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Abstract

In this dissertation we use an upwind finite difference method known as Roe's scheme to solve the flux form of the shallow water equations over a sphere explicitly. The shallow water equations are transformed into a set of conservation laws, together with source terms as a result of the earth's rotation and the presence of mountains on the earth's surface. We incorporate the technique of operator splitting on a regular longitude latitude grid to allow the method to be applied to the two-dimensional shallow water equations. As a result of the first order scheme being highly diffusive we observe the need to use a second order version of Roe's method which incorporates a flux limiter. The method is applied to a standard test case in numerical weather prediction.

Contents

1	Intr	roduction	1
2	Roe	e's method	4
	2.1	Roe's scheme	4
	2.2	Why use Roe's scheme?	6
	2.3	Application to 2-d problems	7
3	$\operatorname{Th}\epsilon$	e equations	ξ
	3.1	Standard form of the shallow water equations	ē
	3.2	Obtaining the conservation form	10
	3.3	Application of Roe's method	12
		3.3.1 Applying to the longitudinal (ϕ) direction	12
		3.3.2 Applying to the latitudinal (θ) direction	19
4	$\operatorname{Th}\epsilon$	e test case and computational considerations	22
	4.1	The test case	22
	4.2	The grid	23
	4.3	Stability analysis	24
5	Res	ults	25

Bi	Bibliography				
6	Con	clusion	30		
	5.2	Translation of the bell over the poles	29		
	5.1	Translation of the bell around the equator	25		

Ch pter 1

Introduction

The weather is a very important factor in many peoples lives, so it is of great interest to know what the weather will be like in advance. Thus there is a need to be able to predict the weather, as it has a great influence on society.

In order to make predictions, mathematical models must be formed. In forming any model, it must be decided exactly what it is that is being modelled, and what factors are involved. Once the factors have been decided, a way must be found to quantify them. Only then may the laws and relationships be postulated which characterise the model. Often the result is a set of complicated differential equations for which there is no analytical solution. If this is the case then numerical techniques must be applied to find a solution to the problem.

Within the atmosphere many complex processes take place. ven if it is desired to model a particular process, quite often it is not possible to isolate that process from all of the others, and some combined effect must been considered.

Numerical weather prediction involves solving numerically a large set of differential equations. ven though certain assumptions and simplifications will have been made to derive the equations, the problem is still a complex one, and finding the solution poses many difficulties. Meteorologists are always looking to improve the numerical methods used to study the weather. Applying different numerical mation conveyed from the different latitudinal lines may conflict. One solution to this problem is of course not to have grid points at the poles.

The poles can again present a problem when regular grids are used, as the grid points bunch together in their vicinity. When explicit finite difference methods are used, this imposes the need for a smaller time step at the poles than required elsewhere on the grid, in order to satisfy the Courant-Friedrichs-Lewy (CFL) stability condition.

The purpose of this project is to implement an upwind TVD scheme, namely Roe's scheme, for the shallow water equations on a rotating sphere, which is a standard test case in numerical weather prediction. In the following section Roe's scheme will be discussed and its advantages and disadvantages over other methods considered. Chapter 3 deals with the shallow water equations and the manipulation involved to enable the implementation of the method. A test case is then presented, the results from which are shown. Finally a discussion is given on the conclusions drawn from the project.

Ch pter 2

Roe's method

2.1 Roe's scheme

Roe [8] proposed a method to obtain an approximate solution to a set of conservation laws of the form

$$\mathbf{w}_t + \mathbf{F}_x = 0 \tag{2.1}$$

based on regarding the data as piecewise constant and solving a set of Riemann problems. A Riemann problem is one where the initial data is constant either side of a discontinuity. If the discontinuity lies at the point x = x' then the initial values are

$$\mathbf{w}(0,x) = \begin{cases} \mathbf{w}_L & \text{if } x < x' \\ \mathbf{w}_R & \text{if } x > x' \end{cases}$$
 (2.2)

where \mathbf{w}_L and \mathbf{w}_R denote the left and right states and x' is the interface between them.

The solution to equation (2.1), \mathbf{w}_{j}^{n} , is regarded as an approximation to the average state between two interfaces, where the interfaces are placed at the mid

points of the cells, i.e.

$$_{j}^{n} = \frac{1}{\Delta} \frac{\left(i + \frac{1}{2}\right)\Delta x}{\left(i - \frac{1}{2}\right)dx} \quad (\Delta)$$
 (23)

where Δ is the grid spacing on a regular grid, and Δ is the time step.

If the problem (2.1) is approximated by

$$t + \tilde{x} = 0 \tag{2.4}$$

where $\tilde{}$ is a constant matrix, then an approximate solution to the exact problem (2.1) can be taken to be the exact solution to the approximate problem (2.4).

The matrix $\tilde{}$ which depends on L and R, can be picked in many ways but in Roe's scheme is chosen so that it satisfies the following properties.

- (i) $\tilde{}$ constitutes a linear mapping from the vector space to the vector space .
- (ii) As $_L$ $_R$, $\tilde{}$ $(_L$ $_R)$ $(_)$, where = —.
- (iii) For any L, R, (L R) (L R) = L R
- (iv) The eigenvectors of are linearly independent.

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the interface being

$$\mathbf{F}_{i+\frac{1}{2}}(\mathbf{w}_L, \mathbf{w}_R) = \frac{1}{2}(\mathbf{F}_L + \mathbf{F}_R) - \frac{1}{2} \sum_{k} \tilde{\alpha}_k \left| \tilde{\lambda}_k \right| \tilde{\mathbf{e}}_k.$$
 (2.6)

For linear problems, this is equivalent to the early Courant-Isaacson-Rees (CIR) method [1].

To apply the method to a non-linear problem, the local linearisation described above can be introduced by choosing \tilde{A} having property U which implies that its eigenvalues and eigenvectors not only satisfy equation (2.5) but also

$$\mathbf{w}_R - \mathbf{w}_L = \sum_k \tilde{\alpha}_k \tilde{\mathbf{e}}_k. \tag{2.7}$$

ach $\tilde{\alpha}$ satisfies a scalar scheme and the method of updating is to

add
$$-\frac{\Delta t}{\Delta x}\tilde{\lambda}_k\tilde{\alpha}_k\tilde{\mathbf{e}}_k$$
 to \mathbf{w}_R if $\tilde{\lambda}_k > 0$ (2.8)

and

add
$$-\frac{\Delta t}{\Delta x}\tilde{\lambda}_k\tilde{\alpha}_k\tilde{\mathbf{e}}_k$$
 to \mathbf{w}_L if $\tilde{\lambda}_k < 0$ (2.9)

where λ_k , $\tilde{\alpha}_k$ and $\tilde{\mathbf{e}}_k$ are determined in the calculation using the values of \mathbf{w} and \mathbf{F} from the current time step, and so the method is explicit.

2.2 Why use Roe's scheme?

In this instance the main advantages of using Roe's scheme are that it allows the use of upwinding which preserves monotonicity of the solution, and of course it is also conservative.

When applied to scalar problems, Roe's method is the same as the first order upwind method, yielding only first order accurate results that are free of oscillations. Often in a general problem where greater accuracy is required, central differences would be used to approximate the spatial derivatives. However in the case of hyperbolic conservation laws, using central differences results in a unconditionally unstable scheme, when applied explicitly. Although implicit central difference schemes have no stability limits, they can be much more computationally expensive to use than explicit methods and can also suffer from oscillations.

Another possibility is the Lax-Wendroff method. This scheme is explicit, second order and conditionally stable. However the method is also prone to oscillations which can give non-physical solutions, such as negative depth. Roe's scheme can be made second order by the introduction of a flux limiter [12]. Limiters add an anti diffusive term to the scheme which helps reduce the effects of diffusion, whilst maintaining an oscillation free method.

To enable Roe's scheme to be applied to a system in 2-d, the equations must be decomposed into a set of one dimensional problems. One technique for achieving this is known as 'operator splitting' (see [2] for application to the uler equations). For the system of equations

$$t + x + y = 0 (210)$$

applying the use of operator splitting in its simplest form results in the following

$$t$$
 x

- t y

At each time step, the solution to equations (2.11) and (2.12) are found in turn over the region and superimposed to obtain the approximate solution of (2.10).

Another more complex form of splitting which can be used to obtain the decomposition is 'Strang splitting'. This is similar to the splitting described above in that it involves rewriting the problem as a set of equations which are similar in form to equations (2.11) and (2.12). The difference between the two types lies in the order in which the equations are solved and in the size of the time step used. If we represent the problem in equation (2.10) using the simpler form of splitting as

$$L^{x}_{\frac{1}{2}\Delta t}L^{y}_{\frac{1}{2}\Delta t}$$

where $L_{\frac{1}{2}\Delta t}^x$ represents equation (2.11) and likewise $L_{\frac{1}{2}\Delta t}^y$ is equation (2.12), then the application of Strang splitting to the problem can be denoted by

$$L^{x}_{\frac{1}{4}\Delta t}L^{y}_{\frac{1}{4}\Delta t}L^{y}_{\frac{1}{4}\Delta t}L^{x}_{\frac{1}{4}\Delta t}$$

which is equivalent to

$$L^{x}_{\frac{1}{4}\Delta t}L^{y}_{\frac{1}{2}\Delta t}L^{x}_{\frac{1}{4}\Delta t}$$

so that we solve the problem along the x coordinate direction using a quarter time step, then the y direction using a half time step and then finally solve again along the x direction using a quarter time step.

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Using \mathbf{i} , \mathbf{j} and \mathbf{k} to denote the longitudinal, latitudinal and outward radial unit vectors respectively, the horizontal vector velocity is represented as $\mathbf{v} = u\mathbf{i} + v\mathbf{j}$ and the spherical horizontal gradient and divergence operators become

$$\nabla h \equiv \frac{\mathbf{i}}{a\cos\theta} \frac{\partial h}{\partial \phi} + \frac{\mathbf{j}}{a} \frac{\partial h}{\partial \theta}$$
 (3.3)

and

$$\nabla \cdot \mathbf{v} \equiv \frac{1}{a \cos \theta} \left[\frac{\partial u}{\partial \phi} + \frac{\partial (v \cos \theta)}{\partial \theta} \right]. \tag{3.4}$$

If equation (3.1) is written in terms of the velocity components then after some rearrangement, the resulting equations are

$$\frac{\partial (h^* u)}{\partial t} + \nabla \cdot (h^* u \mathbf{v}) + \frac{g h^*}{a \cos \theta} \frac{\partial h}{\partial \phi} = \left(f + \frac{u}{a} \tan \theta \right) h^* v \tag{3.5}$$

and

$$\frac{\partial (h^*v)}{\partial t} + \nabla .(h^*v\mathbf{v}) + \frac{gh^*}{a}\frac{\partial h}{\partial \theta} = -\left(f + \frac{u}{a}\tan\theta\right)h^*u. \tag{3.6}$$

xpanding the divergence terms then gives

$$\frac{\partial h^*}{\partial t} + \frac{1}{a \cos \theta} \left[\frac{\partial (h^* u)}{\partial \phi} + \frac{\partial (h^* v \cos \theta)}{\partial \theta} \right] = 0$$
 (3.7)

$$\frac{\partial(h^*u)}{\partial t} + \frac{1}{a\cos\theta} \left[\frac{\partial(h^*u^2)}{\partial\phi} + \frac{\partial(h^*uv\cos\theta)}{\partial\theta} \right] + \frac{gh^*}{a\cos\theta} \frac{\partial h}{\partial\phi} = \left(f + \frac{u}{a}\tan\theta \right) h^*v$$
(3.8)

$$\frac{\partial(h^*v)}{\partial t} + \frac{1}{a\cos\theta} \left[\frac{\partial(h^*uv)}{\partial\phi} + \frac{\partial(h^*v^2\cos\theta)}{\partial\theta} \right] + \frac{gh^*}{a} \frac{\partial h}{\partial\theta} = -\left(f + \frac{u}{a}\tan\theta\right)h^*u.$$
(3.9)

3.2 Obtaining the conservation form

As they stand, equations (3.7) to (3.9) are not in the form required to apply Roe's method. We now proceed to derive the equations in conservation law form as far

as possible. Firstly, multiplying the equations through by $\cos \theta$, and introducing a new variable h' where $h' = h^* \cos \theta$, we obtain

$$\frac{\partial h'}{\partial t} + \frac{\partial}{\partial \theta} \quad \frac{h'u}{a\cos\theta} + \frac{\partial}{\partial \theta} \quad \frac{h'v}{a} = 0 \tag{3.10}$$

$$\frac{\partial}{\partial t} (h'u) + \frac{\partial}{\partial \phi} \frac{h'uv}{a\cos\theta} + \frac{\partial}{\partial \theta} \frac{h'v^2}{a} + \frac{gh'}{a\cos\theta} \frac{\partial h}{\partial \phi} = f + \frac{u}{a}\tan\theta \quad h'v \quad (3.11)$$

$$\frac{\partial}{\partial t}(h'v) + \frac{\partial}{\partial \phi} \frac{h'uv}{a\cos\theta} + \frac{\partial}{\partial \theta} \frac{h'v^2}{a} + \frac{gh'}{a}\frac{\partial h}{\partial \theta} = f + \frac{u}{a}\tan\theta \quad h'u. \quad (3.12)$$

By substituting $h = h^* + h_s$ into equations (3.11) and (3.12), the partial derivatives involving h can be split up, and after some manipulation and rearrangement, equations (3.11) and (3.12) become

$$\frac{\partial}{\partial t} (h'u) + \frac{\partial}{\partial \phi} \frac{h'u^2}{a \cos \theta} + \frac{h'^2 g}{2a \cos^2 \theta} + \frac{\partial}{\partial \theta} \frac{h'uv}{a} =$$

$$f + \frac{u \tan \theta}{a} h'v \frac{gh'}{a \cos \theta} \frac{\partial h_s}{\partial \phi} \tag{3.13}$$

$$\frac{\partial}{\partial t}(h'v) + \frac{\partial}{\partial \phi} \frac{h'uv}{a\cos\theta} + \frac{\partial}{\partial \theta} \frac{h'v^2}{a} + \frac{h'^2g}{2a\cos\theta} = f + \frac{u\tan\theta}{a} h'u \frac{gh'}{a} \frac{\partial h_s}{\partial \theta} \frac{h'^2g\tan\theta}{2a\cos\theta}.$$
(3.14)

Together with equation (3.10), these can now be written as the matrix system

$$t + \phi + \theta = + + \tag{3.15}$$

where

$$= (h', h'u, h'R)^T R \partial \theta . h , h R R R d R 3 16 d \gamma R d$$

$$= \frac{7}{\cos^{2}} + \frac{7}{\cos^{2}} + \frac{7}{2 \cos^{2}} - \frac{7}{\cos^{2}}$$
 (3 17)

$$= \frac{\prime}{2} \frac{\prime}{2} \frac{\prime}{2} + \frac{\prime^2}{2 \cos^2}$$
 (3 18)

$$= 0 \quad \frac{'}{\cos} \quad ^{s} \quad 0 \tag{3.19}$$

$$_{t}$$
 $_{\phi}$ $-$

$$_{t}$$
 $_{\phi}$

$$\frac{1}{a\cos\theta}$$

$$\frac{m^2}{ah'^2\cos\theta} \frac{h'g}{a\cos^2\theta} \frac{2m}{ah'\cos\theta}$$

$$\frac{mn}{ah'^2\cos\theta} \frac{n}{\cos\theta} \frac{n}{\cos\theta}$$

where $\Delta(\)=(\)_R$ $\ (\)_L$ denotes the difference between the right and left states. From our definition of $\$ this leads to the following conditions,

$$\Delta' = {}_1 + {}_2 \tag{3.34}$$

$$\Delta(\ '\) = \ _1 \ + \ _2 \ + \ _3 \tag{3.36}$$

Substitution for $\ _1, \ _2$ and $\ _3$ yields the following expressions for the wavestrengths

$$_{1} = \frac{1}{2}\Delta ' + \frac{1}{2 *}[\Delta (') \Delta ']$$
 (3 37)

$$_{2} = \frac{1}{2}\Delta' \frac{1}{2} = [\Delta(') \Delta']$$
 (3.38)

$$_{3}=\Delta (\ '\)\qquad \Delta\ '$$

In applying the method, it is not equation (3.24) that is solved, but its approximation over the interval $\begin{bmatrix} L & R \end{bmatrix}$. The vectors $\begin{bmatrix} L & R \end{bmatrix}$ and $\begin{bmatrix} L & R \end{bmatrix}$ are the approximations to at the points $\begin{bmatrix} L & R \end{bmatrix}$ and $\begin{bmatrix} L & R \end{bmatrix}$ which lie either side of the interface positioned at $\begin{bmatrix} L & R \end{bmatrix}$. The problem to be solved is

$$\frac{\stackrel{n+1}{P} \stackrel{n}{\stackrel{p}{P}} + \stackrel{\sim}{\stackrel{(R L)}{\Delta}} = \stackrel{\sim}{\stackrel{(n)}{\Delta}} = (340)$$

 $0\ 3 ffi) HOh \\ 35 ffiH\ ff \\ 3\ ffi \\ 5 Hh \\ 6399 \\ i \\ S60 \\ ffi \\ 90). \\ +h0\ 3 ffi \\ 70 \\ \Omega h \\ 6N \\ 3NO \\ 90 \\ ff \\ 6(3) \\ b \\ 9ff \\ Ah \\ Affi \\ 2h \\ 22 \\ (ff \\ 6O \\ 30 \\ th \\ 90 \\ th \\ 100 \\$

Rearranging equation (3.40) gives

$$_{P}^{n+1} = {}_{P}^{n} + \Delta t \tilde{} \frac{\Delta t}{\Delta \phi} \tilde{A} \left({}_{R} \qquad {}_{L} \right). \tag{3.42}$$

To find the new value of P at the next time step, (R - L) and \tilde{P} are projected onto the local eigenvectors. If

$$_{R} \qquad _{L} = \tilde{\alpha}_{k} \tilde{\alpha}_{k}$$

$$(3.43)$$

then

$$\tilde{A}\left(\begin{array}{cc} R & L \end{array}\right) = \int_{k=1}^{3} \tilde{\lambda}_{k} \tilde{\alpha}_{k} \tilde{\lambda}_{k}$$

$$(3.44)$$

where

$$\tilde{A}\Delta = \Delta \tag{3.45}$$

Moreover, expanding the source term in terms of the eigenvectors of \tilde{A}

$$\tilde{A}(n) = \frac{1}{\Delta \phi} \int_{k=1}^{3} \tilde{\beta}_{k} \tilde{\lambda}_{k}$$

$$(3.46)$$

enables equation (3.42) to be rewritten as

$$_{P}^{n+1} = _{P}^{n} + \frac{\Delta t}{\Delta \phi} _{k=1}^{3} \tilde{\lambda}_{k} \tilde{\gamma}_{k}^{\tilde{k}}$$
 (3.47)

where $\tilde{\gamma}_k = \tilde{\alpha}_k + \tilde{\beta}_k/\tilde{\lambda}_k$.

To perform the update, each interface (or cell) is considered in turn. For a specific cell, either L or R will be incremented, depending on the sign of the eigenvalue. The result is that we

add
$$\frac{\Delta t}{\Delta} \tilde{k}_k \tilde{k}_k$$
 to R if $k = 0$ (3.48)

or

add
$$\frac{\Delta}{\Delta} \tilde{k}_k \tilde{k}_k$$
 to L if $k = 0$ (3.49)

where \tilde{k} , \tilde{k} and \tilde{k} are to be determined.

We now go on to find approximate values for the variables at the interface states based on the left and right values at the grid points. The approximate values for the eigenvalues, eigenvectors and wavestrengths within a cell are found by substituting the approximate values for u, v and $\overline{h^*g}$ at the interface into the expressions for the exact values of λ_k , k and k. Thus it remains to find approximations for k, k, and k are k and k. Thus it remains to find approximations for k, k, and k are k and k and k are k are k and k are k are k and k are k and k are k and k are k and k are k are k and k are k and k are k are k and k are k are k and k and k are k and k are k are k are k are k and k are k are k and k are k and k are k and k are k are k and k are k are k are k are k are k and k are k and k are k are k are k are k and k are k ar

Let $\tilde{\psi}^2$ denote the approximation to $\overline{h^*g}$, then from equations (3.44) and (3.45) and, after multiplication through by $\cos \theta$, the following expressions are obtained

$$\Delta(h'u) = \tilde{u} + \tilde{\psi} \quad \tilde{\alpha}_1 + \tilde{u} \quad \tilde{\psi} \quad \tilde{\alpha}_2 \tag{3.50}$$

$$\Delta h'u^2 + \frac{h'^2g}{2\cos\theta} = \tilde{u} + \tilde{\psi}^2 \tilde{\alpha}_1 + \tilde{u} \tilde{\psi}^2 \tilde{\alpha}_2$$
 (3.51)

$$\Delta(h'uv) = \tilde{u} + \tilde{\psi} \quad \tilde{v}\tilde{\alpha}_1 + \tilde{u} \quad \tilde{\psi} \quad \tilde{v}\tilde{\alpha}_2 + \tilde{u}\tilde{\alpha}_3, \tag{3.52}$$

where a tilde above a value denotes its approximation.

By expanding the brackets in equation (3.51) and rearranging the right hand side, the equation becomes

$$\Delta h'u^2 + \frac{h'^2g}{2a\cos\theta} = \tilde{u}^2 + \tilde{\psi}^2 (\tilde{\alpha}_1 + \tilde{\alpha}_2) + 2\tilde{u}\tilde{\psi}(\tilde{\alpha}_1 - \tilde{\alpha}_2). \tag{3.53}$$

Further rearrangement gives

$$\tilde{u^2}\Delta h'$$
 $2\tilde{u}\Delta h'u^2 = \tilde{\psi^2}\Delta h'$ $\Delta \frac{h'g}{2\cos\theta}$ (3.54)

One way of satisfying equation (3.54) is to make it identically zero on both sides.

Thus setting by the right hand side to zero, ² is given by

$$\tilde{r}_2 = \frac{1}{2\cos^2(R_1 + R_2')}$$
 (3.55)

Similarly, solving the quadratic for \tilde{u} gives

$$\tilde{u} = \frac{\Delta (h'u)}{\Delta (h'u)^2} \frac{\Delta (h'u)^2 (\Delta h') \Delta (h'u^2)}{\Delta h'}$$
(3.56)

and taking the negative value for the second term (the positive value leads to nothing physical), \tilde{u} becomes

$$\tilde{u} = \frac{\overline{h_R'} u_R + \overline{h_L'} u_L}{\overline{h_R'} + \overline{h_L'}}.$$
(3.57)

From equation (3.56)

$$\Delta(h'u) \quad \tilde{u}\Delta h' = \quad \overline{h'_R h'_L} \Delta u, \tag{3.58}$$

so if \tilde{h}' is defined as $\overline{h_R'h_L'}$ then

$$\Delta(h'u) \quad \tilde{u}\Delta h' = \tilde{h'}\Delta u. \tag{3.59}$$

Having found \tilde{u} , $\tilde{\psi}$ for a given \tilde{h}' we now find \tilde{v} . Rearranging equation (3.52) yields

$$\Delta(') = \tilde{ } (\tilde{ }_1 + \tilde{ }_2) + \tilde{ } (\tilde{ }_1 - \tilde{ }_2) + \tilde{ }_3$$
 (3.60)

The wavestrengths can be eliminated by using equations (3.37) to (3.39) with , replaced by the interface values and * replaced by . After some rearrangement this gives

$$\Delta(\ '\)\quad \tilde{}\Delta(\ '\)=\tilde{}\left[\Delta(\ '\)\quad \tilde{}\Delta\ '\right] \tag{3.61}$$

As with equation (3.54) we consider each side of equation (3.61) in turn and set both sides to be zero. Then substituting for ~ and taking the differences between the states gives

$$\Delta(') \quad \tilde{\Delta}' = \tilde{\Delta}' \Delta \quad \frac{\overline{R} + \overline{L} L}{\overline{R} + \overline{L}}$$
 (3 62)

1 - -

2 – —

3

1 2 3

3 —— =1

______1 2

____ 1 2 3

1 2 3

1 ———

2

In a manner similar to that applied previously, by writing as

$$= \frac{n}{a}, \frac{mn}{ah'} - \frac{2}{r} + \frac{r^2}{2 \cos^2}$$
 (3.76)

its Jacobian, B, is found to be

$$0 \qquad 0 \qquad \frac{1}{a}$$

$$= --- = \qquad \frac{uv}{a} \qquad \frac{v}{a} \qquad \frac{u}{a}$$

$$\frac{v^2}{a} + \frac{h'g}{a\cos\theta} \qquad 0 \qquad \frac{2v}{a}$$

$$(3.77)$$

with eigenvalues

$$a_1 = - + \frac{a_2}{a_1} \qquad a_2 = - \frac{a_3}{a_2} \qquad (3.78)$$

and eigenvectors

As before we can find expressions for the interface values of $\ , \ , \ \overline{\ }^*$ and $\ '$ which lead to definitions of $\ ^\sim_1, \ ^\sim_2$ and $\ ^\sim_3$ that satisfy

$$\Delta = \int_{k=1}^{3} \tilde{k}_{k}$$
 (3.80)

and

$$\Delta = \sum_{k=1}^{3} \tilde{k} \tilde{k} k \tag{3.81}$$

Following the algebra through the resulting expression for $^{\tilde{}}{}^{2}$ is

$$\tilde{z}^{2} = \frac{\Delta \frac{h'^{2}}{\cos \theta}}{\Delta'} \tag{3.82}$$

However unlike last time, cos will be different for the two states. Because must be a real quantity, as it approximates $\overline{}$, it is necessary that 2 be positive. This can be ensured by replacing the left and right values of cos with an average value, for example the value of cos at the interface. Thus 2 becomes

$$\tilde{r}_{2} = \frac{1}{2\cos I} (r_{R} + r_{L}')$$
 (3.83)

where $_{I}=\frac{1}{2}(_{L}+_{R}).$ We find that $\tilde{}$ and $\tilde{}$ are the some as in the previous case, namely

$$\tilde{a} = \frac{\overline{a}_{R} + \overline{a}_{L} + \overline{a}_{L}}{\overline{a}_{R} + \overline{a}_{L}}$$

$$(3.84)$$

and

$$\tilde{a} = \frac{\frac{a}{R} + \frac{a}{L} L}{\frac{a}{R} + \frac{a}{L}}$$

$$(3.85)$$

Setting ' L'

2 3

$$\tilde{\beta}_2 = -\frac{g\tilde{h}'}{a\tilde{\psi}\cos\theta} \tag{3.92}$$

$$\tilde{\beta}_3 = 0. \tag{3.93}$$

Ch pter

he test c se nd comput tion l consider tions

4.1 The test case

The test case to which we shall apply the method is the first problem in a suite of seven cases proposed by Williamson et al [13], developed specifically for the shallow water equations.

The problem entails predicting the motion of a cosine bell over the globe, and tests the advective part of any scheme, by specifying analytic values for the advecting winds. Different wind directions are used, which alter the path of the bell.

The advecting wind is given by

$$u = u_0 (\cos \theta \cos \alpha + \sin \theta \sin \alpha) \tag{4.1}$$

$$v = -u_0 \sin \phi \sin \alpha, \tag{4.2}$$

where α is the angle between the axis of solid body rotation and the polar axis

of the spherical coordinate system.

In the paper it suggests running the code with $\alpha = 0.0, 0.05, \pi/2 - 0.05$ and $\pi/2$. With α set to zero, the profile moves around the equator. For α equal to $\pi/2$, the trajectory is over the poles.

The initial cosine bell is defined by

$$h^*(\phi, \theta) = \begin{cases} \frac{h_0^*}{2} \left(1 + \cos\left(\frac{\pi r}{R}\right) \right) & \text{if } r < R \\ 0 & \text{if } r \ge R \end{cases}$$

$$\tag{4.3}$$

where $h_0^* = 1000m$, R = a/3 and r is the radius of the great circle distance between (ϕ, θ) and the centre of the bell which is initially at $(\phi_c, \theta_c) = (3\pi/2, 0)$. The radius r can be calculated from

$$r = a\cos^{-1}\left[\sin\theta_c\sin\theta + \cos\theta_c\cos(\phi - \phi_c)\right]. \tag{4.4}$$

The parameter values to be used are

$$a = 6.37122 \times 10^{6} m$$

$$\Omega = 7.292 \times 10^{5} s^{-1}$$

$$a = 9.80616 m s^{-2}$$

and u_0 is to be set as $2\pi a/(12 \text{ days})$ which is equivalent to about $40ms^{-1}$. There are no mountains in this problem, corresponding to h_s being zero everywhere.

If the program is run for 12 days then the initial profile should have returned to its starting point, without any change of shape.

4.2 The grid

The grid we shall use has points equally distributed in the longitudinal and latitudinal directions, with grid spacings $\Delta \phi$ and $\Delta \theta$ respectively. In this instance

the angle θ is not the standard polar coordinate but is instead measured from the equator so that θ lies in the interval $[-\pi/2, \pi/2]$ where $-\pi/2$ is the South pole and $\pi/2$ is the North pole.

There are no nodes at the poles, nodes which lie directly opposite one another closest the poles being $\frac{1}{2}\Delta\theta$ from the pole. To avoid nodes at the pole points, it becomes necessary that there are an even number of intervals in the longitudinal direction.

4.3 Stability analysis

For an equation of the form

$$w_t + au_x = 0 (4.5)$$

the stability (CFL) condition is

$$\left| a \frac{\Delta t}{\Delta x} \right| \le 1. \tag{4.6}$$

In this instance, there are two stability conditions [4] which must both be satisfied.

These are

$$\left| \frac{\tilde{\lambda}_{max}}{a \cos \theta} \frac{\Delta t}{\Delta \phi} \right| \le 1 \tag{4.7}$$

for the ϕ direction, and

$$\left| \frac{\tilde{\lambda}_{max}}{a} \frac{\Delta t}{\Delta \theta} \right| \le 1 \tag{4.8}$$

for the θ direction, where $\tilde{\lambda}_{max} = \max(\tilde{\lambda}_1, \tilde{\lambda}_2, \tilde{\lambda}_3)$ for the two problems.

From the definitions of the approximate eigenvalues for the two directions, we see that condition (4.7) imposes a more severe constraint on the time step than condition (4.8) especially as the profile tends towards the poles where $\cos \theta \to 0$.

The first figure shows a contour plot of the initial height of the cosine bell, where

= * with the centre positioned at () = $(\frac{3\pi}{2} \ 0)$ with a peak value of 1000m.

Running the program for 12 days with = 0 on a 96 73 grid using a time step of 1800 produces figure 2. Although the centre of the bell has returned to its initial position and the contour plots show the profile to be almost symmetric, a large amount of diffusion has taken place (as expected), as can be seen by the fact that the profile has spread out and that the peak value has decreased dramatically. It is expected that refining the grid would reduce this problem and as figure 3 shows where the grid spacing has been halved, this turns out to be the case.

To investigate how much the size of the time step contributes to the diffusion, the program is run with $\Delta = 450$, as shown in figure 4. From comparison with

the diffusion problem as figures 2 and 4 are almost identical.

The results so far suggest that it would be wise to introduce a flux limiter

used to decompose the system of equations. Though the squaring effect is still present, the maximum height values are much better than before.

When the program is run using $\alpha = 0.05$, the results produced are very similar to those already seen and so are not included here.

For all the runs we see that the profile has been elongated along the direction of the flow. There are no oscillations or negative heights present, and the solution remains smooth.

There are a multitude of papers on the solutions of the shallow water equations on a sphere which consider a variety of methods and give many results. We shall only draw a comparison with a few other methods here which have also been applied to the first test case in [13].

In [7] Malcolm applies a number of schemes to the shallow water equations using the first test case of Williamson et al. The results presented were obtained using a similar latitude/longitude grid to that used here. Note that in the plots the values of the height have been scaled up by 1000. The results in figure 11 were obtained using the unified model advection scheme. The contour lines lie in the range 700-1700, corresponding to the presence of negative values for the height using the normal scale. This is a result of the oscillations present. Better solutions were obtained using the 4th order Heun scheme. Figure 12 shows the solution on a 96×73 grid. The peak value has only decreased slightly, again negative values have appeared due to the oscillations occurring along the path that the bell took, and the profile steepened on the left hand side. Using a 288×217 grid the results were improved corresponding to the maximum and minimum height values being raised, and the profile being symmetric about the

line $\phi = 3\pi/2$. Further results were obtained from Heun's scheme incorporating the use of a filtering technique and using both 4th and 6th order diffusion terms. Shown also are the results from using a semi-lagrangian code. We quote in table 1 below the maximum and minimum values attained for the height profiles for some different schemes on a 96×73 grid with a 30 minute time step. Included within the table are some results obtained using a similar TVD scheme to that used here.

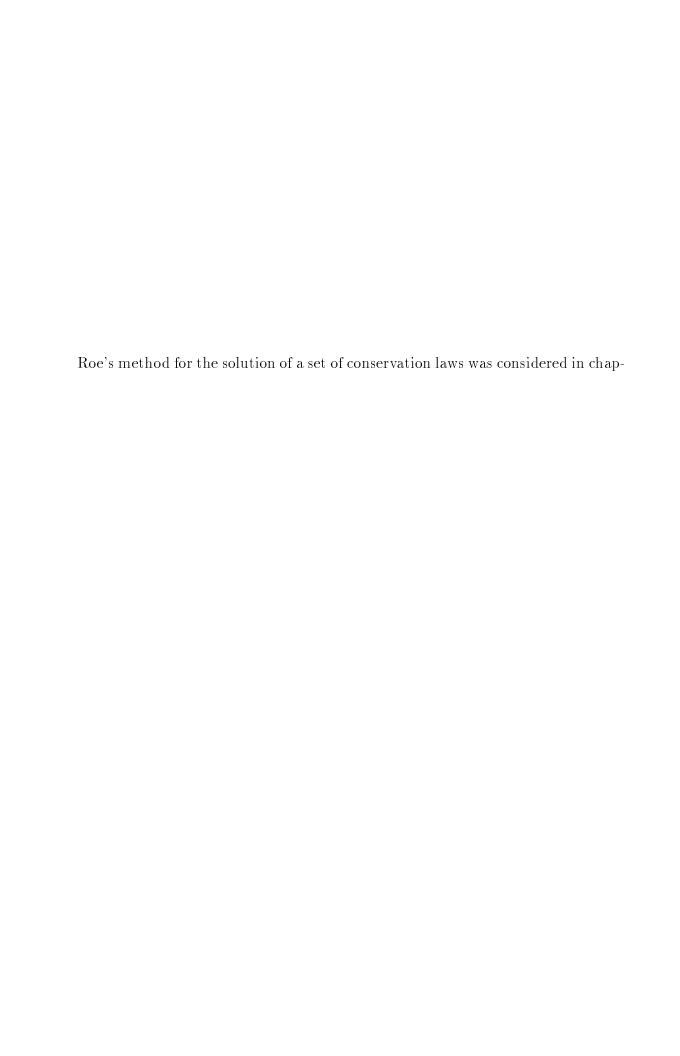
Scheme	Max. height	Min. height
Heun scheme	1937.5	836.4
Heun scheme + 4th order diffusion	1707.3	955.1
TVD (superbee)	1739.5	1000.0
TVD (van leer)	1540.0	1000.0

Table 1:

In a publication by Heikes [5] a new model was presented for the shallow water equations which solves the stream function/velocity potential form of the equations on a new grid. In the paper the results from this model were compared to those obtained using the Arakawa-Lamb and NCAR spectral shallow water models (which both use regular longitudinal/latitudinal grids) on the Williamson et al test cases. Figures 14 and 15 are the results from running the proposed model and the Arakawa-Lamb model on the first test case. In both models elongation of the profiles occurred in the direction of the flow. In this instance the solution obtained using the NCAR model is identical to the initial conditions [6].

The next part of the test is to send the profile over the poles. Because of the need for a much smaller time step than before (in order to satisfy the stability condition), the overall running time is increased, and so only 6 day runs will be

max



solution travelling with different phase speeds. We found that a very small time step had to be used to send the bell over the pole, because of the large wavespeeds in the ϕ direction, so that the solution remained stable. The form of splitting used to decompose the equation affected the quality of the solution, in particular we saw that using Strang splitting gave better results than when the simpler form of operator splitting was used.

be stretched in the direction of the flow corresponding to different parts of the

From a comparison with some other methods we have seen that the problem of the solution being stretched along the flow direction is a common problem to many methods. Although Roe's scheme is only first order and so very diffusive, we were able to reduce this problem by using flux limiters, but not to the extent that we could produce comparable results to some of the more accepted methods. We saw that one of the major problems encounted with some other methods, the presence of oscillations in the solution, was not a problem in this instance.

From the difficulties encountered with the small step size near the pole there is an obvious need to use an implicit scheme to which the size of the time step is not restricted by a stability condition, if a regular latitude/longitude grid is to be used. Alternatively, perhaps the pole problem could be overcome by using a non-regular grid where the grid points do not bunch together at the poles. It might be possible to write an adaptive code which modifies the size of the time step depending on the solution, so that at all times the maximum time step is used that will satisfy the CFL condition. However, although that would produce

a faster code when applied to this particular test case, it would not in practice be more efficient for a general problem. So far only the advective part of the code has been tested. To give a true evaluation of the methods performance further testing is needed.

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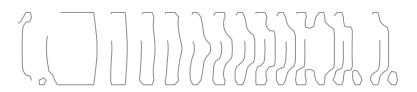
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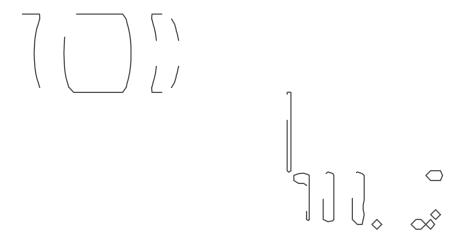


Figure 14: TIG-2562: $\alpha=0,$ a.) Day 3, b.) Day 6, c.) Day 9, d.) Day 12.

Figure 15: Arakawa-Lamb: $\alpha=0,$ a.) Day 3, b.) Day 6, c.) Day 9, d.) Day 12.