



# Supplementary Materials: Comparative Analyses between the Zero-Inertia and Fully Dynamic Models of the Shallow Water Equations for Unsteady Overland Flow Propagation

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## Part A: The numerical procedures

#### A.1 The fractional time step procedure

The two numerical models applied in this paper are based on a time-splitting approach, where a prediction and a correction problem are sequentially solved. For both the numerical procedures, it has been shown that the prediction and the correction steps have the characteristics of a convective and a diffusive problem, respectively [2–5]. For these reasons, we call the prediction system "convective prediction" (CP) system and the correction system "diffusive correction" (DC) system.

Let's assume a general system of balance laws,

$$\frac{\partial \mathbf{U}}{\partial t} + \nabla \cdot \mathbf{F}(\mathbf{U}) = \mathbf{B}(\mathbf{U})$$
 A.1

where **U**, **F**(**U**) and **B**(**U**) are the vector of the unknown variables, the flux vector and the source term respectively. Let  $F^{p}(\mathbf{U})$  and  $\mathbf{B}^{p}(\mathbf{U})$  be a suitable numerical flux and source term respectively, further defined, applying a fractional time-step procedure, we set:

$$\mathbf{F}(\mathbf{U}) = \mathbf{F}^{p}(\mathbf{U}) + \left(\mathbf{F}(\mathbf{U}) - \mathbf{F}^{p}(\mathbf{U})\right)$$
A.2,a

$$\mathbf{B}(\mathbf{U}) = \mathbf{B}^{p}(\mathbf{U}) + (\mathbf{B}(\mathbf{U}) - \mathbf{B}^{p}(\mathbf{U}))$$
A.2,b

Applying integration in time, we split system (A.1) in the two following ones:

$$\mathbf{U}^{k+1/2} - \mathbf{U}^k + \nabla \cdot \int_0^{\Delta t} \mathbf{F}^p \, dt = \int_0^{\Delta t} \mathbf{B}^p \, dt$$
A.3,a

$$\mathbf{U}^{k+1} - \mathbf{U}^{k+1/2} + \nabla \cdot \int_{0}^{\Delta t} \mathbf{F} \, dt - \nabla \cdot \overline{\mathbf{F}}^{p} \Delta t = \int_{0}^{\Delta t} \mathbf{B} \, dt - \overline{\mathbf{B}}^{p} \Delta t \qquad \text{A.3,b}$$

and we call systems (A.3,a) and (A.3,b) prediction and correction systems respectively. In Eqs (A.3), the overbar symbol marks the mean in time values of the numerical flux and source terms computed during the prediction step and apices k, k+1/2 and k+1 mark the values of the unknown variables computed at the beginning of the time step, at the end of the prediction step and at the end of the correction step respectively. The integrals  $\mathbf{\bar{F}}^p \Delta t$  and  $\mathbf{\bar{B}}^p \Delta t$  in Eq. (A.3,b) are computed "a posteriori" after the solution of the prediction problem, as explained in [2,3,5] for the fully dynamic model and in [4] for the diffusive model, respectively. Observe that summing systems (A.3,a) and (A.3,b), the integral of the original system (A.1) is formally obtained. The difference ( $\mathbf{U}^{k+1}-\mathbf{U}^{k+1/2}$ ) in Eq.

(A.3,b) is close to zero as far as the differences of the flux terms integrals  $\left(\nabla \cdot \int_{0}^{\Delta t} \mathbf{F} \, dt - \nabla \cdot \overline{\mathbf{F}}^{p} \Delta t\right)$  and

of the source terms  $\left(\int_{0}^{\Delta t} \mathbf{B} dt - \overline{\mathbf{B}}^{p} \Delta t\right)$ , respectively on the l.h.s. and r.h.s. in the same equation, are

small.

A suitable choice of terms  $\mathbf{F}^{p}(\mathbf{U})$  and  $\mathbf{B}^{p}(\mathbf{U})$  in the prediction step allows a much easier solution of the two steps in Eqs. (A.3) with respect to the original formulation in system (A.1).

In the FDSWEsM, vectors  $\mathbf{U}$ ,  $\mathbf{F}(\mathbf{U})$  and  $\mathbf{B}(\mathbf{U})$  have the following expressions:

$$\mathbf{U} = \begin{pmatrix} h & uh & vh \end{pmatrix}^{T} \qquad \mathbf{F} = \begin{pmatrix} \mathbf{F}_{1} & \mathbf{F}_{2} \end{pmatrix}$$
$$\mathbf{F}_{1} = \begin{pmatrix} uh \\ u^{2}h + \frac{1}{2}gh^{2} \\ uvh \end{pmatrix} \qquad \mathbf{F}_{2} = \begin{pmatrix} vh \\ v^{2}h + \frac{1}{2}gh^{2} \\ uvh \end{pmatrix}$$
A.4,a
$$\mathbf{B} = \begin{pmatrix} p & -gh \left(\frac{\partial z}{\partial x} + \frac{n^{2}u\sqrt{(uh)^{2} + (vh)^{2}}}{h^{7/3}}\right) - gh \left(\frac{\partial z}{\partial y} + \frac{n^{2}v\sqrt{(uh)^{2} + (vh)^{2}}}{h^{7/3}}\right) \end{pmatrix}^{T}$$
$$\mathbf{F}_{1}^{p} = \begin{pmatrix} uh \\ u^{2}h \\ uvh \end{pmatrix} \qquad \mathbf{F}_{2}^{p} = \begin{pmatrix} vh \\ v^{2}h \\ uvh \end{pmatrix}$$

$$\mathbf{F}_{1}^{p} = \begin{bmatrix} u^{2}h \\ uvh \end{bmatrix} \qquad \mathbf{F}_{2}^{p} = \begin{bmatrix} v^{2}h \\ uvh \end{bmatrix}$$
$$\mathbf{B}^{p} = \begin{bmatrix} p & -gh \left( \frac{\partial H^{k}}{\partial x} + \frac{n^{2}u\sqrt{(uh)^{2} + (vh)^{2}}}{h^{7/3}} \right) -gh \left( \frac{\partial H^{k}}{\partial y} + \frac{n^{2}v\sqrt{(uh)^{2} + (vh)^{2}}}{h^{7/3}} \right) \end{bmatrix}^{T} \qquad \text{A.4,b}$$

where  $(.)^T$  is the transposed of vector (.).

In the 0ISWEsM, the same vectors assume the following forms:

$$\mathbf{U} = H \qquad \mathbf{F} = -\frac{h^{5/3}}{n\sqrt{|\nabla H|}}\nabla H \qquad \mathbf{B} = p \qquad A.5,a$$

$$\mathbf{F}^{p} = -\frac{h^{5/3}}{n\sqrt{|\nabla H|^{k}}} (\nabla H)^{k}, \quad \mathbf{B}^{p} = \mathbf{B}$$
A.5,b

In both the FDSWEsM and 0ISWEsM, the gradients of the water level in the prediction step are computed at time level  $t^k$  and are kept constant during the time step (in the source term vector  $\mathbf{B}^p$  and in the flux term  $\mathbf{F}^p$ , respectively for the FDSWEsM and 0ISWEsM. More details in [2-5].

#### A.2. Spatial integration of the governing equations

In both the numerical solvers, the governing equations are integrated over unstructured triangular meshes satisfying the Generalized Delaunay (GD) property [4]. Let  $T_h$  be an unstructured GD triangulation of the 2D bounded domain  $\Omega$ . We call *basic mesh* the triangulation  $T_h$  with  $N_T$  triangles and N nodes and its generic triangle and node are denoted as  $k_T$  and  $P_i$  (i = 1, ..., N), respectively. We construct a *dual mesh*  $E_h$  over the basic mesh and its dual element (or dual finite volume), associated with the node  $P_i$ , is denoted as  $e_i$  (i = 1, ..., N). This is the closed region obtained by merging the sub-triangles given by subdividing each triangle  $k_T$  sharing node  $P_i$ , by means of its axes (see figure 1).  $e_i$  is the Voronoi region (or Voronoi polygon) ([4] and cited references).



Figure S1. The basic and the dual mesh.

### A.3. The numerical solution of the OISWEsM

In the 0ISWEsM [4], the computational cell is the Voronoi polygon  $e_i$  and the storage capacity is concentrated in the node  $P_i$  in the measure of 1/3 of the area of all the triangles sharing  $P_i$ . The authors in [4] assume a linear variation of the water level H inside each triangle, according to the three nodal values.

In the 0ISWEs physical problem, the flow field has an exact scalar potential, that is the water level. For the application of the MAST procedure, at the beginning of each time step, the cells are ordered on the base of their potential values, computed at the end of the previous time step or, for the first time iteration, given by the initial condition.

According to Eqs. (4), (A.2)-(A.3) and (A.5), after integration in space and application of the Green theorem, the integral form of the CP system for the generic computational cell *i* is [4]

$$A_{i}\frac{dH_{i}}{dt} + \sum_{j}Fl_{i,j}^{out} = \sum_{m}Fl_{i,m}^{in} + A_{i}p_{i}, \text{ with, } Fl_{i,j}^{out} = K_{i,j}^{k}h_{i}^{5/3}, A_{i} = \frac{1}{3}\sum_{n=1,N_{T}}|k_{T_{n}}|\delta_{i,n}, A_{i} = \frac{1}{3}\sum_{n=1,N_{T}}|k_{T_{n}}|\delta_{i,$$

where  $A_i$  is the area of cell i,  $|k_{T_n}|$  is the area of triangle n,  $_{i,n}$  is equal to 1 if triangle n shares node i, 0 otherwise,  $Fl_{i,j}^{out}$  is the flux leaving cell i to the any neighbouring downstream (in the potential scale) cell j (with  $H_j^k \leq H_i^k$ ),  $K_{i,j}^k$  is the flux coefficient, further defined,  $Fl_{i,m}^{in}$  is the flux entering cell i from any neighbouring upstream (in the potential scale) cell m with  $H_i^k \leq H_m^k$  and  $p_i$  is source term in node i. The flux coefficient is defined as (see Eqs.(19)-(22) in [4])

$$K_{i,j}^{k} = \left(c_{i,j}^{1} E_{1}^{k} + c_{i,j}^{2} E_{2}^{k}\right) \frac{\left(H_{i}^{k} - H_{j}^{k}\right)}{d_{ij}}$$
A.7

where  $d_{ij}$  is the distance between nodes *i* and *j*, indices 1 and 2 mark the two triangles sharing side *ij*, the coefficient  $E_{1(2)}^k$  is

$$E_m^k = \frac{1}{n_m \sqrt{|\nabla H_m^k|}}, m = 1, 2$$
 A.8

where the sub-index *m* marks all the parameters of the triangle *m* sharing the same side and  $c_{i,j}^{l(2)}$  is the distance between the midpoint of side *ij* and the circumcentre of triangle *m*. If side *ij* shares only one triangle,  $c_{i,j}^2$  and  $E_2^k$  are zero. The distance  $c_{i,j}^{l(2)}$  is computed as in Eqs. (19) in [4], where the

authors prove that the proposed formulation guarantees the consistency of the flux between cells *i* and *j* and the difference of the corresponding potentials  $(H_i^k - H_i^k)$  for an unstructured GD mesh.

If the r.h.s. of system (A.6) is approximated with its mean value in the time step, the solution of the same system is disentangled in the sequential solution of N Ordinary Differential Equations (ODEs) [4],

$$A_{i}\frac{dH_{i}}{dt} + \sum_{j}Fl_{i,j}^{out} = \sum_{m}\overline{Fl}_{i,m}^{in} + A_{i}\overline{p_{i}}$$
A.9

one for each cell, going from the cell with highest to the cell with the lowest potential value. In Eq. (A.9),  $\overline{Fl}_{i,m}^{in}$  is the mean in time value of the flux entering from the upstream (in the potential scale) cell *m*, previously solved, and  $\overline{p_i}$  is the mean (in time) value of  $p_i$ . A very fast semi-analytical solution of the ODEs (A.9) is proposed in [4], which allows to save a lot of computational time.

Call  $h_i^k$  the water depth at the beginning of the time step and  $h_i^{kf}$  its asymptotic steady-state value (i. e. when  $dH_i/dt$  =0), computed according to Eqs. (A.9),

$$h_i^{kf} = \left(\frac{\overline{Fl}_i^{in}}{\sum_j K_{i,j}^k}\right)^{3/5}, \text{ with } \overline{Fl}_i^{in} = \overline{Fl}_i^{in} + A_i \overline{p_i}$$
 A.10

Eq. (A.9) can be written in dimensionless form as:

$$\frac{d\xi}{d\tau} = 1 - \xi^{5/3}, \quad \xi = \frac{h_i}{h_i^{kf}}, \quad , \quad \tau = \frac{dt \; Fl_i^{vin}}{A_i h_i^{kf}}, \text{ if } \qquad h_i^{kf} > h_i^k \qquad A.11,a$$

$$\frac{d\xi}{d\tau} = \xi_{f}^{5/3} - \xi^{5/3}, \quad \xi = \frac{h_{i}}{h_{i}^{k}}, \quad \tau = \frac{dt \ \overline{Fl}_{i}^{vin}}{A_{i}h_{i}^{k}} \left(\frac{h_{i}^{k}}{h_{i}^{kf}}\right)^{5/3}, \text{ if } \qquad h_{i}^{kf} < h_{i}^{k} \qquad A.11, b$$

The proposed semi-analytical solution proposed in [4] is

$$\xi = \frac{\exp(c_1 \tau) + c_2}{\exp(c_1 \tau) + c_3} \qquad \text{if} \quad h_i^{kf} > h_i^k \qquad \text{A.12,a}$$

$$\xi = 1 + (\xi_f - 1) \frac{\exp(c_1 \tau) + c_2}{\exp(c_1 \tau) + c_3} \quad \text{if} \quad h_i^{kf} < h_i^k \quad A.12, b$$

with a proper choice of the  $c_1$ ,  $c_2$  and  $c_3$  coefficients. Using any  $c_3$  value it is possible to match the initial value  $\xi_0$  and its first derivative  $\xi'_0$  by setting:

$$c_2 = \xi_0 (1 + c_3) - 1,$$
  $c_1 = \xi_0 \frac{(1 + c_3)^2}{(c_3 - c_2)}$  if  $h_i^{kf} > h_i^k$  A.13,a

$$c_2 = -1$$
,  $c_1 = \xi_0' \frac{(1+c_3)}{(\xi_f - 1)}$  if  $h_i^{kf} < h_i^k$  A.13,b

The  $c_3$  coefficient affects the maximum error that is obtained according to functions (A.12) using different time step sizes. This optimum depends on  $\xi_0$ , if  $h_i^{kf} > h_i^k$  and on  $\xi_f$ , if  $h_i^{kf} < h_i^k$ . The optimum coefficients have been computed numerically for different possible  $\xi_0$  and  $\xi_f$  values by comparing functions (A.12) with a numerical solution computed using a very small time step. See in table 1 and in figure 7 in [4] the computed optimum  $c_3$  values. See in figures 8,a and 8,b in [4] the

numerical solution of Eqs. (A.11) in the case of respectively  $\xi_0 = 0$  and  $\xi_f = 0$ , compared with the semianalytical solutions (A.12) corresponding to the optimal  $c_3$  values (respectively 0.7469 and -0.8171). The maximum error computed with the initial conditions  $\xi_0 = 0$ , for  $h_i^{kf} > h_i^k$ , or  $\xi_f = 0$ , for  $h_i^{kf} < h_i^k$ , is the worse one and it is smaller than 10<sup>-3</sup>. More details can be found in [4]

After the solution of each ODEs (A.9), the mean in time total flux going from cell i to the neighbouring downstream (in the potential scale) cells is computed applying the local mass balance for cell i [4]. In the framework of the MAST procedure, due to the sequential solution of the cells and their ordering at the beginning of the time iteration, the mean (in time) entering flux is always known before each solution of the ODEs (A.9) [4].

The same spatial discretization adopted for the CP problem is used in the DC problem and the initial condition of the DC problem is the final state obtained after the solution of the CP step, marked with the index k+1/2. Starting from Eqs. (4), (A.2)-(A.3) and (A.5), after spatial integration of the correction problem inside each Voronoi cell, the following DC system is obtained [4]

$$\frac{A_i}{\Delta t}\eta_i + \sum_{n=1,N_T} D_{i,j}^k \left(\eta_i - \eta_j\right) \delta_{i,n} = \sum_{n=1,N_T} D_{i,j}^k \left(\vartheta_j - \vartheta_i\right) \delta_{i,n} \qquad i = 1, \dots, N \qquad A.14,a$$

where

$$\eta = H - H^{k+1/2}, \quad \vartheta = H^k - H^{k+1/2}, \quad D_{i,j}^k = \sum_{m=1,2} \frac{c_{i,j}^m E_m^k}{d_{ij}} (h_l^{km})^{5/3},$$

$$h^{km} = \frac{h^k + h^{k+1/2}}{2}$$
A.14,b

with initial condition = 0. In Eq. (A.14,a), *in* has been previously specified and the sum in Eq. (A.14,b) is extended to the two triangles *m* sharing side *ij*. l = i if  $H_i^k \ge H_j^k$ , l = j if  $H_i^k < H_j^k$ . The proposed formulation of the coefficients  $D_{i,j}^k$  provides the same flux estimation of the CP step [4] and it is similar to the one of a standard linear (*P*1) Finite Element Galerkin scheme. Analogies and differences between the flux formulation adopted in the present DSWEs model and the one of a *P*1 Galerkin scheme are presented and discussed for a GD triangulation in [4].

The matrix of the linear system resulting from Eqs. (A.14) has order *N* (the number of the nodes) and, in the case of a GD triangulation, it is symmetric, positive-definite, strictly diagonally dominant, with *M*-property and system (A.14) is well-conditioned [4]. After the unknowns are computed by solving the system (A.14), the final values of the water levels are updated as

$$H^{k+1} = H^{k+1/2} + \eta^{2}$$
 A.15

The splitting of the original governing equations in the CP and DC steps, allows this 0ISWEsM solver to easily deal with waves/flooding propagation problems over dry domains, as well as a robust solution of wetting/drying problems [4].

A.3.1. Investigation of the behaviour of the proposed 0ISWEsM over refined meshes and smooth surfaces

If the area  $A_i$  of the *i*<sup>th</sup> Voronoi cell goes approaches zero, the dimensionless variable in Eqs. (A.11) approaches infinity, and the solution of the prediction step is given to the asymptotic values in Eqs. (A.12,a) or (A.12,b), respectively if  $h_i^{kf} > h_i^k$  or  $h_i^{kf} < h_i^k$ , shown in figures 8 in [4].

In the correction step, the value of the capacity term  $A_i$  i/t in system (A.14) becomes negligible compared with the flux terms  $\sum_{n=1,N_T} D_{i,j}^k (\eta_i - \eta_j) \delta_{i,n}$  and  $\sum_{n=1,N_T} D_{i,j}^k (\vartheta_j - \vartheta_i) \delta_{i,n}$ 

If the Manning coefficient *n* approaches zero, coefficient  $E_m^k$  in Eq. (A.8) approaches infinity, as well as the flux coefficient in the prediction step  $K_{i,j}^k$  in Eq. (A.7) and  $\sum_j K_{i,j}^k$  in Eq. (A.10).

This implies that the asymptotic steady-state value  $h_i^{kf}$  in Eq. (A.10) approaches zero and the dimensionless variable in Eqs. (A.11) approaches infinity. In this case, the dimensionless solution  $\xi$  of the semi-analytical procedure of the prediction step is given by the asymptotic value of Eq. (A.12,b).

Coefficient  $D_{i,j}^k$  in the correction system in Eq. (A.14,a) approaches infinity (see the third of Eq. (A.14,b)). By dividing the terms in Eq. (A.14,a) by coefficient  $D_{i,j}^k$ , the capacity term becomes negligible with respect to the other terms in the same equation.

A.3.2. Preservation of the water at rest condition (C-property)

The proposed 0ISWEsM preserves the *C*-property (e.g., [7]). For quiescent water, we have, in the prediction step, zero flux entering in each cell and zero gradient of the water level *H*. This implies that in Eq. (A.6) we have  $\sum_{m} Fl_{i,m}^{in} = 0$  and, from Eqs. (A.7)-(A.8)  $\sum_{j} Fl_{i,j}^{out} = 0$ . The solution of the prediction step gives  $H^{k+1/2} = H^k$ . System (A.14) in the correction step becomes

$$\frac{A_i}{\Delta t}\eta_i + \sum_{n=1,N_T} D_{i,j}^k \left(\eta_i - \eta_j\right) \delta_{i,n} = 0$$
A.16

since  $\vartheta = H^k - H^{k+1/2} = 0$ , and this implies that the correction of the water level is zero.

## A.4. The numerical solution of the FDSWEsM

In the FDSWEsM, the computational cell is the triangle  $k_T$ , with the storage capacity concentrated in the circumcentre of  $k_T$  in the measure of the area of  $k_T$  and a piecewise constant value of the unknown variables h (as well as H), uh ad vh is assumed inside each triangle [3, 5].

According to Eqs. (1)-(2), (A.2)-(A.4), by integrating in space the prediction problem and applying the Green's theorem, the integral form of the CP system is [3, 5]

$$\frac{\partial h}{\partial t} |k_T|_e + \sum_{j=1,3} \delta_{j,e} F_{j,e}^{out} = \sum_{j=1,3} (1 - \delta_{j,e}) F_{j,e}^{in} + |k_T|_e p_e \qquad e = 1, \dots, N_T$$
A.17

$$\frac{\partial q_{s,e}}{\partial t} |k_T|_e + \sum_{j=1,3} \delta_{j,e} M_{j,e}^{s,out} + R_e^s = \sum_{j=1,3} (1 - \delta_{j,e}) M_{j,e}^{s,in}, \quad \text{with} \quad s = x, y,$$

$$q_s = \begin{cases} uh & \text{if } s = x \\ vh & \text{if } s = y \end{cases}$$
A.18

where index *e* marks the values of the variables in triangle  $k_{T,e}$ ,  $F_{j,e}^{out(in)}$  and  $M_{j,e}^{x(y)out(in)}$  are the flux and the *x*(*y*) component of the momentum flux, leaving (entering) cell  $k_{T,e}$  across side *j* (*j* = 1, 2, 3) of  $k_{T,e}$ , respectively,  $_{j,e} = 1$  (0) if the flux across side *j* of  $k_{T,e}$  is oriented outward (inward)  $k_{T,e}$ ,  $p_e$  is the source term in cell  $k_{T,e}$  and  $R_e^{x(y)}$  is the source term defined as [3],

$$R_{e}^{s} = bfr_{e}^{s} + gr_{e}^{s} \qquad \text{with} \quad bfr_{e}^{s} = \left|k_{T}\right|_{e} g \frac{n^{2}q_{s,e}\sqrt{\left(uh\right)_{e}^{2} + \left(vh\right)_{e}^{2}}}{h_{e}^{7/3}} \quad \text{and} \\ gr_{e}^{s} = \left|k_{T}\right|_{e} gh_{e} \frac{\partial H_{e}^{k}}{\partial s} \qquad A.19$$

The first term on the l.h.s. of Eq. (A.18) is the integral form of the local inertia (or local acceleration), the two summations of the same equation, on the l.h.s. and r.h.s., respectively, represent the integral form of the convective inertia (or convective acceleration), computed as the line integral of the momentum fluxes across the sides of the computational cell, the third term on the l.h.s of Eq.

(A.18), is the sum of the integral form of the bottom friction  $bfr_e^s$  and water level gradient  $gr_e^s$  terms.

As for the 0ISWEs problem, the authors in [3,5] disentangle the solution of the prediction system in Eqs. (A.17)–(A.18) in the sequential solution of  $N_T$  ODEs, by approximating the r.h.s. of the same equations with their mean values in time computed during the time step.

To apply the sequential solution of the computational cells, it is necessary to order them [3,5], but, unlike the previous DSWEs problem, an exact scalar potential of the flow field does not exist in the FDSWEs physical problem. For this reason, the authors in [2,3,6] add a convective corrective (CC) step, splitting the original prediction step in a "convective prediction" (CP) system and in a "convective correction" (CC) system [3,6].

The computational cells are ordered by applying an iterative procedure based on the inter-cell flux direction between adjacent cells, proposed in [5] and further modified in [6]. In the following, we call  $no_e^k$  the order number of cell  $k_{T,e}$  at the beginning of the generic time step (time level  $t^k$ ).

After the ordering of the computational cells, the CP step is solved as a sequence of small systems of ODEs, from time level  $t^k$  to time level  $t^k + \Delta t$  [5], proceeding from the cell(s) with the lowest  $n o_e^k$  value to the cell(s) with the highest  $n o_e^k$  value. The same procedure is repeated in the CC step, proceeding in the opposite direction, from the cell(s) with the highest to the cell(s) with the lowest  $n o_e^k$  value [6]. Cell  $k_{T,e}$  with order  $n o_e^k$  is solved only after the solution of the neighbouring  $k_{T,ep}$  cells with  $n o_{ep}^k < n o_e^k$  ( $n o_{ep}^k > n o_e^k$ ), in the CP (CC) step [5]. The ODEs system of the CP problem is [3, 5]

$$\frac{dh_{e}}{dt}|k_{T,e}| + \frac{1}{\Delta t} \int_{\Delta t} \sum_{j=1,3} \delta_{j,e} F_{j,e}^{p,out}(t) dt = \sum_{j=1,3} (1 - \delta_{j,e}) \overline{F}_{j,e}^{p,in} + |k_{T,e}| \overline{p}_{e}, e = 1, ..., NT$$
 A.20

$$\frac{dq_{s,e}}{dt} \left| k_{T,e} \right| + \frac{1}{\Delta t} \int_{\Delta t} R_e^s(t) dt + \frac{1}{\Delta t} \int_{\Delta t} \sum_{j=1,3} \delta_{j,e} M_{j,e}^{p,s,out}(t) dt = \sum_{j=1,3} \left( 1 - \delta_{j,e} \right) \overline{M}_{j,e}^{p,s,in} \qquad s$$

$$= x, y$$
A.21

with the symbols specified as above and the leaving flux and momentum flux  $F_{j,e}^{p,out}$  and  $M_{j,e}^{p,x(y),out}$  defined as follows. If the two neighbouring triangles  $k_{T,e}$  and  $k_{T,ep}$  share side  $\mathbf{r}_{i,ip}$  between nodes *i* and *ip*, where  $\mathbf{r}_{i,ip}$  is the *j*<sup>th</sup> side of  $k_{T,e}$  and the *m*<sup>th</sup> side of  $k_{T,ep}$  (*j*, *m* = 1, 2, 3), we define the flux across side *j* of  $k_{T,e}$  as [3]

$$FL_{j,e} = (uh)_{e} (y_{ip} - y_{i}) - (vh)_{e} (x_{ip} - x_{i})$$
A.22

and the flux and the momentum fluxes between  $k_{T,e}$  and  $k_{T,ep}$  as [3]

$$F_{j,e} = FL_{j,e}$$
 if  $FL_{j,e} > 0$  and  $FL_{j,e} > FL_{m,ep}$  A.23,a

$$F_{j,e} = -FL_{m,ep}$$
 otherwise A.23,b

$$M_{j,e}^{x} = F_{j,e} \frac{(uh)_{e}}{h_{e}}, \qquad M_{j,e}^{y} = F_{j,e} \frac{(vh)_{e}}{h_{e}} \qquad \text{if } F_{j,e} = FL_{j,e} \qquad A.24,a$$

$$M_{j,e}^{x} = F_{j,e} \frac{(uh)_{ep}}{h_{ep}}, \qquad M_{j,e}^{y} = F_{j,e} \frac{(vh)_{ep}}{h_{ep}} \qquad \text{otherwise} \qquad A.24,b$$

Flux/momentum flux continuity is guaranteed at each triangle side by Equations (A.23)–(A.24) and  $F_{j,e} = -F_{m,ep}$  and  $M_{j,e}^{x(y)} = -M_{m,ep}^{x(y)}$ .

The leaving flux/momentum flux  $F_{j,e}^{p,out}$  and  $M_{j,e}^{p,x(y),out}$  in Eqs. (A.20)-(A.21) of the CP step, going from cell  $k_{T,e}$  to the neighbouring cell  $k_{T,ep}$  (with  $nO_e^k < nO_{ep}^k$ ), are defined as [3]

$$F_{j,e}^{p,out} = \max\left(0, F_{j,e}\right), \quad M_{j,e}^{p,x,out} = F_{j,e}^{p,out} \frac{(uh)_e}{h_e}, \quad M_{j,e}^{p,y,out} = F_{j,e}^{p,out} \frac{(vh)_e}{h_e} \quad \text{if} \\ no_e^k < no_{ep}^k \quad (A.25,a), \quad F_{j,e}^{p,out} = M_{j,e}^{p,x(y),out} = 0 \quad \text{if} \quad no_e^k > no_{ep}^k \quad (A.25,b)$$

 $\overline{F}_{j,e}^{p,in}$  and  $\overline{M}_{j,e}^{p,s,in}$  on the r.h.s. of Eqs. (A.20)-(A.21) are the mean in time values of the incoming fluxes and momentum fluxes, respectively (know from the solution of the previously solved neighbouring  $k_{T,ep}$  cells with  $no_e^k > no_{ep}^k$  [2, 3, 5, 6]) and  $\overline{p}_e$  is the mean in time value of the source term in  $k_{T,e}$ .

Once the ODEs system (A.20)-(A.21) is solved for  $k_{T,e}$ , the mean in time value of the total flux  $\overline{F}_{e}^{out}$  leaving from  $k_{T,e}$ , from  $t^{k}$  to  $t^{k} + \Delta t$ , is computed according to the local mass balance for cell  $k_{T,e}$  [3]. The mean in time flux  $\overline{F}_{j,e}^{out}$  and momentum fluxes  $\overline{M}_{j,e}^{x(y),out}$ , leaving from the  $j^{\text{th}}$  side of  $k_{T,e}$  to the  $m^{\text{th}}$  side of the neighbouring cell  $k_{T,ep}$  (with  $nO_{e}^{k} < nO_{ep}^{k}$ ) are estimated as in [2, 3, 5]. After that, the same ODEs system (A.20)-(A.21) is solved for cell  $k_{T,ep}$ , selected among the unsolved ones according to

$$no_{ep}^{k} > no_{e}^{k}$$
 and  $no_{ep}^{k} = \min(no_{el}^{k}), k_{T,el} \subset \text{ set of unsolved cells}$  A.26

The same procedure is applied for the solution of the CC step, whose ODEs system is

$$\frac{dh_e}{dt}|T_e| + \frac{1}{\Delta t} \int_{\Delta t} \sum_{j=1,3} \delta_{j,i} F_{j,e}^{c,out}(t) dt = \sum_{j=1,3} (1 - \delta_{j,e}) \overline{F}_{j,e}^{c,in}$$
A.27

$$\frac{dq_{s,e}}{dt}|T_i| + \frac{1}{\Delta t} \int_{\Delta t} \sum_{j=1,3} \delta_{j,e} M_{j,e}^{c,s,out}(t) dt = \sum_{j=1,3} (1 - \delta_{j,e}) \overline{M}_{j,e}^{c,s,in}, s = x, y$$
 A.28

with  $\delta_{j,e}$  defined as above and the leaving flux/momentum flux  $F_{j,e}^{c,out}$  and  $M_{j,e}^{c,x(y),out}$  in Eq. (A.27)-(A.28) of the CC step, going from cell  $k_{T,e}$  to the neighbouring cell  $k_{T,ep}$  (with  $no_e^k > no_{ep}^k$ ), are defined as [3]

$$F_{j,e}^{c,out} = \max\left(0, F_{j,e}\right), \quad M_{j,e}^{c,x,out} = F_{j,e}^{c,out} \frac{(uh)_e}{h_e}, \qquad M_{j,e}^{c,y,out} = F_{j,e}^{c,out} \frac{(vh)_e}{h_e} \quad \text{if} \\ no_e^k > no_{ep}^k \quad (A.29,a), \qquad F_{j,e}^{c,out} = M_{j,e}^{c,x(y),out} = 0 \quad \text{if} \quad A.29,b$$

As motivated in [2, 3], the source terms are allocated in the CP step.

After the solution of the ODEs system (A.27)–(A.28) for cell  $k_{T,e}$ , the total leaving fluxes/momentum fluxes from side j of  $k_{T,e}$  to the neighbouring cell  $k_{T,ep}$  with  $no_e^k > no_{ep}^k$  are computed as for the previous CP step. More details in [2,3,6].

According to Equations (1)-(2), (A.2)-(A.4), after integration in space, neglecting the inertial terms in the momentum equations (as motivated in [2,3]), substituting the momentum equations in the continuity equation and applying the Green's theorem, we obtain a linearized system in the unknown water level [2,3,5],

$$\int_{k_{T,e}} \frac{\partial H}{\partial t} dk_{T,e} + \sum_{j=1,3} \int_{L_{j,e}} -elem_e \frac{\partial H^{k+1}}{\partial n_{j,e}} dl = \sum_{j=1,3} \int_{L_{j,e}} -elem_e \frac{\partial H^k}{\partial n_{j,e}} dl + \sum_{j=1,3} \int_{L_{j,e}} \left(\overline{\mathbf{q}} - \mathbf{q}^{cc}\right) \cdot \hat{\mathbf{n}}_{j,e} dl \qquad A.30,a$$

with 
$$elem_e = g\overline{h_e}\Delta t / \left( 1 + \Delta t g n^2 \overline{\left(\frac{\sqrt{(uh)_e^2 + (vh)_e^2}}{h_e^{7/3}}\right)} \right)$$
 and  $e = 1, ..., A.30, b$   
 $N_T$ 

where  $L_{j,e}$  is the length of side *j* of triangle  $k_{T,e}$ ,  $\hat{\mathbf{n}}_{j,e}$  is its unit orthogonal vector,  $\overline{\mathbf{q}}$  and  $\mathbf{q}^{cc}$  are the mean in time value of the flow rate vector computed during the (CP + CC) steps and its final value, respectively [Error! Reference source not found.] and symbols  $(\bar{*})$  in Eq. (A.30) is the mean in time value of (\*) computed during the (CP + CC) steps. The initial state for the DC step is the solution computed at the end of the CC step. The matrix of the linear system resulting from Eqs. (A.30) has order  $N_T$  (the number of the triangles) and, in the case of a GD triangulation, is well conditioned, symmetric, positive-definite, strictly diagonally dominant, with M-property and system (A.30) is well-conditioned [5]. After solving the DC step, uh and vh are updated, as well as the spatial gradients of the water levels, as explained in [5].

## Part B. Supplementary material for the presented tests

In this Appendix we plot the figures and tables recalled in the text of the main paper.

OISWEsM							FDSWEsM								
r <sub>c</sub>	L <sub>1,h</sub>	r <sub>c</sub>	L <sub>1,uh</sub>	rc	L <sub>2,h</sub>	r <sub>c</sub>	L <sub>2,uh</sub>	rc	L <sub>1,h</sub>	r <sub>c</sub>	L <sub>1,uh</sub>	r <sub>c</sub>	L <sub>2,h</sub>	r <sub>c</sub>	L <sub>2,uh</sub>
case 1							case 1								
0.323	0.0311	0.350	0.0031	0.384	0.0526	0.348	0.0053	0.949	0.0036	0.987	0.0023	0.996	0.0059	0.907	0.0049
0.520	0.0248	0.464	0.0024	0.481	0.0403	0.496	0.0042	1.055	0.0019	1.059	0.0011	1.030	0.0030	0.927	0.0026
0.615	0.0173	0.660	0.0017	0.597	0.0289	0.577	0.0030	1.065	0.0009	1.026	0.0006	1.012	0.0015	1.077	0.0014
	0.0113		0.0011		0.019		0.0020		0.0004		0.0003		0.0007		0.0006
case 2							case 2								
0.712	0.0070	0.502	0.0028	0.799	0.0126	0.476	0.0041	0.971	0.0036	0.994	0.0023	0.948	0.0060	0.963	0.0049
0.817	0.0043	0.634	0.0020	0.844	0.0073	0.694	0.0029	1.047	0.0019	1.089	0.0011	1.050	0.0031	1.048	0.0025
0.991	0.0024	0.876	0.0013	1.083	0.0040	0.807	0.0018	1.103	0.0009	1.030	0.0005	1.001	0.0015	1.007	0.0012
	0.0012		0.0007		0.0019		0.0010		0.0004		0.0003		0.0007		0.0006
case 3							case 3								
0.967	0.0039	0.520	0.0028	0.890	0.0065	0.554	0.0039	0.940	0.0043	0.940	0.0009	0.932	0.0076	0.893	0.0016
1.012	0.0020	0.657	0.0019	0.959	0.0035	0.777	0.0026	1.059	0.0022	1.056	0.0005	1.000	0.0040	0.981	0.0008
1.032	0.0010	0.836	0.0012	1.016	0.0018	0.895	0.0015	1.051	0.0011	1.070	0.0002	1.029	0.0020	1.104	0.0004
	0.0005		0.0007		0.0009		0.0008		0.0005		0.0001		0.0010		0.0002
case 4							case 4								
0.951	0.0039	0.582	0.0026	0.905	0.0065	0.584	0.0037	0.913	0.0043	1.000	0.0009	0.858	0.0076	0.976	0.0016
0.998	0.0020	0.772	0.0017	0.956	0.0035	0.823	0.0025	0.987	0.0023	1.106	0.0005	1.066	0.0042	1.074	0.0008
1.005	0.0010	1.008	0.0010	0.983	0.0018	0.926	0.0014	1.082	0.0011	1.064	0.0002	1.004	0.0020	1.078	0.0004
	0.0005		0.0005		0.0009		0.0007		0.0005		0.0001		0.0010		0.0002

Test 1. Steady flow in a 1D channel with undulating bottom profile

**Figure S1.** Test 1. *L*<sup>1</sup> and *L*<sup>2</sup> norms of the relative errors for *h* and *uh*. Unstructured meshes.

OISWEsM							FDSWEsM								
rc	L <sub>1,h</sub>	r <sub>c</sub>	L <sub>1,uh</sub>	r <sub>c</sub>	L <sub>2,h</sub>	r <sub>c</sub>	L <sub>2,uh</sub>	r <sub>c</sub>	L <sub>1,h</sub>	r <sub>c</sub>	L <sub>1,uh</sub>	r <sub>c</sub>	L <sub>2,h</sub>	r <sub>c</sub>	L <sub>2,uh</sub>
case 1							case 1								
0.324	0.0311	0.355	0.0031	0.291	0.0526	0.308	0.0052	0.961	0.0036	1.019	0.0023	1.016	0.0059	0.924	0.0049
0.523	0.0248	0.493	0.0024	0.575	0.0430	0.485	0.0042	1.063	0.0019	1.079	0.0011	1.040	0.0029	0.929	0.0026
0.586	0.0173	0.601	0.0017	0.613	0.0289	0.659	0.0030	1.066	0.0009	1.084	0.0005	0.974	0.0014	1.170	0.0014
	0.0115		0.0011		0.0189		0.0019		0.0004		0.0003		0.0007		0.0006
case 2							case 2								
0.726	0.0070	0.493	0.00273	0.795	0.0126	0.463	0.0040	1.078	0.0038	1.006	0.0023	0.971	0.0060	0.911	0.0047
0.790	0.0042	0.657	0.00194	0.846	0.0073	0.693	0.0029	1.016	0.0018	1.107	0.0011	1.033	0.0031	1.059	0.0025
1.012	0.0024	0.855	0.00123	1.625	0.0040	0.845	0.0018	1.101	0.0009	1.115	0.0005	1.010	0.0015	1.005	0.0012
	0.0012		0.00068		0.0013		0.0010		0.0004		0.0002		0.0007		0.0006
case 3							case 3								
0.951	0.0039	0.506	0.0027	0.899	0.0065	0.531	0.0038	0.957	0.0043	0.971	0.0009	0.928	0.0076	0.978	0.0016
1.015	0.0020	0.799	0.0019	0.967	0.0035	0.776	0.0026	1.144	0.0022	1.031	0.0005	1.000	0.0040	0.930	3000.0
1.038	0.0010	0.752	0.0011	1.033	0.0018	0.941	0.0015	1.000	0.0010	1.118	0.0002	1.029	0.0020	1.066	0.0004
	0.0005		0.0006		0.0009		0.0008		0.0005		0.0001		0.0010		0.0002
case 4							case 4								
0.930	0.0038	0.548	0.0025	0.902	0.0064	0.582	0.0037	0.900	0.0042	1.064	0.0009	0.866	0.0076	0.943	0.0015
1.000	0.0020	0.768	0.0017	0.950	0.0034	0.835	0.0025	1.003	0.0023	1.070	0.0004	1.057	0.0042	1.007	3000.0
1.000	0.0010	0.977	0.0010	0.986	0.0018	0.929	0.0014	1.111	0.0011	1.014	0.0002	1.003	0.0020	1.061	0.0004
	0.0005		0.0005		0.0009		0.0007		0.0005		0.0001		0.0010		0.0002

**Figure S2.** Test 1. *L*<sup>1</sup> and *L*<sup>2</sup> norms of the relative errors for *h* and *uh*. Structured meshes.





Figure S3. Test 2. L1 and L2 norms of qout.

Test 3. Rainfall in a 2D catchment



**Figure S4.** Test 3. *L*<sup>1</sup> and *L*<sup>2</sup> norms of *q*<sub>out</sub>.



**Figure S5.** Test 3. Computed water depth and vectors of unitary flow rate at different durations. 22 s (top), 30 s (middle), 52 s (bottom).

# Test 4. The Toce river case

Table S1. Test 4. Values of the L1, L2 Linf norms of the relative errors at the gauges [-]

	FL	DSWEsM	no	01	SWEsM 1	<b>1</b> 0	0ISWEsM nopt			
gauge	$L_1$	L2	Linf	$L_1$	L2	Linf	$L_1$	$L_2$	Linf	
P1	6.65E-04	8.04E-04	3.65E-02	9.07E-04	1.10E-03	5.96E-02	9.07E-04	1.09E-03	5.83E-02	
P2	6.83E-04	9.29E-04	3.97E-02	2.69E-03	2.56E-03	7.34E-02	1.64E-03	1.44E-03	5.25E-02	
Р3	9.53E-04	8.73E-04	2.09E-02	1.33E-03	1.39E-03	6.25E-02	1.27E-03	1.39E-03	5.95E-02	
S4	1.29E-03	1.69E-03	5.59E-02	1.40E-03	1.90E-03	5.08E-02	1.23E-03	1.74E-03	4.60E-02	
P4	5.51E-04	1.04E-03	7.59E-02	9.74E-04	1.32E-03	6.64E-02	9.74E-04	1.32E-03	6.64E-02	
S6S	4.29E-04	5.66E-04	1.44E-02	3.45E-03	3.95E-03	5.46E-02	2.31E-03	2.99E-03	4.29E-02	
S6D	1.12E-03	9.76E-04	2.15E-02	1.49E-03	2.43E-03	3.46E-02	8.78E-04	1.32E-03	3.16E-02	
P5	1.23E-03	1.43E-03	4.07E-02	2.71E-03	3.28E-03	6.39E-02	1.64E-03	1.80E-03	5.29E-02	
P8	1.53E-03	2.87E-03	4.32E-02	2.46E-03	3.84E-03	7.06E-02	1.54E-03	2.09E-03	6.30E-02	
S8D	6.33E-04	8.87E-04	2.30E-02	2.96E-03	4.95E-03	5.42E-02	7.10E-04	9.80E-04	2.70E-02	
Р9	8.31E-04	1.07E-03	3.96E-02	4.68E-03	7.33E-03	7.90E-02	2.46E-03	3.89E-03	7.63E-02	
P10	8.06E-04	1.08E-03	2.69E-02	3.18E-03	6.54E-03	7.89E-02	1.15E-03	1.55E-03	4.45E-02	
P12	5.94E-04	8.53E-04	1.70E-02	1.00E-03	1.60E-03	3.91E-02	8.86E-04	1.33E-03	3.91E-02	
P13	1.42E-03	2.32E-03	6.75E-02	2.97E-03	5.61E-03	6.92E-02	2.18E-03	3.22E-03	6.75E-02	
P18	2.05E-03	3.63E-03	5.80E-02	2.24E-03	3.61E-03	7.97E-02	1.16E-03	1.87E-03	4.87E-02	
P19	3.78E-04	7.50E-04	3.00E-02	1.90E-03	2.40E-03	5.57E-02	1.45E-03	1.45E-03	3.88E-02	
P21	1.19E-03	1.37E-03	4.17E-02	3.63E-03	7.23E-03	8.14E-02	1.57E-03	2.28E-03	5.39E-02	
P23	2.59E-03	2.37E-03	4.95E-02	5.52E-03	8.32E-03	1.35E-01	4.38E-03	5.76E-03	8.41E-02	
P24	2.67E-03	2.22E-03	3.53E-02	7.69E-03	1.63E-02	1.80E-01	5.53E-03	9.97E-03	1.38E-01	
P25	2.76E-03	2.44E-03	3.62E-02	5.86E-03	1.65E-02	1.86E-01	3.83E-03	8.07E-03	1.07E-01	
P26	1.80E-03	2.50E-03	4.53E-02	7.31E-03	8.32E-03	1.69E-01	6.12E-03	6.26E-03	8.35E-02	



Figure S6. Test 4. Measured and computed water levels at gauges S6S, P4 and S6D.

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