^2 \right) 2D \right) g^T,
\]

(5)

The PDE for \( g \) and its right hand side are described in detail in [23]. Here \( a \) is an arbitrary constant with dimensions of velocity. As developed in [30], this fictitious wave speed may be treated as a parameter that affects the partitioning of hyperbolic and elliptic terms. Through proper choice of that parameter, the CFL limiting time step of the hyperbolic partition can be improved by several orders of magnitude in the Newtonian limit (\( \lambda \to 0 \)). Here, for purposes of analysis, \( a \) need only satisfy \( \min_d (\rho a^2 + \tau_{dd}) > 0 \).

All together, the coupled PDEs (1)-(3) may be written in the form

\[
\frac{\partial U}{\partial t} + \frac{\partial F}{\partial x_i} = S_h(U) + S_i(U, \nabla U, \Delta U),
\]

(8)

where the left hand side is a system of conservation laws, and the right hand side contains proper hyperbolic source terms, \( S_h \), and improper (elliptic) source terms, \( S_i \). \( U \) is the vector of conserved quantities:

\[
U = (u, M, ge_0, \ldots, ge_{D-1})^T,
\]

(9)

\[
F_d = \left( u_d u - \frac{1}{\rho} \tau e_d, u_d M, gu\delta_{0,d}, \ldots, gu\delta_{D-1,d} \right)^T,
\]

(10)

\[
S_h = \left( -\frac{1}{\rho} \nabla p, -\frac{1}{\lambda} g \tau g^T, 0, \ldots, 0 \right)^T,
\]

(11)

\[
S_i = \left( \nu D \Delta u, 2\left( \frac{\mu_p}{\lambda} - \rho a^2 \right) D g^T, \left[ u \times (\nabla \times g^T) \right] e_0, \ldots, \left[ u \times (\nabla \times g^T) \right] e_{D-1} \right)^T,
\]

(12)

where \( D = 2 \) is the dimensionality of the problem and \( \nu = \mu / \rho \).

We analyze the hyperbolic subsystem in primitive variables, \( W \). The linearization of (8) in primitive variables gives matrices whose eigenvalues are wave speeds, and whose eigenvectors determine the characteristics. If, in the 1D analysis of these linearized equations for direction \( d \), primitive variable \( u_d \) is included, then wave speeds and characteristics describing compressive wave motion are observed. Yet, omission of \( u_d \) and its corresponding stress \( \tau_{dd} \) is also inaccurate [9] since variation in these quantities is permitted by the multidimensional equations. The approach
to this dilemma, after \[9; 8\] is to block partition the primitive equations, treating dependence on gradients of the variables \(u_d\) and \(\tau_{dd}\) as source terms from the point of view of the remaining variables. We will refer to the variable partition \((u_d, \tau_{dd})\) as inactive (subscript \(I\)), and the remaining variable partition as active (subscript \(A\)). For \(d = 0\),

\[
W_T^T = (W_{A,0}^T | W_{I,0}^T) = (u_1, \tau_{10}, \tau_{11}, g_{00}, g_{10}, g_{01}, g_{11} | u_0, \tau_{00}).
\]  

(13)

The primitive variable \(\tau_{01}\) is omitted because \(\tau\) is symmetric. In these variables, the linearized homogeneous advection equation in direction \(d = 0\) is

\[
\frac{\partial W_0}{\partial t} + A_0 \frac{\partial W_0}{\partial x_0} = 0,
\]  

(14)

\[
A_0 = \begin{bmatrix}
A_{AA,0} & A_{AI,0} \\
A_{IA,0} & A_{II,0}
\end{bmatrix}
\]

\[
= \begin{bmatrix}
u_0 & -1/\rho & 0 & 0 & 0 & 0 & 0 & 0 \\
-\rho c_0^2 & u_0 & 0 & 0 & 0 & 0 & 0 & 0 \\
-2\tau_{10} & 0 & u_0 & 0 & 0 & 0 & 0 & 0 \\
g_{01} & 0 & 0 & u_0 & 0 & 0 & 0 & 0 \\
g_{11} & 0 & 0 & 0 & u_0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & u_0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & u_0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0
\end{bmatrix}
\]

(15)

with \(c_d = \sqrt{a^2 + \tau_{dd}/\rho}\). The diagonal matrix of eigenvalues of partition \(A_{AA,0}\) is

\[
A_0 = \text{diag} (u_0 - c_0, u_0, u_0, u_0, u_0, u_0, u_0 + c_0)^T.
\]  

(16)

The corresponding right eigenvectors are given by the columns of

\[
R_0 = \begin{bmatrix}
-c_0 & 0 & 0 & 0 & 0 & c_0 \\
-\rho c_0^2 & 0 & 0 & 0 & 0 & -\rho c_0^2 \\
-2\tau_{10} & 1 & 0 & 0 & 0 & -2\tau_{10} \\
g_{01} & 0 & 1 & 0 & 0 & g_{01} \\
g_{11} & 0 & 0 & 1 & 0 & g_{11} \\
0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0
\end{bmatrix}.
\]  

(17)

2.0.1. An exact Riemann solver. For the incompressible Euler equations, Bell et al. [2] construct edge-centered time-centered predictor states using Taylor series with upwind derivatives. For those equations, their approach is identical to using a higher-order Godunov predictor because upwinding solves exactly the associated
Riemann problem. In the present system of equations, the wave structure is more complex, but there are no genuinely nonlinear waves, that is,

\[
(\nabla_{W_A} A_{kk}) \cdot R e_k = 0,
\]

for each of the 7 waves \( k \) associated with block \( A_{AA} \) of (15). This condition is guaranteed by the fact that the complete solution for the inactive variables is taken to be the average of the input left and right states [8], and therefore the eigenvalues are constant with respect to each component of \( W_A \).

By analysis of the generalized Riemann invariants,

\[
\frac{\partial(W_A)_0}{e_0^T R e_k} = \frac{\partial(W_A)_1}{e_1^T R e_k} = \cdots = \frac{\partial(W_A)_6}{e_6^T R e_k},
\]

for each wave \( k \), it may be concluded (assuming for convenience \( d = 0 \)) that

(i) \( u_1 \) and \( \tau_{10} \) are constant across the 5 contact (speed \( u_0 \)) waves;

(ii) \( g_{01} \) and \( g_{11} \) are constant across the fast \( u_0 \pm c_0 \) waves;

(iii) the generalized Riemann invariants for the \( \pm \) fast waves include the identities

\[
\frac{\partial u_1}{\pm c_0} = \frac{\partial \tau_{10}}{-\rho c_0^2} = \frac{\partial g_{00}}{g_{01}} = \frac{\partial g_{10}}{g_{11}},
\]

where the denominators of each term are constant across the wave. Thus, across each fast wave the change in \( u_1 \) is proportional to \( c_0 \), etc.;

(iv) across the fast waves, the generalized Riemann invariants contain also

\[
\frac{\partial \tau_{11}}{-2 \tau_{10}}.
\]

So, given the change of \( \tau_{10} \) across the given wave, the change in \( \tau_{11} \) is determined.

Let the constant states in the Riemann fan be labeled \( W_L, W_L^*, W_R^*, \) and \( W_R \) in sequence, and let \( \Psi_L (\Psi_R) \) measure the strength of the left (right) fast waves. From observation (iii) one has

\[
\begin{pmatrix}
\frac{u_1}{\tau_{10}} \\
\frac{\tau_{10}}{L^*}
\end{pmatrix}
= \begin{pmatrix}
\frac{u_1}{\tau_{10}} \\
\frac{\tau_{10}}{L}
\end{pmatrix}
- \Psi_L \begin{pmatrix}
\frac{c_0}{\rho c_0^2}
\end{pmatrix},
\]

\[
\begin{pmatrix}
\frac{u_1}{\tau_{10}} \\
\frac{\tau_{10}}{R^*}
\end{pmatrix}
= \begin{pmatrix}
\frac{u_1}{\tau_{10}} \\
\frac{\tau_{10}}{R}
\end{pmatrix}
+ \Psi_R \begin{pmatrix}
\frac{c_0}{-\rho c_0^2}
\end{pmatrix},
\]

and from observation (i) one has \( (u_1, \tau_{10})_{L^*} = (u_1, \tau_{10})_{R^*} \), which couples the fast waves enabling their strength to be simply determined from

\[
\begin{pmatrix}
c_0 & c_0 \\
-\rho c_0^2 & -\rho c_0^2
\end{pmatrix}
\begin{pmatrix}
\Psi_L \\
\Psi_R
\end{pmatrix}
= \begin{pmatrix}
\frac{u_1}{\tau_{10}} \\
\frac{\tau_{10}}{L}
\end{pmatrix}
- \begin{pmatrix}
\frac{u_1}{\tau_{10}} \\
\frac{\tau_{10}}{R}
\end{pmatrix}.
\]

(23)
With $\tau_{10}$ determined across the wave fan, observation (iv) determines $\tau_{11}$:

$$\int_{(\tau_{11})_L}^{(\tau_{11})_L^*} d\tau_{11} = \frac{2}{\rho c_0^2} \int_{(\tau_{10})_L}^{(\tau_{10})_L^*} \tau_{10} d\tau_{10},$$

(24)

$$(\tau_{11})_L^* = (\tau_{11})_L + \frac{1}{\rho c_0^2} \left[ (\tau_{10})_L^2 - (\tau_{10})_L^2 \right].$$

(25)

The same equation holds across the right fast wave. The determination of other variables is then trivial by application of observation (iii). For example, from (20),

$$\begin{align*}
\frac{(g_{00})_{R^*} - (g_{00})_R}{(g_{01})_{R^*}} &= \frac{(u_1)_{R^*} - (u_1)_R}{c_0},
\end{align*}$$

(26)

The active variable solution to our Riemann problem is given by the constant state ($L$, $L^*$, $R^*$, or $R$) containing the zero wave speed characteristic.

### 3. Predictor-corrector formulation

We discretize time in steps $\Delta t$, with $t^{n+1} = t^n + \Delta t^n$. Space is discretized in square cells of length $h$, and $x = hi$ is the lower left corner of cell $i$. Variables $U_i^n$ are cell-centered.

For each time step $n$, the artificial wave speed $a$ is a global constant determined by the heuristic model:

$$a^2 = \min \left\{ \chi(\lambda) a_{\infty}^2 + [1 - \chi(\lambda)] a_0^2, \frac{\rho}{\lambda} \right\},$$

(27)

$$a_{\infty}^2 = \frac{\rho}{\lambda},$$

(28)

$$a_0^2 = \frac{2}{\rho} \min_{i,d} |(\tau_{dd})_i|,$$

(29)

$$\chi(\lambda) = \frac{\lambda}{t_{\text{adv}}} \left[ 1 - e^{-\lambda/(2t_{\text{adv}})} \right] \left( 1 - e^{-\lambda/t_{\text{adv}}} \right),$$

(30)

$$t_{\text{adv}} = \frac{h}{\max_i |u|},$$

(31)

with limiting values $a^2 = a_{\infty}^2$ as $\lambda \to \infty$, and $a^2 = a_0^2$ as $\lambda \to 0$. Note that the conserved quantity $M$ depends on $a$, so a reevaluation of $a$ necessitates a rescaling of $M$ throughout the domain.

The predictor step of the method uses well-established higher-order Godunov approaches [5; 6] to estimate time-centered edge-centered solution values. These predictor states are made discrete divergence-free ($\nabla \cdot u^{n+\frac{1}{2}} = 0$) on a marker-and-cell (MAC) stencil [15].
Fluxes $F_{i \pm e/2}^{n+1/2} = F(U_{i \pm e/2}^{n+1/2})$ computed from these predictor states enter a conservative update:

$$\tilde{U}_{i}^{n+1} = U_{i}^{n} - \frac{\Delta t}{h} \sum_{d=0}^{D-1} \left[ (F_{d})_{i+e_d/2}^{n+1/2} - (F_{d})_{i-e_d/2}^{n+1/2} \right].$$  \hspace{1cm} (32)

The corrector computes $U^{n+1}$ by adding to $\tilde{U}^{n+1}$ implicit and explicit source term contributions, and by use of an approximate cell-centered projection to make $u^{n+1}$ discrete divergence-free.

### 3.1. Predictor

The predictor in our predictor-corrector method consists of the calculation of time-centered edge-centered states, $W_{i+e_d/2}^{n+1/2}$, which are discrete divergence-free. The predictor state is computed in four steps.

1. **First**, the one-dimensional primitive equations are used to estimate time-centered edge-centered states. For the active partition, characteristic tracing and slope limiting occur as in higher-order Godunov methods. For the inactive partition, Taylor series in space and time with centered differences are used. This first step uses strictly one-dimensional equations with no transverse coupling.

2. **Second**, the edge states so obtained are double-valued, and we resolve these with the Riemann solver described above in Section 2.0.1.

3. **Third**, the transverse coupling omitted in the first step is incorporated using cell-centered gradients of the edge-centered states computed by the Riemann solution. The transverse flux correction is described in [5; 28], but we include the transverse terms in terms of primitive variable differences rather than conservative fluxes. The corrected states so-obtained are again double-valued, and another Riemann problem gives a single final result.

4. **Fourth**, time-centered edge-centered velocity data is made discrete divergence-free, i.e.,

$$u := u - \nabla \left[ \Delta^{-1}(\nabla \cdot u) \right].$$  \hspace{1cm} (33)

With $u$ edge-centered, $\nabla \cdot u$ is cell-centered. This projection is exact, in the sense that $\Delta_h = (\nabla \cdot)_h \nabla_h$, with the discrete Laplacian reducing to the standard 5-point stencil in two dimensions away from boundaries. Then, $\Delta^{-1}(\nabla \cdot u)$ is also cell-centered. The discrete gradient operator uses centered divided differences to give edge-centered corrections. The normal and tangential velocity components are updated at each face even though only the normal velocity contributes to the divergence.

The details of the first step is now given. With the active–inactive partitioning introduced in (13), upwind characteristic tracing for the active primitive variables takes the form
where \( \bar{A} = \text{diag}(\bar{A}_{ij}) \) if \( \Lambda_{ii} \geq 0 \), 0 otherwise) is a projection that sets to zero those terms of the diagonal argument matrix corresponding with eigenvalues whose sign is negative/positive, respectively. The subscript \( L \) (\( R \)) indicates that the result is traced to the left (right) side of the edge \( i + e_d/2 \). Where the stencils support it, the derivatives \( \partial W_i/\partial x \) use van Leer limited [34] fourth-order accurate derivatives [4]. The derivatives \( \partial W_i/\partial x \) use second-order centered divided differences. The tilde denotes that source terms have not yet been accounted for. The inactive variables are extrapolated in time using

\[
(W_{A,d})^{n+\frac{1}{2}}_{i+e_d/2,L} = (W_{A,d})^n_i - R_d \bar{A}_{i+e_d/2,L} \left( \frac{\Delta t}{2} A_{d} - \frac{h}{2} I \right) \left( \frac{\partial W_{A,d}}{\partial x_d} \right)^n_i
-
\frac{\Delta t}{2h} \bar{A}_{i+e_d/2,L} \left( \frac{\partial W_{A,d}}{\partial x_d} \right)^n_{i+e_d}
\]

\[
(W_{A,d})^{n+\frac{1}{2}}_{i+e_d/2,R} = (W_{A,d})^n_i - R_d \bar{A}_{i+e_d/2,R} \left( \frac{\Delta t}{2} A_d + \frac{h}{2} I \right) \left( \frac{\partial W_{A,d}}{\partial x_d} \right)^n_i
-
\frac{\Delta t}{2h} \bar{A}_{i+e_d/2,R} \left( \frac{\partial W_{A,d}}{\partial x_d} \right)^n_{i+e_d}
\]

(34)

The velocity source is computed explicitly via

\[
u^{n+\frac{1}{2}}_{i+e_d/2} = u^{n+\frac{1}{2}}_{i+e_d/2} + \frac{\Delta t}{2} \left( -\frac{1}{\rho} \nabla p^{n+\frac{1}{2}} + \nu_s(\Delta \nu n^{n+\frac{1}{2}}) \right)_i,
\]

(36)

where \( \Delta \nu \) the discrete 5-point Laplacian in regular domains. The time-centered pressure is taken from the previous time step. The calculation of \( \nabla p^{n+\frac{1}{2}} \) occurs as the last step of the corrector, (43).

The source term for viscoelastic stress is computed implicitly to properly recover the Newtonian limit \( (\tau \rightarrow 2\mu \rho D \text{ as } \lambda \rightarrow 0) \):

\[
\tau^{n+\frac{1}{2}}_{i+e_d/2,L} = \tau^{n+\frac{1}{2}}_{i+e_d/2,L} + \frac{\Delta t}{2} \left[ -\frac{1}{\lambda} \tau^{n+\frac{1}{2}}_{i+e_d/2,L} + \left( \frac{\mu \rho}{\lambda} - \rho \lambda^2 \right) 2D^{n+\frac{1}{2}} \right].
\]

(37)

The rate of strain tensor, \( D \), is calculated with centered differences.

The source terms for \( g \) are omitted for the following reason. The material
reference frame $X$ can be defined, at the start of each time step, to be equal to $x$, i.e., $g = l$ identically at the start of each time step. With this choice, the source terms for $g$ are zero if evaluated at $t^n$. Resetting $g$ to $l$ necessitates renormalizing $M$ from time step to time step.

3.2. Corrector. The corrector generates time $n + 1$ cell-centered states that are discrete divergence-free. The basic idea is to generate cell-centered time $t^{n+1}$ estimates, $\tilde{U}^{n+1}$, using the flux differencing quadrature (32). To these estimates source terms are added, as described below, to obtain $U^{n+1}$.

The corrector step for the velocity field is more complicated. We would like to use the following update equation (the superscript * indicates that the velocity field is not yet divergence-free):

$$u^{n+1,*} - u^n = \left[ -\nabla \cdot \left( u \otimes u - \frac{1}{\rho} \tilde{\tau} \right)^{n+\frac{1}{2}} \right] + \left( -\frac{1}{\rho} \nabla p^{n-\frac{1}{2}} + v_s \Delta u \right).$$

However, as in [30], we would like for the velocity update equation to properly capture the Newtonian and elastic limits. We modify the predictor step by not including the source terms for $\tau$ in the edge state prediction to instead obtain $\tilde{\tau}$ at edges. However, extra care must be taken since the transverse correction term is still computed with edge states that have been constructed with the $\tau$ sources.

Combining an equation of the form (37) with (38), we arrive at our new update equation for velocity:

$$u^{n+1,*} - u^n = \left[ \nabla \cdot \left( u \otimes u - \frac{2\lambda}{2\lambda + \Delta t} \tilde{\tau} \right)^{n+\frac{1}{2}} \right] - \frac{1}{\rho} \nabla p^{n-\frac{1}{2}}.$$

These equations are expressible as $D$ scalar discrete Helmholtz equations. This discretization is chosen in order to capture the Newtonian and elastic limits, that is, in the Newtonian limit ($\lambda \to 0$) we recover

$$u^{n+1,*} - u^n = \left[ v_s + v_p \right] \Delta u + \left[ -\nabla \cdot \left( u \otimes u \right)^{n+\frac{1}{2}} - \frac{1}{\rho} \nabla p^{n-\frac{1}{2}} \right].$$

and in the elastic limit ($\lambda \to \infty$), where $a^2$ is given by (28), we recover

$$u^{n+1,*} - u^n = v_s \Delta u + \left[ -\nabla \cdot \left( u \otimes u - \frac{\tilde{\tau}}{\rho} \right)^{n+\frac{1}{2}} - \frac{1}{\rho} \nabla p^{n-\frac{1}{2}} \right].$$

The Helmholtz equations (39) are solved using the Runge–Kutta technique of [32], which yields an $l_0$ stable solution in regular and irregular domains. That method specifies the time centering of the Laplacian term.
The last step of the velocity corrector removes the divergence of $u^*$ and calculates the pressure whose gradient will affect the subsequent time step using a pressure-projection formulation [31]. First, a potential $\phi$ is calculated on cell centers with the discrete Laplacian:

$$\Delta \phi = \left[ \nabla \cdot \text{Avg} \left( u^{n+1,*} + \frac{\Delta t}{\rho} \nabla p^{n+\frac{1}{2}} \right) \right],$$  \hspace{0.5cm} (42)

where Avg is an operator that computes face-centered values by averaging neighboring cell-centered values. Pressure is proportional to $\phi$, and

$$\nabla p^{n+\frac{1}{2}} = \frac{\rho}{\Delta t} \nabla \phi.$$  \hspace{0.5cm} (43)

With this gradient, the discrete-divergence-free velocity is

$$u^{n+1} = u^{n+1,*} + \frac{\Delta t}{\rho} (\nabla p^{n-\frac{1}{2}} - \nabla p^{n+\frac{1}{2}}).$$  \hspace{0.5cm} (44)

This projection is approximate, in the sense that $\Delta h \neq (\nabla \cdot) h \nabla h$. As noted by Lai [20], the approximate projection does not remove certain nonphysical oscillatory modes. These are damped by application of a filter

$$u := u + \zeta \nabla (\nabla \cdot u),$$  \hspace{0.5cm} (45)

using a divergence stencil other than the centered divided difference used in (42). We use $\zeta = h^2/5$ in two dimensions which is stable while always damped monopole modes in the experience of [8; 30].

The corrector step for $g$ and $M$ simply follows the flux differencing quadrature (32) followed by a source term update. The source term for $g$ is computed as in [23] using edge — and time — centered values from the predictor. The viscoelastic stress source term is discretized using Crank–Nicholson:

$$M^{n+1} = \tilde{M}^{n+1} + \frac{\Delta t}{2} \left( g \left[ \left( \frac{\mu_p}{\lambda} - \rho a^2 \right) 2D - \frac{1}{\lambda} \tau \right] g^T \right)^n$$

$$+ \frac{\Delta t}{2} \left( g \left[ \left( \frac{\mu_p}{\lambda} - \rho a^2 \right) 2D - \frac{1}{\lambda} \tau \right] g^T \right)^{n+1},$$

rearranged in the form

$$M^{n+1} = \frac{2\lambda}{2\lambda + \Delta t} \tilde{M}^{n+1} - \frac{\Delta t}{2\lambda + \Delta t} M^n$$

$$+ \frac{\Delta t}{2\lambda + \Delta t} \left( g \left[ (\mu_p - \rho a^2 \lambda) 2D + \rho a^2 I \right] g^T \right)^n$$

$$+ \frac{\Delta t}{2\lambda + \Delta t} \left( g \left[ (\mu_p - \rho a^2 \lambda) 2D + \rho a^2 I \right] g^T \right)^{n+1},$$  \hspace{0.5cm} (46)

which is evaluated pointwise.
4. Irregular domains

We use a Cartesian grid embedded boundary method to discretize the fluid equations in the presence of irregular boundaries [6]. In this approach, the irregular domain is discretized as a collection of control volumes formed by the intersection of the problem domain with the square Cartesian grid cells as in a “cookie cutter”. The various operators — the discrete divergence \( \nabla \cdot \), discrete gradient \( \nabla \), and discrete Laplacian \( \Delta \) — are approximated using finite volume differences on the irregular control volumes. Cells are classified as regular if they do not intersect embedded boundaries, irregular if they intersect boundaries, or covered if they have zero fluid volume fraction. Faces are classified in an analogous way. In problems containing irregular domains, the finite volume treatment of the regular cells follows the description of Section 3.

Throughout, time \( t^n \) data \((U)\) will be centered at cell centers, even if that point lies outside the fluid domain. Time \( t^{n+\frac{1}{2}} \) data (fluxes \( F \)) are centered at the centroid of faces,

\[
\dot{x}_{i \pm e_d/2} = \frac{1}{a_{i \pm e_d/2} h^{D-1}} \int_{A_{i \pm e_d/2}} x \, dA, \tag{47}
\]

where \( a_{i \pm e_d/2} \) is the area fraction of a cell edge \( i \pm e_d/2 \) not covered by the embedded boundary, or

\[
a_{i \pm e_d/2} = \frac{A_{i \pm e_d/2}}{h^{D-1}}, \tag{48}
\]

with \( A_{i \pm e_d} \) the area of cell \( i \) on side \( \pm d \) in contact with the fluid. Other geometric quantities used are the volume fraction, defined as

\[
\kappa_i = \frac{V_i}{h^D}, \tag{49}
\]

the area fraction of the domain boundary intersected with cell \( i \), \( A_{i}^{EB} \), and its associated area fraction, defined as

\[
a_{i}^{EB} = \frac{A_{i}^{EB}}{h^{D-1}}, \tag{50}
\]

and the outward-directed vector normal to the embedded boundary interface in cell \( i \), given by

\[
n_i = \frac{1}{a_{i}^{EB} h^{D-1}} \int_{A_{i}^{EB}} n \, dA. \tag{51}
\]
In irregular cells, the quadrature (32) is not appropriate [6]. A stable but nonconservative update is

$$\tilde{U}_{i}^{n+1} = U_{i}^{n} - \Delta t \left[ \kappa_{i}(\nabla \cdot F)_{i}^{C} + (1 - \kappa_{i})(\nabla \cdot F)_{i}^{NC} \right]^{n + \frac{1}{2}},$$

with conservative and nonconservative flux differences given by

$$(\nabla \cdot F)_{i}^{NC} = \frac{1}{h} \sum_{d=0}^{D-1} \left[ (F_{d,i+\varepsilon_{d}/2} - (F_{d,i-\varepsilon_{d}/2}) \right],$$

$$(\nabla \cdot F)_{i}^{C} = \frac{1}{V_{i}} \int_{V_{i}} (\nabla \cdot F) \, dV \approx \frac{1}{\kappa_{i}h} \left[ \sum_{d=0}^{D-1} \sum_{\pm \varepsilon_{d}} \left\{ \pm \kappa_{i,\pm \varepsilon_{d}/2} F_{d}(\hat{x}_{i,\pm \varepsilon_{d}/2}) \right\} + \kappa_{i}\rho \nabla \cdot (\rho F_{EB}) \right],$$

respectively.

Conservation violation is expressed locally by the generalized mass deficit $\delta m_{i}$,

$$\delta m_{i} = \Delta t (1 - \kappa_{i}) \kappa_{i} \left[ (\nabla \cdot F)_{i}^{NC} - (\nabla \cdot F)_{i}^{C} \right]_{i},$$

which is redistributed in a volume-weighted manner according to

$$\tilde{U}_{i}^{n+1} = \tilde{U}_{i}^{n+1} + \sum_{j=\text{neighbor}(i)}^{\mathcal{D}} \frac{\delta m_{j}}{w_{j}},$$

$$w_{i} = \sum_{j=\text{neighbor}(i)}^{\mathcal{D}} \kappa_{j}.$$

The calculation of fluxes on covered faces, and stencils used to re-center fluxes to centroids, are described in [6; 22; 29]. Additional details are given in [24]. Here we describe differences between the regular and irregular domain calculations that are specific to the present algorithm.

We compute the Poisson equation in divergence form, $\Delta \phi \approx \nabla h \cdot (\nabla h \phi) = f$, with discrete divergence given by the conservative form (54). This means that $\kappa \Delta h \phi$ is directly accessible, and division by $\kappa$ can be unstable. For the Laplacian appearing in the velocity source term (36) we use $\kappa \Delta h \phi$ in place of $\Delta h \phi$, which formally introduces an $O(\Delta t)$ discretization error. However, the results obtained by this approximation are stable and appear to not affect the global error.

In irregular cells the discretization of the divergence term in (39) is computed as follows. Define a velocity flux to be

$$F_{u} = u \otimes u - \frac{2\lambda}{2\lambda + \Delta t \rho} \cdot \tilde{W}.$$
Then, compute the divergence of $\mathbf{F}_u$ using (52) and redistribute according to (56).

Covered face values needed in the nonconservative divergence are obtained by extrapolation from face-centered time-centered values, as described in [6]. Unlike [6], we take this extrapolated edge state to represent the unique face value, so no further Riemann problem is solved.

5. Boundary conditions

In the hyperbolic treatment, boundary conditions enter in two ways:

(1) on embedded boundaries, e.g., the computation of $\mathbf{F}^{EB}$ in (54); and

(2) where the Cartesian cells abut the problem domain.

The conservative flux divergence (54) includes the flux derived from data centered at the centroid of the embedded boundary. Such states are derived from cell-centered data using Taylor series, without upwind projection. If $\hat{x}_i^{EB}$ is the centroid relative to the cell center,

$$W^{n+\frac{1}{2}, EB}_i = W^n_i + \hat{x}_i^{EB} \cdot (\nabla W^n)_i + \frac{\Delta t}{2} \left( \frac{\partial W_i}{\partial t} \right)^n$$

$$= W^n_i + \sum_d \left[ (\hat{x}_i^{EB})_d I - \frac{\Delta t}{2} A_d \right] \left( \frac{\partial W^n_i}{\partial x_d} \right)_i + \frac{\Delta t}{2} S^n_i. \quad (59)$$

This extrapolation is implemented without partitioning of $W$ or $A$. The source terms are implemented as with the predictor Section 3.1. The discrete gradient $\nabla W$ uses central differences where possible, or one-sided differences where necessary.

This one-sided boundary value may be incompatible with physical boundary conditions. The approach to boundary conditions uses the ideas of Ghidaglia and Pascal [10]. Let $W_P$ be an extrapolated edge state, as calculated by (59), and let $W_S$ be the final value used to construct the edge flux. In appropriately rotated coordinates, we are interested in the eigenstructure of the matrix $A_{AA}(W_S)$. For each characteristic pointing into the domain, one degree of freedom at the boundary must be specified. For each characteristic pointing out of the domain, a characteristic condition must be met. Specifically, if characteristic $k$ points out of the domain, a sufficient characteristic condition is

$$l_k \cdot (W_P - W_S) = 0. \quad (60)$$

For solid wall boundaries, including the embedded boundaries, this construction is straightforward. We derive $A_{AA}$ on the boundary using active variables taken from $W_P$, and selecting inactive variables on physical grounds. In the present application, the embedded boundaries are stationary surfaces subject to no-flow conditions. Accordingly, the inactive variable $u_n$ is uniquely determined, $u_n = 0$. 
(here subscript $n$ denotes the interface normal direction; subscript $t$ will denote the tangential direction). There is no a priori reason for $\tau_{nn}$ to be affected by boundary conditions, thus we take $\tau_{nn}$ in state $W_S$ equal to its extrapolated value in $W_P$. With these choices, exactly one characteristic of $A_{AA}$ enters the domain, leaving one degree of freedom to be specified. We use the no-slip boundary condition to zero the tangential velocity component. For the characteristic that points out of the domain, the characteristic condition $l_{u_n-c} \cdot (W_P - W_S) = 0$ (if the wall normal is positive) or $l_{u_n+c} \cdot (W_P - W_S) = 0$ (if the wall normal is negative) uniquely determines the shear stress $\tau_{nt}$ component of $W_S$. Thus, for solid wall boundaries, $W_S = W_P$, except for variables $u$ which are taken to be zero on physical grounds and the shear stress which is determined by the characteristic condition.

For inflow and outflow boundaries this procedure is more involved. Let $n$ point out of the domain, so for inflow we have $u_n < 0$. Inflow conditions are either supersonic, $u_n + c < 0$, or not, with $u_n$ and $c$ given by inactive variables taken from the specified inflow condition. When supersonic, all characteristics flow into the domain, and the state $W_S$ is given exclusively by imposed conditions. If not supersonic, only the characteristic $u_n + c$ flows out of the domain, so only one constraint on $W_S$ comes from $W_P$. In this case we determine the shear stress $\tau_{nt}$ component of $W_S$ by solving $l_{u_n+c} \cdot (W_P - W_S) = 0$, with all other components of $W_S$ being prescribed by the inflow condition.

On outflow, we take the inactive variables from $W_P$, and $u_n > 0$. If supersonic, $u_n - c > 0$, no characteristics flow into the fluid domain, and we take $W_S = W_P$. If subsonic, one degree of freedom of $W_S$ is specified by external conditions. In that case, we choose $u_t = 0$ and determine the remaining values of $W_S$ from $l_k \cdot (W_S - W_P) = 0$, for all $k \neq u_n - c$.

Boundary conditions are also required for the Helmholtz velocity correctors, (39), and the divergence-cleaning projections (33) and (42). The implicit velocity equations (39) use homogeneous Dirichlet conditions on solid wall boundaries, inhomogeneous Dirichlet conditions on inflow boundaries (using prescribed far-field values), and homogeneous Neumann conditions on outflow. The discrete Laplacian operator encountered in divergence-cleaning projections uses homogeneous Dirichlet on outflow, and homogeneous Neumann on inflow and solid walls.

6. Results

Results are presented for three fluids: a Maxwell (highly elastic) fluid, characterized by having no solvent viscosity, a nonzero polymeric viscosity, and a nonzero relaxation time; a Newtonian fluid, characterized by having a nonzero solvent viscosity, no polymeric viscosity, and relaxation time of zero; and a hybrid fluid [30] — a Maxwell fluid with an added solvent viscosity. Two geometries are used
Figure 1. Time-dependent $u_\theta$ profiles of a Maxwell fluid with a vortex initial condition in a rectangle. The domain has 256 × 256 cells with 24 time step increments using $\Delta t = 1.6 \times 10^{-3}$. The range is from −0.5 (red) to 0.5 (blue).

that are nonconforming with Cartesian grids; a rotated rectangular geometry, and a circular domain.

For the rectangular geometry, the computational domain has $l = w = 2.0$. The rectangular box has dimensions $l = 1.7$, and $w = 1.0$, and has been rotated 45° to maximize the amount of fluid in the computational domain. The coarse domain has 128 × 128 cells. We have chosen an initial vortex velocity profile that is sufficiently smooth at the vortex edge, given by the function

$$u_\theta(r) = 2.56[(r/0.45)(1 - r/0.45)]^4 H(0.45 - r),$$

where $r$ is the distance to the center of the box and $H$ is the Heaviside step function. This gives a maximum initial speed of $|u| = 1.0$ at $r = 0.225$ (see Figure 1, top left).

For all images corresponding to the angled box geometry, we have rotated the output so the variables are seen with respect to the normal (lengthwise) and transverse (widthwise) directions. The initial pressure is set to zero. We define the characteristic speed, $U$, as the maximum initial velocity and the characteristic length, $L$, as the width of the box.
Figure 2. Profiles for a Maxwell fluid in a rectangle at $t = 0.4224$ (last image in Figure 1). Clockwise from top left: $u_t$, $-0.5$ (red) to $0.5$ (blue); normal stress $\tau_{nn}$, $-0.21$ (red) to $0.31$ (blue); normal stress $\tau_{tt}$, $-0.21$ (red) to $0.28$ (blue); shear stress $\tau_{tn}$, $-0.46$ (red) to $0.33$ (blue); hydrostatic pressure $p$, $0$ (red) to $0.656$ (blue).

Figure 3. Time-dependent $u_0$ profiles of a Maxwell fluid with a vortex initial condition in a disk. The domain has $128 \times 128$ cells with 24 time step increments using $\Delta t = 1.6 \times 10^{-3}$. The range is from $-0.50$ (red) to $0.50$ (blue).

For the circular geometry, the computational domain has $l = w = 1.0$ and the circle has radius $r = 0.45$ to maximize the amount of fluid in the computational domain. The coarse domain has $64 \times 64$ cells. The initial velocity profile is $u_0(r) = 2.56[(r/0.4)(1-r/0.4)]^4 H(0.4-r)$, which gives a maximum initial speed of $|u| = 1.0$ at $r = 0.2$ (see Figure 3, top left). The initial pressure is set to zero. We define the characteristic speed, $U$, as the maximum initial velocity and the characteristic length, $L$, as the diameter of the circle.
A HIGHER-ORDER UPWIND METHOD FOR VISCOELASTIC FLOW 73

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Table 1. Solution error convergence rates for a Maxwell fluid with a vortex initial condition in a rectangle. Data correspond to Figures 1 and 2.

6.1. Maxwell fluid. For the Maxwell fluid, the rheological parameters are $\mu_s = 0$, $\mu_p = 1.0$, $\lambda = 1.0$, and $\rho = 1.0$. This gives the dimensionless parameters $\text{Re} = 1.0$, $\text{We} = 1.0$, and $\text{Ma} = 1.0$ for the rectangular box geometry. The coarse time step for each geometry is $3.2 \times 10^{-3}$, corresponding to $\text{CFL} \approx 0.5$. The time-dependent normal velocity is shown in Figure 1. The elastic wave propagation and reflection off the walls is clearly visible. The transverse velocity, stress, and pressure corresponding to the final image of normal velocity are shown in Figure 2. The solution error convergence after 400 fine time steps is given in Table 1. We use the same rheological parameters for the circular geometry, leading to dimensionless parameters $\text{Re} = 0.9$, $\text{We} = 0.9$, and $\text{Ma} = 1.0$. The time-dependent $u_0$ profiles are shown in Figure 3. Again, the elastic wave propagation and reflection off the walls is easily visible. The $u_1$ component of velocity, stress, and pressure corresponding to the final image of $u_0$ are shown in Figure 4. The solution error convergence after 400 fine time steps is given in Table 2.

For Maxwell fluids, we have observed that additional cell-centered filtering steps (45) are required to prevent the buildup of divergent modes near cells with small
Figure 4. Profiles for a Maxwell fluid in a disk at $t = 0.2688$ (last image in Figure 3). Clockwise from top left: $u_1$, $-0.50$ (red) to $0.50$ (blue); normal stress $\tau_{00}$, $-0.38$ (red) to $0.67$ (blue); normal stress $\tau_{11}$, $-0.38$ (red) to $0.67$ (blue); shear stress $\tau_{10}$, $-0.53$ (red) to $0.53$ (blue); hydrostatic pressure $p$, $0$ (red) to $0.55$ (blue).

Figure 5. Time-dependent $u_n$ profiles of a Newtonian fluid with a vortex initial condition in a rectangle. The domain has $256 \times 256$ cells with 2 time step increments using $\Delta t = 3.75 \times 10^{-3}$. The range is from $-0.25$ (red) to $0.25$ (blue).

In the other flow regimes, the nonzero solvent viscosity in the diffusion equation solver smooths the velocity and helps eliminate the divergent modes and additional filtering steps are not required. The approach taken here to stabilize the method is to perform 1 filter iteration per time step at the coarse resolution, 2 iterations at the medium resolution, and 4 iterations at the fine resolution. The additional filter steps are not required for the other flow regimes, but are included for consistency.

6.2. Newtonian fluid. For the Newtonian fluid, the rheological parameters are $\mu_s = 1.0$, $\mu_p = 0.0$, $\lambda = 1.0 \times 10^{-11}$, and $\rho = 1.0$ leading to dimensionless parameters $Re = 1.0$ and $We = 0.0$ for the rectangular box geometry. Since $\mu_p = 0$, volume fractions.
Table 2. Solution error convergence rates for a Maxwell fluid with a vortex initial condition in a disk. Data correspond to Figures 3 and 4.

The polymeric stress remains zero at all times. The coarse time step for each geometry is $7.5 \times 10^{-3}$, corresponding to CFL $\approx 0.5$. The time-dependent normal velocity is shown in Figure 5, in which the vortex spreads out to fill the box and decays over time. The transverse velocity and pressure corresponding to the final image of normal velocity are shown in Figure 6. The solution error convergence after 40 fine time steps is given in Table 3. Only a small number of time steps are used because after 40 the fluid velocity has already decayed to less than two percent of its initial value.

We use the same rheological parameters for the circular geometry, leading to dimensionless parameters $Re = 0.9$ and $We = 0$. The time-dependent normal velocity is shown in Figure 7. As with the rectangular box case, the vortex spreads out to fill the circle and decays over time. The transverse velocity and pressure corresponding to the final image of normal velocity are shown in Figure 8. The solution error convergence after 20 fine time steps is given in Table 4.

6.3. Hybrid fluid. For the hybrid fluid, the rheological parameters are $\mu_s = 0.1$, $\mu_p = 0.9$, $\lambda = 1.0$, and $\rho = 1.0$ leading to dimensionless parameters $Re = 1.0$, $We =$

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1.0, and $\text{Ma} = 1.05$ for the rectangular box geometry. The initial stress is set to zero. The coarse time step is $3.2 \times 10^{-3}$ for each geometry, corresponding to $\text{CFL} \approx 0.5$. The time-dependent normal velocity is shown in Figure 9. As is the case with the Newtonian fluid, the vortex spreads out and decays over time, with a different shape than the Newtonian case. The transverse velocity, stress, and pressure corresponding

Figure 6. Profiles for a Newtonian fluid in a rectangle at $t = 2.25 \times 10^{-2}$ (last image in Figure 5). Left: $u_t$, $-0.15$ (red) to 0.15 (blue); right: hydrostatic pressure $p$, 0 (red) to 1.96 (blue).

Figure 7. Time-dependent $u_0$ velocity profiles of a Newtonian fluid with a vortex initial condition in a disk. The domain has $128 \times 128$ cells with 2 time step increments using $\Delta t = 3.75 \times 10^{-3}$. The range is from $-0.25$ (red) to 0.25 (blue).

Figure 8. Profiles for a Newtonian fluid in a disk at $t = 2.25 \times 10^{-2}$ (last image in Figure 7). Left: $u_1$, $-0.15$ (red) to 0.15 (blue); right: hydrostatic pressure $p$, 0 (red) to 0.032 (blue).
A HIGHER-ORDER UPWIND METHOD FOR VISCOELASTIC FLOW

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<tr>
<td>(L_\infty)</td>
<td>(u_0)</td>
<td>2.36\text{e}-03</td>
<td>5.46\text{e}-04</td>
<td>2.11</td>
</tr>
<tr>
<td>(u_1)</td>
<td>2.36\text{e}-03</td>
<td>5.46\text{e}-04</td>
<td>2.11</td>
<td></td>
</tr>
<tr>
<td>(p)</td>
<td>4.12\text{e}-02</td>
<td>1.59\text{e}-02</td>
<td>1.37</td>
<td></td>
</tr>
</tbody>
</table>

Table 3. Solution error convergence rates for a Newtonian fluid with a vortex initial condition in a rectangle. Data correspond to Figures 5 and 6.

<table>
<thead>
<tr>
<th>norm</th>
<th>Variable</th>
<th>Coarse Error</th>
<th>Fine Error</th>
<th>Order</th>
</tr>
</thead>
<tbody>
<tr>
<td>(L_1)</td>
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<td>9.44\text{e}-05</td>
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<tr>
<td>(p)</td>
<td>5.12\text{e}-03</td>
<td>1.58\text{e}-04</td>
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<tr>
<td>(L_2)</td>
<td>(u_0)</td>
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<td>2.07</td>
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<td>1.16\text{e}-04</td>
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<tr>
<td>(p)</td>
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<td>2.48\text{e}-04</td>
<td>5.40</td>
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</tr>
<tr>
<td>(L_\infty)</td>
<td>(u_0)</td>
<td>1.06\text{e}-03</td>
<td>5.51\text{e}-04</td>
<td>0.95</td>
</tr>
<tr>
<td>(u_1)</td>
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<td>5.51\text{e}-04</td>
<td>0.95</td>
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Table 4. Solution error convergence rates for a Newtonian fluid with a vortex initial condition in a disk. Data correspond to Figures 7 and 8.

to the final image of normal velocity are shown in Figure 10. The solution error convergence after 200 fine time steps is given in Table 5.

We use the same rheological parameters for the circular geometry, leading to dimensionless parameters \(\text{Re} = 0.9\), \(\text{We} = 0.9\), and \(\text{Ma} = 1.05\). The time-dependent \(u_0\) component of velocity is shown in Figure 11. As with the rectangular box case, the vortex spreads out to fill the circle and decays over time. The transverse velocity, stress, and pressure corresponding to the final image of \(u_0\) are shown in Figure 12. The solution error convergence after 200 fine time steps is given in Table 6.
Figure 9. Time-dependent $u_n$ profiles of a hybrid fluid with a vortex initial condition in a rectangle. The domain has $256 \times 256$ cells with 30 time step increments using $\Delta t = 1.6 \times 10^{-3}$. The range is from $-0.25$ (red) to $0.25$ (blue).

Figure 10. Profiles for a hybrid fluid in a rectangle at $t = 0.144$ (last image in Figure 9). Clockwise from top left: $u_t$, $-0.15$ (red) to $0.15$ (blue); normal stress $\tau_{nn}$, $-0.25$ (red) to $0.37$ (blue); normal stress $\tau_{tt}$, $-0.25$ (red) to $0.37$ (blue); shear stress $\tau_{tn}$, $-0.30$ (red) to $0.29$ (blue); hydrostatic pressure $p$, $0$ (red) to $0.55$ (blue).

Figure 11. Time-dependent $u_0$ profiles of a hybrid fluid with a vortex initial condition in a disk. The domain has $128 \times 128$ cells with 20 time step increments using $\Delta t = 1.6 \times 10^{-3}$. The range is from $-0.25$ (red) to $0.25$ (blue).
<table>
<thead>
<tr>
<th>norm</th>
<th>Variable</th>
<th>Coarse Error</th>
<th>Fine Error</th>
<th>Order</th>
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<td>$\tau_{10}$</td>
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<tr>
<td></td>
<td>$\tau_{11}$</td>
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<td>$p$</td>
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<td>7.53e-05</td>
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<td>2.02e-04</td>
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<td>$\tau_{00}$</td>
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<td>$\tau_{10}$</td>
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<td></td>
<td>$\tau_{11}$</td>
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<tr>
<td></td>
<td>$p$</td>
<td>3.82e-04</td>
<td>1.28e-04</td>
<td>1.58</td>
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<tr>
<td>$L_\infty$</td>
<td>$u_0$</td>
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<td>3.26e-03</td>
<td>1.05</td>
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<td>$u_1$</td>
<td>6.82e-03</td>
<td>3.29e-03</td>
<td>1.05</td>
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<td>$\tau_{00}$</td>
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<td>$\tau_{11}$</td>
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<td>1.06</td>
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</table>

Table 5. Solution error convergence rates for a hybrid fluid with a vortex initial condition in a rectangle. Data correspond to Figures 9 and 10.

Figure 12. Profiles for a hybrid fluid in a disk at $t = 0.096$ (last image in Figure 11). Clockwise from top left: $u_1$, $-0.15$ (red) to $0.15$ (blue); normal stress $\tau_{00}$, $-0.25$ (red) to $0.35$ (blue); normal stress $\tau_{11}$, $-0.25$ (red) to $0.35$ (blue); shear stress $\tau_{10}$, $-0.30$ (red) to $0.30$ (blue); hydrostatic pressure $p$, $0$ (red) to $0.040$ (blue).
Table 6. Solution error convergence rates for a hybrid fluid with a vortex initial condition in a disk. Data correspond to Figures 11 and 12.

### 7. Conclusions

For each of the test problems, we demonstrate second-order convergence of the solution error in $L^1$ and first-order in $L^\infty$ for velocity and stress with an advective CFL time step constraint of $\text{CFL} \approx 0.5$, as expected. This is an improvement over [30], in which less than second-order convergence was obtained with a smaller time step, and the algorithm did not support arbitrary smooth geometries. The algorithm also exhibits at least first-order convergence in $L^1$ for pressure, as expected. In some cases, such as the Maxwell fluid in the rectangular geometry, the convergence rates in $L^\infty$ exceed first-order. This is due to the fact that given the position and shape of the expanded vortex, the largest magnitude errors occur in the interior of the domain, where the algorithm is second-order.

A feature calling for further study is the apparent need for additional projection filters (45) to smooth out the divergence in the velocity field of Maxwell fluids in irregular cells. Approaches include different filtering stencils, or different covered face state extrapolation algorithms.

The first obvious extension to this work is a three-dimensional discretization of the equations. The upwind method [6] and discretizations for Poisson’s equation and
the heat equation [29] in a three-dimensional embedded boundary framework have already been developed, so the extension is straightforward. The methods in this paper have been developed under the assumption that the geometry is sufficiently smooth. Additional studies are required to determine the robustness of the algorithm in the presence of discontinuous geometries, such as abrupt contractions. This will enable comparisons to standard benchmark problems [1; 26; 27; 30], such as the flow of elastic liquids in hard-cornered planar and axisymmetric contractions. Additional studies are also required to examine the robustness of this algorithm under higher values of We and Ma, and for a variety of operating conditions for experimental comparison [12; 13; 14]. In addition, adaptive numerical algorithms for the incompressible Navier–Stokes equations, in which the grid is locally refined in regions of interest, are being developed [21]. Adaptive techniques have already been used with success for hyperbolic conservation laws [6], so these two methods can be combined to develop a new adaptive projection method for incompressible viscoelasticity. Finally, another possible extension is the discretization of more advanced constitutive models, such as the PTT [25], White–Metzner [35] and Giesekus [11] models. The methods in this paper provide a framework for including the additional terms present in these models.

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References


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