

# Behaviour of the Dam-Break Problem for the Serre Equations

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## ABSTRACT

**Keywords:** dispersive waves, conservation laws, Serre equation, finite volume method, finite difference method

## 1 INTRODUCTION

## 2 SERRE EQUATIONS

The Serre equations can be derived as an approximation to the full Euler equations by depth integration similar to (Su and Gardner 1969). They can also be seen as an asymptotic expansion to the Euler equations as well (Lannes and Bonneton 2009). The former is more consistent with the perspective from which numerical methods will be developed while the latter indicates the appropriate regions in which to use these equations as a model for fluid flow. The set up of the scenario under which the Serre approximation is made consists of a two dimensional  $\mathbf{x} = (x, z)$  fluid over a bottom topography as in Figure 1 acting under gravity. Consider a fluid particle at depth  $\xi(\mathbf{x}, t) = h(x, t) + z_b(x) - z$  below the water surface, see Figure 1. Where the water depth is  $h(x, t)$  and  $z_b(x)$  is the bed elevation. The fluid particle is subject to the pressure,  $p(\mathbf{x}, t)$  and gravitational acceleration,  $\mathbf{g} = (0, g)^T$  and has a velocity  $\mathbf{u} = (u(\mathbf{x}, t), w(\mathbf{x}, t))$ , where  $u(\mathbf{x}, t)$  is the velocity in the  $x$ -coordinate and  $w(\mathbf{x}, t)$  is the velocity in the  $z$ -coordinate and  $t$  is time. Assuming that  $z_b(x)$  is constant the Serre equations read (Li et al. 2014)

$$\frac{\partial h}{\partial t} + \frac{\partial(\bar{u}h)}{\partial x} = 0 \quad (1a)$$

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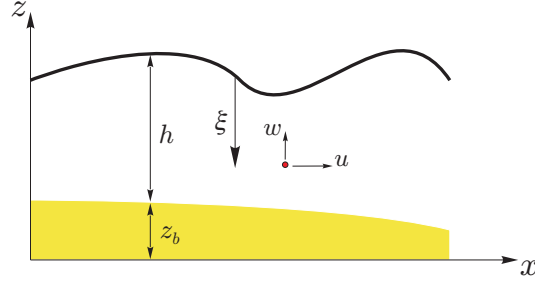


FIG. 1: The notation used for one-dimensional flow governed by the Serre equation.

$$\underbrace{\frac{\partial(\bar{u}h)}{\partial t} + \frac{\partial}{\partial x} \left( \bar{u}^2 h + \frac{gh^2}{2} \right)}_{\text{Shallow Water Wave Equations}} + \underbrace{\frac{\partial}{\partial x} \left( \frac{h^3}{3} \left[ \frac{\partial \bar{u}}{\partial x} \frac{\partial \bar{u}}{\partial x} - \bar{u} \frac{\partial^2 \bar{u}}{\partial x^2} - \frac{\partial^2 \bar{u}}{\partial x \partial t} \right] \right)}_{\text{Dispersion Terms}} = 0. \quad (1b)$$

Serre Equations

Where  $\bar{u}$  is the average of  $u$  over the depth of water.

### Hamiltonian

The Serre equations admits a Hamiltonian which is the same as the energy (Li 2002; Le Métayer et al. 2010). The Hamiltonian is:

$$\mathcal{H} = \frac{1}{2} \int_{-\infty}^{\infty} hu^2 + gh^2 + \frac{h^3}{3} \left( \frac{\partial u}{\partial x} \right)^2 dx \quad (2)$$

We can calculate this numerically by partitioning the total integral into cell-wise integrals. The cell-wise integral can then be calculated by quartic interpolation utilising neighbouring cells and then applying Gaussian quadrature with 3 points over the cell to get a sufficiently high order method to calculate the Hamiltonian for our purposes, in particular our method is at least third order accurate for the  $\partial u / \partial x$  term.

### DIRECT NUMERICAL METHODS

The presence of the mixed spatial temporal derivatives in the momentum equation (1b) makes the Serre equations difficult to solve directly with standard numerical methods. A naive way to avoid this is to approximate (1b) by finite differences and the results of this are presented here. To facilitate this a uniform grid in space will be used with  $\Delta x = x_{i+1} - x_i$  for all  $i$  and quantities evaluated at these grid points will be denoted by subscripts for example  $h_i = h(x_i)$ . The grid in time will be denoted by superscripts for example  $h^n = h(t^n)$ , noting that  $h^n$  is a function in space.

### Finite Difference Approximation to Conservation of Momentum Equation

In [Zoppou thesis/my work] it was demonstrated that an efficient numerical scheme for the Serre equations must be at least second-order accurate thus the derivatives in (1b) will be approximated by second-order finite differences. Firstly (1b) must be expanded, making use of (1a) one obtains

$$h \frac{\partial u}{\partial t} + X - h^2 \frac{\partial^2 u}{\partial x \partial t} - \frac{h^3}{3} \frac{\partial^3 u}{\partial x^2 \partial t} = 0 \quad (3a)$$

where  $X$  contains only spatial derivatives and is

$$X = u h \frac{\partial u}{\partial x} + g h \frac{\partial h}{\partial x} + h^2 \frac{\partial u}{\partial x} \frac{\partial u}{\partial x} + \frac{h^3}{3} \frac{\partial u}{\partial x} \frac{\partial^2 u}{\partial x^2} - h^2 u \frac{\partial^2 u}{\partial x^2} - \frac{h^3}{3} u \frac{\partial^3 u}{\partial x^3} \quad (3b)$$

where the bar over  $u$  has been dropped to simplify notation. Then taking second-order centred finite difference approximation to the time derivatives for (3a) gives

$$h^n \frac{u^{n+1} - u^{n-1}}{2\Delta t} + X^n - (h^n)^2 \frac{\left(\frac{\partial u}{\partial x}\right)^{n+1} - \left(\frac{\partial u}{\partial x}\right)^{n-1}}{2\Delta t} - \frac{(h^n)^3}{3} \frac{\left(\frac{\partial^2 u}{\partial x^2}\right)^{n+1} - \left(\frac{\partial^2 u}{\partial x^2}\right)^{n-1}}{2\Delta t} = 0,$$

$$h^n (u^{n+1} - u^{n-1}) + 2\Delta t X^n - (h^n)^2 \left( \left(\frac{\partial u}{\partial x}\right)^{n+1} - \left(\frac{\partial u}{\partial x}\right)^{n-1} \right) - \frac{(h^n)^3}{3} \left( \left(\frac{\partial^2 u}{\partial x^2}\right)^{n+1} - \left(\frac{\partial^2 u}{\partial x^2}\right)^{n-1} \right) = 0.$$

Introducing

$$Y^n = 2\Delta t X^n - h^n u^{n-1} + (h^n)^2 \left(\frac{\partial u}{\partial x}\right)^{n-1} + \frac{(h^n)^3}{3} \left(\frac{\partial^2 u}{\partial x^2}\right)^{n-1}$$

and rearranging results in

$$h^n u^{n+1} - (h^n)^2 \left(\frac{\partial u}{\partial x}\right)^{n+1} - \frac{(h^n)^3}{3} \left(\frac{\partial^2 u}{\partial x^2}\right)^{n+1} + Y^n = 0.$$

Taking second-order approximations to the spatial derivatives and evaluating the quantities at the correct locations gives

$$h_i^n u_i^{n+1} - (h_i^n)^2 \left( \frac{u_{i+1}^{n+1} - u_{i-1}^{n+1}}{2\Delta x} \right) - \frac{(h_i^n)^3}{3} \left( \frac{u_{i+1}^{n+1} - 2u_i^{n+1} + u_{i-1}^{n+1}}{\Delta x^2} \right) = -Y_i^n \quad (4)$$

66 This can be rearranged into a tri-diagonal matrix that updates  $u$  given its current and pre-  
 67 vious values. So that

$$68 \begin{bmatrix} u_0^{n+1} \\ \vdots \\ u_m^{n+1} \end{bmatrix} = A^{-1} \begin{bmatrix} -Y_0^n \\ \vdots \\ -Y_m^n \end{bmatrix} =: \mathcal{G}_u(\mathbf{u}^n, \mathbf{h}^n, \mathbf{u}^{n-1}, \mathbf{h}^{n-1}, \Delta x, \Delta t). \quad (5)$$

69  
 70 Where

$$71 \quad A = \begin{bmatrix} b_0 & c_0 & & & \\ a_0 & b_1 & c_1 & & \\ & a_1 & b_2 & c_2 & \\ & & \ddots & \ddots & \ddots \\ & & & a_{m-3} & b_{m-2} & c_{m-2} \\ & & & & a_{m-2} & b_{m-1} & c_{m-1} \\ & & & & & a_{m-1} & b_m \end{bmatrix}$$

72  
 with

$$a_{i-1} = \frac{(h_i^n)^2}{2\Delta x} \frac{h_{i+1}^n - h_{i-1}^n}{2\Delta x} - \frac{(h_i^n)^3}{3\Delta x^2}, \quad (6a)$$

$$b_i = h_i^n + \frac{2h_i^n}{3\Delta x^2} \quad (6b)$$

and

$$c_i = -\frac{(h_i^n)^2}{2\Delta x} \frac{h_{i+1}^n - h_{i-1}^n}{2\Delta x} - \frac{(h_i^n)^3}{3\Delta x^2}. \quad (6c)$$

73 Lastly for completeness the final expression for  $Y_i^n$  is given by

$$74 \quad Y_i^n = 2\Delta t \left[ u_i^n h_i^n \frac{u_{i+1}^n - u_{i-1}^n}{2\Delta x} + g h_i^n \frac{h_{i+1}^{n-1} - h_{i-1}^{n-1}}{2\Delta x} + (h_i^n)^2 \left( \frac{u_{i+1}^{n-1} - u_{i-1}^{n-1}}{2\Delta x} \right)^2 \right. \\
+ \frac{(h_i^n)^3}{3} \frac{u_{i+1}^n - u_{i-1}^n}{2\Delta x} \frac{u_{i+1}^n - 2u_i^n + u_{i-1}^n}{\Delta x^2} - (h_i^n)^2 u_i^n \frac{u_{i+1}^n - 2u_i^n + u_{i-1}^n}{\Delta x^2} \\
75 \quad \left. - \frac{(h_i^n)^3}{3} u_i^n \frac{u_{j+2}^n - 2u_{j+1}^n + 2u_{j-1}^n - u_{j-2}^n}{2\Delta x^3} \right] \\
76 \quad - h_i^n u_i^{n-1} + (h_i^n)^2 \frac{u_{i+1}^{n-1} - u_{i-1}^{n-1}}{2\Delta x} + \frac{(h_i^n)^3}{3} \frac{u_{i+1}^{n-1} - 2u_i^{n-1} + u_{i-1}^{n-1}}{\Delta x^2}. \quad (7)$$

76 In particular this is an explicit numerical method for (1b), that requires the current and  
 77 previous values of  $h$  and  $u$ .

## 78 The Lax Wendroff Method for Conservation of Mass Equation

79 Because the conservation of mass equation (1a) has no mixed derivative term standard  
80 techniques of conservation laws can be used. This paper will use two, one using the Lax  
81 Wendroff method as was done by El et al. (2006) and the other will use the same finite  
82 difference approximation process as the above section. Both of these are theoretically  
83 second-order accurate. To make these methods precise they will be presented here in  
84 sufficient replicable detail.

85 Note that (1a) is in conservative law form for  $h$  where the Jacobian is  $u$ . Thus using  
86 the previously defined spatio-temporal discretisation the lax-wendroff update for  $h$  is

$$\begin{aligned} h_i^{n+1} = & h_i^n - \frac{\Delta t}{2\Delta x} ((uh)_{i+1}^n - (uh)_{i-1}^n) \\ & + \frac{\Delta t^2}{2\Delta x^2} \left( \frac{u_{i+1}^n - u_i^n}{2} ((uh)_{i+1}^n - (uh)_i^n) - \frac{u_i^n - u_{i-1}^n}{2} ((uh)_i^n - (uh)_{i-1}^n) \right). \end{aligned} \quad (8)$$

89 Performing this update for all  $i$  will be denoted by  $\mathcal{E}_h(\mathbf{u}^n, \mathbf{h}^n, \Delta x, \Delta t)$ .

## 90 Second Order Finite Difference Method

91 To follow the process above [] to obtain a second-order finite difference approximation  
92 to (1a) the derivatives are first expanded then approximated by second order centered finite  
93 differences to give

$$\frac{h_i^{n+1} - h_i^{n-1}}{2\Delta t} + u_i^n \frac{h_{i+1}^n - h_{i-1}^n}{2\Delta x} + h_i^n \frac{u_{i+1}^n - u_{i-1}^n}{2\Delta x} = 0. \quad (9)$$

96 After rearranging this to give an update formula one obtains

$$h_i^{n+1} = h_i^{n-1} - \Delta t \left( u_i^n \frac{h_{i+1}^n - h_{i-1}^n}{\Delta x} + h_i^n \frac{u_{i+1}^n - u_{i-1}^n}{\Delta x} \right). \quad (10)$$

99 Performing this update for all  $i$  will be denoted by  $\mathcal{G}_h(\mathbf{u}^n, \mathbf{h}^n, \mathbf{h}^{n-1}, \Delta x, \Delta t)$ .

## 100 The Finite Difference Methods

101 To summarise the first numerical method which naively approximates all derivatives  
102 by finite differences has the following update algorithm

$$\left. \begin{aligned} \mathbf{h}^{n+1} &= \mathcal{G}_h(\mathbf{u}^n, \mathbf{h}^n, \Delta x, \Delta t) \\ \mathbf{u}^{n+1} &= \mathcal{G}_u(\mathbf{u}^n, \mathbf{h}^n, \mathbf{u}^{n-1}, \mathbf{h}^{n-1}, \Delta x, \Delta t) \end{aligned} \right\} \mathcal{G}(\mathbf{u}^n, \mathbf{h}^n, \mathbf{u}^{n-1}, \mathbf{h}^{n-1}, \Delta x, \Delta t). \quad (11)$$

105 While the second method which follows from a naive interpretation of the numerical  
106 method described by El et al. (2006) is

$$\left. \begin{aligned} \mathbf{h}^{n+1} &= \mathcal{E}_h(\mathbf{u}^n, \mathbf{h}^n, \Delta x, \Delta t) \\ \mathbf{u}^{n+1} &= \mathcal{G}_u(\mathbf{u}^n, \mathbf{h}^n, \mathbf{u}^{n-1}, \mathbf{h}^{n-1}, \Delta x, \Delta t) \end{aligned} \right\} \mathcal{E}(\mathbf{u}^n, \mathbf{h}^n, \mathbf{u}^{n-1}, \mathbf{h}^{n-1}, \Delta x, \Delta t). \quad (12)$$

## CONSERVATIVE FORM OF THE SERRE EQUATIONS

To overcome the aforementioned difficulty of mixed derivatives the Serre equations (1) can be reformulated into conservative form which has no mixed spatio-temporal derivatives. This is accomplished by the introduction of a new quantity (Le Métayer et al. 2010; Zoppou 2014)

$$G = uh - h^2 \frac{\partial h}{\partial x} \frac{\partial u}{\partial x} - \frac{h^3}{3} \frac{\partial^2 u}{\partial x^2}. \quad (13)$$

Consequently, (1) can be rewritten as

$$\frac{\partial h}{\partial t} + \frac{\partial(uh)}{\partial x} = 0 \quad (14a)$$

and

$$\frac{\partial G}{\partial t} + \frac{\partial}{\partial x} \left( Gu + \frac{gh^2}{2} - \frac{2h^3}{3} \frac{\partial u}{\partial x} \frac{\partial u}{\partial x} \right) = 0. \quad (14b)$$

## A Hybrid Finite Difference-Volume Method for Serre Equations in Conservative Form

The conservative form (14) allows for a wider range of numerical techniques such as finite element methods (Li et al. 2014) and finite volume methods (Le Métayer et al. 2010; Zoppou 2014). In this paper the first-, second- and third-order finite difference-volume methods of [] will be used. These have been well validated and their order of accuracy has been confirmed.

## NUMERICAL SIMULATIONS

In this section the methods introduced in this paper will be validated by using them to approximate an analytic solution of the Serre equations, this will also be used to verify their order of accuracy. Then an in depth comparison of these methods for a smooth approximation to the discontinuous dam break problem will be provided to investigate the behaviour of these equations in the presence of discontinuities. This is a problem that so far has only received a proper treatment in (El et al. 2006), with other research giving only a cursory look into the topic.

## SOLITON

Currently cnoidal waves are the only family of analytic solutions to the Serre equations (Carter and Cienfuegos 2011). Solitons are a particular instance of cnoidal waves

that travel without deformation and have been used to verify the convergence rates of the described methods in this paper.

For the Serre equations the solitons have the following form

$$h(x, t) = a_0 + a_1 \operatorname{sech}^2(\kappa(x - ct)), \quad (15a)$$

$$u(x, t) = c \left( 1 - \frac{a_0}{h(x, t)} \right), \quad (15b)$$

$$\kappa = \frac{\sqrt{3a_1}}{2a_0 \sqrt{a_0 + a_1}} \quad (15c)$$

and

$$c = \sqrt{g(a_0 + a_1)} \quad (15d)$$

where  $a_0$  and  $a_1$  are input parameters that determine the depth of the quiescent water and the maximum height of the soliton above that respectively. In the simulation  $a_0 = 1\text{m}$ ,  $a_1 = 1\text{m}$  for  $x \in [-50\text{m}, 250\text{m}]$  and  $t \in [0\text{s}, 50\text{s}]$ . With  $\Delta t = 0.5C_r\Delta x$  where  $C_r = 1/\sqrt{g(a_0 + a_1)}$  which satisfies the CFL condition.

## Results

This numerical experiment and its results have been reported by []. Thus in this paper only the results for the two methods introduced in this paper are given.

## SMOOTHED DAM-BREAK

The discontinuous dam-break problem can be approximated smoothly using the hyperbolic tangent function. Such an approximation will be called a smoothed dam-break problem and will be defined as such

$$h(x, 0) = h_0 + \frac{h_1 - h_0}{2} (1 + \tanh(\alpha(x_0 - x))), \quad (16a)$$

$$u(x, 0) = 0.0\text{m/s}. \quad (16b)$$

Where  $a$  is given and controls the width of the transition between the two dam-break heights of  $h_0$  and  $h_1$ . For large  $\alpha$  the width is small and vice versa. For a fixed  $\Delta x$  there are large enough  $\alpha$  values such that the transition width is zero. This experiment was run

for both of the methods described in this paper and the 3 different order finite difference-volume methods described in []. In this particular simulation  $h_0 = 1.0m$ ,  $h_1 = 1.8m$  on  $x \in [0m, 1000m]$  for  $t \in [0s, 30s]$  with  $x_0 = 500m$ . The simulations were run changing both  $\Delta x$  and  $\alpha$ . To ensure stability  $\Delta t = 0.01\Delta x$  while for the second-order finite volume method  $\theta = 1.2$ . Since this experiment involves a very large amount of data the analysis will be broken up into three sections: decreasing  $\Delta x$ , increasing  $\alpha$  and finally differences between the methods.

## Changing $\Delta x$

Decreasing  $\Delta x$  allows the numerical method to better approximate the analytic solution to the equations. So for our valid [] numerical methods it would be expected that smaller  $\Delta x$ 's provide a closer approximation to the analytic solution. This was demonstrated for smooth problems [] above.

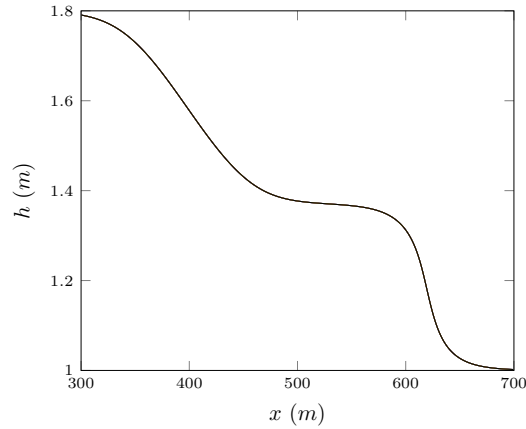
In this comparison we pick an  $\alpha$  and a method and investigate the result of decreasing  $\Delta x$ . Because the smoothness of the initial conditions depends on both  $\Delta x$  and  $\alpha$  one must be careful that the initial conditions do not change from discontinuous to smooth as  $\Delta x$  is altered as then we are no longer comparing smooth problems. This is of particular importance for the two finite difference methods as they do not correctly handle discontinuous initial conditions.

The first and most important observation is that there are four types of behaviour as  $\Delta x \rightarrow 0$  depending on the  $\alpha$  and the numerical method. It was found that the second- and third-order methods had similar  $\alpha$  ranges determining the behaviour while the first-order had very different ranges, because of this large difference the term higher-order will be used to refer to all second- and third- order methods. Also for the purposes of simplicity these scenario's will be demonstrated by solutions of the FDVM as they are better for illustrative purposes. The four scenarios are identified by the behaviour of the solutions when  $\Delta x$  is small and they correspond to different results in the literature.

The first behaviour which will be referred to as the non-oscillatory scenario has such smooth initial conditions that there are no introduced oscillations. This scenario has a highest  $\alpha$  of 0.025 with lowest unknown, in these ranges the smoothed dam-break problem is a very poor approximation to the dam-break problem. This behaviour was observed for all methods when  $\alpha = 0.025$  and an example case for the third-order method is plotted in Figure 2. This example demonstrates rapid convergence with all the solutions being graphically identical. This scenario resembles the solution of the shallow water wave equations in that it contains only a rarefaction and a shock with no dispersive waves.

For the results in Figure 2 the Hamiltonian was calculated numerically to be 10335.9175488(units) for the initial conditions and 10335.9175485 (units) for the final time of the third-order method with  $\Delta x = 10/2^{10}$  was  $7.3744593423e - 06 \times 10^{-6}$ . This gives a relative error of  $3.55675452797 \times 10^{-11}$  for the conservation of the Hamiltonian. This low relative error





(a)

FIG. 2: Smooth dam break problem for  $\alpha = 0.025$  for  $\Delta x = 10/2^{10}$  (blue),  $\Delta x = 10/2^9$  (green),  $\Delta x = 10/2^8$  (red),  $\Delta x = 10/2^7$  (cyan),  $\Delta x = 10/2^6$  (magenta),  $\Delta x = 10/2^5$  (yellow),  $\Delta x = 10/2^4$  (black)

suggests that our numerical method is solving the equations appropriately validating our results.

The second will be referred to as the flat scenario due to the presence of a constant height state between the oscillations at the shock and rarefaction fan. This scenario occurs between at least  $\alpha = 0.05$  and  $\alpha = 1$  for the higher-order methods and occurs from at least  $\alpha = 0.05$  to  $\alpha = 1000$  for the first-order method (so far). This scenario corresponds to the results presented by Le Métayer et al. (2010) and Mitsotakis et al. (2014).

An example plot demonstrating this scenario for the third-order method with  $\alpha = 0.5$  can be seen in Figure 3. As  $\Delta x$  decreases the solutions converge which is sensible since for the  $\Delta x$  in Figure [] the initial conditions are smooth as can be seen in Figure [] and these methods have been verified for smooth problems. So that by  $\Delta x = 10/2^8$  the solutions for higher  $\Delta x$  are visually identical.

For the results in Figure 2 the Hamiltonian of the initial conditions was calculated numerically to be 10395.5623489(units). The relative error for the conservation of the Hamiltonian by the third order method with  $\Delta x = 10/2^{10}$  was  $6.12566742634 \times 10^{-9}$ . This low relative error suggests that our numerical method is solving the equations appropriately validating our results. We are however seeing the trend that will continue further on that our relative error is increasing as we better approximate the discontinuous initial conditions of the dam break problem.

The third scenario will be referred to as the contact discontinuity scenario due to the

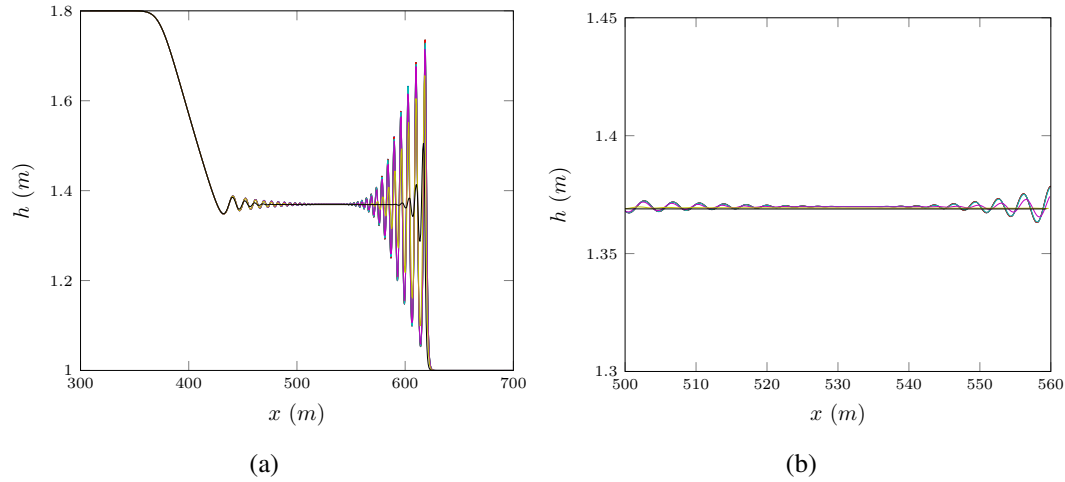


FIG. 3: Smooth dam break problem for  $\alpha = 0.5$  for  $\Delta x = 10/2^{10}$  (blue),  $\Delta x = 10/2^9$  (green),  $\Delta x = 10/2^8$  (red),  $\Delta x = 10/2^7$  (cyan),  $\Delta x = 10/2^6$  (magenta),  $\Delta x = 10/2^5$  (yellow),  $\Delta x = 10/2^4$  (black)

use of that term to describe it by El et al. (2006). For the higher-order methods it occurs at  $\alpha = 2.5$  and so far has not occurred for the first order method. The contact discontinuity scenarios main feature is that the oscillations from the rarefaction fan and the shock decay and appear to meet at a point as can be seen in Figure 4. For the experiments performed this doesn't appear to be an actual centre point but rather that the oscillations decay so quickly around the 'contact discontinuity' that it appears to be the case. All the higher order methods so far have not shown a converged solution as  $\Delta x$  decreases. However it does appear that convergence is likely with the solutions getting closer together.

For the results in Figure 4 the Hamiltonian of the initial conditions was calculated numerically to be 10398.0737089(units). The relative error for the conservation of the Hamiltonian by the third order method with  $\Delta x = 10/2^{10}$  was  $2.01012400037 \times 10^{-8}$ . This shows that we are still accurately capturing the behaviour of the equations validating the of El et al. (2006).

The fourth scenario will be referred to as the bump scenario due to the oscillations no longer decaying down towards a point but rather growing around where the contact discontinuity was in the previous scenario as can be seen in Figure 5. This behaviour has hitherto not been published and is certainly not an expected result. There are some important observations. Firstly changing  $\alpha$  does increase the height of the bump for the lowest resolution methods although after increasing  $\alpha$  has no effect. The behaviour of these solutions in Figure 5 do not clearly show convergence.

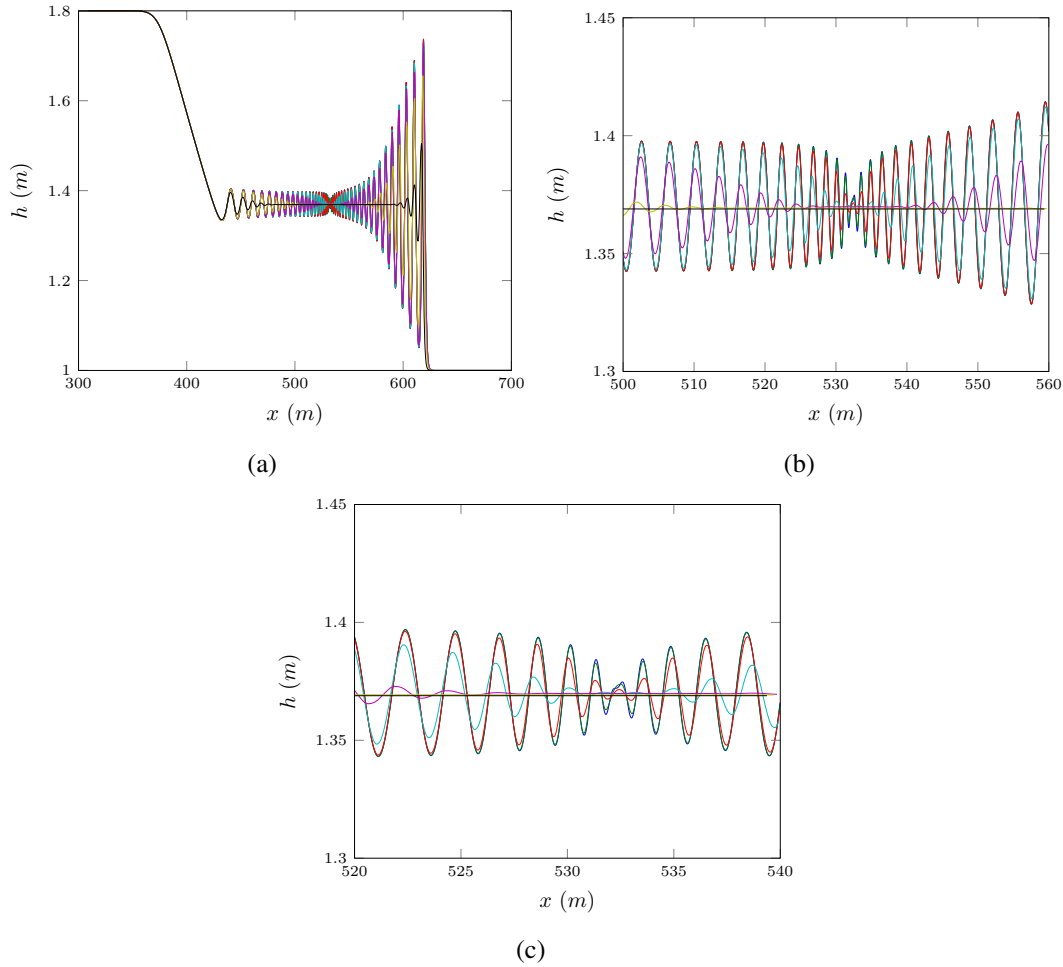


FIG. 4: Smooth dam break problem for o3 [] with  $\alpha = 2.5$  for  $\Delta x = 10/2^{10}$  (blue),  $\Delta x = 10/2^9$  (green),  $\Delta x = 10/2^8$  (red),  $\Delta x = 10/2^7$  (cyan),  $\Delta x = 10/2^6$  (magenta),  $\Delta x = 10/2^5$  (yellow),  $\Delta x = 10/2^4$  (black)

For the results in Figure 4 the Hamiltonian of the initial conditions was calculated numerically to be 10398.6937378(units). The relative error for the conservation of the Hamiltonian by the third order method with  $\Delta x = 10/2^{10}$  was  $7.3744593423e - 06 \times 10^{-6}$ . This shows that we are still accurately capturing the behaviour of the equations validating the of El et al. (2006).

All of the scenarios described above and displayed using the higher-order FDVM also occur for the FDM, however because finite differences cannot properly handle discontinuities this is a little more subtle. Firstly, since for each  $\alpha$  there is a  $\Delta x$  such that for larger

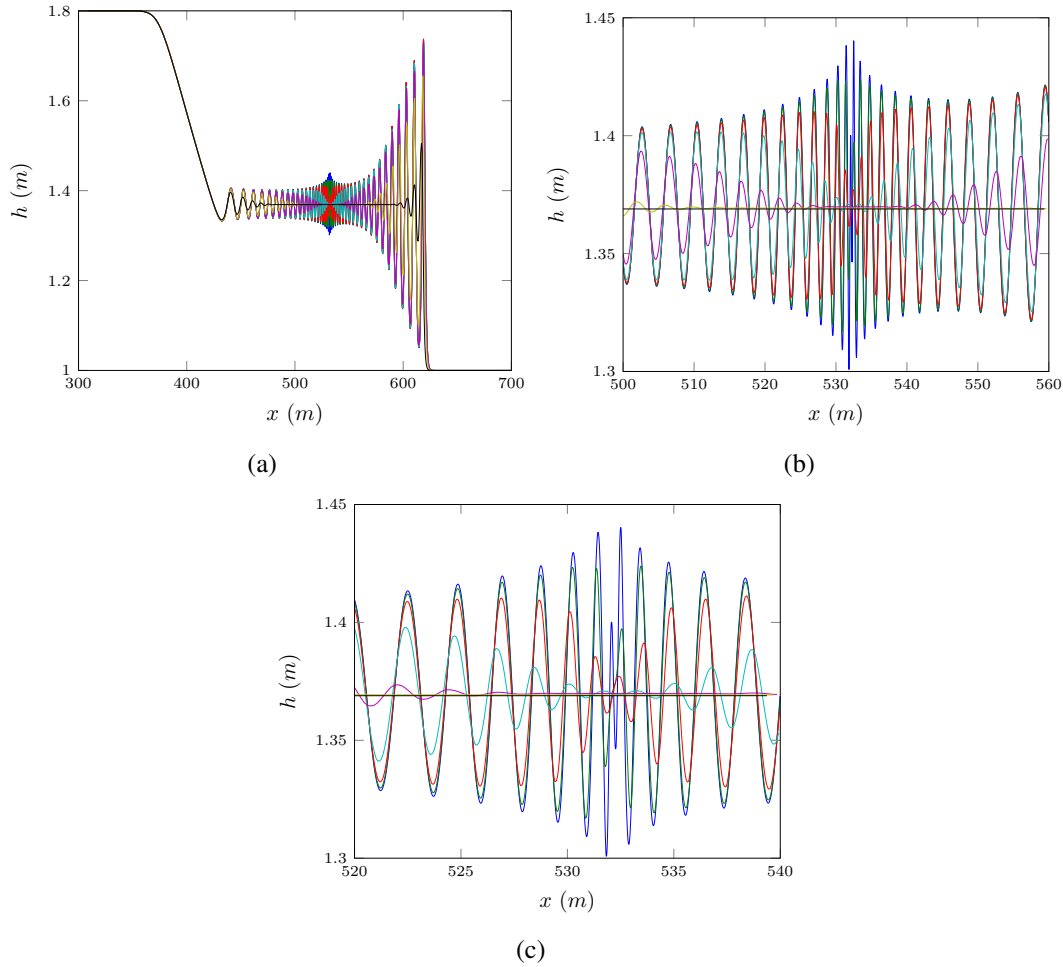


FIG. 5: Smooth dam break problem for o3 [] with  $\alpha = 1000.0$  for  $\Delta x = 10/2^{10}$  (blue),  $\Delta x = 10/2^9$  (green),  $\Delta x = 10/2^8$  (red),  $\Delta x = 10/2^7$  (cyan),  $\Delta x = 10/2^6$  (magenta),  $\Delta x = 10/2^7$  (yellow),  $\Delta x = 10/2^8$  (black)

$\Delta x$  the smooth dam break problem is no longer smooth enough for a finite difference approximation to be appropriate. This becomes a problem for the contact discontinuity and bump scenarios since they require higher  $\alpha$  and are thus more discontinuous to begin with. The result of this are non-physical looking oscillations for large  $\Delta x$  values that were not replicated by the FDVM and thus can be attributed to this flaw of FDM as in Figure [].

Overall there were two types of trending behaviours as  $\Delta x$  was decreased one for the FDM and another for the FDVM. FDM decreased the number of oscillations in the solution as in Figure [], while FDVM increased the number of oscillations in the solution as can

be seen in Figure []. This is explained by Zoppou and Roberts (1996) as the FDM are second order finite difference approximations their errors are dissipative thus introducing oscillatory errors which are most prominent when  $\Delta x$  and therefore the errors are large. While the behaviour of the FDVM is explained by a series of effects [] [TVD, treating things as cell averages, thus flattening things in cells,].

[(verify convergence rates near discontinuities?)

### Changing $\alpha$

Increasing  $\alpha$  allows the initial conditions (7) to approach the dam break problem with  $h_1$  to the left and  $h_0$  to the right centred around  $x_0$ . So it would be expected that as  $\alpha \rightarrow \infty$  that the solution of the smooth dam break problem would approach the corresponding dam break problem. This is the case for numerical methods because for a fixed  $\Delta x$   $\alpha$  can be chosen large enough that (7) is precisely the dam break problem. This can be seen in Figure [] with  $\Delta x =$  where the required  $\alpha$  for this to occur is below 1000 which was the maximum  $\alpha$  value used in these experiments. However, only the FDVM were able to handle such large  $\alpha$ 's because the initial conditions are not smooth enough to allow for stability in the FDM as can be seen in Figure []. While the FDVM handled this quite well and for all  $\Delta x$  tested as  $\alpha$  increased the solutions converged, even though for higher  $\Delta x$  []  $\alpha$  was not large enough to make (7) a jump discontinuity.

This confirms the superiority of the FDVM to handle non smooth initial conditions and the inability of FDM to handle them. Even near discontinuous initial conditions caused problems for the FDM with the introduction of oscillations that were not replicated by the FDVM and appeared to be non-physical. An example of these transitional solutions between the properly smooth initial conditions and the unstable discontinuous ones can be seen in Figure []. [(only compare the models when FD started smooth enough)

For the range of  $\alpha$ 's which are smooth enough for the FDM to be appropriate then as  $\alpha$  increases the number of oscillations increases as well for both the FDM and the FDVM. So that the smoothness of the initial conditions controls the oscillations but this depends on  $\Delta x$  since for a fixed  $\alpha$  the smoothness of the discretised initial conditions depends on  $\Delta x$ . [] (relative smoothness, more universal number)

It was observed that  $\Delta x$  can be chosen large enough such that increasing  $\alpha$  does not resolve some of the more complex structure observed for smaller  $\Delta x$  values. This  $\Delta x$  depends on the model most notably for the first-order finite difference-volume scheme this  $\Delta x$  is very small. An example of this for the third-order FDVM scheme can be seen in Figure [].

### Comparison of Models

The first-order FDVM was too diffuse and

## CONCLUSIONS

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