

# **A new version of transfer matrix method for multibody systems**

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**Abstract** In order to avoid the global dynamics equations and increase the computational efficiency for multibody system dynamics (MSD), the transfer matrix method of multibody system (MSTMM) has been developed and applied very widely in research and engineering in recent 20 years. It differs from ordinary methods in multibody system dynamics with respect to the feature that there is no need for a global dynamics equation, and it uses loworder matrices for high computational efficiency. For linear systems, MSTMM is exact even if continuous elements like beams are involved. The discrete time MSTMM, however, has to use local linearization. In order to release the method from such approximations, a new version of MSTMM is presented in this paper where translational and angular accelerations, on the one hand, and internal forces and moments, on the other hand, are used as state variables. Already linear relationships among these quantities are utilized, which results in new element transfer matrices and algorithms making the study of multibody systems as simple as the study of single bodies. The proposed approach also allows combining MSTMM with any general numerical integration procedure. Some numerical examples of MSD are given to demonstrate the proposed method.

**Keywords** Multibody system dynamics · Transfer matrix method · Transfer equation · Computational efficiency · Computational precision

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# **1 Introduction**

Multibody system dynamics methods (MSDM) have developed rapidly in the last 50 years [[1](#page-19-0)[–3](#page-19-1)] and provide powerful tools for studying dynamics of various mechanical systems. Although the various ordinary MSDMs have different styles, at the same time almost all of them share the following characteristics: firstly, it is necessary to develop the global dynamics equations of the system, and they have to be deduced again if the system's topological structure is changed; secondly, the order of the global dynamics equations is not less than the number of degrees of freedom of the system, which may become very high for complex systems resulting in rather long computational time.

The approach proposed in the paper is different from this and has a different origin. In 1986, Kumar and Sankar developed a discrete time transfer matrix method (DTTMM) for structural dynamics of time-variant systems by combining the transfer matrix method with a numerical integration procedure [\[4\]](#page-19-2). In 1989, Rui and others extended the transfer matrix method to multibody systems (MSTMM) for vibration analysis of linear multi-rigidflexible-body systems by developing new transfer matrices [[5](#page-19-3)], where eigenvalues of linear multi-rigid-flexible-body systems can be computed easily and with high precision [\[6](#page-19-4)]. For general nonlinear multi-rigid-body and multi-rigid-flexible-body systems, the discrete time transfer matrix method of multibody systems (MSDTTMM) was developed by combining MSTMM with a numerical integration procedure  $[7–10]$  $[7–10]$  $[7–10]$ . In this case, however, linearization is mandatory to deduce the transfer equation of elements, which essentially describes the governing equation of motion represented by a linear relationship between the state vectors of inboard and outboard ends of an element. In [\[7](#page-19-5)], the Newmark-Beta numerical integration procedure is introduced, leading to only second-order computational precision. Although a higher order numerical integration procedure could be adopted, the deduction process is rather tedious.

This can be avoided with a new version of MSTMM presented in this paper. Translational and angular accelerations, together with internal forces and moments, are taken as new state variables instead of position coordinates as in the original MSDTTMM. This results in totally different transfer matrices of elements and algorithms compared with the original ones described in  $[7-10]$  $[7-10]$  $[7-10]$ . The proposed method expands the advantages of MSTMM by allowing more sophisticated numerical integration procedures such as any Runge–Kutta method to be used. Global dynamics equations of the system are still avoided. Instead, involved matrices have low order and the setup of the global transfer equation is highly programmable. The proposed method is simple, straightforward, efficient, practical, and provides a powerful tool for MSD. Numerical examples in Sect. [4](#page-9-0) will show good agreement with simulation results obtained by an ordinary dynamics method.

### **2 General theorems and steps of the new version of MSTMM**

Any complex multibody system may be divided into various elements including bodies (rigid bodies, elastic bodies, lumped masses, etc.) and hinges (joints, ball-and-socket, pins, linear and rotary springs, linear and rotary dampers, etc.), which are connected at intermediate points. The general idea of MSTMM is to use kinematic and kinetic quantities at these connection points as state variables which are related by the characteristics of the element in between. The characteristics are described by linear transfer equations from one connection point to the other, which are called input and output points *I* and *O* of the element, see, e.g., Fig. [1](#page-2-0). For a specific multibody system, the element transfer equations are combined

<span id="page-2-0"></span>

<span id="page-2-1"></span>by simple matrix operations to an overall transfer equation which may be solved for given boundary states at the endpoints of the system. These steps will be explained in more detail in the following.

#### **2.1 Coordinate systems and sign convention**

<span id="page-2-2"></span>In order to describe the position and orientation of connecting points, we assign a righthanded Cartesian reference frame *K* to each of them, see, e.g., input point *I* in Fig. [1.](#page-2-0) Its spatial position in the inertial Cartesian frame  $\{oxyz\}$  is then given by coordinates x, y, and *z*, and its orientation by three space-angles  $\theta_1$ ,  $\theta_2$ ,  $\theta_3$  [[10](#page-19-6), [12](#page-19-7)] which describe successive rotations about the inertial *x*-, *y*-, and *z*-axes. The latter define the direction cosine matrix *A* for coordinate transformations from *K* to  $\{oxyz\}$ , which can be built from the elementary rotation matrices  $A_{x,\theta_1}$ ,  $A_{y,\theta_2}$ ,  $A_{z,\theta_3}$  as

$$
A = A_{z,\theta_3} A_{y,\theta_2} A_{x,\theta_1} = \begin{bmatrix} c_2 c_3 & s_1 s_2 c_3 - c_1 s_3 & c_1 s_2 c_3 + s_1 s_3 \\ c_2 s_3 & s_1 s_2 s_3 + c_1 c_3 & c_1 s_2 s_3 - s_1 c_3 \\ -s_2 & s_1 c_2 & c_1 c_2 \end{bmatrix}
$$
 (1)

where  $s_i = \sin \theta_i$ ,  $c_i = \cos \theta_i$ ,  $i = 1, 2, 3$ .

By differentiating *A* with respect to time, we may deduce the skew-symmetric matrix of the angular velocity vector  $\mathbf{\Omega} = [\Omega_x, \Omega_y, \Omega_z]^T$  of frame *K* about the inertial frame *oxyz* as  $\tilde{\mathbf{\Omega}} = \dot{\mathbf{A}} \mathbf{A}^{\mathrm{T}}$ , which can be rearranged as  $\mathbf{\Omega} = \mathbf{A} \mathbf{H} \dot{\mathbf{\theta}}$ . By further differentiation we find [\[12\]](#page-19-7)

$$
\ddot{\theta} = H^{-1}A^{\mathrm{T}}\dot{\Omega} - H^{-1}\dot{H}\dot{\theta}
$$

where

$$
\boldsymbol{H} = \begin{bmatrix} 1 & 0 & -s_2 \\ 0 & c_1 & s_1 c_2 \\ 0 & -s_1 & c_1 c_2 \end{bmatrix}, \qquad \boldsymbol{\dot{H}} = \begin{bmatrix} 0 & 0 & -c_2 \dot{\theta}_2 \\ 0 & -s_1 \dot{\theta}_1 & c_1 c_2 \dot{\theta}_1 - s_1 s_2 \dot{\theta}_2 \\ 0 & -c_1 \dot{\theta}_1 & -s_1 c_2 \dot{\theta}_1 - c_1 s_2 \dot{\theta}_2 \end{bmatrix},
$$

and  $\theta = [\theta_1, \theta_2, \theta_3]^T$ . It should be pointed out that any kind of Euler/Bryan angle formulation has its inherent singularity, which is why the angular acceleration  $\dot{\mathbf{z}}$  rather than  $\ddot{\theta}$  is used in the state vector hereinafter. Once approaching a singular configuration, one can resort to another kind of Euler/Bryan angle to avoid the singularity problem. Alternatively, many other methods, such as Rodriguez parameters, Euler parameters, etc., can be also adopted to describe the orientation and avoid singular orientation. If not mentioned explicitly, all vectors are described w.r.t. the inertial frame.

In contrast to the sign conventions of the original MSDTTMM [\[10\]](#page-19-6), the following definition is used: positive translational and angular accelerations coincide with positive directions of the coordinate system; positive directions of inboard forces and outboard torques acting on the element coincide with positive directions, whereas positive outboard forces and inboard torques coincide with negative directions of the coordinate system.

#### **2.2 New state vectors, transfer equations and transfer matrices of elements**

In the original MSDTTMM [\[10\]](#page-19-6), the position and orientation coordinates of the connection points have been used as state variables. For example, for spatial motion the state vector was summarized from coordinates of the connecting point, rotation angles, internal torques and forces, and an artificial one as

$$
z = \begin{bmatrix} x & y & z & \theta_x & \theta_y & \theta_z & m_x & m_y & m_z & q_x & q_y & q_z & 1 \end{bmatrix}^T. \tag{2}
$$

<span id="page-3-0"></span>However, since kinematics typically is nonlinear, linearization had to be involved in the derivation of the transfer equations and also an integration scheme had to be part of the final transfer relations.

<span id="page-3-3"></span>In order to avoid all these drawbacks, the proposed approach combines kinematics and kinetics equations on the acceleration level. Therefore, the new state vectors summarize acceleration variables of connection points and forces and moments between any two multibody system elements moving in space:

$$
z = \begin{bmatrix} \ddot{x} & \ddot{y} & \ddot{z} & \dot{\Omega}_x & \dot{\Omega}_y & \dot{\Omega}_z & m_x & m_y & m_z & q_x & q_y & q_z & 1 \end{bmatrix}^T, \tag{3}
$$

or

$$
z = \begin{bmatrix} \ddot{r}^T & \dot{\Omega}^T & m^T & q^T & 1 \end{bmatrix}^T, \tag{4}
$$

where

<span id="page-3-1"></span>
$$
\ddot{\mathbf{r}} = \begin{bmatrix} \ddot{x} \\ \ddot{y} \\ \ddot{z} \end{bmatrix}, \quad \dot{\mathbf{\Omega}} = \begin{bmatrix} \dot{\Omega}_x \\ \dot{\Omega}_y \\ \dot{\Omega}_z \end{bmatrix}, \quad m = \begin{bmatrix} m_x \\ m_y \\ m_z \end{bmatrix}, \quad q = \begin{bmatrix} q_x \\ q_y \\ q_z \end{bmatrix}.
$$
 (5)

Here  $\ddot{r}$  and **Ω** are accelerations and angular accelerations being correlated to internal forces *q* and torques *m*, respectively, where all quantities are described in the inertial frame. The "1" at the end of the state vector  $(3)$  is used to account for external forces, as well as centrifugal and Coriolis forces, as will become clear in the following.

<span id="page-3-2"></span>For a system moving in the  $x$ ,  $y$ -plane, the state vectors reduce to

$$
z = \begin{bmatrix} \ddot{x} & \ddot{y} & \dot{\Omega}_z & m_z & q_x & q_y & 1 \end{bmatrix}^\mathrm{T},\tag{6}
$$

and for a one-dimensional motion the reduced state vector reads as

$$
z = \begin{bmatrix} \ddot{x} & q_x & 1 \end{bmatrix}^\mathrm{T} . \tag{7}
$$

<span id="page-4-0"></span>As mentioned above, relations between translational and angular accelerations, on the one hand, and resulting forces and torques, on the other hand, are always related linearly by Newton's second law and Euler's theorem. Also relations between accelerations of different points resulting from rigid body kinematics are linear. Therefore, both types of relations qualify for expressing interrelations between state vectors  $z_{i,l}$  and  $z_{i,O}$  of the input end *I* and output end *O* of an element *j* by linear algebraic equations in matrix form

$$
z_{j,O} = \mathbf{U}_j z_{j,I} \tag{8}
$$

<span id="page-4-2"></span>where the element transfer matrix  $U_j$  has to summarize all physically relevant characteristics of the considered element. For spatial motion the transfer matrix is typically a  $13 \times 13$ matrix according to the state vectors [\(3\)](#page-3-0), whereas for planar and one-dimensional problems the size reduces to  $7 \times 7$  $7 \times 7$  and  $3 \times 3$ , respectively, according to ([6](#page-3-1)) and(7).

### **2.3 Transfer equation and transfer matrix of overall system**

A multibody system is built up from elements by connecting output *O* of an element *j* with input *I* of another element *k*. According to kinematics and Newton's third law, and the specific sign conventions in Sect. [2.1](#page-2-1), the state vectors of these coinciding output and input ends are equal. This allows substituting input states  $z_{k,I}$  by output states  $z_{i,O}$  and the associ-ated transfer equation [\(8\)](#page-4-0), i.e.,  $z_{k,O} = U_k z_{k,I} \equiv U_k z_{i,O} = U_k U_j z_{j,I}$ , which finally leads to products of element transfer matrices. Thus, element transfer matrices may be considered as building blocks which are assembled together according to the topology of the considered system by simple matrix multiplication. The transfer equations [\(8](#page-4-0)) exist or may be derived for rigid bodies, flexible bodies, gas, fluid, and all kind of joints and hinges. Some of them will be derived in Sect. [3](#page-5-0).

<span id="page-4-3"></span><span id="page-4-1"></span>For chain systems consisting of *n* elements, the overall transfer equation can be obtained until the final substitution of transfer equations relates the output state  $z_{n,Q}$  of the *n*th element to the input state  $z_{1,I}$  of the first element by

$$
z_{n,O} = \boldsymbol{U}_{1-n} z_{1,I} \tag{9}
$$

where the overall transfer matrix is given by the product

<span id="page-4-4"></span>
$$
\boldsymbol{U}_{1-n} = \boldsymbol{U}_n \cdots \boldsymbol{U}_2 \boldsymbol{U}_1 \tag{10}
$$

of element transfer matrices.

By applying the boundary conditions, to eliminate the known boundary variables of state vectors from  $z_{1,I}$  and  $z_{n,O}$ , summarizing the remaining unknown boundary states and constant scalar "1" in a vector  $\bar{z}$ , and re-ordering the equations, ([9](#page-4-1)) can be re-written as

$$
\bar{U}\bar{z} = 0 \tag{11}
$$

in order to compute the unknown boundary states. As a result, the boundary state vectors are known completely and allow computing the remaining state vectors in between by successively applying the element transfer equations [\(8](#page-4-0)).

#### **2.4 Simulation of system motion**

It is maybe worthwhile mentioning that in general the transfer matrices depend on position (especially orientation) and/or velocity coordinates. Thus, the procedure in Sect. [2.3](#page-4-2) has to be performed at any time instant  $t_i$  as follows:

- (i) For a time instant  $t_i$ , position and velocity coordinates are given (especially from initial conditions in the beginning).
- (ii) With this information, transfer matrices in ([8](#page-4-0)) can be computed for all elements of the system.
- (iii) The overall transfer matrix may be build up from element matrices according to the system topology; for chain systems by simple multiplication ([10](#page-4-3)).
- (iv) Reduction and solution of  $(11)$  yields unknown boundary states.
- (v) Successive application of [\(8\)](#page-4-0) yields all accelerations as part of state vectors ([3\)](#page-3-0).
- (vi) Any integration scheme may proceed to next time instant.

This solution procedure of the new version of MSTMM is somehow similar to the recursive solution schemes of multibody systems. However, the recursive method has two sweeps through the system, a backward sweep from terminal bodies to the root body to eliminate the bodies and constraints one by one, and a forward sweep from the root body to the terminal bodies to compute the unknown quantities (including the accelerations of bodies, the generalized acceleration of joints and the constraint forces) successively; whereas only the forward sweep is necessary in the proposed method. The backward sweep in the recursive method herein is replaced by the overall transfer equation [\(9](#page-4-1)), in which the overall transfer matrix ([10](#page-4-3)) could be computed by a parallel algorithm, not necessarily by serial processing. If using the proposed approach, any chain system of any length with any boundary conditions reduces to small scale equations  $(9)$  $(9)$ ,  $(10)$ ; closed-loop systems can be handled like chain systems as is done in [[11](#page-19-8)], and the difficulty and computational scale will not increase distinctly compared with the corresponding chain system; the state vector of any point in a system may be directly "transferred" from state vector of any boundary point or any point inside the system.

### <span id="page-5-0"></span>**3 Transfer matrices of some typical elements**

In the following, transfer matrices of some body and hinge elements will be developed with respect to the inertial reference frame respectively for the new version of MSTMM.

### **3.1 Transfer matrix of a rigid body moving in space**

The transfer equation of a rigid body can be deduced directly from its kinematics and dynamics equations. It is taken as an example to sketch the idea of the proposed method for bodies. Transfer equations of flexible bodies can be deduced in a similar way. A rigid body moving in space with single input end *I* and single output end *O* is shown in Fig. [1,](#page-2-0) where *C* denotes the mass center; the subscript *I* denotes the body-fixed coordinate system with origin *I*, while  $\{oxyz\}$  is the inertial coordinate system.

The inertial position coordinates of output point *O* are given by

$$
r_0 = r_I + r_{IO} = r_I + A_I l_{IO}
$$
 (12)

where  $r_I$  is the position of the input point w.r.t. the inertial frame and  $l_{IO}$  describes vector  $\vec{r}_{IO}$  in the body-fixed coordinate system. This formula also applies to point *C*, where  $A_I$ is the direction cosine matrix [\(1\)](#page-2-2) of the body-fixed frame. By differentiation w.r.t. time, we find

$$
\dot{\boldsymbol{r}}_0 = \dot{\boldsymbol{r}}_I + \dot{\boldsymbol{A}}_I \boldsymbol{l}_{IO} = \dot{\boldsymbol{r}}_I + \dot{\boldsymbol{A}}_I \boldsymbol{A}_I^{\mathrm{T}} \boldsymbol{A}_I \boldsymbol{l}_{IO} = \dot{\boldsymbol{r}}_I + \tilde{\boldsymbol{A}}_I \boldsymbol{r}_{IO}
$$

<span id="page-6-2"></span><span id="page-6-0"></span>and by further differentiation

<span id="page-6-6"></span>
$$
\ddot{\boldsymbol{r}}_{O} = \ddot{\boldsymbol{r}}_{I} - \tilde{\boldsymbol{r}}_{IO} \dot{\boldsymbol{\Omega}}_{I} + \tilde{\boldsymbol{\Omega}}_{I} \tilde{\boldsymbol{\Omega}}_{IVIO}.
$$
\n(13)

Using Newton's law and Euler's theorem and considering the sign convention in Sect. [2.1,](#page-2-1) the dynamics equations of the rigid body can be gained as

$$
m\ddot{\boldsymbol{r}}_C = \boldsymbol{q}_I - \boldsymbol{q}_O + \boldsymbol{f}_C,\tag{14}
$$

<span id="page-6-3"></span>
$$
\dot{G}_I = -\tilde{r}_{IO}q_O + m_O - m_I + m_C - m\tilde{r}_{IC}\ddot{r}_I + \tilde{r}_{IC}f_C
$$
\n(15)

where *m* is the body mass,  $q_l$  and  $q_o$  are the forces acting on *I* and *O*,  $f_c$  and  $m_c$  denote external force and torque acting on the mass center of the rigid body,  $m<sub>I</sub>$  and  $m<sub>O</sub>$  are torques acting on the points *I* and *O*,  $r_{IO}$  and  $r_{IC}$  are position vectors from *I* to *O* and *C*, respectively, and  $\ddot{\mathbf{r}}_I$  is the absolute acceleration of point *I*. All quantities are described in the inertial frame. The acceleration of the mass center  $\ddot{r}_C$  can be substituted analogously to ([13](#page-6-0)), where  $O \rightarrow C$ . The moment of momentum is given as  $G_I = A_I J_I A_I^T \mathbf{\Omega}_I$  where the inertia tensor w.r.t. *I* is described in the body-fixed frame. Later on, also the angular velocity  $\omega_I = A_I^T \Omega_I$  in the body-fixed frame will be used, see ([20](#page-6-1)).

Since the angular velocities and thus accelerations of body-fixed reference frames located at any point of the same rigid body are equal, we find

<span id="page-6-4"></span>
$$
\dot{\Omega}_O = \dot{\Omega}_I. \tag{16}
$$

Equations  $(13)$ – $(15)$  $(15)$  $(15)$  can be rewritten as

$$
\ddot{\boldsymbol{r}}_O = \ddot{\boldsymbol{r}}_I + \boldsymbol{E}_1 \dot{\boldsymbol{\Omega}}_I + \boldsymbol{E}_2, \tag{17}
$$

$$
\boldsymbol{q}_{O} = \boldsymbol{q}_{I} + \boldsymbol{E}_{3}\ddot{\boldsymbol{r}}_{I} + \boldsymbol{E}_{4}\dot{\boldsymbol{\Omega}}_{I} + \boldsymbol{E}_{5},
$$
\n(18)

$$
\boldsymbol{m}_O = \boldsymbol{m}_I + \boldsymbol{E}_6 \boldsymbol{q}_O + \boldsymbol{E}_7 \ddot{\boldsymbol{r}}_I + \boldsymbol{E}_8 \dot{\boldsymbol{\Omega}}_I + \boldsymbol{E}_9 \tag{19}
$$

<span id="page-6-1"></span>where

$$
E_1 = -\tilde{r}_{IO}, \qquad E_2 = \tilde{\Omega}_I \tilde{\Omega}_I r_{IO}, \qquad E_3 = -mI_3,
$$
  
\n
$$
E_6 = \tilde{r}_{IO}, \qquad E_7 = m\tilde{r}_{IC}, \qquad E_8 = A_I J_I A_I^{\mathrm{T}},
$$
  
\n
$$
E_4 = m\tilde{r}_{IC}, \qquad E_5 = f_C - m\tilde{\Omega}_I \tilde{\Omega}_I r_{IC},
$$
  
\n
$$
E_9 = -m_C - \tilde{r}_{IC} f_C + A_I \tilde{\omega}_I J_I \omega_I.
$$
\n(20)

<span id="page-6-5"></span>With state vector ([4\)](#page-3-3), [\(16\)](#page-6-3)–([19](#page-6-4)) can be combined in a transfer equation [\(8\)](#page-4-0) with the  $13 \times 13$ transfer matrix of a rigid body

$$
U_{j} = \begin{bmatrix} I_{3} & E_{1} & O_{3\times 3} & O_{3\times 3} & E_{2} \\ O_{3\times 3} & I_{3} & O_{3\times 3} & O_{3\times 3} & O_{3\times 1} \\ E_{6}E_{3} + E_{7} & E_{6}E_{4} + E_{8} & I_{3} & E_{6} & E_{6}E_{5} + E_{9} \\ E_{3} & E_{4} & O_{3\times 3} & I_{3} & E_{5} \\ O_{1\times 3} & O_{1\times 3} & O_{1\times 3} & O_{1\times 3} & 1 \end{bmatrix}_{j} . \tag{21}
$$

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#### <span id="page-7-5"></span><span id="page-7-2"></span><span id="page-7-1"></span>**3.2 Transfer equation of a smooth ball-and-socket hinge**

For a smooth ball-and-socket hinge *j* , its mass and size are neglected. Therefore, the position coordinates, as well as the corresponding time derivatives, and internal forces and torques are equal. This results in the following equations:

<span id="page-7-0"></span>
$$
\ddot{\mathbf{r}}_{j,O} = \ddot{\mathbf{r}}_{j,I},\tag{22}
$$

$$
\boldsymbol{q}_{j,O} = \boldsymbol{q}_{j,I},\tag{23}
$$

$$
\mathbf{m}_{j,O} = \mathbf{m}_{j,I}.\tag{24}
$$

If its outboard body's outboard hinge is also a smooth ball-and-socket hinge  $j + 2$ , the internal torques of  $j + 2$  vanish since the friction is neglected. This results in

$$
\boldsymbol{m}_{j+1,O} = \mathbf{0} \tag{25}
$$

where  $j + 1$  is the corresponding outboard body of the smooth ball-and-socket hinge  $j$ .

The transfer equation for the outboard rigid body of the acceleration hinge *j* is

$$
z_{j+1,0} = U_{j+1} z_{j+1,I}
$$
 (26)

<span id="page-7-6"></span>where the transfer matrix  $(21)$  $(21)$  $(21)$  may be partitioned as

$$
U_{j+1} = \begin{bmatrix} u_{1,1} & u_{1,2} & u_{1,3} & u_{1,4} & u_{1,5} \\ u_{2,1} & u_{2,2} & u_{2,3} & u_{2,4} & u_{2,5} \\ u_{3,1} & u_{3,2} & u_{3,3} & u_{3,4} & u_{3,5} \\ u_{4,1} & u_{4,2} & u_{4,3} & u_{4,4} & u_{4,5} \\ 0_{1\times3} & 0_{1\times3} & 0_{1\times3} & 0_{1\times3} & 1 \end{bmatrix}_{j+1}
$$
 (27)

<span id="page-7-3"></span>Then the third row yields a relation of the input and output state variables of this rigid body as

$$
\boldsymbol{m}_{j+1,O} = \boldsymbol{u}_{3,1}\ddot{\boldsymbol{r}}_{j+1,I} + \boldsymbol{u}_{3,2}\dot{\boldsymbol{\Omega}}_{j+1,I} + \boldsymbol{u}_{3,3}\boldsymbol{m}_{j+1,I} + \boldsymbol{u}_{3,4}\boldsymbol{q}_{j+1,I} + \boldsymbol{u}_{3,5}.
$$
 (28)

<span id="page-7-4"></span>Solving the above equation for  $\dot{Q}_{i+1,I}$  with relation [\(25\)](#page-7-0) yields

$$
\dot{\Omega}_{j+1,I} = -u_{3,2}^{-1}u_{3,1}\ddot{r}_{j+1,I} - u_{3,2}^{-1}u_{3,3}m_{j+1,I} - u_{3,2}^{-1}u_{3,4}q_{j+1,I} - u_{3,2}^{-1}u_{3,5}.
$$
 (29)

<span id="page-7-7"></span>With  $z_{i+1} = z_{i,0}$  and [\(22\)](#page-7-1)–([24](#page-7-2)), [\(29\)](#page-7-3) may be reformulated as

$$
\dot{\Omega}_{j,O} = -u_{3,2}^{-1}u_{3,1}\ddot{r}_{j,I} - u_{3,2}^{-1}u_{3,3}m_{j,I} - u_{3,2}^{-1}u_{3,4}q_{j,I} - u_{3,2}^{-1}u_{3,5}.
$$
 (30)

By combining ([22](#page-7-1))–[\(24\)](#page-7-2) and ([30](#page-7-4)), the transfer matrix of smooth ball-and-socket hinge *j* can be obtained as

$$
U_{j} = \begin{bmatrix} I_{3} & O_{3\times 3} & O_{3\times 3} & O_{3\times 3} & O_{3\times 1} \\ -u_{3,2}^{-1}u_{3,1} & O_{3\times 3} & -u_{3,2}^{-1}u_{3,3} & -u_{3,2}^{-1}u_{3,4} & -u_{3,2}^{-1}u_{3,5} \\ O_{3\times 3} & O_{3\times 3} & I_{3} & O_{3\times 3} & O_{3\times 1} \\ O_{3\times 3} & O_{3\times 3} & O_{3\times 3} & I_{3} & O_{3\times 1} \\ O_{1\times 3} & O_{1\times 3} & O_{1\times 3} & O_{1\times 3} & 1 \end{bmatrix},
$$
(31)

<span id="page-8-8"></span><span id="page-8-0"></span>



#### <span id="page-8-6"></span><span id="page-8-4"></span><span id="page-8-3"></span>**3.3 Transfer equation of smooth pin hinge moving in space**

<span id="page-8-1"></span>A smooth pin hinge *j* with inboard rigid body *j* −1 and outboard rigid body *j* +1 is shown in Fig. [2.](#page-8-0) The position coordinates coincide as well as the corresponding time derivatives, and internal forces and torques are equal:

$$
\ddot{\mathbf{r}}_{j,O} = \ddot{\mathbf{r}}_{j,I},\tag{32}
$$

$$
\boldsymbol{q}_{j,O} = \boldsymbol{q}_{j,I},\tag{33}
$$

$$
\mathbf{m}_{j,O} = \mathbf{m}_{j,I}.\tag{34}
$$

<span id="page-8-7"></span>In the direction of the rotation axis  $\zeta_{i,O}$ , the torque vanishes. By transforming  $m_{j,I}$  to the outboard frame with  $A_{j,O}^T$  and projecting it with  $H_2 = [0 \ 0 \ 1]$  onto the *ζ* -axis, this may be expressed by

$$
\boldsymbol{H}_2 \boldsymbol{A}_{j,O}^\mathrm{T} \boldsymbol{m}_{j,I} = 0. \tag{35}
$$

<span id="page-8-5"></span>From kinematics we find for the angular velocities  $\mathbf{\Omega}_{i,O} = \mathbf{\Omega}_{i,I} + A_{i,O}\omega_r$  where  $\omega_r =$  $[0 0 \dot{\Theta}_j]^{T}$ . By differentiation we get

$$
\dot{\mathbf{\Omega}}_{j,O} = \dot{\mathbf{\Omega}}_{j,I} + \tilde{\mathbf{\Omega}}_{j,O} A_{j,O} \omega_r + A_{j,O} \dot{\omega}_r.
$$
 (36)

The last term of the above equation can be eliminated by premultiplication with  $H_1A_{j,O}^T$ which is firstly a coordinate transformation to the outboard reference frame and then a projection onto directions perpendicular to the rotation axis. By this we finally get

<span id="page-8-2"></span>
$$
\boldsymbol{H}_1 \boldsymbol{A}_{j,O}^{\mathrm{T}} \boldsymbol{\dot{\Omega}}_{j,O} = \boldsymbol{H}_1 \boldsymbol{A}_{j,O}^{\mathrm{T}} \boldsymbol{\dot{\Omega}}_{j,I} + \boldsymbol{H}_1 \boldsymbol{A}_{j,O}^{\mathrm{T}} \boldsymbol{\tilde{\Omega}}_{j,O} \boldsymbol{A}_{j,O} \boldsymbol{\omega}_r
$$
(37)

where

$$
\boldsymbol{H}_1 = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \end{bmatrix}.
$$

Similar to Sect. [3.2](#page-7-5) we now have to take into account outboard body  $j + 1$ . We assume that its outboard element is also a pin hinge, where we have according to  $(35)$ 

$$
H_2 A_{j+1,O}^{\mathrm{T}} m_{j+1,O} = 0.
$$
 (38)

By substituting  $(28)$  $(28)$  $(28)$  into  $(38)$ , we obtain

$$
H_2A_{j+1,0}^{\mathrm{T}}u_{3,1}\ddot{\mathbf{r}}_{j+1,l} + H_2A_{j+1,0}^{\mathrm{T}}u_{3,2}\dot{\mathbf{\Omega}}_{j+1,l} + H_2A_{j+1,0}^{\mathrm{T}}u_{3,3}\mathbf{m}_{j+1,l} + H_2A_{j+1,0}^{\mathrm{T}}u_{3,4}\dot{\mathbf{q}}_{j+1,l} + H_2A_{j+1,0}^{\mathrm{T}}u_{3,5} = 0.
$$
 (39)

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With  $z_{i+1,I} = z_{i,O}$  and [\(32\)](#page-8-3)–([34](#page-8-4)), this may be reformulated as

$$
-H_2A_{j+1,0}^{\mathrm{T}}u_{3,2}\dot{\Omega}_{j,0} = H_2A_{j+1,0}^{\mathrm{T}}u_{3,1}\ddot{r}_{j,I} + H_2A_{j+1,0}^{\mathrm{T}}u_{3,3}m_{j,I} + H_2A_{j+1,0}^{\mathrm{T}}u_{3,4}q_{j,I} + H_2A_{j+1,0}^{\mathrm{T}}u_{3,5}.
$$
 (40)

<span id="page-9-1"></span>In combination with  $(37)$ , we get

$$
E_0 \dot{\Omega}_{j,O} = E_1 \ddot{r}_{j,I} + E_2 \dot{\Omega}_{j,I} + E_3 m_{j,I} + E_4 q_{j,I} + E_5,
$$
\n(41)

or

$$
\dot{\mathbf{\Omega}}_{j,O} = \mathbf{E}_0^{-1} \mathbf{E}_1 \ddot{\mathbf{r}}_{j,I} + \mathbf{E}_0^{-1} \mathbf{E}_2 \dot{\mathbf{\Omega}}_{j,I} + \mathbf{E}_0^{-1} \mathbf{E}_3 \mathbf{m}_{j,I} + \mathbf{E}_0^{-1} \mathbf{E}_4 \mathbf{q}_{j,I} + \mathbf{E}_0^{-1} \mathbf{E}_5, \qquad (42)
$$

where

$$
E_0 = \begin{bmatrix} H_1 A_{j,0}^{\mathrm{T}} \\ -H_2 A_{j+1,0}^{\mathrm{T}} u_{3,2} \end{bmatrix}, \qquad E_1 = \begin{bmatrix} O_{2\times 3} \\ H_2 A_{j+1,0}^{\mathrm{T}} u_{3,1} \end{bmatrix}, \qquad E_2 = \begin{bmatrix} H_1 A_{j,0}^{\mathrm{T}} \\ O_{1\times 3} \end{bmatrix},
$$
  

$$
E_3 = \begin{bmatrix} O_{2\times 3} \\ H_2 A_{j+1,0}^{\mathrm{T}} u_{3,3} \end{bmatrix}, \qquad E_4 = \begin{bmatrix} O_{2\times 3} \\ H_2 A_{j+1,0}^{\mathrm{T}} u_{3,4} \end{bmatrix}, \qquad E_5 = \begin{bmatrix} H_1 A_{j,0}^{\mathrm{T}} \tilde{\Omega}_{j,0} A_{j,0} \omega_r \\ H_2 A_{j+1,0}^{\mathrm{T}} u_{3,5} \end{bmatrix}.
$$

<span id="page-9-2"></span>Combining [\(32\)](#page-8-3), ([33](#page-8-6)), ([34](#page-8-4)), and ([42](#page-9-1)) and arranging them according to the configuration of state vector ([4\)](#page-3-3) yields the transfer matrix of a smooth pin hinge *j* :

$$
U_{j} = \begin{bmatrix} I_{3} & O_{3\times 3} & O_{3\times 3} & O_{3\times 3} & O_{3\times 1} \\ E_{0}^{-1}E_{1} & E_{0}^{-1}E_{2} & E_{0}^{-1}E_{3} & E_{0}^{-1}E_{4} & E_{0}^{-1}E_{5} \\ O_{3\times 3} & O_{3\times 3} & I_{3} & O_{3\times 3} & O_{3\times 1} \\ O_{3\times 3} & O_{3\times 3} & O_{3\times 3} & I_{3} & O_{3\times 1} \\ O_{1\times 3} & O_{1\times 3} & O_{1\times 3} & O_{1\times 3} & 1 \end{bmatrix}_{j}
$$
(43)

<span id="page-9-3"></span><span id="page-9-0"></span>It should be noted the second time derivative of joint coordinate  $\ddot{\theta}_i$  can be obtained if the state vectors of the pin hinge's input and output point are already known. This is achieved by premultiplying [\(36\)](#page-8-7) with  $H_2A_{j,O}^T$ , resulting in

$$
\ddot{\Theta}_j = H_2 A_{j,O}^{\mathrm{T}} \dot{\Omega}_{j,O} - H_2 A_{j,O}^{\mathrm{T}} \dot{\Omega}_{j,I} - H_2 A_{j,O}^{\mathrm{T}} \tilde{\Omega}_{j,O} A_{j,O} \omega_r.
$$
(44)

### **4 Numerical examples**

In order to verify the proposed approach, the dynamics of several spatial pendulum systems and chain systems are simulated numerically by both the proposed method and the classical Newton–Euler method, respectively, where the latter sometimes is also called Lagrange method. For convenience only chain systems are considered in the following. However, the method is also applicable to closed-loop systems, tree systems, network systems, etc., analogously to the original MSDTTMM [\[10\]](#page-19-6).

#### **4.1 Dynamics of a spatial multibody system with ball-and-socket hinges**

Let us firstly consider a pendulum system with three rigid bodies moving in space and connected by three smooth ball-and-socket hinges under the effect of gravity as shown in Fig. [3](#page-11-0). The body and hinge elements are numbered successively, where 1, 3, 5 denote joints, 2, 4, 6 denote bodies, and 0 denotes the boundary ends of the system. The parameters of the three rigid bodies are identical, where the coordinates of the mass center and the output point and the moment of inertia relative to the input-end in the body-fixed frame are as follows:

<span id="page-10-0"></span>
$$
\mathbf{r}_{IC} = [0.5 \ 0 \ 0]^T \text{ m}, \quad \mathbf{r}_{IO} = [1 \ 0 \ 0]^T \text{ m}, \quad \mathbf{J}_I = \text{diag}\left(\frac{1}{6}, \frac{5}{12}, \frac{5}{12}\right) \text{ kg m}^2.
$$

The space-angles given in Sect. [2.1](#page-2-1) of bodies 2, 4, 6 are collected as the generalized coordinates of this system and the initial conditions of the system are

$$
\hat{\boldsymbol{\theta}}_{2,4,6} = \mathbf{0}, \qquad \dot{\boldsymbol{\theta}}_{2,4} = \mathbf{0}, \qquad \dot{\boldsymbol{\theta}}_6 = \begin{bmatrix} 0 & 0.1 & 0 \end{bmatrix}^T \text{ rad/s}.
$$

The overall transfer equation of the system is

$$
z_{6,0} = U_{2-6}z_{2,1} \quad \text{where } U_{2-6} = U_6U_5U_4U_3U_2. \tag{45}
$$

The transfer matrices  $U_2$ ,  $U_4$ ,  $U_6$  referring to bodies correspond to ([21](#page-6-5)), whereas  $U_3$ ,  $U_5$ result from [\(31\)](#page-7-7). The ball-socket-hinge left to body 2 determines the boundary state

$$
z_{2,1} = \begin{bmatrix} 0 & 0 & 0 & \dot{\Omega}_x & \dot{\Omega}_y & \dot{\Omega}_z & 0 & 0 & 0 & q_x & q_y & q_z & 1 \end{bmatrix}_{2,1}^T
$$

whereas the output state of body 6 is free with zero forces and zero torques, but unknown accelerations:

$$
z_{6,0} = \begin{bmatrix} \ddot{x} & \ddot{y} & \ddot{z} & \dot{\Omega}_x & \dot{\Omega}_y & \dot{\Omega}_z & 0 & 0 & 0 & 0 & 0 & 0 & 1 \end{bmatrix}_{6,0}^T.
$$

Thus, the reduced vector of unknown boundaries in [\(11](#page-4-4)) reads as

$$
\bar{z} = \begin{bmatrix} \dot{\Omega}_{x,2,1} & \dot{\Omega}_{y,2,1} & \dot{\Omega}_{z,2,1} & q_{x,2,1} & q_{y,2,1} & q_{z,2,1} \end{bmatrix}
$$

$$
\ddot{x}_{6,0} \quad \ddot{y}_{6,0} \quad \ddot{z}_{6,0} \quad \dot{\Omega}_{x,6,0} \quad \dot{\Omega}_{y,6,0} \quad \dot{\Omega}_{z,6,0} \quad 1 \end{bmatrix}^{\mathrm{T}}.
$$

The time histories of the rotation angles of rigid body 2 obtained by the proposed method and by Newton–Euler method are shown in Fig. [4](#page-11-1) represented by 'line' and ' $\times$ ', respectively. It can be seen clearly that the computational results of the two methods have good agreement, validating the correctness of the new formulation of the transfer matrix method for multibody systems.

#### **4.2 Dynamics of a spatial multibody system with pin hinges**

In the second example, the ball-and-socked hinges in Fig. [3](#page-11-0) are substituted by three smooth pin hinges, leading to the spatial arm system in Fig. [5.](#page-11-2) The overall transfer equation is identical to [\(45\)](#page-10-0), except that  $U_3$ ,  $U_5$  now have to be taken from ([43](#page-9-2)) and the boundary conditions of the system are

$$
z_{2,1} = \begin{bmatrix} 0 & 0 & 0 & 0 & 0 & \dot{\Omega}_z & m_x & m_y & 0 & q_x & q_y & q_z & 1 \end{bmatrix}^T_{2,1},
$$

<span id="page-11-1"></span><span id="page-11-0"></span>

<span id="page-11-2"></span>The joint coordinates [\(44\)](#page-9-3) of pin hinges 1, 3, 5 are collected as the generalized coordinates of this system. Time histories of the relative rotation angles of these pin hinges with initial conditions

$$
\Theta_1 = \Theta_3 = \Theta_5 = 0
$$
 rad,  $\dot{\Theta}_1 = \dot{\Theta}_5 = 0$  rad/s,  $\dot{\Theta}_3 = 1$  rad/s

again coincide for the proposed and Newton–Euler method as shown in Fig. [6](#page-12-0).

# **4.3 Dynamics of a planer pendulum system**

The dynamics of a single pendulum moving in plane is considered to illustrate the increased computational precision of the proposed approach compared to the original MSDTTMM. In Fig. [7,](#page-12-1) the  $\dot{\theta}$  reference trajectory is computed by solving the dynamics equation  $J\ddot{\theta}$  +

<span id="page-12-1"></span><span id="page-12-0"></span>

 $mgl \cos \theta/2 = 0$  with a 4th order Runge–Kutta (R–K) method with a constant step size of  $h = 1.0E-3$  s for initial conditions  $\theta(0) = -\pi/2$  rad and  $\dot{\theta}(0) = -5.4$  rad/s. Dynamics parameters are given as  $J = ml^2/3$ ,  $m = 1$  kg,  $l = 1$  m.

In comparison, the simulation results of the proposed approach combined with 4th order R–K integration method and the original MSDTTMM with constant step size  $h = 1.0E-1$  s are drawn. Obviously, the error of the proposed method is smaller than that of the MS-DTTMM even in the first three integration time steps, where computational errors at each time instant  $t$  are shown in Table [1.](#page-13-0) This demonstrates the higher precision of the proposed approach using exact acceleration computation over the original approach using local linearization.

For computational speed assessment, we consider a planar pendulum system formed by homogeneous rods with length 1 m and mass 1 kg, which are connected by smooth hinges. The CPU used is an i7-3610QM processor with max frequency of 3.3 GHz. Firstly, both the proposed approach and Lagrange equations are computed by using a single thread. The

	h(s)	$e(0.1)$ rad/s	$e(0.2)$ rad/s	$e(0.3)$ rad/s
Reference trajectory	1.0E-3	$\qquad \qquad -$		-
Proposed approach	$1.0E-1$	$-1.21E-4$	$-1.84E-4$	$-1.95E-4$
<b>MSDTTMM</b>	1.0E-1	$-1.05E-2$	$-2.44E-2$	$-3.59E-2$

<span id="page-13-0"></span>**Table 1** Computational error  $e(t)$  at each time instant

global dynamics equation established by Lagrange method can be obtained as

$$
A\begin{bmatrix} \ddot{\theta}_{1} \\ \ddot{\theta}_{2} \\ \vdots \\ \ddot{\theta}_{n} \end{bmatrix} = -B\begin{bmatrix} \dot{\theta}_{1}^{2} \\ \dot{\theta}_{2}^{2} \\ \vdots \\ \dot{\theta}_{n}^{2} \end{bmatrix} - \begin{bmatrix} gl_{1}(m_{1}/2 + m_{2} + \cdots + m_{n}) \cos \theta_{1} \\ gl_{2}(m_{2}/2 + m_{3} + \cdots + m_{n}) \cos \theta_{2} \\ \vdots \\ gl_{n}(m_{n}/2) \cos \theta_{n} \end{bmatrix}
$$
(46)

where  $A = [a_{ij}]$ ,  $B = [b_{ij}]$  with

$$
a_{ij} = \begin{cases} \left(\frac{m_j}{4} + m_{j+1} + \dots + m_n\right)l_j^2 + J_{C_j} & \text{for } i = j, \\ l_i l_j \left(\frac{m_j}{2} + m_{j+1} + \dots + m_n\right) \cos(\theta_i - \theta_j) & \text{for } i \neq j, \end{cases}
$$
  
\n
$$
b_{ij} = \begin{cases} 0 & \text{for } i = j, \\ l_i l_j \left(\frac{m_j}{2} + m_{j+1} + \dots + m_n\right) \sin(\theta_i - \theta_j) & \text{for } i < j, \\ -b_{ji} & \text{for } i > j, \end{cases}
$$

and  $l_j$ ,  $m_j$ ,  $\theta_j$ ,  $J_{C_j}$  are the length, mass, rotation angle and moment of inertial w.r.t. its mass center of rod *j*, respectively.

As shown in Fig. [8](#page-14-0)(a), the time consumption of the proposed approach is much less than that of a classical Lagrange method, which demonstrates the high computational speed of the new version of MSTMM. Under the same computational condition, Fig. [8\(](#page-14-0)b) shows a comparison with the order (n) recursive method [[13](#page-19-9)] which has a similar speed.

# **4.4 Dynamics of a planar multibody system with spring and damping hinges modeled by acceleration hinges**

Figure [9](#page-14-1) shows a chain multibody system composed of masses moving along the *x*-axis. The interactions among the body elements are due to spring and damper elements. The system parameters are

$$
m_i = 0.1 \text{ kg}
$$
  $(i = 2, 4, 6, 8, 10),$   $k_j = 1000 \text{ N/m},$   
 $c_j = 0 \text{ N/(ms}^{-1})$   $(j = 1, 3, 5, 7, 9).$ 

The displacements  $x_{i,I}$   $(i = 2, 4, \ldots, 10)$  of all the bodies are collected as generalized coordinates of the system and the initial conditions are

$$
x_{i,I}(0) = \frac{i}{2} \cdot 0.1 \text{ m}, \qquad \dot{x}_{i,I}(0) = 0 \quad (i = 2, 4, \dots, 10).
$$

The external forces are

$$
f_2 = 100 \text{ N}, \qquad f_k = 0 \text{ N} \quad (k = 4, 6, 8, 10).
$$



<span id="page-14-0"></span>**Fig. 8** Time cost of the new version of MSTMM compared to (**a**) Lagrange method and (**b**) recursive scheme

<span id="page-14-1"></span>The overall transfer equation reads as

$$
z_{10,0} = U_{2-10}z_{2,1}
$$
 where  $U_{2-10} = U_{10}U_9U_8U_7U_6U_5U_4U_3U_2$ 

with  $U_i$  ( $i = 2, 4, 6, 8, 10$ ) denoting the transfer matrix of body elements and  $U_j$  ( $j =$ 3*,* 5*,* 7*,* 9*)* denoting the transfer matrix of hinge elements. The state vector *z* in this example may be reduced according to [\(7](#page-3-2)). The transfer matrix of body elements  $U_i$  ( $i = 2, 4, 6, 8, 10$ ) can be obtained by reducing transfer matrix  $(21)$  $(21)$  $(21)$  of a spatial rigid body to a one dimensional

<span id="page-15-0"></span>



problem, namely

$$
U_i = \begin{bmatrix} 1 & 0 & 0 \\ -m_i & 1 & F_i \\ 0 & 0 & 1 \end{bmatrix} \quad (i = 2, 4, ..., 10)
$$
 (47)

where  $F_i$  is the summation of the external force and all the elastic and damping forces acting on body *i*, i.e.,

$$
F_i = f_i + k_{i-1}(x_{i-2} - x_i) + c_{i-1}(\dot{x}_{i-2} - \dot{x}_i)
$$
  
+  $k_{i+1}(x_{i+2} - x_i) + c_{i+1}(\dot{x}_{i+2} - \dot{x}_i)$   $(i = 2, 4, ..., 8),$  (48)

$$
F_{10} = f_{10} + k_9(x_8 - x_{10}) + c_9(\dot{x}_8 - \dot{x}_{10}).
$$
\n(49)

The two adjacent rigid bodies of a hinge  $j$  ( $j = 3, 5, 7, 9$ ) in this example are relatively free, and such joints are called accelerations hinges in this paper. For such a hinge *j* , we get

$$
(\ddot{x})_{j,O} \equiv (\ddot{x})_{j+1,I} = F_{j+1}/m_{j+1},\tag{50}
$$

$$
(q_x)_{j,O} = (q_x)_{j,I}.
$$
\n(51)

Collecting these two equations according to the state vector ([7](#page-3-2)) yields the transfer matrix

$$
\boldsymbol{U}_{j} = \begin{bmatrix} 0 & 0 & F_{j+1}/m_{j+1} \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix} \quad (j = 3, 5, 7, 9) \tag{52}
$$

of this hinge element. The formulas of a generalized spatial elastic hinge and acceleration hinge are provided in the Appendix.

The boundary conditions of the system are

$$
z_{2,1} = [\ddot{x}, 0, 1]_{2,1}^{T}
$$
,  $z_{10,0} = [\ddot{x}, 0, 1]_{10,0}^{T}$ .

The time histories in Fig. [10](#page-15-0) of the position coordinates of body element 2 and body element 10 again show good agreement between the proposed method and Newton–Euler method.

## **5 Conclusions**

The new version of MSTMM presented in the paper improves the computational precision of the transfer matrix method to the same level as ordinary multibody system approaches based on global dynamics equations. This is achieved by formulating transfer equations between accelerations and forces instead of position coordinates and forces as in the original MSDTTMM. The new formulation avoids linearization and allows using any integration algorithm. Simultaneously it simplifies the transfer matrices compared to the original MSDTTMM, while keeping the features of transfer matrix method of avoiding global dynamics equations of the multibody system. For long chains the computational speed is comparable to that of recursive schemes.

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# **Appendix: Elastic hinge and acceleration hinge**

### **A.1 Elastic hinge**

An elastic hinge [\[14\]](#page-19-10) is sometimes also referred to as bushing element in ADAMS software, coupling element in  $[15]$  $[15]$  $[15]$ , or force element in  $[16]$ , whose mass and size are neglected. It models springs and dampers acting between two bodies so that force and torque magnitudes can be expressed by six Cartesian components, which are linear functions of translational and rotational displacements between two adjacent bodies. For an elastic hinge *j* shown in Fig. [11,](#page-17-0) the inboard and outboard body elements are connected with this hinge. The action and reaction between the two connected body elements  $j - 1$  and  $j + 1$  are represented by the elastic forces due to relative deformations of the hinge. Generally speaking, these elastic forces can be treated as part of the external forces acting on the connected body elements. Thus, the first step is to find the elastic forces.

In Fig. [11,](#page-17-0)  $I \xi_1 \eta_1 \zeta_1$  is the inboard coordinate system of hinge *j*,  $O \xi_2 \eta_2 \zeta_2$  is the outboard frame, and  $I\xi_3\eta_3\zeta_3$  is an intermediate coordinate system used to formulate large relative rotation. The origin point and *ζ* -axis of  $I\xi_3η_3\zeta_3$  coincide with those of  $I\xi_1η_1\zeta_1$ , respectively. *Iξ*<sub>3</sub>*η*<sub>3</sub>*ζ*<sub>3</sub> can be obtained by rotating *Iξ*<sub>1</sub>*η*<sub>1</sub>*ζ*<sub>1</sub> about the *ζ* -axis by angle *θ<sub>r</sub>*, where the value of  $\theta_r$  is achieved such that the relative rotation between  $O\xi_2\eta_2\zeta_2$  and  $I\xi_3\eta_3\zeta_3$  is minimized. *ζ*<sup>1</sup> is the axis of large relative rotation of hinge *j* .

There are various ways to choose an appropriate value of  $\theta_r$ , and the following is just a suggestion. Firstly, one obtains the direction cosine matrices  $A_{0,1}$  and  $A_{0,2}$  of  $I\xi_1\eta_1\zeta_1$  and  $O_{2n/2}^2$ <sub>*ζ*2</sub> with respect to *oxyz*. The relative direction cosine matrix of  $O_{2n/2}^2$ <sub>*ζ*2</sub> with respect to  $I\xi_1\eta_1\zeta_1$  is then given as

$$
A_{1,2} = A_{0,1}^{\mathrm{T}} A_{0,2}.
$$
 (53)

Assuming that  $A_{1,2}$  is obtained by three continuous rotations according to Sect. [2.1](#page-2-1), one can deduce three corresponding rotation angles  $\theta_1$ ,  $\theta_2$ ,  $\theta_3$ . Then one can treat the third rotation angle  $\theta_3$  as the unknown variable  $\theta_r$ . In a similar way,  $\dot{\theta}_r$ , the time derivative of  $\theta_r$ , can be obtained. Then the orientation and angular velocity of  $I\xi_3\eta_3\zeta_3$  are totally obtained.

#### <span id="page-17-0"></span>**Fig. 11** Spatial elastic hinge

In a similar way, the relative orientation cosine matrix of  $O_{52} \eta_2 \zeta_2$  with respect to  $I_{53} \eta_3 \zeta_3$ and its time derivative with respect to time can be obtained, denoted as  $A_{3,2}$  and  $A_{3,2}$ , respectively. Then one can obtain the three corresponding infinitesimal rotation angles and their time derivatives with respect to time, which are denoted as  $\theta_{3,2} = [\theta_x \theta_y \theta_z]^T$  and  $\dot{\theta}_{3,2} = [\dot{\theta}_x \ \dot{\theta}_y \ \dot{\theta}_z]^{\mathrm{T}}$ , respectively.

The elastic torques and forces of hinge *j* decomposed in  $I\xi_3\eta_3\zeta_3$  can be obtained as follows:

$$
\mathbf{m} = [k'_x \theta_x + c'_x \dot{\theta}_x \quad k'_y \theta_y + c'_y \dot{\theta}_y \quad k'_z \theta_r + c'_z \dot{\theta}_r]^T, \n\mathbf{f} = [k_x \Delta x + c_x \Delta \dot{x} \quad k_y \Delta y + c_y \Delta \dot{y} \quad k_y \Delta z + c_y \Delta \dot{z}]^T
$$
\n(54)

where  $\Delta r = [\Delta x \Delta y \Delta z]^T$  is the position vector from the origin of  $O \xi_2 \eta_2 \zeta_2$  to that of  $I\xi_3\eta_3\zeta_3$  decomposed in  $I\xi_3\eta_3\zeta_3$ .  $k_x$ ,  $k_y$ ,  $k_z$ ,  $c_x$ ,  $c_y$  and  $c_z$  are the stiffness and damping coefficients.  $k'_x$ ,  $k'_y$ ,  $k'_z$ ,  $c'_x$ ,  $c'_y$ , and  $c'_z$  are the rotational stiffness and damping coefficients, respectively.

The elastic forces [\(54\)](#page-17-1) will be treated as part of the external forces in [\(14\)](#page-6-6) and ([15](#page-6-2)) acting on the connected body elements and the elastic hinge may be treated as an acceleration hinge whose transfer equation is deduced in the following section.

# **A.2 Transfer equations of acceleration hinges moving in space**

The acceleration hinge may be considered as a dummy hinge, but without interaction between the two adjacent bodies. The translational and angular acceleration of the output end of an acceleration hinge can be obtained from the transfer matrix of its outboard body. The processes to deduce this transfer matrix are similar to those for smooth ball-and-socket hinges and pin hinges in Sects. [3.2](#page-7-5) and [3.3.](#page-8-8) The internal forces and torques of the elastic hinge will be treated as external forces and torques acting on its inboard and outboard body, respectively.

For a massless hinge *j* moving in space, the following force and moment equilibria can be obtained:

<span id="page-17-3"></span><span id="page-17-2"></span>
$$
\boldsymbol{q}_{j,O} = \boldsymbol{q}_{j,I},\tag{55}
$$

$$
\mathbf{m}_{j,O} = \mathbf{m}_{j,I}.\tag{56}
$$

<span id="page-17-1"></span>

If its outboard body's outboard hinge is also an acceleration hinge  $j + 2$ , the internal forces and torques of  $j + 2$  vanish since they are treated as part of the external forces acting on its two adjacent bodies. This results in

$$
\boldsymbol{q}_{j+1,O} = \boldsymbol{0},\tag{57}
$$

$$
\boldsymbol{m}_{j+1,O} = \mathbf{0} \tag{58}
$$

where  $j + 1$  is the corresponding outboard body of the acceleration hinge  $j$ .

The transfer equation for the outboard rigid body of the acceleration hinge *j* is

$$
z_{j+1,O} = U_{j+1} z_{j+1,I}
$$
\n(59)

<span id="page-18-2"></span>where the transfer matrix  $(21)$  $(21)$  $(21)$  may be partitioned as

$$
U_{j+1} = \begin{bmatrix} u_{1,1} & u_{1,2} & u_{1,3} & u_{1,4} & u_{1,5} \\ u_{2,1} & u_{2,2} & u_{2,3} & u_{2,4} & u_{2,5} \\ u_{3,1} & u_{3,2} & u_{3,3} & u_{3,4} & u_{3,5} \\ u_{4,1} & u_{4,2} & u_{4,3} & u_{4,4} & u_{4,5} \\ 0_{1\times 3} & 0_{1\times 3} & 0_{1\times 3} & 0_{1\times 3} & 1 \end{bmatrix}_{j+1}
$$
(60)

<span id="page-18-4"></span><span id="page-18-3"></span>Then the third and fourth rows yield a relation of internal moments and forces between input and output of this rigid body as

$$
\boldsymbol{m}_{j+1,O} = \boldsymbol{u}_{3,1}\ddot{\boldsymbol{r}}_{j+1,I} + \boldsymbol{u}_{3,2}\dot{\boldsymbol{\Omega}}_{j+1,I} + \boldsymbol{u}_{3,3}\boldsymbol{m}_{j+1,I} + \boldsymbol{u}_{3,4}\boldsymbol{q}_{j+1,I} + \boldsymbol{u}_{3,5},
$$
(61)

$$
\boldsymbol{q}_{j+1,O} = \boldsymbol{u}_{4,1} \ddot{\boldsymbol{r}}_{j+1,I} + \boldsymbol{u}_{4,2} \dot{\boldsymbol{\Omega}}_{j+1,I} + \boldsymbol{u}_{4,3} \boldsymbol{m}_{j+1,I} + \boldsymbol{u}_{4,4} \boldsymbol{q}_{j+1,I} + \boldsymbol{u}_{4,5}.
$$
 (62)

Substituting  $(57)$  and  $(58)$  into  $(61)$  and  $(62)$  $(62)$  $(62)$  yields

$$
\begin{bmatrix} u_{3,1} & u_{3,2} \\ u_{4,1} & u_{4,2} \end{bmatrix} \begin{bmatrix} \ddot{r}_{j+1,I} \\ \dot{\Omega}_{j+1,I} \end{bmatrix} + \begin{bmatrix} u_{3,3} & u_{3,4} \\ u_{4,3} & u_{4,4} \end{bmatrix} \begin{bmatrix} m_{j+1,I} \\ q_{j+1,I} \end{bmatrix} + \begin{bmatrix} u_{3,5} \\ u_{4,5} \end{bmatrix} = 0.
$$
 (63)

<span id="page-18-5"></span>With  $z_{i+1, I} = z_{i, O}$  and [\(55\)](#page-17-2)–([56](#page-17-3)), [\(63\)](#page-18-4) may be reformulated as

$$
\begin{bmatrix} \ddot{\mathbf{r}}_{j,O} \\ \dot{\mathbf{\Omega}}_{j,O} \end{bmatrix} = -\hat{\mathbf{u}}_1^{-1} \hat{\mathbf{u}}_2 \begin{bmatrix} \mathbf{m}_{j,I} \\ \mathbf{q}_{j,I} \end{bmatrix} - \hat{\mathbf{u}}_1^{-1} \hat{\mathbf{u}}_3
$$
(64)

where

$$
\hat{u}_1 = \begin{bmatrix} u_{3,1} & u_{3,2} \\ u_{4,1} & u_{4,2} \end{bmatrix}, \qquad \hat{u}_2 = \begin{bmatrix} u_{3,3} & u_{3,4} \\ u_{4,3} & u_{4,4} \end{bmatrix}, \qquad \hat{u}_3 = \begin{bmatrix} u_{3,5} \\ u_{4,5} \end{bmatrix}.
$$
 (65)

By combining  $(55)$  $(55)$ ,  $(56)$  and  $(65)$  $(65)$  $(65)$ , the transfer matrix of acceleration hinge *j* can be obtained as

$$
U_{j} = \begin{bmatrix} \mathbf{0}_{6\times6} & -\hat{\mathbf{u}}_{1}^{-1}\hat{\mathbf{u}}_{2} & -\hat{\mathbf{u}}_{1}^{-1}\hat{\mathbf{u}}_{3} \\ \mathbf{0}_{6\times6} & \mathbf{I}_{6} & \mathbf{0}_{6\times1} \\ \mathbf{0}_{1\times6} & \mathbf{0}_{1\times6} & 1 \end{bmatrix}_{j}.
$$
 (66)

<span id="page-18-1"></span><span id="page-18-0"></span>

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# <span id="page-19-5"></span><span id="page-19-4"></span><span id="page-19-3"></span><span id="page-19-2"></span><span id="page-19-1"></span><span id="page-19-0"></span>**References**

- 1. Wittenburg, J.: Dynamics of Systems of Rigid Bodies. Teubner, Stuttgart (1977)
- 2. Schiehlen, W.: Multibody Systems Handbook. Springer, Berlin (1990)
- 3. Shabana, A.A.: Dynamics of Multibody Systems. Cambridge University Press, Cambridge (2009)
- 4. Kumar, A.S., Sankar, T.S.: A new transfer matrix method for response analysis of large dynamic systems. Comput. Struct. **23**, 545–552 (1986)
- <span id="page-19-6"></span>5. Rui, X., Wang, G., Lu, Y.: Transfer matrix method for linear multibody system. Multibody Syst. Dyn. **19**(3), 179–207 (2008)
- <span id="page-19-8"></span>6. Bestle, D., Abbas, L., Rui, X.: Recursive eigenvalue search algorithm for transfer matrix method of linear flexible multibody systems. Multibody Syst. Dyn. **32**, 429–444 (2014)
- <span id="page-19-7"></span>7. Rui, X., He, B., Lu, Y., et al.: Discrete time transfer matrix method for multibody system dynamics. Multibody Syst. Dyn. **14**(4), 317–344 (2005)
- <span id="page-19-9"></span>8. Rui, X., He, B., Rong, B., et al.: Discrete time transfer matrix method for multi-rigid-flexible-body system moving in plane. Multibody Syst. Dyn. **223**(K1), 23–42 (2009)
- 9. Rui, X., Rong, B., Wang, G., et al.: Discrete time transfer matrix method for dynamics analysis of complex weapon systems. Sci. China, Technol. Sci. **54**(5), 1061–1071 (2011)
- <span id="page-19-12"></span><span id="page-19-11"></span><span id="page-19-10"></span>10. Rui, X., Yun, L., Lu, Y., et al.: Transfer Matrix Method for Multibody System and Its Applications. Science Press, Beijing (2008)
- 11. Rui, X., Zhang, J., Zhou, Q.: Automatic deduction theorem of overall transfer equation of multibody systems. Adv. Mech. Eng. **6**, 378047 (2014)
- 12. Kane, T.R., Likine, P.W., Levinson, D.A.: Spacecraft Dynamics. McGraw-Hill, New York (1983)
- 13. Brandl, H., Johanni, R., Otter, M.: An algorithm for the simulation of multibody systems with kinematic loops. In: Proceedings of the IFToMM Seventh World Congress on the Theory of Machines and Mechanisms, pp. 407–411 (1987)
- 14. Pestel, E.C., Leckie, F.A.: Matrix Method in Elastomechanics. McGraw-Hill, New York (1963)
- 15. Popp, K., Schielen, W.: Ground Vehicle Dynamics. Springer, Berlin (2010)
- 16. Shabana, A.: Computational Dynamics. Wiley, Chichester (2010)