Nonlinear finite volume method with discrete maximum principle for the two-phase flow model.

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Abstract: The discrete maximum principle is a meaningful requirement for numerical schemes used in multiphase flow models. It eliminates numerical pressure overshoots and undershoots, which may cause unnatural Darcy velocities and wrong numerical saturations. In this paper we study the application of the nonlinear finite volume method with discrete maximum principle [1] to the two-phase flow model. The method satisfies the discrete maximum principle for numerical pressures of incompressible fluids with neglected capillary pressure. For non-zero capillary pressure and constant phase viscosities the discrete maximum principle holds for numerical global pressure.

Introduction

The discrete maximum principle is a very desirable property of numerical schemes used for discretization of numerical models dealing with concentrations, densities, absolute temperatures. It provides non-negativity of solution an eliminates artificial sources and sinks. In multiphase flow models pressure gradient appears to be the main source of transport. Thus erroneous local extremum leads to wrong transport terms and consequently to wrong numerical solutions.

In this article we adapt nonlinear multi-point flux discretization scheme presented in [1, 2] for two-phase flow model. In the first section we describe the discretization scheme in application to diffusion equation. Two-phase flow models and fully implicit scheme for it are described in the third and fourth sections respectively. Then we prove that nonlinear multi-point flux discretization for the two-phase flow model leads to solution satisfying the discrete maximum principle. Then the Jacobian of the model residuals is computed and Newton iterations are described. Finally we present numerical experiments validating the discrete maximum principle of the solution. The solution is compared with solutions obtained using two-point linear and nonlinear schemes [4].

1 Nonlinear multi-point finite volume flux discretization

In this section we describe nonlinear multi-point flux discretization scheme (see [1, 2]) in application to diffusion problem.

Let Ω be a three-dimensional polyhedral domain with the Lipschitz boundary $\Gamma = \Gamma_N \cup \Gamma_D$. The diffusion equation for unknown variable p with the Dirichlet or Neumann boundary conditions written in a mixed form reads:

$$\mathbf{q} = -\mathbb{K}\nabla p, \qquad \text{div } \mathbf{q} = g \qquad \text{in } \Omega, \\ p = g_D \qquad \text{on } \Gamma_D, \\ \mathbf{n} \cdot \mathbf{q} = g_N \qquad \text{on } \Gamma_N.$$
(1)

Here $\mathbb{K}(\mathbf{x})$ is a symmetric positive definite (possibly anisotropic) diffusion tensor, $f(\mathbf{x})$ is a source term, $g_D(\mathbf{x})$ and $g_N(\mathbf{x})$ are given Dirichlet and Neumann boundary conditions for Γ_D and Γ_N parts of the boundary, consequently.

The weak maximum principle [6] for elliptic problems implies the maximum principle for diffusion equation solution. The last one states that for $f \leq 0$ the variable $p(\mathbf{x})$ satisfies:

$$\max_{\mathbf{x}\in\bar{\Omega}} p(\mathbf{x}) \le \max_{\mathbf{x}\in\Gamma_D\cup\Gamma_N} p(\mathbf{x}).$$

The minimum principle is formulated accordingly: for $f \ge 0$ the pressure $p(\mathbf{x})$ satisfies:

$$\min_{\mathbf{x}\in\bar{\Omega}} p(\mathbf{x}) \geq \min_{\mathbf{x}\in\Gamma_D\cup\Gamma_N} p(\mathbf{x}).$$

Note, that if $g_N \ge 0$ then maximum principle states additionally, that p cannot attain maximum on Γ_N . If $g_n \le 0$ then p cannot attain minimum on Γ_N .

1.1 Nonlinear FV discretization scheme

The FV scheme uses one degree of freedom per cell T, p_T , collocated at cell barycenter \mathbf{x}_T . Integrating the mass balance equation (1) over T and using the divergence theorem, we obtain:

$$\sum_{f \in \partial T} \sigma_{T,f} \, \mathbf{q}_f \cdot \mathbf{n}_f = \int_T g \, \mathrm{d}x, \qquad \mathbf{q}_f = \frac{1}{|f|} \int_f \mathbf{q} \, \mathrm{d}s, \tag{2}$$

where $\mathbf{q}_f \cdot \mathbf{n}_f$ is the total flux across face f, and $\sigma_{T,f}$ is either 1 or -1 depending on the mutual orientation of the normal vectors \mathbf{n}_f and \mathbf{n}_T (\mathbf{n}_T denotes the outward normal vector), $|n_f| = |f|$ and |f| denotes area of face f.

Both nonlinear discretization schemes (multi-point and two-point) exploit the same idea. The monotone two-point scheme was described in details in [3, 4]. Here we describe only the multi-point discretization.

1.2 Multi-point flux discretization



Figure 1: Two representations of co-normal vector $l_f = \mathbb{K} \cdot \mathbf{n}_f$ (2D example).

First for each cell-face pair we need to find a triplet (see Fig. 1 for a 2D example), a set of three vectors \mathbf{t}_* such that for the co-normal vector $l_f = \mathbb{K} \cdot \mathbf{n}_f$ we have

$$l_f = \alpha \mathbf{t}_1 + \beta \mathbf{t}_2 + \gamma \mathbf{t}_3, \tag{3}$$

where coefficients α , β and γ are non-negative.

Since the flux normal component is in fact the directional derivative along the co-normal vector l_f , it can also be represented as the sum of three derivatives along \mathbf{t}_* which are approximated by central differences:

$$q_{+} = \alpha'_{+} (p_{+} - p_{+,1}) + \beta'_{+} (p_{+} - p_{+,2}) + \gamma'_{+} (p_{+} - p_{+,3}),$$
(4)

where coefficients $\alpha'_{+}, \beta'_{+}, \gamma'_{+}$ are normalized by $|t_{+,i}|/|l_f|$ coefficients (3) for cell T_{+} .

For the opposite co-normal vector $-l_f$ we have the similar representation with non-negative coefficients:

$$q_{-} = \alpha'_{-} (p_{-} - p_{-,1}) + \beta'_{-} (p_{-} - p_{-,2}) + \gamma'_{-} (p_{-} - p_{-,3}).$$
(5)

Now we can take a linear combination of (4) and (5) with non-negative coefficients μ_+ and μ_- :

$$\mathbf{q}_f \cdot \mathbf{n}_f = \mu_+ q_+ + \mu_-(-q_-) \tag{6}$$

Flux approximation requires the linear combination of μ_+ , μ_- to be convex:

$$\mu_{+} + \mu_{-} = 1. \tag{7}$$

To construct the *multi-point* nonlinear discretization, we set equal two representation of the flux:

$$\mu_{+}q_{+} = -\mu_{-}q_{-}.$$
(8)

The solution of equations (7) and (8) for 2D case is considered in detail in [1]. Here we use these results in 3D case.

If $|q_+| = |q_-| = 0$ then the solution of (7),(8) in not unique. In this case we choose $\mu_+ = \mu_- = 1/2$. Otherwise consider two cases. The first case is for $q_+q_- \leq 0$. The solution is

$$\mu_{+} = \frac{q_{-}}{q_{-} - q_{+}}, \qquad \mu_{-} = \frac{q_{+}}{q_{+} - q_{-}},$$

and μ_{\pm} are non-negative. Thus,

$$\mathbf{q}_f \cdot \mathbf{n}_f = \frac{2q_+q_-}{q_- - q_+} = -\frac{2q_-q_+}{q_+ - q_-},$$

and numerical flux has two equivalent algebraic representations:

$$\mathbf{q}_f \cdot \mathbf{n}_f = 2\mu_+ \Big(\alpha'_+(p_+ - p_{+,1}) + \beta'_+(p_+ - p_{+,2}) + \gamma'_+(p_+ - p_{+,3}) \Big) = (9)$$

$$= A_{+,1}(p_+ - p_{+,1}) + A_{+,2}(p_+ - p_{+,2}) + A_{+,3}(p_+ - p_{+,3}),$$

and

$$-\mathbf{q}_{f} \cdot \mathbf{n}_{f} = 2\mu_{-} \left(\alpha'_{-}(p_{-} - p_{-,1}) + \beta'_{-}(p_{-} - p_{-,2}) + \gamma'_{-}(p_{-} - p_{-,3}) \right) = (10)$$
$$= A_{-,1}(p_{-} - p_{-,1}) + A_{-,2}(p_{-} - p_{-,2}) + A_{-,3}(p_{-} - p_{-,3}),$$

with non-negative coefficients $A_{\pm,k}$, k = 1, 2, 3. Note that these coefficients depend on fluxes and hence on values of p at neighboring cells. Thus, the resulting multi-point flux approximation is *nonlinear* and its stencil is embedded into the set of the closest neighboring cells.

The second case, $q_+q_- > 0$, leads to a potentially degenerate diffusive flux. In order to avoid this degeneracy, the authors in [5] re-group the terms in equation (6) as follows:

$$\mathbf{q}_f \cdot \mathbf{n}_f = \mu_+ \widetilde{q}_+ + \mu_- (-\widetilde{q}_-) + (\mu_+ \alpha'_+ + \mu_- \alpha'_-)(p_+ - p_-),$$

where $\tilde{q}_{+} = \beta'_{+}(p_{+}-p_{+,2}) + \gamma'_{+}(p_{+}-p_{+,3}), \quad \tilde{q}_{-} = \beta'_{-}(p_{-}-p_{-,2}) + \gamma'_{-}(p_{-}-p_{-,3}).$ We find μ_{+} and μ_{-} as a solution of

$$\widetilde{q}_+\mu_+ + \widetilde{q}_-\mu_- = 0,$$

$$\mu_+ + \mu_- = 1.$$

Again if the solution is not unique, we choose $\mu_+ = \mu_- = 1/2$. Otherwise, for $\tilde{q}_+ \tilde{q}_- \leq 0$ we get

$$\mathbf{q}_{f} \cdot \mathbf{n}_{f} = 2\mu_{+}\widetilde{q}_{+} + (\mu_{+}\alpha'_{+} + \mu_{-}\alpha'_{-})(p_{+} - p_{-}) = A_{+,1}(p_{+} - p_{+,1}) + A_{+,2}(p_{+} - p_{+,2}) + A_{+,3}(p_{+} - p_{+,3}) = -2\mu_{-}\widetilde{q}_{-} - (\mu_{+}\alpha'_{+} + \mu_{-}\alpha'_{-})(p_{-} - p_{+}) = -A_{-,1}(p_{-} - p_{-,1}) - A_{-,2}(p_{-} - p_{-,2}) - A_{-,3}(p_{-} - p_{-,3}),$$
(11)

where $A_{+,1} = A_{-,1} = \mu_+ \alpha_{+,1} + \mu_- \alpha_{-,1}$. If $\tilde{q}_+ \tilde{q}_- > 0$

$$q_f = (\mu_+ \alpha'_+ + \mu_- \alpha'_-)(p_+ - p_-) = A_{+,1}(p_+ - p_-).$$

Calculating $A_{\pm,k}$ and inserting the diffusive fluxes $\mathbf{q}_f \cdot \mathbf{n}_f$ in the mass balance equation, we get an algebraic problem. For a logically cubic mesh, the considered finite volume scheme results in the conventional seven-point stencil.

2 Two-phase black oil model

We consider a two-phase flow of immiscible fluids in a porous medium [7, 8]. The phase, that wets the medium more than the other, is called wetting phase and is indicated by subscript w. The other phase is the non-wetting phase and indicated by o.

The basic equations for the two-phase flow are the following:

• Mass conservation for each phase:

$$\frac{\partial \phi \rho_{\alpha} S_{\alpha}}{\partial t} = -\nabla \cdot (\rho_{\alpha} \mathbf{u}_{\alpha}) + q_{\alpha}, \quad \alpha = w, o.$$
(12)

• Darcy's law:

$$\mathbf{u}_{\alpha} = -\frac{k_{r\alpha}}{\mu_{\alpha}} \mathbb{K} \left(\nabla p_{\alpha} - \rho_{\alpha} g \nabla z \right), \quad \alpha = w, o.$$
(13)

• Two fluids fill the voids:

$$S_w + S_o = 1, \tag{14}$$

• Pressure difference between phases is given by capillary pressure:

$$p_o - p_w = p_c(S_w),\tag{15}$$

where \mathbb{K} is the absolute permeability tensor, ρ_{α} is the phase density, μ_{α} is the viscosity, $k_{r\alpha}$ is the relative phase permeability, ϕ is the porosity, g is the gravity term and q_{α} is the source/sink well term.

On the reservoir boundary we consider no-flow (homogeneous Neumann) condition.

Wells are introduced through one of the following conditions: either given bottom hole pressure p_{bh} , or given component flux q_{α} , or given total flux $q_T = q_w + q_o$ for producer wells.

All the test cases in this article were performed for vertical perfect wells. The formula for the well term was suggested by Peaceman [9]. For a cell T with center \mathbf{x}_T connected to the well we have:

$$q_{\alpha} = \frac{k_{r\alpha}}{\mu_{\alpha}} WI \Big(p_{bh} - p_{\alpha} - \rho_{\alpha} g(z_{bh} - z) \Big) \delta(\mathbf{x} - \mathbf{x}_{T}), \tag{16}$$

WI is the *well index* which does not depend on the properties of fluids but depends on properties of the media, $\delta(\mathbf{x} - \mathbf{x}_T)$ is the Dirac function.

Choosing oil pressure $P = p_o$ and water saturation $S = S_w$ as primary unknowns, defining a formation volume factor $B_{\alpha} = \rho_{\alpha,0}/\rho_{\alpha}$, and mobility $\lambda_{\alpha} = k_{r\alpha}/(\mu_{\alpha}B_{\alpha})$ and using (12)-(15) we can get the equivalent formulation:

$$\frac{\partial}{\partial t}\frac{\phi S}{B_w} + \nabla \cdot \mathbb{K}\lambda_w (\nabla P - \nabla P_c(S) - \frac{\rho_{w,0}}{B_w}g\nabla z) = \frac{q_w}{\rho_{w,0}},\tag{17}$$

$$\frac{\partial}{\partial t}\frac{\phi(1-S)}{B_o} + \nabla \cdot \mathbb{K}\lambda_o(\nabla P - \frac{\rho_{o,0}}{B_o}g\nabla z) = \frac{q_o}{\rho_{o,0}}.$$
(18)

In the discrete counterparts of (12)-(13) the mobilities $\lambda_{\alpha}(S, p)$ on the face f_{ij} are taken upwinded:

$$\lambda_{\alpha}(S) = \begin{cases} \lambda_{\alpha}(S_i, p_i) & \text{if flow is directed from cell } i \text{ to cell } j, \\ \lambda_{\alpha}(S_j, p_j) & \text{if flow is directed from cell } j \text{ to cell } i. \end{cases}$$

The phase mobilities for well-producer are taken upwinded from the cell. For well-injector we have only water injected and thus take the downstream mobility from the cell with the well: $\lambda_{inj} = (\frac{k_{rw}}{\mu_w B_w} + \frac{k_{ro}}{\mu_o B_o})_{cell}$. We also assume that there is no capillary pressure in wells, so all the well fluxes depend on the same (oil) pressure.

3 Fully implicit scheme for two-phase flow equations

First we apply the implicit scheme to the mass conservation equations (12):

$$\frac{\left(\frac{\phi S_{\alpha}}{B_{\alpha}}\right)^{n+1} - \left(\frac{\phi S_{\alpha}}{B_{\alpha}}\right)^{n}}{\Delta t^{n+1}} = -\operatorname{div}(\mathbf{u}_{\alpha}^{n+1}) + \left(\frac{q_{\alpha}}{\rho_{\alpha,0}}\right)^{n+1}, \quad \alpha = w, o.$$
(19)

Now we can write down the nonlinear residual equations for the l^{th} approximation to a quantity evaluated at time step n + 1 inside grid cell T_i :

$$R_{\alpha,i}^{l} = \int_{T_{i}} \left[\left(\frac{\phi S_{\alpha}}{B_{\alpha}} \right)_{i}^{l} - \left(\frac{\phi S_{\alpha}}{B_{\alpha}} \right)_{i}^{n} + \Delta t^{n+1} \left(\operatorname{div} \mathbf{u}_{\alpha} - \frac{q_{\alpha}}{\rho_{\alpha,0}} \right)_{i}^{l} \right] \, \mathrm{d}x, \quad \alpha = w, o.$$

$$\tag{20}$$

The discrete counterpart of (19) can be written as:

$$R_{\alpha,i} = 0, \quad \alpha = w, o \tag{21}$$

for all grid cells at every time step.

4 Discrete maximum principle

In order to derive discrete maximum principle we introduce assumptions on two-phase flow model coefficients:

- (a1) constant water and oil densities: $B_{\alpha} = 1, \ \alpha = w, o,$
- (a2) constant porosity: $\phi = const$,
- (a3) capillary pressure is neglected: $p_c \equiv 0$,
- (a4) gravitational term is neglected: $u_{\alpha} = -\frac{k_{r\alpha}}{\mu_{\alpha}}\mathbb{K}(\nabla p_{\alpha}), \ \alpha = w, o.$

Using summation over α in (21) and assumptions (a1–a4) we get

$$-\int_{T_i} \operatorname{div} \left(\mathbb{K}\lambda \nabla p_o\right)^l dx = \int_{T_i} \left(\frac{q_w}{\rho_w} + \frac{q_o}{\rho_o}\right)^l dx,$$
(22)

where $\lambda = \frac{k_{rw}}{B_w \mu_w} + \frac{k_{ro}}{B_o \mu_o}$.

Using divergence theorem we rewrite (22):

$$\sum_{f \in \partial T} \lambda_f^l \boldsymbol{q}_f^l \cdot \boldsymbol{n}_f = \int_{T_i} \left(\frac{q_w}{\rho_w} + \frac{q_o}{\rho_o} \right)^l dx \tag{23}$$

where n_f is the exterior normal vector to face f, $|n_f| = |f|$, |f| is the area of face f, and q_f is the flux over face f. The discrete flux normal component $q_f \cdot n_f$ may be defined via the pressure differences according to section 1.2:

$$\sum_{f \in \partial T} \sum_{i=1}^{3} \lambda_{f}^{l} A_{i,f}(p_{T}^{l} - p_{T,i}^{l}) = \int_{T_{i}} \left(\frac{q_{w}}{\rho_{w}} + \frac{q_{o}}{\rho_{o}}\right)^{l} dx.$$
(24)

Let \mathcal{T} stand for the set of all cells and \mathcal{T}_D and \mathcal{T}_N stand for the set of cells with Dirichlet boundary faces and Neumann boundary faces, respectively. We also denote all boundary cells by $\mathcal{T}_B = \mathcal{T}_D \cup \mathcal{T}_N$. **Theorem 1** (Discrete maximum principle). Let the solution p to (24) exist. Let \mathcal{T}_{inj} be the set of cells where $\frac{q_w}{\rho_w} + \frac{q_o}{\rho_o} > 0$. Then

$$\max_{T \in \mathcal{T} \setminus (\mathcal{T}_{inj} \cup \mathcal{T}_B)} p_T \le p_{max} = \max_{\mathcal{T}_{inj} \cup \mathcal{T}_B} p_T.$$
(25)

Proof. Assume that p_T in a cell $T \in \mathcal{T} \setminus (\mathcal{T}_{inj} \cup \mathcal{T}_B)$ has the maximum value.

Coefficients $A_{i,f}$ in (24) are nonnegative by construction. As p_T has the maximum value, from non-negativity of coefficients $A_{i,f}$ and the form of equation (24) one derives that $\boldsymbol{q}_f^l \cdot \boldsymbol{n}_f$ are non-negative. λ is a positive coefficient. Thus the left hand side of equation (24) is non-negative, and the right of (24) is non-positive in $\mathcal{T} \setminus (\mathcal{T}_{inj} \cup \mathcal{T}_B)$. Then both sides of the equation are equal to zero.

Since $\lambda_f \boldsymbol{q}_f^l \cdot \boldsymbol{n}_f$ are non-negative and their sum is zero, we get $\boldsymbol{q}_f^l \cdot \boldsymbol{n}_f = 0$ $\forall f \in \partial T$. Since the discrete normal flux on face f contains the term $p_T - p_{T'}$ where T' is the cell neighbouring T through f, we conclude that $p_T = p_{T'}$.

Assuming that the mesh $\mathcal{T}/\mathcal{T}_{inj}$ is face-connected, we get that p is constant in $\mathcal{T} \setminus (\mathcal{T}_{inj} \cup \mathcal{T}_B)$.

Remark 1. If we have no-flow Neumann condition for p on face f or positive Neumann condition (outflux), flow over this face $q_f^l \cdot n_f$ is non-negative and the proof of the theorem 1 remains correct for this case. Thus p also cannot attain maximum on Neuman boundary (otherwise $p \equiv \text{const}$ in the whole domain).

Remark 2. The authors of [1] also construct nonlinear multi-point flux discretization for discontinious diffusion tensor \mathbb{K} . Under assumptions (a1-a4) the discrete maximum principle for solution of the two-phase flow model also holds for discontinious case. The proof is similar to the proof of the theorem 1.

5 Nonlinear coefficients variation

If we use a Jacobian-based nonlinear solver, we need a variation of coefficients $A_{\pm,i}$ in flux discretizations.

First we write variations for q_{\pm} and \tilde{q}_{\pm} :

$$\Delta q_{\pm} = (\alpha'_{\pm} + \beta'_{\pm} + \gamma'_{\pm})p_{\pm} - \alpha'_{\pm}\Delta p_{\pm,1} - \beta'_{\pm}\Delta p_{\pm,2} - \gamma'_{\pm}\Delta p_{\pm,3},$$

$$\Delta \widetilde{q}_{\pm} = (\beta'_{\pm} + \gamma'_{\pm})p_{\pm} - \beta'_{\pm}\Delta p_{\pm,2} - \gamma'_{\pm}\Delta p_{\pm,3}.$$

Variations for μ_{\pm} and $A_{\pm,i}$ depend on the q_+, q_- . If $|q_+| + |q_-| > 0$ and $q_+q_- \le 0$, then

$$\Delta \mu_{\pm} = \frac{\Delta q_{\mp}}{q_{\mp} - q_{\pm}} - (\Delta q_{\mp} - \Delta q_{\pm}) \frac{q_{\mp}}{(q_{\mp} - q_{\pm})^2},$$
$$\Delta A_{\pm,1} = 2\alpha'_{\pm} \Delta \mu_{\pm}, \qquad \Delta A_{\pm,2} = 2\beta'_{\pm} \Delta \mu_{\pm}, \qquad \Delta A_{\pm,3} = 2\gamma'_{\pm} \Delta \mu_{\pm}.$$

Otherwise variation for μ_{\pm} depends on the \tilde{q}_+, \tilde{q}_- . If $|\tilde{q}_+| + |\tilde{q}_-| > 0$ and $\tilde{q}_+\tilde{q}_- \leq 0$, then

$$\begin{split} \Delta \mu_{\pm} &= \frac{\Delta \widetilde{q}_{\mp}}{\widetilde{q}_{\mp} - \widetilde{q}_{\pm}} - (\Delta \widetilde{q}_{\mp} - \Delta \widetilde{q}_{\pm}) \frac{\widetilde{q}_{\mp}}{(\widetilde{q}_{\mp} - \widetilde{q}_{\pm})^2}, \\ \Delta A_{\pm,1} &= \alpha'_{+} \Delta \mu_{+} + \alpha'_{-} \Delta \mu_{-}, \qquad \Delta A_{\pm,2} = 2\beta'_{\pm} \Delta \mu_{\pm}, \qquad \Delta A_{\pm,3} = 2\gamma'_{\pm} \Delta \mu_{\pm}. \\ \text{If } |\widetilde{q}_{+}| + |\widetilde{q}_{-}| > 0 \text{ and } \widetilde{q}_{+} \widetilde{q}_{-} > 0, \text{ then} \end{split}$$

$$\Delta \mu_{\pm} = \frac{\Delta \widetilde{q}_{\mp}}{\widetilde{q}_{\mp} - \widetilde{q}_{\pm}} - (\Delta \widetilde{q}_{\mp} - \Delta \widetilde{q}_{\pm}) \frac{\widetilde{q}_{\mp}}{(\widetilde{q}_{\mp} - \widetilde{q}_{\pm})^2},$$
$$\Delta A_{\pm,1} = \alpha'_{+} \Delta \mu_{+} + \alpha'_{-} \Delta \mu_{-}, \qquad \Delta A_{\pm,2} = 0, \qquad \Delta A_{\pm,3} = 0.$$

In the case of $|q_+| = |q_-| = 0$ or $|\tilde{q}_+| = |\tilde{q}_-| = 0$ we refer to remark 3. The final variation for q_{\pm} is

$$\Delta q_{\pm} = (A_{\pm,1} + A_{\pm,2} + A_{\pm,3}) \Delta p_{\pm} - A_{\pm,1} \Delta p_{\pm,1} - A_{\pm,2} \Delta p_{\pm,2} - A_{\pm,3} \Delta p_{\pm,3} + \Delta A_{\pm,1} (p_{\pm} - p_{\pm,1}) + \Delta A_{\pm,2} (p_{\pm} - p_{\pm,2}) + \Delta A_{\pm,3} (p_{\pm} - p_{\pm,3}).$$

Remark 3. Coefficients μ_{\pm} in case of zero fluxes $|q_{+}| = |q_{-}| = 0$ or zero modified fluxes $|\tilde{q}_{+}| = |\tilde{q}_{-}| = 0$ have discontinuities with respect to $p_{\pm,i}$. These cases pose a technical problem for the Newton solver. Possible solution of this problem is to induce small perturbation in the pressure field. Alternatively one can use a solver without usage of Jacobian for one timestep.

6 Newton method for two-phase flow model

We suggest to use Newton's method to solve nonlinear system (21) with Darcy velocities (13):

$$J(x^l)\delta x^l = -R(x^l), \tag{26}$$

$$x^{l+1} = x^l + \delta x^l, \tag{27}$$

where l is the l^{th} Newton step, x is a vector of primary unknowns in all grid cells,

$$x = \left(\begin{array}{c} p_o \\ S_w \end{array}\right),$$

R is the vector of nonlinear residuals in all grid cells,

$$R(x) = \left(\begin{array}{c} R_w(x) \\ R_o(x) \end{array}\right),$$

and J is the Jacobian matrix:

$$J(x) = \begin{pmatrix} \frac{\partial R_w}{\partial p_o}(x) & \frac{\partial R_w}{\partial S_w}(x) \\ \frac{\partial R_o}{\partial p_o}(x) & \frac{\partial R_o}{\partial S_w}(x) \end{pmatrix}.$$

We terminate Newton's method when the norm of the residual vector drops below ε_{nwt} .

Below we consider the construction of Jacobian matrix. We divide the residuals into two parts: accumulation (including well terms) and transport, $R_{\alpha,i} = R_{\alpha,i}^{acc} + R_{\alpha,i}^{trans}$, where:

$$\begin{split} R_{\alpha,i}^{acc} &= V_i \left[\left(\frac{\phi S_\alpha}{B_\alpha} \right)_i^l - \left(\frac{\phi S_\alpha}{B_\alpha} \right)_i^n \right] - \Delta t^{n+1} \left(\frac{q_\alpha}{\rho_{\alpha,0}} \right)_i^l, \quad \alpha = w, o, \\ R_{\alpha,i}^{trans} &= \Delta t^{n+1} \int_{T_i} (\operatorname{div} \mathbf{u}_\alpha) \, \mathrm{d}x, \quad \alpha = w, o. \end{split}$$

We also take advantage of the following dependencies:

- $S_o = 1 S_w$, (see eq. (14)),
- $p_w = p_o p_c(S_w)$, with piecewise linear function $p_c(S_w)$ given by table data (see eq. (15)),
- $k_{r\alpha} = k_{r\alpha}(S_w)$ are piecewise linear functions given by table data,
- $\mu_{\alpha} = \mu_{\alpha}(p_o)$ are piecewise linear functions given by table data,
- $B_{\alpha} = B_{\alpha}(p_o)$ are piecewise linear functions given by table data,

•
$$\phi = \phi (1 + c_R (p_o - p_o^0)).$$

6.0.1 Accumulation term

The variation of the accumulation term is as follows:

$$\Delta R_{w,i}^{acc} = V_i \left[\Delta \left(\frac{\phi S_w}{B_w} \right) \right] - \Delta t^{n+1} \Delta \left(\frac{q_w}{\rho_{w,0}} \right),$$
$$\Delta R_{o,i}^{acc} = V_i \left[\Delta \left(\frac{\phi S_o}{B_o} \right) \right] - \Delta t^{n+1} \Delta \left(\frac{q_o}{\rho_{o,0}} \right),$$

where

$$\Delta\left(\frac{\phi S_w}{B_w}\right) = \frac{\phi}{B_w}\Delta S_w + S_w\left(\frac{c_R}{B_w} - \frac{\phi}{B_w^2}\frac{dB_w}{dp_o}\right)\Delta p_o,$$

$$\Delta\left(\frac{\phi S_o}{B_o}\right) = -\frac{\phi}{B_o}\Delta S_w + (1 - S_w)\left(\frac{c_R}{B_o} - \frac{\phi}{B_o^2}\frac{dB_o}{dp_o}\right)\Delta p_o.$$

6.0.2 Transport term

Now we consider the transport term composed of Darcy fluxes

$$R_{\alpha,i}^{trans} = \Delta t^{n+1} \int_{\underline{\Gamma}_i} (\mathbf{u}_{\alpha} \cdot \mathbf{n}) \, \mathrm{d}s \approx \Delta t^{n+1} \sum_{f \in \underline{\Gamma}_i} \mathbf{u}_{\alpha,f} \cdot \mathbf{n}_f.$$

If we use the multi-point discretization of the flux, then we have two different representations of flux over face f for two neighbouring cells T_{\pm} :

$$(\mathbf{u}_{w,f}^{h} \cdot \mathbf{n}_{f})_{\pm} = - \left(\frac{k_{rw}}{\mu_{w}B_{w}}\right)_{f} \left(A_{\pm,1}^{p_{o}}(p_{o,\pm} - p_{o,\pm,1}) + A_{\pm,2}^{p_{o}}(p_{o,\pm} - p_{o,\pm,2}) + A_{\pm,3}^{p_{o}}(p_{o,\pm} - p_{o,\pm,3})\right) \\ + \left(\frac{k_{rw}}{\mu_{w}B_{w}}\right)_{f} \left(A_{\pm,1}^{p_{c}}(p_{c,\pm} - p_{c,\pm,1}) + A_{\pm,2}^{p_{c}}(p_{c,\pm} - p_{c,\pm,2}) + A_{\pm,3}^{p_{c}}(p_{c,\pm} - p_{c,\pm,3})\right) \\ + \left(\frac{k_{rw}}{\mu_{w}B_{w}^{2}}\right)_{f} \left(\rho_{w,0}g(A_{\pm,1}^{z}(z_{\pm} - z_{\pm,1}) + A_{\pm,2}^{z}(z_{\pm} - z_{\pm,2}) + A_{\pm,3}^{z}(z_{\pm} - z_{\pm,3})\right)),$$

$$(28)$$

$$(\mathbf{u}_{o,f}^{h} \cdot \mathbf{n}_{f})_{\pm} = - \left(\frac{k_{ro}}{\mu_{o}B_{o}}\right)_{f} \left(A_{\pm,1}^{p_{o}}(p_{o,\pm} - p_{o,\pm,1}) + A_{\pm,2}^{p_{o}}(p_{o,\pm} - p_{o,\pm,2}) + A_{\pm,3}^{p_{o}}(p_{o,\pm} - p_{o,\pm,3})\right) \\ + \left(\frac{k_{ro}}{\mu_{o}B_{o}^{2}}\right)_{f} \left(\rho_{o,0}g\left(A_{\pm,1}^{z}(z_{\pm} - z_{\pm,1}) + A_{\pm,2}^{z}(z_{\pm} - z_{\pm,2}) + A_{\pm,3}^{z}(z_{\pm} - z_{\pm,3})\right)\right).$$

$$(29)$$

Here $k_{r\alpha} = k_{r\alpha}(\tilde{S})$, \tilde{S} is the upwinded value of water saturation on face fand $B_{\alpha} = B_{\alpha}(\tilde{p})$, $\mu_{\alpha} = \mu_{\alpha}(\tilde{p})$, \tilde{p} is the upwinded value of oil pressure on face f and coefficients $A_{\pm,i}$ depend on variable value in the neighbouring cells:

$$A_{\pm,i}^{p_{\alpha}} = A_{\pm,i}^{p_{\alpha}}(p_{\alpha,+,1}, p_{\alpha,+,2}, p_{\alpha,+,3}, p_{\alpha,-,1}, p_{\alpha,-,2}, p_{\alpha,-,3}),$$
$$A_{\pm,i}^{z} = A_{\pm,i}^{z}(z_{+,1}, z_{+,2}, z_{+,3}, z_{z-,1}, z_{z-,2}, z_{-,3}).$$

We define auxiliary variables and derivatives:

$$\begin{split} \lambda_{g,\alpha} &= \frac{k_{r\alpha}}{\mu_w B_w^2}, \\ &\frac{d\lambda_{g,\alpha}}{d\widetilde{S}} = \frac{d\lambda_{\alpha}}{d\widetilde{S}} / B_w, \\ &\frac{d\lambda_{g,\alpha}}{d\widetilde{p}} = \left(\frac{d\lambda_{\alpha}}{d\widetilde{p}} B_w - \lambda_{\alpha} \frac{dB_w}{d\widetilde{p}}\right) / B_w^2, \\ \mathcal{D}_1 &= A_{\pm,1}^{p_o}(p_{o,\pm} - p_{o,\pm,1}) + A_{\pm,2}^{p_o}(p_{o,\pm} - p_{o,\pm,2}) + A_{\pm,3}^{p_o}(p_{o,\pm} - p_{o,\pm,3}), \\ \mathcal{D}_2 &= A_{\pm,1}^{p_c}(p_{c,\pm} - p_{c,\pm,1}) + A_{\pm,2}^{p_c}(p_{c,\pm} - p_{c,\pm,2}) + A_{\pm,3}^{p_c}(p_{c,\pm} - p_{c,\pm,3}), \\ \mathcal{D}_{3,\alpha} &= \rho_{\alpha,0}g(A_{\pm,1}^z(z_{\pm} - z_{\pm,1}) + A_{\pm,2}^z(z_{\pm} - z_{\pm,2}) + A_{\pm,3}^z(z_{\pm} - z_{\pm,3})). \end{split}$$

Using (28) and (29) we get the following representation for the flux variation for each of two phases:

$$\Delta(\mathbf{u}_{w,f}^{h} \cdot \mathbf{n}_{f}) = \begin{bmatrix} \left(\frac{d\lambda_{w}}{d\tilde{S}_{w}}\right)\left(-\mathcal{D}_{1}+\mathcal{D}_{2}\right) + \frac{d\lambda_{g,w}}{d\tilde{S}_{w}}\mathcal{D}_{3,w} \end{bmatrix} \Delta \widetilde{S}_{w} + \\ \left[\left(\frac{d\lambda_{w}}{d\tilde{p}_{o}}\right)\left(-\mathcal{D}_{1}+\mathcal{D}_{2}\right) + \frac{d\lambda_{g,w}}{d\tilde{p}_{o}}\mathcal{D}_{3,w} \right] \Delta \widetilde{p}_{o} - \\ \lambda_{w}\left(\Delta q_{\pm}^{p_{o}}\right) + \\ \lambda_{w}\left(\Delta q_{\pm}^{p_{c}}\frac{dP_{c}}{dS_{w}}\Delta S_{w}\right) + \\ \lambda_{g,w}\rho_{w,0}g(\Delta q_{\pm}^{z}), \end{bmatrix} \Delta \widetilde{S}_{w} + \\ \left[\left(\frac{d\lambda_{o}}{d\tilde{S}_{w}}\right)\left(-\mathcal{D}_{1}+\mathcal{D}_{2}\right) + \frac{d\lambda_{g,o}}{d\tilde{S}_{w}}\mathcal{D}_{3,\alpha}\right] \Delta \widetilde{S}_{w} + \\ \left[\left(\frac{d\lambda_{o}}{d\tilde{p}_{o}}\right)\left(-\mathcal{D}_{1}+\mathcal{D}_{2}\right) + \frac{d\lambda_{g,o}}{d\tilde{p}_{o}}\mathcal{D}_{3,\alpha}\right] \Delta \widetilde{p}_{o} - \\ \lambda_{o}\left(\Delta q_{\pm}^{p_{o}}\right) + \\ \lambda_{g,o}\rho_{o,0}g(\Delta q_{\pm}^{z}), \end{bmatrix}$$
(30)

where variation for the flux q_{\pm}^{y} of the variable y is defined in section 5.

7 Numerical experiments

We consider two pseudo-2D numerical experiments, each uses three different flux discretization schemes: nonlinear multi-point, nonlinear two-point and linear two-point. Both test cases use the uniform $32 \times 32 \times 1$ grid for the computational domain $[-50; 50] \times [-50; 50] \times [4000; 4010]$. The first experiment is for one injecting and one producing well, and the second one is for three injecting and one producing well.



Figure 2: Schemes of the numerical experiments.

Well locations are presented on Fig. 2. Injecting well pressure is equal to 4100 and producing well pressure is 3900.

Both experiments have been conducted for incompressible phases with constant viscosities $\mu_w = 1$, $\mu_o = 50$, constant porosity $\phi = 0.2$, zero capillary pressure and zero gravity terms, as required by the derivation of the discrete maximum principle theorem 1. Relative permeabilities are shown of Fig. 3 and the absolute permeability tensor is:

$$k = R_z(-\theta_z)diag(k_1, k_2, k_3)R_z(\theta_z),$$

where $k_1 = k_3 = 100, k_2 = 0.1, \theta_z = 112.5^\circ$, and $R_z(\alpha)$ is the matrix of the rotation in xy-plane.

Pressure for the two wells problem (Fig. 2, left) is presented on Fig. 4. The nonlinear multi-point scheme and the linear two-point scheme satisfy the discrete maximum principle for pressure as expected, while the nonlinear two-point flux discretization violates the discrete maximum principle. It was demonstrated in [3] that the mesh refinement reduces the magnitude of the



Figure 3: Relative permeabilities.

overshoots and undershoots for the nonlinear two-point scheme, yet the DMP is still violated. In contrast to two nonlinear schemes, the linear two-point scheme does not provide approximation for the fluxes which results in nonphysical pressure field: the solution if not aligned with the main anisotropy direction. Therefore, only the nonlinear multi-point scheme is consistent and satisfies the discrete maximum principle.



Figure 4: Pressure for t = 2000 for different flux discretization schemes (numerical experiment with 2 wells). Overshoot (p > 4100) and undershoot (p < 3900) domains are denoted by O and U respectively.

The four wells test case (Fig. 2, right) demonstrates the similar pressure solution violating the discrete maximum principle and the consequences of pressure DMP violation for the saturation solution. Fig. 5 shows the pressure solution for the time t = 100. Again, the discrete maximum principle is violated by the two-point nonlinear scheme in the vast region. The top left injecting well cell has the overshoot and the pressure is higher than the well bottom hole pressure 4100. As a result the top left injecting well starts to pump out water instaed of injecting. The nonphysical flow is magnified by the use of the downstream discretization of well mobility $\lambda_{inj} = (\lambda_w + \lambda_{inj})$ λ_o _{cell} > 0, and water saturation of the well cell decreases even after it becomes zero. This leads to the nonphysical negative saturation on the top left well cell (see Fig. 6, middle), water saturation on this cell drops up to -0.800. Water saturations obtained by the multi-point nonlinear scheme, the nonlinear two-point scheme and the two-point linear schemes are compared in Fig. 6. Following the lack of the flux approximation of the linear two-point scheme the water flow is not alighed with the main anisotropy direction.



Figure 5: Pressure for t = 100 for different flux discretization schemes (numerical experiment with 4 wells). Overshoot (p > 4100) and undershoot (p < 3900) domains are denoted by O and U respectively.

8 Conclusion



Figure 6: Water saturation for t = 100 for different flux discretization schemes (numerical experiment with 4 wells). Bottom panel is colored according to extended saturation legend. Initial saturation is s(0) = 0.15.

It was proved that under certain assumptions the nonlinear multi-point flux discretization scheme satisfies the discrete maximum principle for twophase flow model. The scheme was implemented and the numerical results demonstrate that the solution for the pressure satisfies the discrete maximum principle in the cases where the monotone nonlinear two-point scheme fails. Unlike the conventional two-point linear scheme the nonlinear multi-point scheme preserves the flux approximation for the non-orthogonal grids and anisotropic permeability tensors. It was shown that the new scheme may be used with the Newton solver in fully implicit time discretization.

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