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ANALYSIS OF THACKER'S METHOD FOR SOLVING THE LINEARIZED SHALLOW WATER EQUATIONS

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1. INTRODUCTION.

In their simplest form, the shallow water equations read

$$\partial_t \vec{U}(x,t) = -b(x)\vec{\nabla}H(x,t) + fR\vec{U}(x,t), \quad x \in \Omega, \quad t \geq 0, \quad (1.1)$$

$$\partial_t H(x,t) = -\vec{\nabla} \cdot \vec{U}(x,t), \quad x \in \Omega, \quad t \geq 0, \quad (1.2)$$

$$\vec{U}(x,t) \cdot \vec{n}(x) = 0, \quad x \in \partial\Omega, \quad t \geq 0, \quad (1.3)$$

$$\vec{U}(x,0) = \vec{U}_0(x), \quad H(x,0) = H_0(x), \quad x \in \Omega. \quad (1.4)$$

Here $\Omega \subset \mathbb{R}^2$ is a bounded open domain with C^∞ boundary $\partial\Omega$ and closure $\bar{\Omega}$. \vec{n} is the outwards unit normal to $\partial\Omega$. $\vec{U}(x,t) = (\vec{U}_1(x,t), U_2(x,t))$ is a two components vector related to the average horizontal velocity. $H(x,t)$ is the height of the surface of the basin. Up to a constant factor, $b(x)$ is the depth of the basin. We shall assume that $b(\cdot)$ is a $C^\infty(\bar{\Omega})$ strictly positive function. f which represents the intensity of the Coriolis forces is taken to be constant. \vec{U}_0 and H_0 are given initial conditions. R is the $(-\pi/2)$ rotation operator acting in \mathbb{R}^2 , i.e. $R(x_1, x_2) = (x_2, -x_1)$. The tangential vector $\vec{\tau}$ at $\partial\Omega$ is given by $\vec{\tau} = -R\vec{n}$.

Equations (1.1),(1.2) can be easily set in the framework of the theory of semigroups. Let $\mathcal{H} = (L_2(\Omega))^3$ be the Hilbert space with scalar product

$$((\vec{u},g),(\vec{v},h))_{\mathcal{H}} = \int_{\Omega} \frac{1}{b(x)} \vec{u}(x) \cdot \vec{v}(x) + g(x)h(x) \quad (1.5)$$

and associate norm $\|\cdot\|_{\mathcal{H}}$ where $\vec{u} = (u_1, u_2)$, $\vec{v} = (v_1, v_2)$. We define the operator L with domain $\mathcal{D}(L)$ by the relations:

$$\mathcal{D}(L) = \{(\vec{u},g) \in \mathcal{H} \mid \vec{\nabla} \cdot \vec{u} \in L^2(\Omega), \vec{u} \cdot \vec{n} = 0 \text{ on } \partial\Omega, g \in H^1(\Omega)\} \quad (1.6)$$

$$L(\vec{u},g) = (-b\vec{\nabla}g, -\vec{\nabla} \cdot \vec{u}), \quad (1.7)$$

where $H^k(\Omega)$ is the classical Sobolev space of order k . One can verify that L is a skewadjoint operator, i.e. the adjoint of L is $-L$ which implies in particular for (\vec{u},g) and (\vec{v},h) in $\mathcal{D}(L)$ the following relation:

$$(L(\vec{u},g),(\vec{v},h))_{\mathcal{H}} = -((\vec{u},g),L(\vec{v},h))_{\mathcal{H}} \quad (1.8)$$

It follows that L is the infinitesimal generator of a conservative group so that Problem (1.1)-(1.4) possesses a unique solution with the property:

$$\|(\vec{U}(\cdot, t), H(\cdot, t))\|_{\mathcal{H}}^2 = \text{constant}; \quad (1.9)$$

furthermore, from (1.2), (1.3), one deduces immediately the law of mass conservation

$$\int_{\Omega} H(x, t) = \text{constant}. \quad (1.10)$$

Remarque 1.1: In [4], we give some results concerning the regularity of solutions of Problems (1.1)-(1.4).

The purpose of this paper is to analyse a numerical method proposed by Thacker [1], [2] for solving Problems (1.1)-(1.4); more exactly, we shall consider in fact two variants of Thacker's scheme.

2. DISCRETIZATION.

We consider a sequence $\{\mathcal{D}_h\}$ of standard triangulations of Ω , as shown in the figure. h denotes the maximum length of the sides of the triangles of \mathcal{D}_h . We assume that all angles of all triangles of all triangularizations are bounded from below by some positive constant. For a particular \mathcal{D}_h , let N be equal to three times the total number of nodes minus the number of nodes belonging to the boundary $\partial\Omega$, Ω_h will denote the interior of the union of the triangles of \mathcal{D}_h .

Let us consider a fixed triangularization \mathcal{D}_h . For each node P_k , let Λ_k be the polygon formed by the triangles containing P_k and let μ_k be the measure of Λ_k . Z_k will denote the set of indices j such that $P_j \in \partial\Lambda_k$; in the figure, $Z_2 = \{2, 3, 4, 5\}$, $Z_6 = \{7, 8, 9, 10, 11\}$; clearly, $k \in Z_k$ if and only if $P_k \in \partial\Omega$. For some node P_k , let $P_j \in \partial\Lambda_k$, i.e. $j \in Z_k$; let $P_\alpha \in \partial\Lambda_k$ be the node preceding P_j with respect to the trigonometric orientation and let $P_\beta \in \partial\Lambda_k$ be the node following P_j ; we define the vector:

$$\vec{Q}_{jk} = P_\alpha \vec{P}_\beta; \quad (2.1)$$

in the figure, we have, for example, $\vec{Q}_{8,6} = P_7 \vec{P}_5$, $\vec{Q}_{5,2} = P_4 \vec{P}_2$, $\vec{Q}_{22} = P_5 \vec{P}_3$. Furthermore, we introduce at $P_k \in \partial\Omega$ the approximate tangent and normal vectors:

$$\vec{T}_k = \frac{1}{|\vec{Q}_{kk}|} \vec{Q}_{kk}, \quad N_k = R \vec{T}_k. \quad (2.2)$$

For a function ϕ defined on Ω , let ϕ_0 be the continuous, piecewise linear (with respect to \mathcal{D}_h) function defined on Ω_h and equal to ϕ at the nodes; for a vector function $\vec{\psi}$, we define, in the same way, componentwise its interpolant $\vec{\psi}_0$. By Greens's formula, we have the identities:

$$\frac{1}{\mu_k} \int_{\Lambda_k} \vec{\nabla} \phi_0 = \frac{1}{2\mu_k} \sum_{j \in Z_k} \phi(P_j) (R\vec{Q}_{jk}), \quad (2.3)$$

$$\frac{1}{\mu_k} \int_{\Lambda_k} \vec{\nabla} \cdot \vec{\phi}_0 = \frac{1}{2\mu_k} \sum_{j \in Z_k} \vec{\phi}(P_j) \cdot (R\vec{Q}_{jk}); \quad (2.4)$$

clearly, the right members of (2.3), (2.4) define a natural approximation of $\vec{\nabla} \phi(P_k)$ and $\vec{\nabla} \cdot \vec{\phi}(P_k)$ respectively.

With the help of (2.3), (2.4), we now define a space semi-discretization of Problem (1.1)-(1.4). For all nodes P_k , $H_k(t)$ is an approximation of $H(P_k, t)$; for interior nodes $P_k \in \Omega$, $\vec{U}_k(t) = (U_{k1}(t), U_{k2}(t))$ is an approximation of $\vec{U}(P_k, t)$; for boundary nodes $P_k \in \partial\Omega$, $U_{Tk}(t)$ is an approximation of the tangential component of $\vec{U}(P_k, t)$, i.e. of $\vec{U}(P_k, t) \cdot \vec{\tau}(P_k)$. Method IS is then defined by the relations:

$$\frac{\mu_k}{b_k} \dot{\vec{U}}_k(t) = -\frac{1}{2} \sum_{j \in Z_k} H_j(t) (R\vec{Q}_{jk}) + \frac{\mu_k}{b_k} f R\vec{U}_k(t), \quad P_k \in \Omega; \quad (2.5)$$

$$\frac{\mu_k}{b_k} \dot{U}_{Tk}(t) = -\frac{1}{2} \sum_{j \in Z_k} H_j(t) (R\vec{Q}_{jk}) \cdot \vec{\tau}_k, \quad P_k \in \partial\Omega; \quad (2.6)$$

$$\mu_k \dot{H}_k(t) = -\frac{1}{2} \sum_{j \in Z_k} \vec{U}_j(t) \cdot (R\vec{Q}_{jk}), \quad P_k \in \bar{\Omega}; \quad (2.7)$$

$$\vec{U}_k(0) = \vec{U}_0(P_k), \quad P_k \in \Omega; \quad U_{Tk}(0) = \vec{U}_0(P_k) \cdot \vec{\tau}(P_k), \quad P_k \in \partial\Omega; \quad (2.8)$$

$$H_k(0) = H_0(P_k), \quad P_k \in \bar{\Omega}; \quad (2.9)$$

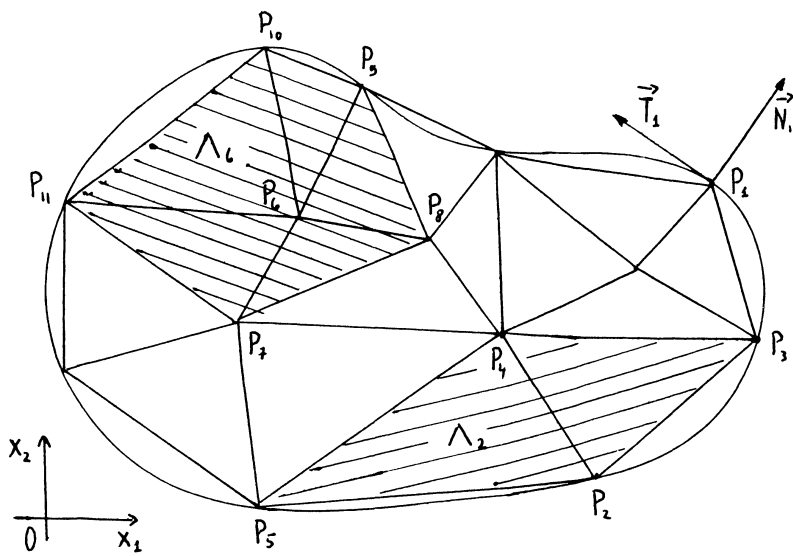
here the "dot" represents the time derivative, $b_k = b(P_k)$ and in (2.7),

$$\vec{U}_j(t) = U_{Tj}(t) \vec{\tau}_j \quad \text{if } P_j \in \partial\Omega.$$

By choosing any fixed order, all the unknown function $U_{k1}(t)$, $U_{k2}(t)$, $U_{Tk}(t)$ and $H_k(t)$ can be set in a single vector $w(t)$ of dimension N . Then Problem (2.5)-(2.9) can be written in the compact form

$$D\dot{w}(t) = Aw(t); \quad w(0) = w_0; \quad (2.10)$$

where D is a diagonal matrix with diagonal elements of the form μ_k or μ_k/b_k . Because of property (1.8) for L , one could expect that A is an antisymmetric matrix. Due to difficulties at the boundary, which seem inherent to the problems and impossible to overcome in a natural way, A is only "almost" antisymmetric. By inspection of the figure, one can show:



Lemma 2.1: Let P_k and P_j be two different nodes belonging to a same triangle. We suppose that at most one of them belong to the boundary $\partial\Omega$. Then

$$\vec{Q}_{jk} + \vec{Q}_{kj} = \vec{0}. \tag{2.11}$$

As easily seen from (2.5)-(2.7), A would be exactly antisymmetric if (2.11) would hold when if both P_k and P_j belong to $\partial\Omega$. There are several ways to modify Scheme (2.5)-(2.7) for obtaining the desired property of antisymmetry. One of them consists in replacing in (2.5), (2.6) \vec{Q}_{jk} by $-\vec{Q}_{kj}$; by Lemma 2.1, (2.5) is not modified whereas (2.6) becomes

$$\frac{u_k}{b_k} \dot{U}_{TK}(t) = \frac{1}{2} \sum_{j \in Z_k} H_j(t) (R\vec{Q}_{jk}) \cdot \vec{r}_k, \quad P_k \in \partial\Omega. \tag{2.12}$$

Method IIS is then defined by Relations (2.5), (2.12), (2.7), (2.8), (2.9) and can be written as

$$D\dot{w}(t) = Bw(t), \quad w(0) = w_0 \tag{2.13}$$

where B is an antisymmetric matrix of order N .

For a vector $v \in \mathbb{R}^N$ and a matrix G of order N , let

$$\|v\| = \left(\sum_{i=1}^N |v_i|^2 \right)^{1/2}, \quad \|G\| = \sup_{v \in \mathbb{R}^N} \frac{\|Gv\|}{\|v\|}.$$

By using the smoothness of $\partial\Omega$ and the angle property of the sequence $\{\mathcal{A}_h\}$, one can verify:

Lemma 2.2: There exist three constants c_1, c_2, c_3 , independent of h such that

$$\text{a) } \|D^{-1/2} (A-B)D^{-1/2}\| \leq c_1, \quad \text{b) } c_2 h^{-1} \leq \|D^{-1/2} B D^{-1/2}\| \leq c_3 h^{-1}.$$

By using Lemma 2.2a, the antisymmetry of B , Relations (2.7), (2.9) in connection with the definition (2.2) of \tilde{T}_k , one can deduce the following properties of Methods IS and IIS.

Proposition 2.1:

- a) If w is solution of (2.13), then $\|D^{1/2} w(t)\|^2 = \|D^{1/2} w_0\|^2$.
- b) There is a constant c , independent of t and of h , such that if w is solution of (2.10), then $\|D^{1/2} w(t)\|^2 \leq e^{ct} \|D^{1/2} w_0\|^2$.
- c) If w is solution of (2.10) or of (2.13), then $\sum_{P_k \in \bar{\Omega}} \mu_k H_k(t) = \sum_{P_k \in \bar{\Omega}} \mu_k H_0(P_k)$.

Remark 2.1: Since the sum of all μ_k 's is equal to three times the area of Ω_k , Proposition 2.1 appears to be, up to a factor 3, the discrete counterpart of Properties (1.9), (1.10) of the exact solution.

We now turn to the time discretisation. Let $\tau > 0$ be the time increment and set $t_n = n\tau$. We shall apply to (2.10) and (2.13) the two-step method, sometimes called "leap-frog" scheme. Method I is then defined by the relations:

$$w_{1/2} = w_0 + \frac{\tau}{2} D^{-1} A w_0 \quad (2.14)$$

$$w_1 = w_0 + \tau D^{-1} A w_{1/2} \quad (2.15)$$

$$w_{n+1} = w_{n-1} + 2\tau D^{-1} A w_n, \quad n = 1, 2, 3, \dots \quad (2.16)$$

Method II is defined by replacing in (2.14)-(2.16). A by B .

The stability analysis of Method II is trivial since $D^{-1} B$ or $D^{-1/2} B D^{-1/2}$ have a pure imaginary spectrum. By using Strang's Lemma [3] in connection with Lemma 2.2a, we easily deduce the stability of Method I from that of Method II.

Proposition 2.2: There exist a function $\alpha: (0,1) \rightarrow \mathbb{R}$ and a constant c , both independent of h, τ and n such that if $\tau < 1/\|D^{-1/2} B D^{-1/2}\|$, we have:

$$\text{a) } \|D^{-1/2} w_n\|^2 \leq \alpha(\tau \|D^{-1/2} B D^{-1/2}\|) \|D^{1/2} w_0\|^2 \text{ for Method II,}$$

$$\text{b) } \|D^{1/2} w_n\|^2 \leq \alpha(\tau \|D^{-1/2} B D^{-1/2}\|) e^{ctn} \|D^{1/2} w_0\|^2 \text{ for Method I.}$$

Remark 2.2: Methods I and II satisfy a law of mass conservation as Methods IS and IIS (see Proposition 2.1c).

Remark 2.3: Proposition 2.2a,b prove the stability of both Method I and II; however Method II appears to be "more" stable than Method I.

Remark 2.4: w_n can be written as a vector of order N with components U_{kn1} , U_{kn2} , U_{Tkn} and H_{kn} . If the Coriolis term $f = 0$, then for Methods I and II, it is possible to compute U_{kn1} , U_{kn2} , U_{Tkn} only at even values of n and H_{kn} at odd values of n which reduces the computer time and the storage requirements by a factor 2; real Thacker's scheme, which is somewhat more difficult to analyse, keeps this property even for $f \neq 0$. In fact, if $f = 0$, Method I is identical to Thacker's scheme (10),(11),(11') in [2] p.683.

Remark 2.5: Thacker [2] has remarked that his scheme can be considered, to some extent, as a lumped version of a Galerkin method. In [4], we briefly analyse the effect of "lumping" on stability.

Remark 2.6: The stability condition $\tau < 1/\|D^{-1/2}BD^{-1/2}\|$ in Proposition 2.2, implies, by Lemma 2.2b that $\tau = O(h)$.

3. ERROR ESTIMATES

Our estimates will be based on the standard consistency+stability argument. Stability has already been analyzed in Section 2.

We begin with a classical study of consistency by assuming that the components of the solution of Problem (1.1)-(1.4) belong to $C^0([0,T]; C^3(\bar{\Omega}))$. We first associate to this solution a vector $u(t) \in \mathbb{R}^N$ in the following way; let $w(t)$ the solution of the semi-discretized problem (2.10); we set: $u_i(t) = U_\ell(P_k, t)$ if $w_i(t) = U_{k\ell}$, $\ell = 1, 2$;

$$u_i(t) = \tilde{U}(P_k, t) \cdot \tilde{z}(P_k) \text{ if } w_i(t) = U_{TK}(t); \quad u_i(t) = H(P_k, t) \text{ if } w_i(t) = H_k(t).$$

Clearly, the time discretization which is of order two, will induce errors of size $O(\tau^2)$, which, by Remark 2.6, can be written as $O(h^2)$. Let us define for $T > 0$:

$$\epsilon_I(T) = \max_{0 \leq t \leq T} \|D^{1/2}(\dot{u}(t) - D^{-1}Au(t))\|, \quad \epsilon_{II}(T) = \max_{0 \leq t \leq T} \|D^{1/2}(\dot{u}(t) - D^{-1}Bu(t))\|,$$

$$R_{I(II)}(T) = \max_{0 \leq t_n \leq T} \|D^{1/2}(u(t_n) - w_n)\| \text{ if } w_n \text{ is obtained by Method I(II).}$$

For $i = I, II$, $R_i(T)$ is the error of Method i , whereas, by using (1.1),(1.2), $\epsilon_i(T)$ is the space consistency error.

In the following, we shall say that for a node P_k , Λ_k is symmetric, if for each $P_j \in \partial\Omega_k$, there exists $P_\ell \in \partial\Lambda_k$ which is symmetric to P_j with respect to P_k . Clearly if $P_k \in \partial\Omega$, Λ_k cannot be symmetric. The basic difference schemes defined by (2.3), (2.4) are of order 2, with respect to h , if Λ_k is symmetric; otherwise there are only of order 1. We shall say that the sequence of triangularizations $\{\mathcal{D}_h\}$ possesses Property G if there exists a constant c , independent of h such that for all $P_k \in \partial\Omega$ one has $|\vec{T}_k - \vec{\zeta}(P_k)| \leq ch^2$; Property G implies a certain regularity in the distribution of the nodes on the boundary. Elementary but tedious calculations allow to establish:

Lemma 3.1: For any fixed $T > 0$ and $i = I, II$, we have: a) $\epsilon_i(T) = O(h^{1/2})$; b) $\epsilon_i(T) = O(h)$ if Property G is satisfied; c) $\epsilon_i(T) = O(h^{3/2})$ if Λ_k is symmetric for all $P_k \in \Omega$ and if Property G is satisfied.

From Lemma 3.1 follows immediately.

Proposition 3.1: Let $T > 0$ and $\xi \in (0,1)$ be fixed numbers. For each triangularization \mathcal{D}_h , τ is chosen in such a way that $0 < \tau < \xi / \|\mathcal{D}^{-1/2} B \mathcal{D}^{-1/2}\|$. Then for $i = I, II$:
a) $R_i(T) = O(h^{1/2})$; b) $R_i(T) = O(h)$ if Property G is satisfied; c) $R_i(T) = O(h^{3/2})$ if Λ_k is symmetric for all $P_k \in \Omega$ and if Property G is satisfied.

Remark 3.1: Suppose, that, instead of (2.2), we set $\vec{T}_k = \vec{\zeta}(P_k)$ (exact tangent vector). Then: a) We lose the exact mass conservation property for both Methods I and II /see Remark 2.2); b) Proposition 3.1 remains valid for Method I; Proposition 3.1a remains valid for Method II.

We now turn to an error analysis under weaker regularity assumptions. We shall suppose that the components of the solution of Problem (1.1)-(1.4) belong to $C^0([0,T]; H^2(\Omega))$. For simplicity, we shall furthermore assume that Ω is convex so that $\Omega_h \subset \Omega$. The difficulty here comes from the fact that the time derivative of the solution is not a continuous function of the space variable. Let $\vec{U}_{kn} = (U_{kn1}, U_{kn2})$, H_{kn} be the approximate solution obtained by Method I or II corresponding to the exact solution $\vec{U}(P_k, t_n)$, $H(P_k, t_n)$; here we set $\vec{U}_{kn} = U_{Tkn} \vec{T}_k$ if $P_k \in \partial\Omega$. Let V_h be the space of continuous piecewise linear functions on Ω_h corresponding to \mathcal{D}_h . We define the function $\vec{U}_h(x, t) = (U_{h1}(x, t), U_{h2}(x, t))$, $H_h(x, t)$, for $x \in \Omega_h$ and $t = t_n$, in the following way: $U_{h\beta}(\cdot, t_n) \in V_h$, $\beta = 1, 2$, $H_h(\cdot, t_n) \in V_h$; $\vec{U}_h(P_k, t_n) = \vec{U}_{kn}$, $H_h(P_k, t_n) = H_{kn}$ for all $P_k \in \bar{\Omega}$.

The main trick will consist in introducing functions $Y_h(x, t) = (Y_{h1}(x, t), Y_{h2}(x, t))$, $Z_h(x, t)$ belonging to V_h for fixed t and which are, for fixed t , Clément's approximations of the exact solution's components; for the notion of Clément's approximation, see [7]. Corresponding to the exact equation (1.2), we have for all $P_k \in \bar{\Omega}$ the following identity:

$$\mu_k Z_h(P_k, t_{n+1}) - \mu_k Z_h(P_k, t_{n-1}) + \tau \sum_{j \in \tilde{Z}_k} \vec{\nabla}(P_j, t_n) \cdot (R\vec{Q}_{jk}) \quad (3.1)$$

$$= \{\mu_k Z_h(P_k, t_{n+1}) - \mu_k Z_h(P_k, t_{n-1}) - 2\tau \int_{\Lambda_k} \dot{H}(x, t_n)\} \quad (3.2)$$

$$+ \tau \left\{ \sum_{j \in \tilde{Z}_k} \vec{\nabla}(P_j, t_n) \cdot (R\vec{Q}_{jk}) - 2 \int_{\Lambda_k} \vec{\nabla} \cdot \vec{U}(x, t_n) \right\}. \quad (3.3)$$

The "time" error term (3.2) can be easily estimated by using the fact that the operations of time derivative and Clément's approximation commute. The "space" error term (3.3) can be handled by remarking the following identity which is a direct consequence of (2.4):

$$\sum_{j \in \tilde{Z}_k} \vec{\nabla}(P_j, t_n) \cdot (R\vec{Q}_{jk}) - 2 \int_{\Lambda_k} \vec{\nabla} \cdot \vec{\nabla}(x, t_n) = 0.$$

Similarly to (3.1)-(3.3), we can write an equation corresponding to (1.1); it is slightly more complicated to handle because of the boundary condition (1.3) and of the presence of the function $b(x)$. With Proposition 2.2, this allows to get error estimates between Clément's approximation and the solution of Method I or II. The final result is contained in the following proposition:

Proposition 3.2: Let $T > 0$ and $\xi \in (0, 1)$ be fixed numbers. We suppose: a) each component of the exact solution belong to $C^0([0, T]; H^2(\Omega))$; b) Property G is satisfied; c) for each triangularization, τ is chosen such that $0 < \tau < \xi / \|\mathcal{D}^{-1/2} B \mathcal{D}^{-1/2}\|$. Then for both Methods I and II we have:

$$\max_{0 \leq t_n \leq T} \int_{\Omega_h} \left\{ \frac{1}{b(x)} |\vec{U}(x, t_n) - \vec{U}_h(x, t_n)|^2 + |H(x, t_n) - H_h(x, t_n)|^2 \right\}^{1/2} = O(h).$$

Remark 3.2: In [4], we give some numerical results.

Remark 3.3: In order to compute the spectrum of the operator L defined in Section 1, one could think of using the same space discretization as in Method I or II; however this generates spurious eigenvalues. For a proper treatment of this problem, see [5], [6].

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