

# ON MULTISYMPLECTICITY OF PARTITIONED RUNGE–KUTTA METHODS

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**Abstract.** Previously, it has been shown that discretising a multi-Hamiltonian PDE in space and time with partitioned Runge–Kutta methods gives rise to a system of equations that formally satisfy a discrete multisymplectic conservation law. However, these studies use the same partitioning of the variables into two partitions in both space and time. This gives rise to a large number of cases to be considered, each with its own set of conditions to be satisfied. We present here a much simpler set of conditions, covering all of these cases, where the variables are partitioned independently in space and time into an arbitrary number of partitions.

In general, it is not known when such a discretisation of a multi-Hamiltonian PDE will give rise to a well defined numerical integrator. However, a numerical integrator that is explicit will typically be well defined. In this paper, we give sufficient conditions on a multi-Hamiltonian PDE for a Lobatto IIIA–IIIB discretisation in space to give rise to explicit ODEs and an algorithm for constructing these ODEs.

**1. Introduction.** A multi-Hamiltonian PDE in one time and one space dimension is a PDE which can be written as a first order system in the form

$$\mathbf{K}\mathbf{z}_t + \mathbf{L}\mathbf{z}_x = \nabla_{\mathbf{z}}S(\mathbf{z}), \quad (1.1)$$

where  $\mathbf{z} \in \mathbb{R}^n$ ,  $\mathbf{K}$  and  $\mathbf{L}$  are non-zero skew-symmetric matrices and  $S(\mathbf{z})$  is a smooth function [4].

Along solutions,  $\mathbf{z}(t, x)$ , to Eq. (1.1) the multisymplectic conservation law,

$$\omega_t + \kappa_x = 0, \quad (1.2)$$

holds, where  $\omega = \frac{1}{2}\mathbf{K}d\mathbf{z} \wedge d\mathbf{z}$  and  $\kappa = \frac{1}{2}\mathbf{L}d\mathbf{z} \wedge d\mathbf{z}$  are 2-forms and  $d\mathbf{z}$  satisfies the first variation of the PDE,

$$\mathbf{K}d\mathbf{z}_t + \mathbf{L}d\mathbf{z}_x = \mathbf{D}_{\mathbf{z}\mathbf{z}}S(\mathbf{z})d\mathbf{z} \quad (1.3)$$

where  $\mathbf{D}_{\mathbf{z}\mathbf{z}}S(\mathbf{z})$  is a symmetric matrix.

One definition of a multisymplectic integrator is a numerical method that exactly preserves a discrete analogue of Eq. (1.2) (a so-called discrete multisymplectic conservation law) by applying a symplectic one-step method in space and time [9]. An important fact here is that multisymplectic integrators do not conserve Eq. (1.2) exactly, but rather different multisymplectic integrators preserve different discrete multisymplectic conservation laws, i.e. different discretisations of Eq. (1.2). This is in contrast to symplectic integrators for ODEs, which conserve symplecticity exactly. Some of the consequences of preserving a discrete multisymplectic conservation law are:

- (i) exact preservation of some integrals, e.g. potential vorticity [12];
- (ii) both energy and momentum are approximately locally conserved [16, 17, 6];
- (iii) quasi-periodic orbits and chaotic regions are preserved (KAM theory) [21];
- (iv) the ability to take comparatively large time-steps and retain long-time stability [10].

In the past several authors [6, 11, 15, 17] have given discretisations of Eq. (1.1) which they have shown to formally satisfy a discrete multisymplectic conservation law. What these authors typically fail to consider is whether the resulting system of equations forms a well defined numerical integrator. Some problems that may occur in such discretisations are [19]:

- (i) there may be no obvious choice of dependent variables;
- (ii) the discrete equations may not be well defined locally (i.e. there may not be one equation per dependent variable per cell);
- (iii) the discrete equations may not be well defined globally (i.e. there may not be one equation per dependent variable across all spatial grid points when boundary conditions are imposed);
- (iv) the discrete equations may not have a solution, or may not have a unique solution or isolated solutions.

Difficulties due to these problems already occur for the most popular multisymplectic integrator, the Preissman box scheme. With periodic boundary conditions in one space dimension, the discrete equations typically only have solutions with an odd number of grid points, while with an even number of grid points they have no solution (nonlinear problems) or an infinite number of solutions (linear problems). With higher order Runge–Kutta (RK) methods these problems are even worse [18].

Problems (iii) and (iv) will, in general, be avoided if a discretisation method is used which gives rise to explicit multisymplectic integrators. In order to construct an explicit multisymplectic integrator, it is necessary for the discretisation in each dimension to be explicit and symplectic. For PDEs in one space and one time dimension, this condition means that a symplectic spatial discretisation must give rise to explicit ODEs in time (or vice-versa since space and time are treated on an equal footing). This rules out discretisation by symplectic RK methods. However, for some partitioned Runge–Kutta (PRK) methods this is possible, e.g. the well-known 5-point method obtained by applying leapfrog in space and time to the nonlinear wave equation,  $u_{tt} - u_{xx} = -V'(u)$ , gives the explicit multisymplectic integrator [6]:

$$\frac{1}{(\Delta t)^2} \begin{bmatrix} 1 & & \\ & -2 & \\ & & 1 \end{bmatrix} u - \frac{1}{(\Delta x)^2} \begin{bmatrix} 1 & -2 & 1 \end{bmatrix} u = -V'(u), \quad (1.4)$$

where we have used the notation of centred stencils.

Thus, in this paper we will be concerned with applying a PRK discretisation in space to obtain explicit ODEs in time. In particular we will consider the Lobatto IIIA–IIIB class of PRK discretisation, which, under certain requirements on the PDE, avoids problems (i) and (ii) and allows explicit ODEs to be obtained.

The remainder of this paper consists of four sections. In Section 2 we will describe a PRK discretisation with an arbitrary number of partitions and show that such a discretisation in time and space gives rise to a natural discrete multisymplectic conservation law which is formally satisfied. In Section 3 we give the conditions on the coefficients of a PRK discretisation to be of Lobatto IIIA–IIIB type and specify our reasons for considering the Lobatto IIIA–IIIB class of PRK discretisation. In Section 4 we give the conditions on a multi-Hamiltonian PDE such that the application of a Lobatto IIIA–IIIB discretisation in space allows one to construct explicit ODEs and then present an algorithm for constructing these ODEs. We follow this with several examples of PDEs that satisfy these conditions (such as the nonlinear wave equation and the nonlinear Schrödinger equation) and some examples of PDEs that do not. In Section 5 we will discuss some properties of the ODEs formed through our construction algorithm and give a shortcut for constructing these ODEs. We will also discuss the discretisation of these ODEs in time and their behaviour with respect to boundary conditions.

**2. Partitioned Runge–Kutta Discretisation.** When a differential equation,

$$\mathbf{z}_t = f(\mathbf{z}), \quad (2.1)$$

is discretised with a PRK discretisation, the vector of dependent variables  $\mathbf{z} \in \mathbb{R}^n$  is partitioned into several partitions,  $\mathbf{z}^{(\gamma)} \in \mathbb{R}^{n_\gamma}$  with  $\sum_\gamma n_\gamma = n$ . Typically, the number of partitions is two but it is possible for the number of partitions to be as high as  $n$ . A grid is then introduced where we take the grid points (or *nodes*) (for convenience only) to have equal spacing  $\Delta t$  and we adopt the following notation: let cell  $i$  be the region in the domain defined by  $t \in [i\Delta t, (i+1)\Delta t)$ , let  $z^\gamma$  be the entry  $\gamma$  in  $\mathbf{z}$ , let  $\mathbf{z}_i^{(\gamma)} \in \mathbb{R}^{n_\gamma}$  be the vector of variables in partition  $\gamma$  at the node in cell  $i$ , let  $\mathbf{Z}_{i,j}^{(\gamma)} \in \mathbb{R}^{n_\gamma}$  be the vector of variables in partition  $\gamma$  at stage  $j$  in cell  $i$  and let the lack of a raised index ( $\gamma$ ) indicate the unpartitioned variable.

For an  $r$ -stage PRK discretisation of Eq. (2.1) one obtains a set of equations coupling the node values  $\mathbf{z}_i$  to the stage values  $\mathbf{Z}_{i,j}$  at  $r$  internal stages given by

$$\begin{aligned} \mathbf{Z}_{i,j}^{(\gamma)} &= \mathbf{z}_i^{(\gamma)} + \Delta t \sum_{k=1}^r a_{jk}^{(\gamma)} \partial_t \mathbf{Z}_{i,k}^{(\gamma)}, \quad j = 1, \dots, r, \\ \mathbf{z}_{i+1}^{(\gamma)} &= \mathbf{z}_i^{(\gamma)} + \Delta t \sum_{j=1}^r b_j^{(\gamma)} \partial_t \mathbf{Z}_{i,j}^{(\gamma)}, \end{aligned} \quad (2.2)$$

for each  $\gamma$ , where the new variables  $\partial_t \mathbf{Z}_{i,j}$  satisfy Eq. (2.1), i.e.

$$\partial_t \mathbf{Z}_{i,j} = f(\mathbf{Z}_{i,j}), \quad (2.3)$$

and the coefficients  $b_j^{(\gamma)}$  and  $a_{jk}^{(\gamma)}$  are chosen to satisfy certain order conditions.

The conditions for a two-partition PRK discretisation of a canonical Hamiltonian ODE, with partitioning  $\mathbf{z}^{(1)} = \mathbf{q}$  and  $\mathbf{z}^{(2)} = \mathbf{p}$ , to be symplectic are [1]

$$-a_{kj}^{(1)} b_k^{(2)} - b_j^{(1)} a_{jk}^{(2)} + b_j^{(1)} b_k^{(2)} = 0 \quad \text{for all } j, k, \quad (2.4)$$

while the conditions for an RK discretisation (i.e. a one-partition PRK discretisation with  $\mathbf{z}^{(1)} = \mathbf{z}$ ,  $n_1 = n$ ) of the same ODE to be symplectic are [20]

$$-a_{kj}^{(1)} b_k^{(1)} - b_j^{(1)} a_{jk}^{(1)} + b_j^{(1)} b_k^{(1)} = 0 \quad \text{for all } j, k. \quad (2.5)$$

Generally, for a PRK discretisation with coefficients satisfying Eq. (2.4), the coefficients will not satisfy Eq. (2.5).

When the PDE (1.1) is discretised in space with an  $r$ -stage PRK discretisation, the set of equations that one obtains are given by

$$\begin{aligned} \mathbf{Z}_{i,j}^{(\gamma)} &= \mathbf{z}_i^{(\gamma)} + \Delta x \sum_{k=1}^r a_{jk}^{(\gamma)} \partial_x \mathbf{Z}_{i,k}^{(\gamma)}, \quad j = 1, \dots, r, \\ \mathbf{z}_{i+1}^{(\gamma)} &= \mathbf{z}_i^{(\gamma)} + \Delta x \sum_{j=1}^r b_j^{(\gamma)} \partial_x \mathbf{Z}_{i,j}^{(\gamma)}, \end{aligned} \quad (2.6)$$

for each  $\gamma$ , where the new variables  $\partial_x \mathbf{Z}_{i,j}$  satisfy Eq. (1.1), i.e.

$$\mathbf{K} \partial_t \mathbf{Z}_{i,j} + \mathbf{L} \partial_x \mathbf{Z}_{i,j} = \nabla_{\mathbf{z}} S(\mathbf{Z}_{i,j}). \quad (2.7)$$

Eqs. (2.6) and (2.7) form a differential-algebraic equation (DAE) for  $\mathbf{Z}_{i,j}$  and  $\mathbf{z}_i$ . However, in this DAE there are no ODEs for the node values and the constraints only apply to  $\mathbf{L}\mathbf{Z}_{i,j}$ , not  $\mathbf{Z}_{i,j}$ . Furthermore,  $\mathbf{L}$  may not have full rank, which may prevent one from obtaining a system of explicit ODEs for the  $\mathbf{Z}_{i,j}$ .

Previous studies of the PDE (1.1) discretised in space and time with PRK methods have concluded that such discretisations satisfy a natural discrete approximation of the multisymplectic conservation law (1.2) [11]. However, these studies use the same partitioning of the variables for both the space and time discretisations, which leads to a large number of cases to be considered, each with its own set of conditions to be satisfied. This choice of partitioning in each dimension is important as the conditions for the discretised equations to satisfy the discrete multisymplectic conservation law depend upon  $\mathbf{K}$  and  $\mathbf{L}$ .

For example, given a multi-Hamiltonian PDE and a two-partition PRK discretisation in time with coefficients satisfying Eq. (2.4), if the PDE has no time derivatives of the variables in the second partition, then the discretisation is in fact an RK discretisation with the same coefficients as the first of the PRK pair, which will not in general satisfy Eq. (2.5).

To consider the most general case, we will now assume the finest possible partitioning of the variables, namely  $n$  partitions where, for each entry  $\gamma$  in  $\mathbf{z}$  we have that  $n_\gamma = 1$  and the partition  $\mathbf{z}^{(\gamma)}$  consists simply of the variable  $z^\gamma$ . We will use the notation  $dZ_{i,j}^{\gamma,n,m}$  to represent the entry  $\gamma$  in  $\mathbf{z}$  at stage  $j$  of cell  $i$  in space and stage  $m$  of cell  $n$  in time, where a lack of either the index  $j$  or  $m$  indicates the node variable of cell  $i$  in space or cell  $n$  in time respectively. Also, let  $b_j^{(\gamma)}$  and  $a_{ij}^{(\gamma)}$  be the coefficients of the spatial PRK discretisation for the partitioning associated with the variable  $z^\gamma$  and let  $B_m^{(\gamma)}$  and  $A_{nm}^{(\gamma)}$  be the coefficients of the temporal PRK discretisation for the partitioning associated with the variable  $z^\gamma$ .

The following theorem gives a much simpler set of conditions for PRK discretisations of Eq. (1.1) in space and time to satisfy a discrete multisymplectic conservation law. Since it immediately applies to any other partitioning of the variables by simply equating the  $b_j^{(\gamma)}$  and  $a_{ij}^{(\gamma)}$  coefficients of the appropriate partitions in space or time, this set of conditions encompasses all of the cases considered in previous studies.

**THEOREM 2.1.** *A multi-Hamiltonian PDE (1.1) discretised by a PRK method in space and another PRK method in time has a discrete multisymplectic conservation law, given by*

$$\Delta x \sum_j b_j (\omega_{i,j}^{n+1} - \omega_{i,j}^n) + \Delta t \sum_m B_m (\kappa_{i+1}^{n,m} - \kappa_i^{n,m}) = 0, \quad (2.8)$$

where  $\omega_{i,j}^n = \frac{1}{2} \sum_{\beta,\gamma} \mathbf{K}_{\beta\gamma} dZ_{i,j}^{\gamma,n} \wedge dZ_{i,j}^{\beta,n}$  and  $\kappa_i^{n,m} = \frac{1}{2} \sum_{\beta,\gamma} \mathbf{L}_{\beta\gamma} dZ_i^{\gamma,n,m} \wedge dZ_i^{\beta,n,m}$  when the following conditions hold:

$$\begin{aligned} b_j^{(\gamma)} &= b_j, \\ -a_{kj}^{(\gamma)} b_k^{(\beta)} - b_j^{(\gamma)} a_{jk}^{(\beta)} + b_j^{(\gamma)} b_k^{(\beta)} &= 0 \end{aligned} \quad (2.9)$$

for all  $j, k$  and pairs  $(\beta, \gamma)$  such that  $\mathbf{L}_{\beta\gamma} \neq 0$  and

$$\begin{aligned} B_m^{(\gamma)} &= B_m, \\ -A_{nm}^{(\gamma)} B_n^{(\beta)} - B_m^{(\gamma)} A_{mn}^{(\beta)} + B_m^{(\gamma)} B_n^{(\beta)} &= 0 \end{aligned} \quad (2.10)$$

for all  $m, n$  and pairs  $(\beta, \gamma)$  such that  $\mathbf{K}_{\beta\gamma} \neq 0$ .

*Proof.*

$$\begin{aligned}
& (\kappa_{i+1}^{n,m} - \kappa_i^{n,m}) \\
&= \frac{1}{2} \sum_{\beta,\gamma} \left( \mathbf{L}_{\beta\gamma} dZ_{i+1}^{\gamma,n,m} \wedge dZ_{i+1}^{\beta,n,m} - \mathbf{L}_{\beta\gamma} dZ_i^{\gamma,n,m} \wedge dZ_i^{\beta,n,m} \right) \\
&= \frac{1}{2} \sum_{\beta,\gamma} \mathbf{L}_{\beta\gamma} \left( (dZ_i^{\gamma,n,m} + \Delta x \sum_j b_j^{(\gamma)} \partial_x dZ_{i,j}^{\gamma,n,m}) \wedge (dZ_i^{\beta,n,m} \right. \\
&\quad \left. + \Delta x \sum_k b_k^{(\beta)} \partial_x dZ_{i,k}^{\beta,n,m}) - dZ_i^{\gamma,n,m} \wedge dZ_i^{\beta,n,m} \right) \\
&= \frac{1}{2} \sum_{\beta,\gamma} \mathbf{L}_{\beta\gamma} \left( \Delta x (dZ_i^{\gamma,n,m} \wedge \sum_k b_k^{(\beta)} \partial_x dZ_{i,k}^{\beta,n,m} \right. \\
&\quad \left. + \sum_j b_j^{(\gamma)} \partial_x dZ_{i,j}^{\gamma,n,m} \wedge dZ_i^{\beta,n,m}) + (\Delta x)^2 \sum_{j,k} b_j^{(\gamma)} b_k^{(\beta)} \partial_x dZ_{i,j}^{\gamma,n,m} \wedge \partial_x dZ_{i,k}^{\beta,n,m} \right) \\
&= \frac{1}{2} \sum_{\beta,\gamma} \mathbf{L}_{\beta\gamma} \left( \Delta x \sum_k (dZ_{i,k}^{\gamma,n,m} - \Delta x \sum_j a_{kj}^{(\gamma)} \partial_x dZ_{i,j}^{\gamma,n,m}) \wedge b_k^{(\beta)} \partial_x dZ_{i,k}^{\beta,n,m} \right. \\
&\quad \left. + \Delta x \sum_j b_j^{(\gamma)} \partial_x dZ_{i,j}^{\gamma,n,m} \wedge (dZ_{i,j}^{\beta,n,m} - \Delta x \sum_k a_{jk}^{(\beta)} \partial_x dZ_{i,k}^{\beta,n,m}) \right. \\
&\quad \left. + (\Delta x)^2 \sum_{j,k} b_j^{(\gamma)} b_k^{(\beta)} \partial_x dZ_{i,j}^{\gamma,n,m} \wedge \partial_x dZ_{i,k}^{\beta,n,m} \right) \\
&= \frac{1}{2} \sum_{\beta,\gamma} \mathbf{L}_{\beta\gamma} \left( \Delta x \left( \sum_k b_k^{(\beta)} dZ_{i,k}^{\gamma,n,m} \wedge \partial_x dZ_{i,k}^{\beta,n,m} + \sum_j b_j^{(\gamma)} \partial_x dZ_{i,j}^{\gamma,n,m} \wedge dZ_{i,j}^{\beta,n,m} \right) \right. \\
&\quad \left. + (\Delta x)^2 \sum_{j,k} (-a_{kj}^{(\gamma)} b_k^{(\beta)} - b_j^{(\gamma)} a_{jk}^{(\beta)} + b_j^{(\gamma)} b_k^{(\beta)}) \partial_x dZ_{i,j}^{\gamma,n,m} \wedge \partial_x dZ_{i,k}^{\beta,n,m} \right) \\
&= \Delta x \sum_{\beta,\gamma,j} b_j^{(\gamma)} \mathbf{L}_{\beta\gamma} \partial_x dZ_{i,j}^{\gamma,n,m} \wedge dZ_{i,j}^{\beta,n,m} \\
&\quad + \frac{1}{2} (\Delta x)^2 \sum_{\beta,\gamma} \mathbf{L}_{\beta\gamma} \sum_{j,k} (-a_{kj}^{(\gamma)} b_k^{(\beta)} - b_j^{(\gamma)} a_{jk}^{(\beta)} + b_j^{(\gamma)} b_k^{(\beta)}) \partial_x dZ_{i,j}^{\gamma,n,m} \wedge \partial_x dZ_{i,k}^{\beta,n,m}.
\end{aligned} \tag{2.11}$$

When  $\mathbf{L}_{\beta\gamma}$  is non-zero, the  $(\Delta x)^2$  term above is zero if

$$-a_{kj}^{(\gamma)} b_k^{(\beta)} - b_j^{(\gamma)} a_{jk}^{(\beta)} + b_j^{(\gamma)} b_k^{(\beta)} = 0 \quad \text{for all } j, k. \tag{2.12}$$

Similarly,

$$\begin{aligned}
& (\omega_{i,j}^{n+1} - \omega_{i,j}^n) = \Delta t \sum_{\beta,\gamma,m} B_m^{(\gamma)} \mathbf{K}_{\beta\gamma} \partial_t dZ_{i,j}^{\gamma,n,m} \wedge dZ_{i,j}^{\beta,n,m} \\
&+ \frac{1}{2} (\Delta t)^2 \sum_{\beta,\gamma} \mathbf{K}_{\beta\gamma} \sum_{m,l} (-A_{lm}^{(\gamma)} B_l^{(\beta)} - B_m^{(\gamma)} A_{ml}^{(\beta)} + B_m^{(\gamma)} B_l^{(\beta)}) \partial_t dZ_{i,j}^{\gamma,n,m} \wedge \partial_t dZ_{i,k}^{\beta,n,m}
\end{aligned} \tag{2.13}$$

and when  $\mathbf{K}_{\beta\gamma}$  is non-zero, the  $(\Delta t)^2$  term is zero if

$$-A_{nm}^{(\gamma)} B_n^{(\beta)} - B_m^{(\gamma)} A_{mn}^{(\beta)} + B_m^{(\gamma)} B_n^{(\beta)} = 0 \quad \text{for all } m, n. \tag{2.14}$$

Now, writing Eq. (1.3) in components and taking its wedge product with  $dz^\beta$  gives

$$\sum_{\gamma} (\mathbf{K}_{\beta\gamma} \partial_t dz^\gamma \wedge dz^\beta + \mathbf{L}_{\beta\gamma} \partial_x dz^\gamma \wedge dz^\beta) = 0 \quad \text{for all } \beta \quad (2.15)$$

since  $\mathbf{D}_{\mathbf{z}\mathbf{z}}S(\mathbf{z})$  is symmetric. Thus, in general

$$\sum_{\gamma,j,m} b_j^{(\gamma)} B_m^{(\gamma)} \mathbf{L}_{\beta\gamma} (\partial_x dZ_{i,j}^{\gamma,n,m} \wedge dZ_{i,j}^{\beta,n,m}) = - \sum_{\gamma,j,m} b_j^{(\gamma)} B_m^{(\gamma)} \mathbf{K}_{\beta\gamma} (\partial_t dZ_{i,j}^{\gamma,n,m} \wedge dZ_{i,j}^{\beta,n,m}) \quad (2.16)$$

when  $b_j^{(\gamma)} = b_j$  and  $B_m^{(\gamma)} = B_m$  for all  $j, m$  and  $\gamma$ .

Therefore, if Eqs. (2.9) and (2.10) hold then we can see from Eqs. (2.11) and (2.13) that the discrete multisymplectic conservation law (2.8) holds.  $\square$

The discrete multisymplectic conservation law (Eq. (2.8)) is an approximation to the integral

$$\int_{i\Delta x}^{(i+1)\Delta x} (\omega(x, (n+1)\Delta t) - \omega(x, n\Delta t)) dx + \int_{n\Delta t}^{(n+1)\Delta t} (\kappa((i+1)\Delta x, t) - \kappa(i\Delta x, t)) dt = 0, \quad (2.17)$$

which is the integral of Eq. (1.2) over the cell with lower left corner at  $(i\Delta x, n\Delta t)$ .

Now, suppose we have a two-partition PRK discretisation in space where the coefficients satisfy Eq. (2.4) but not Eq. (2.5), then for Eq. (2.12) to be satisfied the partitioning of the variables in space must be chosen such that  $\kappa$  only has terms of the form  $dz^{(1)} \wedge dz^{(2)}$ . Similarly, given a two-partition PRK discretisation in time where the coefficients satisfy Eq. (2.4) but not Eq. (2.5), for Eq. (2.14) to be satisfied the partitioning of the variables in time must be chosen such that  $\omega$  only has terms of the form  $dz^{(1)} \wedge dz^{(2)}$ .

Theorem 2.1 shows that if the partitioning in space and time is chosen appropriately, then a PRK discretisation in space and time with coefficients satisfying Eq. (2.4) will result in an integrator that formally satisfies a multisymplectic conservation law given by Eq. (2.8). However, this does not guarantee that the integrator is well defined. The approach we take to obtaining a well defined multisymplectic integrator is to apply an explicit symplectic PRK discretisation in each dimension.

We define an *explicit discretisation* in space as a discretisation for which the time derivatives of the dependent variables may be written explicitly in terms of the dependent variables. Their derivation may involve solving linear systems, but these must be independent of the PDE. An *explicit local discretisation* is an explicit discretisation for which these ODEs depend only on nearby values of the dependent variables.

In Section 4 we will give the conditions on a multi-Hamiltonian PDE such that one can obtain an explicit local symplectic PRK discretisation in space based on Lobatto IIIA–IIIB and we will give an algorithm for obtaining the explicit ODEs in time.

**3. Lobatto IIIA–IIIB.** The particular class of PRK discretisation that we consider in this paper is a two-partition discretisation known as Lobatto IIIA–IIIB. For

these methods, the coefficients  $a_{ij}^{(1)}$ ,  $a_{ij}^{(2)}$  and  $b_j^{(1)} = b_j^{(2)} = b_j$  are determined by [7]:

$$\begin{aligned} B(r) : \quad & \sum_{i=1}^r b_i c_i^{k-1} = \frac{1}{k} \quad \text{for } k \leq r \\ C(r) : \quad & \sum_{j=1}^r a_{ij}^{(1)} c_j^{k-1} = \frac{1}{k} c_i^k \quad \text{for } i = 1, \dots, r \text{ and } k \leq r, \\ D(r) : \quad & \sum_{i=1}^r b_i c_i^{k-1} a_{ij}^{(2)} = \frac{1}{k} b_j (1 - c_j^k) \quad \text{for } j = 1, \dots, r \text{ and } k \leq r, \end{aligned} \quad (3.1)$$

where the  $c_i$  are zeros of the Lobatto quadrature polynomial

$$\frac{d^{r-2}}{dx^{r-2}} (x^{r-1}(x-1)^{r-1}). \quad (3.2)$$

While the Lobatto IIIA and Lobatto IIIB classes of RK methods have each been known since the mid 1960s, their coefficients do not satisfy Eq. (2.5) and it was only discovered relatively recently that the Lobatto IIIA–IIIB class of PRK methods formed by combining Lobatto IIIA and Lobatto IIIB has coefficients that satisfy Eq. (2.4) [14, 22]. Thus for a discretisation of Eq. (1.1), if the partitioning of the variables in each of the space and time dimensions can be chosen such that the two-form associated with each dimension only has terms of the form  $dz^{(1)} \wedge dz^{(2)}$ , then the resulting integrator will satisfy a discrete multisymplectic conservation law.

The reason we consider the Lobatto IIIA–IIIB class of PRK discretisations is because their coefficients are related in the following way:

$$a_{1j}^{(1)} = 0, \quad a_{rj}^{(1)} = b_j \quad \text{for all } j, \quad (3.3)$$

$$a_{ir}^{(2)} = 0, \quad a_{i1}^{(2)} = b_1 \quad \text{for all } i, \quad (3.4)$$

and the  $(r-2) \times (r-2)$  matrix  $\mathbf{C}$  with entries

$$\mathbf{C}_{i-1, j-1} = \sum_{k,l} a_{ik}^{(1)} (b_l - \delta_{kl}) a_{lj}^{(2)} \quad \text{for } 2 \leq i, j \leq r-1 \quad (3.5)$$

is invertible.

The relations given in Eqs. (3.3) and (3.4) are a direct consequence of Eqs. (3.1) and (3.2) and give us three properties which will be required in our algorithm for constructing explicit ODEs in the next section. Firstly, from Eq. (3.3) we can see that for  $\gamma = 1$ , a node value is equal to the first stage value associated with that node and also equal to the last stage value associated with the previous node. Secondly, Eq. (3.4) gives us that both  $\sum_j b_j a_{jr}^{(2)}$  and  $b_1 - \sum_j b_j a_{j1}^{(2)}$  are zero. Lastly, Eqs. (3.3) and (3.4) together give

$$\sum_{k,l} a_{ik}^{(1)} (b_l - \delta_{kl}) a_{lj}^{(2)} = 0 \quad \text{if either } i \in \{1, r\} \text{ or } j \in \{1, r\}, \quad (3.6)$$

where  $\delta_{kl}$  is the Kronecker delta. Eq. (3.5) will be used in the construction algorithm directly.

The coefficients for Lobatto IIIA–IIIB methods can be written succinctly as pairs of Butcher tableaux, we give below the coefficients for  $r = 2, 3$  and 4.

$$r = 2 : \quad \text{IIIA: } \begin{array}{c|cc} 0 & 0 & 0 \\ 1 & \frac{1}{2} & \frac{1}{2} \\ \hline & \frac{1}{2} & \frac{1}{2} \end{array}, \quad \text{IIIB: } \begin{array}{c|cc} 0 & \frac{1}{2} & 0 \\ 1 & \frac{1}{2} & 0 \\ \hline & \frac{1}{2} & \frac{1}{2} \end{array}. \quad (3.7)$$

Second order Lobatto IIIA–IIIB is often referred to as generalised leapfrog.

$$r = 3 : \quad \text{IIIA: } \begin{array}{c|ccc} 0 & 0 & 0 & 0 \\ \frac{1}{2} & \frac{5}{24} & \frac{1}{3} & -\frac{1}{24} \\ 1 & \frac{1}{6} & \frac{2}{3} & \frac{1}{6} \\ \hline & \frac{1}{6} & \frac{2}{3} & \frac{1}{6} \end{array}, \quad \text{IIIB: } \begin{array}{c|ccc} 0 & \frac{1}{6} & -\frac{1}{6} & 0 \\ \frac{1}{2} & \frac{1}{6} & \frac{1}{3} & 0 \\ 1 & \frac{1}{6} & \frac{5}{6} & 0 \\ \hline & \frac{1}{6} & \frac{2}{3} & \frac{1}{6} \end{array}. \quad (3.8)$$

$$r = 4 : \quad \text{IIIA: } \begin{array}{c|ccccc} 0 & 0 & 0 & 0 & 0 \\ \frac{5-\sqrt{5}}{10} & \frac{11+\sqrt{5}}{120} & \frac{25-\sqrt{5}}{120} & \frac{25-13\sqrt{5}}{120} & \frac{-1+\sqrt{5}}{120} \\ \frac{5+\sqrt{5}}{10} & \frac{11-\sqrt{5}}{120} & \frac{25+13\sqrt{5}}{120} & \frac{25+\sqrt{5}}{120} & \frac{-1-\sqrt{5}}{120} \\ 1 & \frac{1}{12} & \frac{5}{12} & \frac{5}{12} & \frac{1}{12} \\ \hline & \frac{1}{12} & \frac{5}{12} & \frac{5}{12} & \frac{1}{12} \end{array}, \quad (3.9)$$

$$\text{IIIB: } \begin{array}{c|ccccc} 0 & \frac{1}{12} & \frac{-1-\sqrt{5}}{24} & \frac{-1+\sqrt{5}}{24} & 0 \\ \frac{5-\sqrt{5}}{10} & \frac{1}{12} & \frac{25+\sqrt{5}}{120} & \frac{25-13\sqrt{5}}{120} & 0 \\ \frac{5+\sqrt{5}}{10} & \frac{1}{12} & \frac{25+13\sqrt{5}}{120} & \frac{25-\sqrt{5}}{120} & 0 \\ 1 & \frac{1}{12} & \frac{11-\sqrt{5}}{24} & \frac{11+\sqrt{5}}{24} & 0 \\ \hline & \frac{1}{12} & \frac{5}{12} & \frac{5}{12} & \frac{1}{12} \end{array}.$$

**4. Explicit ODEs.** In the one dimensional situation (i.e. time integration), the dependent variables are the  $\mathbf{z}_i$ ; Eq. (2.2) determines the stage variables  $\mathbf{Z}_{i,j}$  and defines a map from  $\mathbf{z}_i$  to  $\mathbf{z}_{i+1}$ . In contrast, for situations where the dimension is greater than one (e.g. for PDEs of the form of Eq. (1.1)), if one applies a PRK discretisation in space, then the dependent variables will typically be the stage variables  $\mathbf{Z}_{i,j}$ , while the node variables  $\mathbf{z}_i$  and the new variables  $\partial_x \mathbf{Z}_{i,j}$  will be eliminated using the PDE to yield a set of ODEs in time for the  $\mathbf{Z}_{i,j}$ . As we shall see in the following theorem this elimination depends upon the structure of not only  $\mathbf{K}$  and  $\mathbf{L}$ , but also of  $S(\mathbf{z})$ .

**THEOREM 4.1.** *Consider a multi-Hamiltonian PDE (1.1) where the  $\mathbf{K}$  and  $\mathbf{L}$  matrices have the following structure:*

$$\mathbf{K} = \begin{bmatrix} & -I_{\frac{1}{2}(d_1+d_2)} & \\ I_{\frac{1}{2}(d_1+d_2)} & & \\ & & 0_{d_1} \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} & & I_{d_1} \\ & 0_{d_2} & \\ -I_{d_1} & & \end{bmatrix} \quad (4.1)$$

where  $d_1 = n - \text{rank}(\mathbf{K})$ ,  $d_2 = n - 2d_1 \leq d_1$ ,  $I_d$  is the  $d \times d$  identity matrix,  $0_d$  is the  $d \times d$  zero matrix.

Let the variables  $\mathbf{z}$  be partitioned into two partitions  $\mathbf{z}^{(1)} \in \mathbb{R}^{d_1+d_2}$  and  $\mathbf{z}^{(2)} \in \mathbb{R}^{d_1}$ , where we denote the first  $d_1$  components of  $\mathbf{z}^{(1)}$  by  $\mathbf{q}$ , the last  $d_2$  components of  $\mathbf{z}^{(1)}$



by  $\mathbf{v}$  and the components of  $\mathbf{z}^{(2)}$  by  $\mathbf{p}$  such that the PDE may be written as

$$\begin{bmatrix} & -\mathbf{I}_{\frac{1}{2}(d_1+d_2)} \\ \mathbf{I}_{\frac{1}{2}(d_1+d_2)} & \\ & \mathbf{0}_{d_1} \end{bmatrix} \begin{bmatrix} \mathbf{q} \\ \mathbf{v} \\ \mathbf{p} \end{bmatrix}_t + \begin{bmatrix} & & I_{d_1} \\ & \mathbf{0}_{d_2} & \\ -I_{d_1} & & \end{bmatrix} \begin{bmatrix} \mathbf{q} \\ \mathbf{v} \\ \mathbf{p} \end{bmatrix}_x = \begin{bmatrix} \nabla_{\mathbf{q}} S(\mathbf{z}) \\ \nabla_{\mathbf{v}} S(\mathbf{z}) \\ \nabla_{\mathbf{p}} S(\mathbf{z}) \end{bmatrix}. \quad (4.2)$$

If the function  $S(\mathbf{z})$  can be written in the form

$$S(\mathbf{z}) = T(\mathbf{p}) + V(\mathbf{q}) + \widehat{V}(\mathbf{v}) \quad (4.3)$$

where  $T(\mathbf{p}) = \frac{1}{2}\mathbf{p}^t \beta \mathbf{p}$  and  $\widehat{V}(\mathbf{v}) = \frac{1}{2}\mathbf{v}^T \alpha \mathbf{v}$  such that  $|\beta| \neq 0$  and  $|\alpha| \neq 0$ , then applying an  $r$ -stage Lobatto IIIA–IIIB PRK discretisation in space to the PDE leads to a set of explicit local ODEs in time in the stage variables associated with  $\mathbf{q}$ .

*Proof.* Due to the form of  $S(\mathbf{z})$ , the central  $d_2$  rows of Eq. (4.2) allow us to write entry  $i$  in  $\mathbf{v}$  as

$$v_i = \sum_{j=1}^{d_2} (\alpha^{-1})_{i,j} \partial_t q_{j+\frac{1}{2}(d_1-d_2)} \quad (4.4)$$

and hence

$$\partial_t v_i = \sum_{j=1}^{d_2} (\alpha^{-1})_{i,j} \partial_t^2 q_{j+\frac{1}{2}(d_1-d_2)}. \quad (4.5)$$

Substituting Eq. (4.5) into Eq. (4.2), we can eliminate the  $\mathbf{v}$  variables in favour of higher order derivatives in time of the  $\mathbf{q}$  variables. This lets us write Eq. (4.2) as

$$\mathbf{Kz}_t + \mathbf{Lz}_x - \mathcal{Ez}_{tt} = \nabla_{\mathbf{z}} S(\mathbf{z}), \quad (4.6)$$

where  $\mathbf{z}$ ,  $\mathbf{K}$ ,  $\mathbf{L}$ ,  $\mathcal{E}$  and  $S(\mathbf{z})$  are the new vectors, matrices and functions given below:

$$\mathbf{z} = \begin{bmatrix} \mathbf{q} \\ \mathbf{p} \end{bmatrix}, \quad \mathbf{K} = \begin{bmatrix} & & -\mathbf{I}_{\frac{1}{2}(d_1-d_2)} \\ & \mathbf{0}_{d_2} & \\ \mathbf{I}_{\frac{1}{2}(d_1-d_2)} & & \\ & & & \mathbf{0}_{d_1} \end{bmatrix}, \quad (4.7)$$

$$\mathbf{L} = \begin{bmatrix} & \mathbf{I}_{d_1} \\ -\mathbf{I}_{d_1} & \end{bmatrix}, \quad \mathcal{E} = \begin{bmatrix} \mathbf{0}_{\frac{1}{2}(d_1-d_2)} & & & \\ & \alpha^{-1} & & \\ & & \mathbf{0}_{\frac{1}{2}(d_1-d_2)} & \\ & & & \mathbf{0}_{d_1} \end{bmatrix}$$

and  $S(\mathbf{z}) = T(\mathbf{p}) + V(\mathbf{q})$ .

Note that if  $d_2 = 0$ , then Eq. (4.2) and Eq. (4.6) are identical; i.e.  $\widehat{V}(\mathbf{v}) \equiv 0$  and  $\mathcal{E}$  is a  $d_1 \times d_1$  matrix of zeros.

We shall now give a five step algorithm for constructing explicit local ODEs in time from an  $r$ -stage Lobatto IIIA–IIIB PRK discretisation of Eq. (4.6). However, before we begin, it is necessary to introduce the following notation which will be used throughout the remainder of this text.

- (i)  $Z_{i,j}^\eta$  is the stage variable at stage  $j$  in cell  $i$  for the entry  $\eta$  in  $\mathbf{z}$ ,
- (ii)  $\mathbf{Z}_i^\eta$  is the vector of stage variables in cell  $i$  for the entry  $\eta$  in  $\mathbf{z}$ ,
- (iii)  $Z_{i,j}$  is  $Z_{i,j}^\eta$  for all values of  $\eta$ ,
- (iv)  $\partial_t^n Z_{i,j}^\eta$  is a variable representing the first ( $n = 1$ ) and second ( $n = 2$ ) time derivatives of  $Z_{i,j}^\eta$ ,
- (v)  $z_i^\eta$  is the node variable in cell  $i$  for the entry  $\eta$  in  $\mathbf{z}$ ,
- (vi)  $\partial_{z^\eta} S(\mathbf{Z}_i)$  is the vector of stage values at cell  $i$  obtained by taking the derivative of the function  $S(\mathbf{z})$  with respect to the entry  $\eta$  in  $\mathbf{z}$ ,
- (vii)  $\mathbf{A}^{(1)}$  is the  $r \times r$  matrix of  $a_{ij}$  values for Lobatto IIIA,
- (viii)  $\mathbf{A}^{(2)}$  is the  $r \times r$  matrix of  $a_{ij}$  values for Lobatto IIIB,
- (ix)  $\mathbf{b}$  is the common vector of length  $r$  of  $b_j$  values for Lobatto IIIA and IIIB,
- (x)  $\mathbf{1}$  is a vector of length  $r$  with all entries equal to 1.

Now, Eq. (4.6) discretised in space by an  $r$ -stage Lobatto IIIA–IIIB PRK discretisation results in the following system of implicit ODEs:

$$\mathbf{Q}_i^\eta = q_i^\eta \mathbf{1} + \Delta x \mathbf{A}^{(1)} (-\partial_{p^\eta} T(\mathbf{P}_i)), \quad (4.8)$$

$$q_{i+1}^\eta = q_i^\eta + \Delta x \mathbf{b}^T (-\partial_{p^\eta} T(\mathbf{P}_i)), \quad (4.9)$$

$$\mathbf{P}_i^\eta = p_i^\eta \mathbf{1} + \Delta x \mathbf{A}^{(2)} (\partial_{q^\eta} V(\mathbf{Q}_i) + g_i^\eta), \quad (4.10)$$

$$p_{i+1}^\eta = p_i^\eta + \Delta x \mathbf{b}^T (\partial_{q^\eta} V(\mathbf{Q}_i) + g_i^\eta), \quad (4.11)$$

for  $1 \leq \eta \leq d_1$ , where

$$g_i^\eta = \begin{cases} \partial_t \mathbf{Q}_i^{\eta + \frac{1}{2}(d_1 + d_2)}, & 1 \leq \eta \leq \frac{1}{2}(d_1 - d_2), \\ -\partial_t \mathbf{Q}_i^{\eta - \frac{1}{2}(d_1 + d_2)}, & \frac{1}{2}(d_1 + d_2) < \eta \leq d_1, \\ \sum_{\theta=1}^{d_2} (\alpha^{-1})_{\eta - \frac{1}{2}(d_1 - d_2), \theta} \partial_t^2 \mathbf{Q}_i^{\theta + \frac{1}{2}(d_1 - d_2)}, & \frac{1}{2}(d_1 - d_2) < \eta \leq \frac{1}{2}(d_1 + d_2). \end{cases} \quad (4.12)$$

It should be noted that for the simpler case where  $d_2 = 0$ , the third option for  $g_i^\eta$  vanishes.

### Construction Algorithm:

#### Step 1:

A special property of the Lobatto IIIA discretisation is that the first row of the coefficient matrix  $\mathbf{A}^{(1)}$  is zero and the last row of  $\mathbf{A}^{(1)}$  is  $\mathbf{b}^T$ .

Due to this property, we can see that the first row of Eq. (4.8) gives  $q_i^\eta = Q_{i,1}^\eta$  and comparing the last row of Eq. (4.8) with Eq. (4.9) gives  $q_{i+1}^\eta = Q_{i,r}^\eta$ . Furthermore, from these two identities we can conclude that  $Q_{i,r}^\eta = Q_{i+1,1}^\eta$ ,  $\partial_t Q_{i,r}^\eta = \partial_t Q_{i+1,1}^\eta$  and  $\partial_t^2 Q_{i,r}^\eta = \partial_t^2 Q_{i+1,1}^\eta$ .

#### Step 2:

Since  $T(\mathbf{p}) = \frac{1}{2} \mathbf{p}^T \beta \mathbf{p}$  and  $|\beta| \neq 0$  we have that  $\mathbf{P}_i = \beta^{-1} \nabla_{\mathbf{p}} T(\mathbf{P}_i)$ . Also a property of all RK and PRK discretisations is that  $\mathbf{b}^T \mathbf{1} = 1$ . Therefore we can substitute  $\mathbf{P}_i^\eta$  from Eq. (4.10) into Eq. (4.9) and rearrange to get

$$p_i^\eta = -\frac{1}{\Delta x} \sum_{\zeta=1}^{d_1} \left( (\beta^{-1})_{\eta, \zeta} (Q_{i+1,1}^\zeta - Q_{i,1}^\zeta) \right) - \Delta x \mathbf{b}^T \mathbf{A}^{(2)} (\partial_{q^\eta} V(\mathbf{Q}_i) + g_i^\eta). \quad (4.13)$$

Note that this rearrangement is possible since  $\beta$  operates on the index  $\eta$ , while  $\mathbf{b}$  and  $\mathbf{A}$  operate on the index  $j$  as given in the notation scheme.

Step 3:

Substituting  $\mathbf{P}_i^\eta$  from Eq. (4.10) into Eq. (4.8) and then substituting  $p_i^\eta$  from Eq. (4.13) into the resulting equation gives

$$\begin{aligned}
\mathbf{Q}_i^\eta &= Q_{i,1}^\eta \mathbf{1} - \Delta x \mathbf{A}^{(1)} \left( \sum_{\zeta=1}^{d_1} \beta_{\eta,\zeta} (\mathbf{P}_i^\zeta) \right) \\
&= Q_{i,1}^\eta \mathbf{1} - \Delta x \mathbf{A}^{(1)} \left( \sum_{\zeta=1}^{d_1} \beta_{\eta,\zeta} (p_i^\zeta \mathbf{1} + \Delta x \mathbf{A}^{(2)} (\partial_{q^\zeta} V(\mathbf{Q}_i) + g_i^\zeta)) \right) \\
&= Q_{i,1}^\eta \mathbf{1} - \Delta x \mathbf{A}^{(1)} \left( \sum_{\zeta=1}^{d_1} \beta_{\eta,\zeta} \left( \left[ -\frac{1}{\Delta x} \sum_{\xi=1}^{d_1} ((\beta^{-1})_{\zeta,\xi} (Q_{i,r}^\xi - Q_{i,1}^\xi)) \right. \right. \right. \\
&\quad \left. \left. \left. - \Delta x \mathbf{b}^T \mathbf{A}^{(2)} (\partial_{q^\zeta} V(\mathbf{Q}_i) + g_i^\zeta) \right] \mathbf{1} + \Delta x \mathbf{A}^{(2)} (\partial_{q^\zeta} V(\mathbf{Q}_i) + g_i^\zeta) \right) \right).
\end{aligned} \tag{4.14}$$

Rearranging and applying  $\beta^{-1}$  gives

$$\begin{aligned}
&\frac{1}{(\Delta x)^2} \sum_{\zeta=1}^{d_1} (\beta^{-1})_{\eta,\zeta} \left[ \mathbf{Q}_i^\zeta - Q_{i,1}^\zeta \mathbf{1} - \mathbf{A}^{(1)} (Q_{i,r}^\zeta - Q_{i,1}^\zeta) \mathbf{1} \right] \\
&= \mathbf{A}^{(1)} \left[ (\mathbf{b}^T \mathbf{A}^{(2)} (\partial_{q^\eta} V(\mathbf{Q}_i) + g_i^\eta)) \mathbf{1} - \mathbf{A}^{(2)} (\partial_{q^\eta} V(\mathbf{Q}_i) + g_i^\eta) \right], \\
&= \mathbf{A}^{(1)} (\mathbf{1b}^T - \mathbf{I}) \mathbf{A}^{(2)} (\partial_{q^\eta} V(\mathbf{Q}_i) + g_i^\eta).
\end{aligned} \tag{4.15}$$

Now, the first and last rows of the left-hand side of Eq. (4.15) are zero as are the first and last rows and columns of  $\mathbf{A}^{(1)} (\mathbf{1b}^T - \mathbf{I}) \mathbf{A}^{(2)}$ . Therefore, we denote rows 2 to  $r-1$  of  $\left[ \mathbf{Q}_i^\zeta - Q_{i,1}^\zeta \mathbf{1} - \mathbf{A}^{(1)} (Q_{i,r}^\zeta - Q_{i,1}^\zeta) \mathbf{1} \right]$  by  $\mathbf{d}_i^\zeta$ , the block of  $\mathbf{A}^{(1)} (\mathbf{1b}^T - \mathbf{I}) \mathbf{A}^{(2)}$  from (2,2) to  $(r-1, r-1)$  by  $\mathbf{C}$  and rows 2 to  $r-1$  of  $\partial_{q^\eta} V(\mathbf{Q}_i) + g_i^\eta$  by  $\mathbf{e}_i^\eta$ .

Then, noting that  $\mathbf{C}$  has full rank due to Eq. (3.5), we can write

$$\frac{1}{(\Delta x)^2} \sum_{\zeta=1}^{d_1} (\beta^{-1})_{\eta,\zeta} \mathbf{C}^{-1} \mathbf{d}_i^\zeta = \mathbf{e}_i^\eta. \tag{4.16}$$

Recalling the definition of  $g_i^\eta$ , Eq. (4.16) immediately allows us to write down explicit formulas for  $\partial_t Q_{i,k}^\eta$  in terms of  $\mathbf{Q}_i$  for  $1 < k < r$  and  $1 \leq \eta \leq \frac{1}{2}(d_1 - d_2)$  or  $\frac{1}{2}(d_1 + d_2) < \eta \leq d_1$  and for  $\partial_t^2 Q_{i,k}^\eta$  in terms of  $\mathbf{Q}_i$  for  $1 < k < r$  and  $\frac{1}{2}(d_1 - d_2) < \eta \leq \frac{1}{2}(d_1 + d_2)$ .

Step 4:

Substituting  $p_i^\eta$  from Eq. (4.13) into Eq. (4.11) for both  $p_i^\eta$  and  $p_{i+1}^\eta$  gives

$$\begin{aligned}
&-\frac{1}{(\Delta x)^2} \sum_{\zeta=1}^{d_1} (\beta^{-1})_{\eta,\zeta} (Q_{i+2,1}^\zeta - 2Q_{i+1,1}^\zeta + Q_{i,1}^\zeta) = \\
&\quad \mathbf{b}^T \mathbf{A}^{(2)} (\partial_{q^\eta} V(\mathbf{Q}_{i+1}) + g_{i+1}^\eta) + (\mathbf{b}^T - \mathbf{b}^T \mathbf{A}^{(2)}) (\partial_{q^\eta} V(\mathbf{Q}_i) + g_i^\eta)
\end{aligned} \tag{4.17}$$

for each  $\eta$ .

Of importance here, is that Eq. (4.17) does not involve the variables  $\partial_t Q_{i+1,r}^\eta$  or  $\partial_t^2 Q_{i+1,r}^\eta$  since the last entry of  $\mathbf{b}^T \mathbf{A}^{(2)}$  is zero. Neither does it involve the variables  $\partial_t Q_{i,1}^\eta$  or  $\partial_t^2 Q_{i,1}^\eta$  since the first entry of  $\mathbf{b}^T - \mathbf{b}^T \mathbf{A}^{(2)}$  is also zero.

Step 5:

Substituting the formulas for  $\partial_t Q_{i,k}^\eta$  and  $\partial_t^2 Q_{i,k}^\eta$  found in Step 3 into Eq. (4.17) and recalling that  $\partial_t Q_{i,r}^\eta = \partial_t Q_{i+1,1}^\eta$  and  $\partial_t^2 Q_{i,r}^\eta = \partial_t^2 Q_{i+1,1}^\eta$ , we can obtain explicit formulas for  $\partial_t Q_{i+1,1}^\eta$  in terms of  $\mathbf{Q}_i$  and  $\mathbf{Q}_{i+1}$  for  $1 \leq \eta \leq \frac{1}{2}(d_1 - d_2)$  and  $\frac{1}{2}(d_1 + d_2) < \eta \leq d_1$  and for  $\partial_t^2 Q_{i+1,1}^\eta$  in terms of  $\mathbf{Q}_i$  and  $\mathbf{Q}_{i+1}$  for  $\frac{1}{2}(d_1 - d_2) < \eta \leq \frac{1}{2}(d_1 + d_2)$ .

Thus, for each cell  $i$  in our grid, we have a system of explicit ODEs for either the first or second time derivatives of the stage variables  $\mathbf{Q}_i$  in terms of local values of  $\mathbf{Q}_i$ .  $\square$

While the conditions on  $\mathbf{K}$ ,  $\mathbf{L}$  and  $S(\mathbf{z})$  in the above theorem may at first appear restrictive, they allow several important equations such as the nonlinear wave and nonlinear Schrödinger equations. A notable exception is the Korteweg-de Vries equation for which  $S(\mathbf{z})$  is not separable. It is also worth noting that the conditions on  $\mathbf{K}$ ,  $\mathbf{L}$  and  $S(\mathbf{z})$  are the same as those required for the continuous system to be written as a system of PDEs in the variables  $\mathbf{q}$  and are similar to those required for a separable Hamiltonian system to be written as a system of second order ODEs.

The structure of  $\mathbf{K}$  is known as the ‘‘Darboux normal form’’ of  $\mathbf{K}$  and a change of coordinates will allow any skew-symmetric matrix to be written this way. If putting  $\mathbf{K}$  in Darboux normal form gives  $\mathbf{L}$  the following structure

$$\mathbf{L} = \begin{bmatrix} & & \Lambda \\ & 0_{d_2} & \\ -\Lambda^T & & \end{bmatrix} \quad (4.18)$$

for some  $d_1 \times d_1$  matrix  $\Lambda$  with  $|\Lambda| \neq 0$ , then the following change of coordinates in the  $\mathbf{p}$  variables can put  $\mathbf{L}$  in the form given in Eq. (4.1). Let  $\hat{\mathbf{p}} = \Lambda \mathbf{p}$  and  $\hat{T}(\hat{\mathbf{p}}) = T(\Lambda^{-1} \hat{\mathbf{p}}) = T(\mathbf{p})$ , then  $\nabla_{\hat{\mathbf{p}}} \hat{T}(\hat{\mathbf{p}}) = \Lambda \nabla_{\mathbf{p}} T(\mathbf{p}) = \Lambda \beta \mathbf{p} = \Lambda \beta \Lambda^{-1} \hat{\mathbf{p}}$  and  $S(\mathbf{z})$  still has the desired structure  $S(\mathbf{z}) = V(\mathbf{q}) + \frac{1}{2} \hat{\mathbf{p}}^T (\Lambda \beta \Lambda^{-1}) \hat{\mathbf{p}}$ .

The upper left  $(d_1 + d_2) \times (d_1 + d_2)$  block of  $\mathbf{L}$  being all zeros is fulfilled for PDEs which, when written as a first order system with  $\mathbf{K}$  in Darboux normal form, have no equations involving both a time and space derivative of the same variable; i.e.  $z_t^\eta + z_x^\eta = f(\mathbf{z})$  does not appear for any  $\eta$ .

**4.1. Examples.** Here we give several examples of common multi-Hamiltonian PDEs. For the PDEs that satisfy the requirements of Theorem 4.1 we give the ODEs that one obtains by applying the construction algorithm to those PDEs. For PDEs that do not satisfy the requirements of Theorem 4.1 we show why they fail and where the construction algorithm breaks down. We also give a PDE constructed so as to require the full use of Theorem 4.1.

**4.1.1. Nonlinear wave equation.** Our first example is the nonlinear wave equation,

$$u_{tt} = u_{xx} - V'(u), \quad (4.19)$$

which can be written as a multi-Hamiltonian PDE in the form of Eq. (4.2) with [4]

$$\mathbf{z} = \begin{bmatrix} u \\ v \\ w \end{bmatrix}, \quad \mathbf{K} = \begin{bmatrix} 0 & -1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} 0 & 0 & 1 \\ 0 & 0 & 0 \\ -1 & 0 & 0 \end{bmatrix} \quad (4.20)$$

and  $S(\mathbf{z}) = V(u) + \frac{1}{2}v^2 - \frac{1}{2}w^2$ .

Here,  $d_1 = d_2 = 1$  with  $\mathbf{z}^{(1)} = \{u, v\}$  and  $\mathbf{z}^{(2)} = \{w\}$ . We also have  $\alpha = -\beta = 1$ , thus we can see that  $\mathbf{K}$ ,  $\mathbf{L}$  and  $S(\mathbf{z})$  satisfy the requirements of Theorem 4.1. Upon eliminating the variable  $v$ , we obtain the PDE (4.6) with

$$\mathbf{z} = \begin{bmatrix} u \\ w \end{bmatrix}, \quad \mathbf{K} = \begin{bmatrix} 0 & 0 \\ 0 & 0 \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} 0 & 1 \\ -1 & 0 \end{bmatrix}, \quad \mathcal{E} = \begin{bmatrix} 1 & 0 \\ 0 & 0 \end{bmatrix} \quad (4.21)$$

and  $S = V(u) - \frac{1}{2}w^2$ .

Applying the construction algorithm for  $r = 2$  gives the following pair of ODEs for each cell  $i$ :

$$\begin{aligned} \partial_t^2 U_{i,1} &= \frac{1}{(\Delta x)^2} (U_{i-1,1} - 2U_{i,1} + U_{i+1,1}) - V'(U_{i,1}), \\ \partial_t^2 U_{i,2} &= \partial_t^2 U_{i+1,1}. \end{aligned} \quad (4.22)$$

Recalling from Step 1 that  $\mathbf{q}_i = \mathbf{Q}_{i,1}$  and noting that the last ODE is simply the first ODE of the next cell, it is convenient to drop the second ODE and rewrite the first ODE in terms of the node variable  $u_i$ :

$$\partial_t^2 u_i = \frac{1}{(\Delta x)^2} (u_{i-1} - 2u_i + u_{i+1}) - V'(u_i). \quad (4.23)$$

Applying the construction algorithm for  $r = 3$  gives the following triplet of ODEs for each cell  $i$ ,

$$\begin{aligned} \partial_t^2 U_{i,1} &= \frac{1}{(\Delta x)^2} (-U_{i-1,1} + 8U_{i-1,2} - 14U_{i,1} + 8U_{i,2} - U_{i+1,1}) - V'(U_{i,1}), \\ \partial_t^2 U_{i,2} &= \frac{1}{(\Delta x)^2} (4U_{i,1} - 8U_{i,2} + 4U_{i+1,1}) - V'(U_{i,2}), \\ \partial_t^2 U_{i,3} &= \partial_t^2 U_{i+1,1}, \end{aligned} \quad (4.24)$$

which cannot be written in terms of the node variables alone.

**4.1.2. NLS equation.** Our second example is the famous cubic-potential non-linear Schrödinger (NLS) equation,

$$i\psi_t + \psi_{xx} + 2|\psi|^2\psi = 0, \quad (4.25)$$

where  $\psi \in \mathbb{C}$ . Taking  $\psi = p + iq$  and separating the real and imaginary components of NLS allows the PDE to be written in the form of Eq. (4.2) with [13]

$$\mathbf{z} = \begin{bmatrix} p \\ q \\ v \\ w \end{bmatrix}, \quad \mathbf{K} = \begin{bmatrix} 0 & -1 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \\ -1 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 \end{bmatrix} \quad (4.26)$$

and  $S = -\frac{1}{2}(p^2 + q^2)^2 - \frac{1}{2}(v^2 + w^2)$ .

Here we have  $d_1 = 2$  and  $d_2 = 0$  with  $\mathbf{z}^{(1)} = \{p, q\}$  and  $\mathbf{z}^{(2)} = \{v, w\}$ .  $S(\mathbf{z})$  can be written as Eq. (4.3) with  $V(\mathbf{q}) = -\frac{1}{2}(p^2 + q^2)$  and  $T(\mathbf{p}) = \frac{1}{2}\mathbf{p}^T\beta\mathbf{p}$  where

$$\beta = \begin{bmatrix} -1 & 0 \\ 0 & -1 \end{bmatrix} \text{ and } \mathbf{p} = \begin{bmatrix} v \\ w \end{bmatrix}, \quad (4.27)$$

thus the NLS equation also satisfies the requirements of Theorem 4.1.

Applying the construction algorithm for an  $r$ -stage discretisation gives  $r$  ODEs for each element of  $\mathbf{z}^{(1)}$  at cell  $i$ . As with the nonlinear wave equation, if we use the 2-stage discretisation then for each element of  $\mathbf{z}^{(1)}$  at cell  $i$  we can drop the ODE for the second stage variable and write the ODE for the first stage variable in terms of the node variables. The resulting ODEs are

$$\begin{aligned} \partial_t p_i &= -\frac{1}{(\Delta x)^2}(q_{i-1} - 2q_i + q_{i+1}) - 2(p_i^2 + q_i^2)q_i, \\ \partial_t q_i &= \frac{1}{(\Delta x)^2}(p_{i-1} - 2p_i + p_{i+1}) + 2(p_i^2 + q_i^2)p_i. \end{aligned} \quad (4.28)$$

These are precisely the ODEs one obtains by applying second order finite differences in space to Eq. (4.25). The same statement applies for other PDEs that satisfy the conditions of Theorem 4.1, thus we note that 2-stage Lobatto IIIA–IIIB discretisation in space is equivalent to second order finite differences in space up to second order differences when applied to such a PDE.

For  $r = 3$  we obtain a triplet of ODEs for each element of  $\mathbf{z}^{(1)}$  at cell  $i$ :

$$\begin{aligned} \partial_t P_{i,1} &= -\frac{1}{(\Delta x)^2}(-Q_{i-1,1} + 8Q_{i-1,2} - 14Q_{i,1} + 8Q_{i,2} - Q_{i+1,1}) \\ &\quad - 2(P_{i,1}^2 + Q_{i,1}^2)Q_{i,1}, \\ \partial_t P_{i,2} &= -\frac{1}{(\Delta x)^2}(4Q_{i,1} - 8Q_{i,2} + 4Q_{i+1,1}) - 2(P_{i,2}^2 + Q_{i,2}^2)Q_{i,2}, \\ \partial_t P_{i,3} &= \partial_t P_{i+1,1}, \\ \partial_t Q_{i,1} &= \frac{1}{(\Delta x)^2}(-P_{i-1,1} + 8P_{i-1,2} - 14P_{i,1} + 8P_{i,2} - P_{i+1,1}) \\ &\quad + 2(P_{i,1}^2 + Q_{i,1}^2)P_{i,1}, \\ \partial_t Q_{i,2} &= \frac{1}{(\Delta x)^2}(4P_{i,1} - 8P_{i,2} + 4P_{i+1,1}) + 2(P_{i,2}^2 + Q_{i,2}^2)P_{i,2}, \\ \partial_t Q_{i,3} &= \partial_t Q_{i+1,1}. \end{aligned} \quad (4.29)$$

**4.1.3. Boussinesq equation.** Our third example is the Boussinesq equation,

$$p_{tt} = (\varepsilon p_{xx} + V'(p))_{xx}, \quad (4.30)$$

which, when written as a multi-Hamiltonian PDE, shares the same  $\mathbf{z}$ ,  $\mathbf{z}^{(1)}$ ,  $\mathbf{z}^{(2)}$ ,  $\mathbf{K}$  and  $\mathbf{L}$  as the NLS equation above [8]. The only difference is the function  $S(\mathbf{z})$  which is given by  $S(\mathbf{z}) = -V(p) + \frac{1}{2}(w^2 - \frac{1}{\varepsilon}v^2)$ .

As before, the requirements of Theorem 4.1 are satisfied and applying the construction algorithm gives  $r$  ODEs for each element of  $\mathbf{z}^{(1)}$  at cell  $i$ . For  $r = 2$ , we

once again drop the ODEs for the second stage variables and write the first ODEs in terms of the node variables as

$$\begin{aligned}\partial_t p_i &= \frac{1}{(\Delta x)^2}(q_{i-1} - 2q_i + q_{i+1}), \\ \partial_t q_i &= \frac{\varepsilon}{(\Delta x)^2}(p_{i-1} - 2p_i + p_{i+1}) + V'(p).\end{aligned}\tag{4.31}$$

For  $r = 3$  we get

$$\begin{aligned}\partial_t P_{i,1} &= \frac{1}{(\Delta x)^2}(-Q_{i-1,1} + 8Q_{i-1,2} - 14Q_{i,1} + 8Q_{i,2} - Q_{i+1,1}), \\ \partial_t P_{i,2} &= \frac{1}{(\Delta x)^2}(4Q_{i,1} - 8Q_{i,2} + 4Q_{i+1,1}), \\ \partial_t P_{i,3} &= \partial_t P_{i+1,1}, \\ \partial_t Q_{i,1} &= \frac{\varepsilon}{(\Delta x)^2}(-P_{i-1,1} + 8P_{i-1,2} - 14P_{i,1} + 8P_{i,2} - P_{i+1,1}) + V'(P_{i,1}), \\ \partial_t Q_{i,2} &= \frac{\varepsilon}{(\Delta x)^2}(4P_{i,1} - 8P_{i,2} + 4P_{i+1,1}) + V'(P_{i,2}), \\ \partial_t Q_{i,3} &= \partial_t Q_{i+1,1}.\end{aligned}\tag{4.32}$$

**4.1.4. Korteweg-de Vries (KdV) equation.** Our fourth example is the KdV equation,

$$u_t = V'(u)_x + \nu u_{xxx}\tag{4.33}$$

and can be written in the form of Eq. (4.2) with [5]

$$\mathbf{z} = \begin{bmatrix} u \\ \phi \\ v \\ w \end{bmatrix}, \quad \mathbf{K} = \begin{bmatrix} 0 & -1 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \\ -1 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 \end{bmatrix}\tag{4.34}$$

and with  $S(\mathbf{z}) = -\frac{1}{2}uw - V(u) - \frac{1}{2\nu}v^2$ . Here,  $d_1 = 2$ ,  $d_2 = 0$  and  $\mathbf{z}$  is partitioned into  $\mathbf{z}^{(1)} = \{u, \phi\}$  and  $\mathbf{z}^{(2)} = \{v, w\}$ .

While the  $\mathbf{K}$  and  $\mathbf{L}$  matrices have the required structure for Theorem 4.1, the function  $S(\mathbf{z})$  does not. Specifically, the  $-uw$  term in  $S(\mathbf{z})$  prevents us from writing  $T(\mathbf{p}) = \frac{1}{2}\mathbf{p}^T \beta \mathbf{p}$  and so Step 2 of the construction algorithm cannot be carried out.

For example, discretising the KdV equation with two-stage Lobatto IIIA–IIIB gives

$$\begin{aligned}v_{i+\frac{1}{2}} &= v_{i-\frac{1}{2}} + \Delta x(\partial_t \phi_i - V'(u_i) - \frac{1}{4}(w_{i+\frac{1}{2}} + w_{i-\frac{1}{2}})), \\ w_{i+\frac{1}{2}} &= w_{i-\frac{1}{2}} - \Delta x \partial_t u_i, \\ -u_{i+1} &= -u_i - \Delta x \frac{1}{\nu} v_{i+\frac{1}{2}}, \\ -\phi_{i+1} &= -\phi_i - \Delta x \frac{1}{4}(u_i + u_{i+1}),\end{aligned}\tag{4.35}$$

where  $u_i = U_{i,1}$ ,  $u_{i+1} = U_{i,2}$ ,  $\phi_i = \Phi_{i,1}$ ,  $\phi_{i+1} = \Phi_{i,2}$ ,  $v_{i+\frac{1}{2}} = V_{i,1} = V_{i,2}$  and  $w_{i+\frac{1}{2}} = W_{i,1} = W_{i,2}$ .

Introducing the operators  $D$  and  $M$ , where  $Du_i = \frac{1}{\Delta x}(u_{i+1} - u_i)$  and  $Mu_i = \frac{1}{2}(u_{i+1} + u_i)$ , allows us to write this system as

$$\begin{aligned} Dv_{i-\frac{1}{2}} &= \partial_t \phi_i - V'(u_i) - \frac{1}{2}Mw_{i-\frac{1}{2}}, \\ Dw_{i-\frac{1}{2}} &= -\partial_t u_i, \\ -Du_i &= -\frac{1}{\nu}v_{i+\frac{1}{2}}, \\ -D\phi_i &= -\frac{1}{2}Mu_i. \end{aligned} \tag{4.36}$$

Eliminating all the variables other than the original variable  $u$  gives the implicit ODE

$$M\partial_t u_i = DV'(u_i) + \nu D^3 u_{i-1}. \tag{4.37}$$

In general,  $M$  is not invertible, thus further conditions are required (e.g. periodic boundary conditions with an odd number of grid points) to form a well defined integrator from this implicit ODE.

This is none other than the *narrow box scheme*, introduced in [2] and derived as a finite volume scheme (and shown to be more accurate than the box scheme) in [3]. Thus, we have shown that the narrow box scheme is multisymplectic.

**4.1.5. Camassa-Holm (CH) equation.** Our fifth example is the CH equation,

$$u_t - u_{xxt} = V'(u)_x + \alpha u_{xxx}. \tag{4.38}$$

This equation can be written in the form of Eq. (1.1) with  $\mathbf{z} = [u, \nu, \phi, w, \rho, v]^T$ ,

$$\mathbf{K} = \begin{bmatrix} 0 & \frac{1}{2} & -\frac{1}{2} & 0 & 0 & 0 \\ -\frac{1}{2} & 0 & 0 & 0 & 0 & 0 \\ \frac{1}{2} & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} 0 & 0 & 0 & 0 & \frac{1}{2} & \alpha \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ -\frac{1}{2} & 0 & 0 & 0 & 0 & 0 \\ -\alpha & 0 & 0 & 0 & 0 & 0 \end{bmatrix} \tag{4.39}$$

and  $S(\mathbf{z}) = uw - V(u) - \frac{\alpha}{2}v^2 - \frac{1}{2}\nu\rho$ . To the best of our knowledge, this is the first occurrence of the multi-Hamiltonian structure of the CH equation in the literature.

Putting  $\mathbf{K}$  into its Darboux normal form puts  $\mathbf{L}$  into the form

$$\mathbf{L} = \begin{bmatrix} 0_3 & \Lambda \\ -\Lambda^T & 0_3 \end{bmatrix} \tag{4.40}$$

where  $\Lambda$  is a  $3 \times 3$  matrix with  $\text{rank}(\Lambda) = 2$ . Thus, we cannot write  $\mathbf{L}$  in the form of Eq. (4.1) and so the Camassa-Holm equation does not satisfy the requirements of Theorem 4.1.

However, partitioning  $\mathbf{z}$  into  $\mathbf{z}^{(1)} = \{u, \nu, \phi\}$  and  $\mathbf{z}^{(2)} = \{w, \rho, v\}$ , then discretising



the CH equation with two-stage Lobatto IIIA–IIIB using the  $D$  and  $M$  notation gives

$$\begin{aligned}
\frac{1}{2}D\rho_{i-\frac{1}{2}} + \alpha Dv_{i-\frac{1}{2}} &= Mw_{i-\frac{1}{2}} - V'(u_i) - \frac{1}{2}\partial_t \nu_i + \frac{1}{2}\partial_t \phi_i, \\
0 &= -\frac{1}{2}M\rho_{i-\frac{1}{2}} + \frac{1}{2}\partial_t u_i, \\
-Dw_{i-\frac{1}{2}} &= -\frac{1}{2}\partial_t u_i, \\
D\phi_i &= Mu_i, \\
-\frac{1}{2}Du_i &= -\frac{1}{2}M\nu_i, \\
-\alpha Du_i &= -\alpha v_{i-\frac{1}{2}}.
\end{aligned} \tag{4.41}$$

Eliminating  $\nu$ ,  $\phi$ ,  $w$ ,  $\rho$  and  $v$  gives the implicit ODE

$$(M^2 - D^2)\partial_t u_i = MDV'(u_i) + \alpha MD^3 u_{i-1}. \tag{4.42}$$

As with the KdV equation, the operator on the left hand side cannot be locally inverted, although it is at least typically invertible.

**4.1.6. Benjamin-Bona-Mahony (BBM) equation.** Our sixth example is the BBM equation,

$$u_t - \alpha u_{xxt} = V'(u)_x. \tag{4.43}$$

Writing this equation in the form of Eq. (1.1) with  $\mathbf{z} = [u, \nu, \phi, w, \rho]^T$  gives

$$\mathbf{K} = \begin{bmatrix} 0 & \frac{\alpha}{2} & -\frac{1}{2} & 0 & 0 \\ -\frac{\alpha}{2} & 0 & 0 & 0 & 0 \\ \frac{1}{2} & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} 0 & 0 & 0 & 0 & \frac{\alpha}{2} \\ 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 & 0 \\ 0 & 0 & 1 & 0 & 0 \\ -\frac{\alpha}{2} & 0 & 0 & 0 & 0 \end{bmatrix} \tag{4.44}$$

and  $S(\mathbf{z}) = uw - V(u) - \frac{\alpha}{2}\nu\rho$ . The multi-Hamiltonian structure of the BBM equation presented here also appears to be the first such occurrence in the literature.

As with the Camassa-Holm equation in the previous example, putting  $\mathbf{K}$  into its Darboux normal form results in an  $\mathbf{L}$  which cannot be written in the form of Eq. (4.1) and so the BBM equation does not satisfy the requirements of Theorem 4.1.

Partitioning  $\mathbf{z}$  into  $\mathbf{z}^{(1)} = \{u, \nu, \phi\}$  and  $\mathbf{z}^{(2)} = \{w, \rho\}$ , then discretising the BBM equation with two-stage Lobatto IIIA–IIIB using the  $D$  and  $M$  notation gives

$$\begin{aligned}
\frac{\alpha}{2}D\rho_{i-\frac{1}{2}} &= Mw_{i-\frac{1}{2}} - V'(u_i) - \frac{\alpha}{2}\partial_t \nu_i + \frac{1}{2}\partial_t \phi_i, \\
0 &= -\frac{\alpha}{2}M\rho_{i-\frac{1}{2}} + \frac{\alpha}{2}\partial_t u_i, \\
-Dw_{i-\frac{1}{2}} &= -\frac{1}{2}\partial_t u_i, \\
D\phi_i &= Mu_i, \\
-\frac{\alpha}{2}Du_i &= -\frac{\alpha}{2}M\nu_i.
\end{aligned} \tag{4.45}$$

Eliminating  $\nu$ ,  $\phi$ ,  $w$  and  $\rho$  gives the implicit ODE

$$(M^2 - \alpha D^2)\partial_t u_i = MDV'(u_i). \tag{4.46}$$

**4.1.7. A made up example.** Our last example is contrived to satisfy the requirements of Theorem 4.1 and demonstrates the case when  $d_2 \neq d_1$  and  $d_2 \neq 0$ . We have chosen  $d_1 = 3$ ,  $d_2 = 1$  and a multi-Hamiltonian PDE (1.1) with  $\mathbf{z} = [q^1, q^2, q^3, v, p^1, p^2, p^3]^T$ ,

$$\mathbf{K} = \begin{bmatrix} 0 & 0 & -1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} 0 & 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ -1 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & -1 & 0 & 0 & 0 & 0 \end{bmatrix} \quad (4.47)$$

and  $S(\mathbf{z}) = V(\mathbf{q}) + \frac{1}{2}\mathbf{p}^T\beta\mathbf{p} + \frac{\alpha}{2}v^2$ , where  $\alpha$  is a constant and

$$\beta = \begin{bmatrix} 1 & 1 & -\frac{1}{2} \\ 1 & 1 & 0 \\ -\frac{1}{2} & 0 & 1 \end{bmatrix}. \quad (4.48)$$

This corresponds to the PDE

$$\begin{aligned} \partial_t q^1 &= -2q_{xx}^1 + 2q_{xx}^2 + \partial_{q^3}V(\mathbf{q}) \\ \frac{1}{\alpha}\partial_t^2 q^2 &= -4q_{xx}^1 + 3q_{xx}^2 - 2q_{xx}^3 - \partial_{q^2}V(\mathbf{q}) \\ \partial_t q^3 &= 4q_{xx}^1 - 4q_{xx}^2 + 2q_{xx}^3 - \partial_{q^1}V(\mathbf{q}) \end{aligned} \quad (4.49)$$

Eliminating the variable  $v$  in favour of higher derivatives in time of  $q^2$  gives Eq. (4.6) with

$$\mathbf{z} = \begin{bmatrix} q^1 \\ q^2 \\ q^3 \\ p^1 \\ p^2 \\ p^3 \end{bmatrix}, \quad \mathbf{K} = \begin{bmatrix} 0 & 0 & -1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \end{bmatrix}, \quad \mathbf{L} = \begin{bmatrix} 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 \\ -1 & 0 & 0 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 & 0 & 0 \\ 0 & 0 & -1 & 0 & 0 & 0 \end{bmatrix}, \quad (4.50)$$

$S(\mathbf{z}) = V(\mathbf{q}) + \frac{1}{2}\mathbf{p}^T\beta\mathbf{p}$ , and the only non-zero entry of  $\mathcal{E}$  given by  $\mathcal{E}_{2,2} = \frac{1}{\alpha}$ .

If we applying the construction algorithm for  $r = 2$  then once again we can drop the ODEs for the second stage variables and write the ODEs for the first stage variables in terms of the node variables giving the following ODEs at cell  $i$ :

$$\begin{aligned} \partial_t q_i^1 &= \frac{1}{(\Delta x)^2}(-2q_{i-1}^1 + 2q_{i-1}^2 + 4q_i^1 - 4q_i^2 - 2q_{i+1}^1 + 2q_{i+1}^2) + \partial_{q^3}V(\mathbf{q}_i), \\ \partial_t^2 q_i^2 &= \frac{\alpha}{(\Delta x)^2}(-4q_{i-1}^1 + 3q_{i-1}^2 - 2q_{i-1}^3 + 8q_i^1 - 6q_i^2 + 4q_i^3 - 4q_{i+1}^1 + 3q_{i+1}^2 - 2q_{i+1}^3) \\ &\quad - \alpha\partial_{q^2}V(\mathbf{q}_i), \\ \partial_t q_i^3 &= \frac{1}{(\Delta x)^2}(4q_{i-1}^1 - 4q_{i-1}^2 + 2q_{i-1}^3 - 8q_i^1 + 8q_i^2 - 4q_i^3 \\ &\quad + 4q_{i+1}^1 - 4q_{i+1}^2 + 2q_{i+1}^3) - \partial_{q^1}V(\mathbf{q}_i). \end{aligned} \quad (4.51)$$

**5. Discussion.** We would like to point out that the discretisation in space by Lobatto IIIA–IIIB in the above examples only modifies the linear component of the multi-Hamiltonian PDE, i.e. the discrete approximation of  $\mathbf{Lz}_x$ . The reason for this is that throughout the construction algorithm, the nonlinear components of the multi-Hamiltonian PDE always appear coupled to the time derivatives as the expression  $\partial_q^\eta V(\mathbf{Q}_i) + g_i^\eta$ .

Furthermore, we note that in the examples above, the same pattern of coefficients arises from discretising different PDEs with the same order Lobatto IIIA–IIIB discretisation. For example, with  $r = 2$  the coefficients in the approximation of  $\mathbf{q}_{xx}$  have a weighting proportional to  $[1, -2, 1]$ , while for  $r = 3$  these coefficients are proportional to  $[-1, 8, -14, 8, -1]$  for the first ODE and  $[4, -8, 4]$  for the second ODE. This behaviour continues for higher values of  $r$ , e.g. for  $r = 4$  the approximation of  $\mathbf{q}_{xx}$  in the first ODE has coefficients proportional to  $[1, \frac{1}{2}(25 - 15\sqrt{5}), \frac{1}{2}(25 + 15\sqrt{5}), -52, \frac{1}{2}(25 + 15\sqrt{5}), \frac{1}{2}(25 - 15\sqrt{5}), 1]$ , the second ODE has coefficients proportional to  $[5 + 3\sqrt{5}, -20, 10, 5 - 3\sqrt{5}]$  and the third ODE has coefficients proportional to  $[5 - 3\sqrt{5}, 10, -20, 5 + 3\sqrt{5}]$ . For higher values of  $r$  these patterns of the coefficients in the approximation of  $\mathbf{q}_{xx}$  become increasingly complicated, yet for a given value of  $r$ , these patterns remain the same regardless of the PDE under consideration.

The reason these patterns of coefficients occur for different PDEs is due to Eqs. (4.16) and (4.17). For a given value of  $r$ ,  $\mathbf{C}$  and  $\mathbf{d}_i^\zeta$  are fixed regardless of the PDE. Similarly, the coefficients  $\mathbf{b}^T \mathbf{A}^{(2)}$  and  $\mathbf{b}^T - \mathbf{b}^T \mathbf{A}^{(2)}$  in Eq. (4.17) are completely determined by  $r$ . Thus, when solving Eqs. (4.16) and (4.17) for  $g_i^\eta$ , the same weighting of the nearby stage variables occurs for  $\mathbf{q}_{xx}$  for different PDEs.

For an  $r$  stage discretisation, the approximation to  $\mathbf{q}_{xx}$  at stage  $j$  is given by

$$-\frac{1}{(\Delta x)^2} (\mathbf{C}^{-1} \mathbf{d}_i^\zeta)_{j-1} = \frac{1}{(\Delta x)^2} \sum_{k=2}^{r-1} (\mathbf{C}^{-1})_{j-1, k-1} ((1 - c_k) Q_{i,1}^\zeta - Q_{i,k}^\zeta + c_k Q_{i,r}^\zeta) \quad (5.1)$$

for  $2 \leq j \leq r - 1$  and  $1 \leq \eta \leq r$ , where  $\mathbf{C}_{i-1, k-1}$  and  $c_k$  are given by Eqs. (3.5) and (3.2) respectively, and by

$$\frac{1}{2b_1(\Delta x)^2} \left( \sum_{k=2}^{r-1} \left( (\mathbf{b}^T \mathbf{A}^{(2)})_k (\mathbf{C}^{-1} \mathbf{d}_i^\zeta)_{k-1} + (\mathbf{b}^T - \mathbf{b}^T \mathbf{A}^{(2)})_k (\mathbf{C}^{-1} \mathbf{d}_{i+1}^\zeta)_{k-1} \right) + (Q_{i+2,1}^\zeta - 2Q_{i+1,1}^\zeta + Q_{i,1}^\zeta) \right), \quad (5.2)$$

for  $j = 1$ , where  $b_1$  is the first entry in  $\mathbf{b}$ .

This suggests the following shortcut:

1. Write the PDE with only terms of the form  $\mathbf{z}_{xx}$  (no  $\mathbf{z}_x$ ).
2. Replace the  $\mathbf{z}_{xx}$  terms with the PRK finite differences of the desired order.

Now, the system of ODEs that one obtains from applying Theorem 4.1 to an appropriate PDE can be written as a Hamiltonian system; e.g. for the Boussinesq equation and  $r = 2$ , the system of ODEs at node  $i$  can be written as

$$\partial_t \mathbf{z}_i = \mathbf{J}^{-1} \nabla_{\mathbf{z}_i} H_i \quad (5.3)$$

where

$$\mathbf{z}_i = \begin{bmatrix} q_i \\ p_i \end{bmatrix}, \quad \mathbf{J}^{-1} = \begin{bmatrix} 0 & 1 \\ -1 & 0 \end{bmatrix} \quad (5.4)$$

and

$$H_i = \frac{1}{(\Delta x)^2}(-q_{i-1}q_i + q_i^2 - q_iq_{i+1} + \varepsilon p_{i-1}p_i - \varepsilon p_i^2 + \varepsilon p_i p_{i+1}) + V(p_i). \quad (5.5)$$

If the nonlinear terms in such a Hamiltonian system are separable, then one can apply an explicit symplectic PRK discretisation in time to obtain an explicit (and hence well defined) high-order local multisymplectic integrator. If the nonlinear terms are not separable then other explicit time integrators may be applied, e.g. symplectic splitting methods [19], which may give superior performance (in terms of speed and stability) over implicit integrators. Even if no explicit time integrator can be applied to the Hamiltonian system, there may be some benefits to having a spatial discretisation that gives rise to explicit ODEs, e.g. the ODEs may be less stiff than those obtained from an implicit discretisation.

In the examples in the previous section, the systems of ODEs arising from the nonlinear wave equation and the Boussinesq equation both have separable Hamiltonians and thus allow for a high-order explicit symplectic PRK discretisation to be applied in time. The NLS equation is not so fortunate, however its non-linearity is only quadratic and thus for an  $r$ -stage Lobatto IIIA–IIIB discretisation in time it is necessary to solve a system of  $r - 1$  coupled quadratic equations for each update of  $\mathbf{P}_i$  or  $\mathbf{Q}_i$ . For  $r = 2$ , this quadratic equation can be solved explicitly (in particular, the same root of the quadratic is always taken) and an explicit (and hence well defined), local, high-order in space, multisymplectic integrator can be formed.

Another point that we would like to make is about how the ODEs that one obtains from our construction algorithm handle boundary conditions. Many other discretisation schemes (e.g. implicit midpoint, higher order Gaussian Runge–Kutta) either do not remain well defined or they require extra conditions to be so [2, 18]. However, our ODEs remain well defined under periodic, Dirichlet and Neumann boundary conditions without any further restrictions. For example, 3-stage Lobatto IIIA–IIIB applied to the NLS equation with Neumann boundary conditions,  $\psi_x = 0$ , applied to the left boundary as  $v_1 = w_1 = 0$  leads to the following ODEs:

$$\begin{aligned} \partial_t P_{1,1} &= -\frac{1}{(\Delta x)^2}(-14Q_{1,1} + 16Q_{1,2} - 2Q_{2,1}) - 2(P_{1,1}^2 + Q_{1,1}^2)Q_{1,1}, \\ \partial_t Q_{1,1} &= \frac{1}{(\Delta x)^2}(-14P_{1,1} + 16P_{1,2} - 2P_{2,1}) + 2(P_{1,1}^2 + Q_{1,1}^2)P_{1,1}, \end{aligned} \quad (5.6)$$

which are equivalent to the first and fourth lines of Eq. (4.29), where the points outside the domain are treated as phantom points, i.e.  $Q_{0,1} = Q_{2,1}$  and  $Q_{0,2} = Q_{1,2}$ .

Finally, we would like to point out that although Theorem 4.1 is stated for the Lobatto IIIA–IIIB class of PRK discretisations, it applies equally well to any PRK discretisation satisfying Eqs. (3.3), (3.4) and (3.5). We leave it as an open question as to whether there are any other PRK discretisations that satisfy Eqs. (3.3), (3.4) and (3.5).

In this paper we have deliberately restricted our attention to the structural properties of PRK discretisation, namely its multisymplecticity and explicitness. Results on its dynamical properties, such as order and dispersion, will be reported elsewhere.

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