

Equations of Motion for Multiple-Rigid-Body Systems via Symbolic Manipulation

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Conventional general-purpose multibody simulation programs, although easily accessible to users, fail to provide explicit equations of motion for given spacecraft configurations. In addition, they require more computer time for performing simulations than do corresponding special-purpose programs. Nevertheless, the derivation of the equations of motion for a spacecraft consisting of two or more rigid bodies is, in general, a formidable task. In this paper, it is shown how the symbolic manipulation language FORMAC can be used in conjunction with Kane's equations to generate on a digital computer explicit equations of motion for given multiple-rigid-body systems.

Introduction

IN recent years, spacecraft designers have become increasingly involved with the dynamics of multiple-rigid-body systems. When a spacecraft is modeled as a collection of two or more rigid bodies, the derivation of the associated equations of motion becomes laborious and subject to careless errors if performed by hand. Accordingly, much attention has been directed toward the development of techniques for generating equations of motion by digital computer.

One of the earliest attempts at formulating equations of motion for multiple-rigid-body systems is attributable to Hooker and Margulies,¹ who derived dynamical equations for the attitude motion of an arbitrary number n of point-connected rigid bodies, the only restriction on the configurations included in this treatment being that chains of connected bodies must not form closed loops. The dynamical equations of Newton and Euler were written for each body, and a procedure was described for the elimination of "constraint torques" that arise at connection points having associated with them only one or two rotational degrees of freedom, thus yielding a set of $3n$ second-order scalar differential equations of motion for the system. A similar set of equations was obtained by Roberson and Wittenburg.² Later, Hooker³ devised a method for eliminating the constraint torques in such a way as to reduce the total number of equations from $3n$ to the number of rotational degrees of freedom of the system. Velman⁴ and Russell,⁵ using a different approach, chose not to isolate each of the n point-connected rigid bodies individually, but instead constructed vector equations of motion for n different subsets of bodies, including, ultimately, a complete set of equations for the entire system. Features of both the discrete-body and nested-body approaches were combined by Frisch⁶ in deriving a set of equations which forms the basis for a digital computer program. Other simulation programs have been developed by Fleischer⁷ and Fleischer and Likins⁸ based upon Refs. 1 and 3. More recently, Huston and Passerello⁹ have made use of Lagrange's form of D'Alembert's principle as developed by Kane¹⁰ and have, consequently, avoided the necessity of considering nonworking interbody constraint forces and torques in their derivation of the equations of motion and corresponding computer program.

All of the aforementioned formulations share certain disadvantages. First, the equations of motion for any particular multiple-rigid-body configuration exist only as indicated operations in computer programs. That is, the analyst never sees explicit terms in the equations as he would if the equations had been derived by hand and programmed term-by-term. Also, the equations must be completely constructed, using a substantial number of DO loops, at every step of a numerical integration and are not, therefore, cast in a form that can be used advantageously with certain computers, such as the CDC 6600, which have the capability of performing several simultaneous arithmetic operations on explicit expressions. Hence, there is sufficient motivation to seek a way of employing the speed and accuracy of a computer in the construction of explicit equations of motion for systems too large to be dealt with by hand. To this end, one can make use of non-numeric computation, that is, symbolic manipulation.

Any symbolic manipulation language being considered for use in formulating dynamical equations of motion must be able to carry out analytical differentiation. Among the more common such languages are SYMBAL,¹¹ TRIGMAN,¹² and FORMAC.¹³ SYMBAL, however, being an ALGOL-based language, produces output that is not, in general, easily read or directly usable with fast FORTRAN numerical integrators. Although TRIGMAN is written in FORTRAN, its use is limited to problems having no more than 29 polynomial or 10 trigonometric variables. FORMAC, based on PL/I, is implemented easily by users familiar only with FORTRAN and contains a powerful array of manipulative functions, including provisions for releasing computer storage space that no longer is needed. Furthermore, FORMAC possesses a variety of output formats, including punched-card options available in a newer version of the language.¹⁴ Hence, FORMAC is ideally suited for the task at hand and is, in fact, employed in the present paper.

The next choice which must be made is that of a method of analytical dynamics to be used with FORMAC. Adopted in the present work is Kane's development of Lagrange's form of D'Alembert's principle, which, as previously pointed out in connection with the work of Huston and Passerello,⁹ provides for the automatic elimination of nonworking interbody constraint forces and torques. In addition, this method is applied in conjunction with Kane's u 's taken as dependent variables (see Ref. 10, pp. 44-48, 188-192), these being linear combinations of time derivatives of system-generalized coordinates (for example, angular velocity measure numbers), the use of which produces equations having less complexity than those derived using independent generalized coordinates themselves as dependent variables.

This paper contains a brief account of the elements of the FORMAC language which are directly applicable to the

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formulation of equations of motion for multiple-rigid-body systems. An example configuration consisting of two point-connected rigid bodies then is analyzed in order to illustrate the use of FORMAC with Kane's equations. For this case, steps are given for generating the explicit equations of motion and for subsequently putting them into a form appropriate for numerical integration. Finally, simulation results obtained from these equations are presented as a practical application of the method.

Elements of the FORMAC Language

In order to bring to light the basic capabilities of FORMAC which are applicable to the derivation of equations of motion, a short summary of the relevant portions of the language will now be given. A complete description of the FORMAC language may be found in Refs. 13 and 14.

To begin with, FORMAC assignment statements are designed by the word LET. For example, the statement $\text{LET}(X1 = C*(A+B)**3)$; assigns to the FORMAC variable $X1$ the value $C(A+B)^3$. More than one assignment statement may appear within the range of a single LET. If one wishes to express $X1$ in its fully expanded form using both the binomial theorem and the distributive law, one can invoke the EXPAND routine. Execution of the statement $\text{LET}(X2 = \text{EXPAND}(X1))$; results in the assignment $X2 = 3C A^2 B + 3C A B^2 + C A^3 + C B^3$.

Another useful routine for manipulating expressions is COEFF. Written in the general format $\text{LET}(E3 = \text{COEFF}(E1, E2))$; COEFF assigns to $E3$ the coefficient of the subexpression $E2$ in the expression $E1$. For example, with $X2$ as previously defined, the statement $\text{LET}(X3 = \text{COEFF}(X2, C))$; results in $X3 = 3 A^2 B + 3 A B^2 + A^3 + B^3$.

One of the most important attributes of FORMAC is its ability to deal with unspecified functions. For instance, $\text{LET}(X4 = D*T*\text{SIN}(T) + Q.(T))$; assigns to $X4$ the sum of two functions of a variable T , the first depending on T explicitly and the second, designated by a decimal point, an implicit function of T . This feature is found to be most useful in connection with DERIV, one of FORMAC's analytical differentiation routines. In this instance, the statement $\text{LET}(X5 = \text{DERIV}(X4, T))$; produces $X5 = D T \cos(T) + D \sin(T) + Q^{(1)}(T)$, where the superscript 1 enclosed in parentheses signifies the first derivative of the function Q with respect to its argument T .

If one desires to assign a name to the first derivative of an unspecified function, one can apply the FORMAC routine pair DIFF, CHAIN. With $X4$ retaining its previous value, the sequence of statements $\text{LET}(\text{DIFF}(Q) = \text{CHAIN}(\text{QDOT}))$; $X6 = \text{DERIV}(X4, T)$; leads to $X6 = \text{QDOT} + D T \cos(T) + D \sin(T)$, where $X6$ differs from $X5$ by the fact that, in $X6$, $Q(1) \cdot (T)$ has been replaced with QDOT .

One can avail oneself of FORMAC's substitution capability by employing the routine REPLACE, which, in its general format, is written as $\text{LET}(E = \text{REPLACE}(E1, A1, B1, A2, B2, \dots, AN, BN))$; where $E1, A1, B1, \dots, AN, BN$ are arbitrary FORMAC expressions. Execution is accomplished as follows: Each occurrence of $A1$ in $E1$ is replaced with $B1$ to yield an expression $E2$. Then each occurrence of $A2$ in $E2$ is replaced with $B2$ to produce an expression $E3$. The process is continued until each occurrence of AN in EN is replaced with BN , at which time the resulting expression is returned as the value of E . As an illustration, $\text{LET}(X7 = \text{REPLACE}(X6, \text{SIN}(T), \text{ST}, \text{COS}(T), \text{CT}))$; results in $X7 = \text{QDOT} + \text{ST} D + \text{CT} D T$, where $X6$ is defined as in the preceding example.

Printed output in FORMAC is handled by the PRINT _OUT statement. For example, $\text{PRINT_OUT}(X7)$; produces the printed line $X7 = \text{QDOT} + \text{ST} D + \text{CT} D T$. Note that the expression has been automatically edited to afford facility in reading. In the latest version of the language¹⁴ FORMAC expressions also may be punched on cards. The routine PUNCH_F, used in the same manner as PRINT_OUT, results

in the punched output of syntactically correct FORTRAN statements. In addition, there are similar punching routines for PL/I and double-precision FORTRAN.

The processing of long FORMAC expressions requires a significant amount of computer core. FORMAC provides two routines for circumventing major storage difficulties, SAVE and ATOMIZE. $\text{SAVE}(E1; E2; \dots; EN)$; copies the expressions named by $E1, \dots, EN$ into secondary storage while releasing the space in core that they previously occupied. A subsequent reference to any Ei ($i=1, \dots, N$) results in the automatic return to core of the expression that Ei names. $\text{ATOMIZE}(E1; E2; \dots; EN)$; causes the space occupied by the expressions named by $E1, \dots, EN$ to be released. The value of any FORMAC variable that has been "atomized" is represented by a null string. For example, $\text{ATOMIZE}(X7)$; yields, upon output, $X7 =$.

One now can begin to comprehend the potential utility of the FORMAC capability, of which the preceding discussion covers but a small part. The idea behind a FORMAC-based scheme for deriving equations of motion is this: rather than create large general-purpose multibody programs that numerically construct and integrate equations of motion for any spacecraft configuration, one can proceed from the concept of treating each multiple-rigid-body system as a separate case and using FORMAC to derive its explicit equations of motion once and for all. The example that follows is intended to demonstrate the feasibility of such an approach.

Application

The system to be considered is a spacecraft (see Fig. 1) consisting of two rigid bodies: a main body B , and a rotor R that is connected to B at point P by a hinge whose axis ζ maintains a fixed orientation relative to both B and R . Lines ξ_i, η_i ($i=1, 2, 3$) are, respectively, principal axes of inertia of B and R for their mass centers B^* and R^* , and ζ is presumed to be parallel to ξ_1 (and hence to η_1). Thus, if one introduces a dextral set of unit vectors b_1, b_2, b_3 , respectively parallel to ξ_1, ξ_2, ξ_3 , and unit vectors r_i parallel to η_i ($i=2, 3$) such that b_1, r_2, r_3 also form a dextral set, then R can be brought into a general orientation with respect to B by initially aligning r_i with b_i ($i=2, 3$) and then subjecting R to a rotation about ζ relative to B characterized by $q b_1$. The mass of B is m_B , and the moments of inertia of B about ξ_i are β_i ($i=1, 2, 3$). Similarly, R has mass m_R , and its moments of inertia about η_i are denoted by ρ_i ($i=1, 2, 3$). The spacecraft is assumed to be moving in a Newtonian reference frame N .

The first step in the process is to record the vector and dyadic quantities from which Kane's equations of motion for this system can be constructed. To begin with, the unit vectors r_2, r_3 can be written as

$$r_2 = \cos q b_2 + \sin q b_3 \quad (1)$$

$$r_3 = -\sin q b_2 + \cos q b_3 \quad (2)$$

so that all vectors and dyadics herein ultimately can be exhibited in the basis b_1, b_2, b_3 for the sake of uniformity. Next, the angular velocity ${}^N\omega^B$ of B in N always can be expressed in terms of functions u_1, u_2, u_3 of time t as

$${}^N\omega^B = u_1 b_1 + u_2 b_2 + u_3 b_3 \quad (3)$$

so that the angular velocity ${}^N\omega^R$ of R in N then is given by

$${}^N\omega^R = (u_1 + u_4) b_1 + u_2 b_2 + u_3 b_3 \quad (4)$$

where

$$u_4 \triangleq \dot{q} \quad (5)$$

The angular acceleration ${}^N\alpha^B$ of B in N can be written as

$${}^N\alpha^B = \dot{u}_1 b_1 + \dot{u}_2 b_2 + \dot{u}_3 b_3 \quad (6)$$

whereas ${}^N\alpha^R$, the angular acceleration of R in N , is furnished by the equation

$${}^N\alpha^R = {}^B d/dt({}^N\omega^R) + {}^N\omega^B \times {}^N\omega^R \quad (7)$$

where ${}^B d/dt$ denotes differentiation with respect to t in B .

One can express the position vectors p^{P/B^*} and p^{P/R^*} of P relative to B^* and R^* as

$$p^{P/B^*} = b_1 b_1 + b_2 b_2 + b_3 b_3 \quad (8)$$

$$p^{P/R^*} = r_1 b_1 + r_2 r_2 + r_3 r_3 \quad (9)$$

where b_i, r_i ($i=1,2,3$) are constants, and p^{R^*/B^*} , the position vector of R^* relative to B^* , as

$$p^{R^*/B^*} = p^{P/B^*} - p^{P/R^*} \quad (10)$$

The velocity ${}^N v^{B^*}$ of B^* in N can be written in terms of functions u_5, u_6, u_7 of t as

$${}^N v^{B^*} = u_5 b_1 + u_6 b_2 + u_7 b_3 \quad (11)$$

so that ${}^N v^{R^*}$, the velocity of R^* in N , is obtained from

$${}^N v^{R^*} = {}^N v^{B^*} + \frac{{}^B d}{dt} (p^{R^*/B^*}) + {}^N \omega^B \times p^{R^*/B^*} \quad (12)$$

The accelerations ${}^N a^{B^*}$, ${}^N a^{R^*}$ of B^* and R^* in N then are given by

$${}^N a^{B^*} = \frac{{}^B d}{dt} ({}^N v^{B^*}) + {}^N \omega^B \times {}^N v^{B^*} \quad (13)$$

$${}^N a^{R^*} = \frac{{}^B d}{dt} ({}^N v^{R^*}) + {}^N \omega^B \times {}^N v^{R^*} \quad (14)$$

and inertia dyadics I^{B/B^*} and I^{R/R^*} of B for B^* and R for R^* can be written as

$$I^{B/B^*} = \beta_1 b_1 b_1 + \beta_2 b_2 b_2 + \beta_3 b_3 b_3 \quad (15)$$

$$I^{R/R^*} = \rho_1 b_1 b_1 + \rho_2 r_2 r_2 + \rho_3 r_3 r_3 \quad (16)$$

thus providing the quantities necessary for the construction of inertia forces $(F^*)_B$, $(F^*)_R$ and inertia torques $(T^*)_B$, $(T^*)_R$ for B and R in N , given by (see Ref. 10, pp. 91, 115-116)

$$(F^*)_B = -m_B {}^N a^{B^*} \quad (17)$$

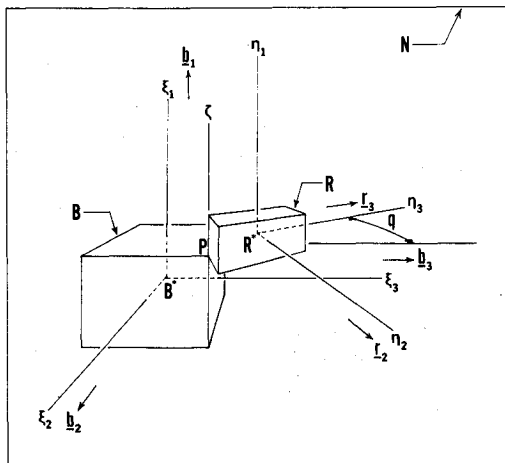


Fig. 1 Schematic representation of spacecraft.

$$(F^*)_R = -m_R {}^N a^{R^*} \quad (18)$$

$$(T^*)_B = (I^{B/B^*} \cdot {}^N \omega^B) \times {}^N \omega^B - I^{B^*/B^*} \cdot {}^N \alpha^B \quad (19)$$

$$(T^*)_R = (I^{R/R^*} \cdot {}^N \omega^R) \times {}^N \omega^R - I^{R^*/R^*} \cdot {}^N \alpha^R \quad (20)$$

The system of contact and body forces acting on B is assumed to be equivalent to a force $(F)_B$, whose line of action passes through B^* , given by

$$(F)_B = F_1 b_1 + F_2 b_2 + F_3 b_3 \quad (21)$$

and a couple of torque $(T)_B$ given by

$$(T)_B = T_1 b_1 + T_2 b_2 + T_3 b_3 \quad (22)$$

whereas R is assumed to be free of such forces.

Generalized active forces F_r and generalized inertia forces F_r^* associated with u_r ($r=1, \dots, 7$) then can be obtained from the expressions

$$F_r = \frac{\partial {}^N v^{B^*}}{\partial u_r} \cdot (F)_B + \frac{\partial {}^N \omega^B}{\partial u_r} \cdot (T)_B \quad (r=1, \dots, 7) \quad (23)$$

$$F_r^* = \frac{\partial {}^N v^{B^*}}{\partial u_r} \cdot (F^*)_B + \frac{\partial {}^N v^{R^*}}{\partial u_r} \cdot (F^*)_R + \frac{\partial {}^N \omega^B}{\partial u_r} \cdot (T^*)_B + \frac{\partial {}^N \omega^R}{\partial u_r} \cdot (T^*)_R \quad (r=1, \dots, 7) \quad (24)$$

so that, finally, one arrives at Kane's equations, given by

$$F_r + F_r^* = 0 \quad (r=1, \dots, 7) \quad (25)$$

(Details of Kane's theory may be found throughout Ref. 10.)

Equation (25) represents a set of first-order ordinary differential equations of the form

$$X \dot{u} = Y \quad (26)$$

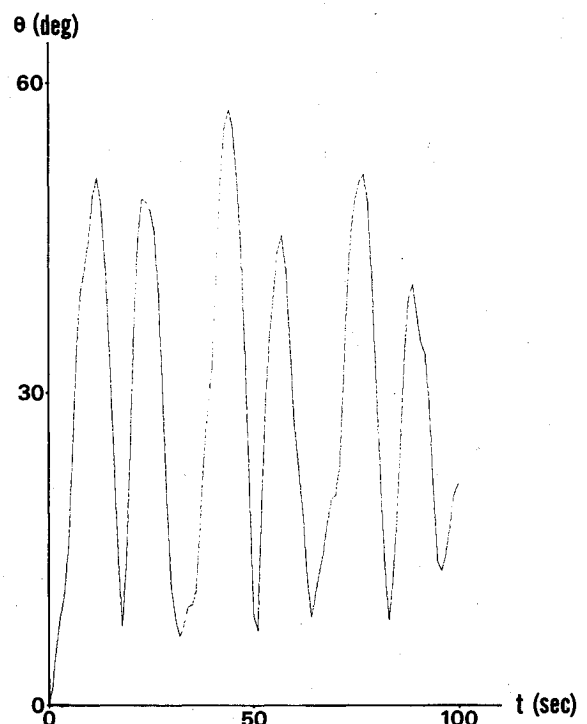


Fig. 2 Simulation results.

where X is a 7×7 matrix of functions of q , Y is a 7×1 matrix of functions of q and u_r ($r=1, \dots, 7$), and \dot{u} is the 7×1 matrix containing $\dot{u}_1, \dots, \dot{u}_7$. Note that one cannot obtain a complete description of the motion of the spacecraft from Eq. (26) alone. One must supply, in addition, a set of kinematical equations that relate u_1, \dots, u_7 to variables characterizing the orientation of R relative to B , the orientation of B in N , and the position of B^* in N . The first of these equations is furnished by Eq. (5), and the remainder may be chosen at one's discretion. For example, if the relationship between b_1, b_2, b_3 and a dextral set of mutually perpendicular unit vectors n_1, n_2, n_3 fixed in N is specified by the Euler parameters, $\epsilon_1, \epsilon_2, \epsilon_3, \epsilon_4$, and if the position vector $p^{B^*/O}$ of B^* relative to a point O fixed in N is written in the form $p^{B^*/O} = x_1 n_1 + x_2 n_2 + x_3 n_3$, where x_i ($i=1,2,3$) are functions of t , then the associated kinematical equations are given by¹⁵

$$\dot{\epsilon}_1 = \frac{1}{2}(u_1 \epsilon_4 - u_2 \epsilon_3 + u_3 \epsilon_2) \quad (27)$$

$$\dot{\epsilon}_2 = \frac{1}{2}(u_1 \epsilon_3 + u_2 \epsilon_4 - u_3 \epsilon_1) \quad (28)$$

$$\dot{\epsilon}_3 = \frac{1}{2}(-u_1 \epsilon_2 + u_2 \epsilon_1 + u_3 \epsilon_4) \quad (29)$$

$$\dot{\epsilon}_4 = \frac{1}{2}(-u_1 \epsilon_1 - u_2 \epsilon_2 - u_3 \epsilon_3) \quad (30)$$

$$\dot{x}_1 = u_5 C_{11} + u_6 C_{12} + u_7 C_{13} \quad (31)$$

$$\dot{x}_2 = u_5 C_{21} + u_6 C_{22} + u_7 C_{23} \quad (32)$$

$$\dot{x}_3 = u_5 C_{31} + u_6 C_{32} + u_7 C_{33} \quad (33)$$

where the direction cosines $C_{ij} = n_i \cdot b_j$ ($i, j=1,2,3$) are given by

$$C_{11} = 1 - 2\epsilon_2^2 - 2\epsilon_3^2 \quad (34)$$

$$C_{12} = 2(\epsilon_1 \epsilon_2 - \epsilon_3 \epsilon_4) \quad (35)$$

$$C_{13} = 2(\epsilon_3 \epsilon_1 + \epsilon_2 \epsilon_4) \quad (36)$$

$$C_{21} = 2(\epsilon_1 \epsilon_2 + \epsilon_3 \epsilon_4) \quad (37)$$

$$C_{22} = 1 - 2\epsilon_3^2 - 2\epsilon_1^2 \quad (38)$$

$$C_{23} = 2(\epsilon_2 \epsilon_3 - \epsilon_1 \epsilon_4) \quad (39)$$

$$C_{31} = 2(\epsilon_3 \epsilon_1 - \epsilon_2 \epsilon_4) \quad (40)$$

$$C_{32} = 2(\epsilon_2 \epsilon_3 + \epsilon_1 \epsilon_4) \quad (41)$$

$$C_{33} = 1 - 2\epsilon_1^2 - 2\epsilon_2^2 \quad (42)$$

Although Eqs. (1-4) and (6-24) are comparatively simple expressions when written in vector-dyadic form, complications arise when one attempts to use them to construct Eq. (26) by hand. In the first place, the process of generating Eq. (25) requires the evaluation of numerous derivatives, cross products, and dot products. To make matters worse, obtaining Eq. (26) from Eq. (25) involves the identification and isolation of 56 rather complicated subexpressions. Appendix A contains the listing of a FORMAC program that performs all of the aforementioned manipulations and produces as output the elements of the matrices X and Y

punched on cards as FORTRAN statements. The program works with the b_i ($i=1,2,3$) measure numbers of the vectors appearing in Eqs. (1-4) and (6-24). A listing of the punched output, produced by the Amdahl 470 computer at Texas A&M University via a remote terminal located at The University of Texas at Austin, may be found in Appendix B. Similar punched decks can be obtained for other spacecraft configurations by writing and implementing the corresponding FORMAC programs. In this way, one can create a "library" consisting of decks containing equations of motion for particular spacecraft of interest. Such decks can be employed directly with a linear algebraic equation solver routine and a numerical integrator to perform simulations on a digital computer.

Simulation

As a check on the correctness of the equations listed in Appendix B, a test simulation was performed using values of the system parameters chosen as follows:

m_B	= 500 kg	m_R	= 10 kg
β_1	= 110 kg-m ²	ρ_1	= 50 kg-m ²
β_2	= 100 kg-m ²	ρ_2	= 45 kg-m ²
β_3	= 90 kg-m ²	ρ_3	= 10 kg-m ²
b_1	= 3 m	r_1	= -0.1 m
b_2	= 4 m	r_2	= -0.2 m
b_3	= 5 m	r_3	= -3.5 m

The quantities F_i, T_i ($i=1,2,3$) were set equal to zero, and initial conditions for Eqs. (5) and (26-33) were taken to be

u_i	= 0	($i=1,2,3,5,6,7$)
u_4	= 1 rad/sec	
q	= 0	
x_i	= 0	($i=1,2,3$)
ϵ_i	= 0	($i=1,2,3$)
ϵ_4	= 1	

A sample of the results obtained from the simulation is presented in Fig. 2 in the form of a 100-sec time-plot of the angle θ between n_1 and b_1 . Values of θ were computed at each time step of the numerical integration from the expression [see Eq. (34)].

$$\theta = \cos^{-1}(C_{11}) \quad (43)$$

One can see from Fig. 2 that the spacecraft is performing motions in N which are oscillatory in nature.

During the simulation, computations also were made of the kinetic energy K of the spacecraft and the magnitude H of the system's central angular momentum vector. It was found that K and H retained their respective initial values of 86.4500 N-m and 365.7642 kg-m²/sec throughout the numerical integration. Since conservation of both of these quantities must follow from the conditions $F_i=0, T_i=0$ ($i=1,2,3$), the claim that the results given in Appendix B (and in Fig. 2) are correct is well founded.

Conclusion

A successful attempt has been made to derive the explicit equations of motion for a multiple-rigid-body spacecraft using symbolic manipulation. The particular configuration considered herein was chosen because its equations of motion are of small enough size to be accommodated in an appendix to the paper, and yet possess sufficient complexity to warrant the use of a computer in their derivation. Although, in this investigation, FORMAC was not applied to the analysis of systems composed of more than two rigid bodies, the prospects for doing so appear quite good, particularly since the program listed in Appendix A consumed only 27 sec of CPU time while requiring 196K bytes of computer memory in producing the equations tabulated in Appendix B.

Appendix A: FORMAC Program

```

ROTOR1: PROCEDURE OPTIONS(MAIN);
FORMAC_OPTIONS;
F_FUNCTION(SQUISH);

/*
/*
/****** FUNCTION PROCEDURE "SQUISH" *****/
/*
SQUISH: F_PROCEDURE(S);
LET(S=REPLACE(S,SIN(Q.(T)),SQ,
COS(Q.(T)),CQ));
DO JSQ=1 TO 7; LET(JSQ="JSQ");
LET(S=REPLACE(S,U(JSQ).(T),U(JSQ)));
END;
F_RETURN(S);
F_END SQUISH;

/*
/*
/****** SUBSCRIPTS FOR CROSS PRODUCTS *****/
/*
LET(P1(1)=2; P1(2)=3; P1(3)=1;
P2(1)=3; P2(2)=1; P2(3)=2);

/*
/*
/****** ASSIGNMENT OF NAMES TO *****/
/****** TIME DERIVATIVES OF U'S *****/
/****** AND TIME DERIVATIVE OF Q *****/
/*
DO J=1 TO 7; LET(J="J");
LET(DIFF(U(J))=CHAIN(U(J)));
END;
LET(DIFF(Q)=CHAIN(U(4).(T)));

/*
/*
/****** ANGULAR VELOCITY MEASURE NUMBERS *****/
/*
DO J=1 TO 3; LET(J="J");
LET(OB(J)=U(J).(T); OR(J)=U(J).(T));
END;
LET(OR(1)=OR(1)+U(4).(T));

/*
/*
/****** ANGULAR ACCELERATION MEASURE NUMBERS FOR R *****/
/*
DO J=1 TO 3; LET(J="J");
LET(ALR(J)=DERIV(OR(J),T)+OB(P1(J))*OR(P2(J))
-OB(P2(J))*OR(P1(J)));
END;

/*
/*
/****** DIRECTION COSINES *****/
/****** OF R RELATIVE TO B *****/
/*
LET(CR(1,1)=1; CR(1,2)=0; CR(1,3)=0;
CR(2,1)=0; CR(2,2)=COS(Q.(T)); CR(2,3)=-SIN(Q.(T));
CR(3,1)=0; CR(3,2)=SIN(Q.(T)); CR(3,3)=COS(Q.(T));

/*
/*
/****** MEASURE NUMBERS OF POSITION VECTOR *****/
/****** OF R* RELATIVE TO B* *****/
/*
DO J=1 TO 3; LET(J="J");
LET(P(J)=B(J)-(R(1)*CR(J,1)+R(2)*CR(J,2)+R(3)*CR(J,3)));
END;

/*
/*
/****** MEASURE NUMBERS OF VELOCITY VECTORS *****/
/****** OF MASS CENTERS OF BODIES *****/
/*
DO J=1 TO 3; LET(J="J");
LET(VB(J)=U(J+4).(T);
VR(J)=VB(J)+DERIV(P(J),T)+OB(P1(J))*P(P2(J))
-OB(P2(J))*P(P1(J)));
END;

/*
/*

```

```

/***** MEASURE NUMBERS OF ACCELERATION VECTORS *****/
/***** OF MASS CENTERS OF BODIES *****/
/*
DO J=1 TO 3; LET(J="J");
LET(AB(J)=DERIV(VB(J),T)+OB(P1(J))*VB(P2(J))
      -OB(P2(J))*VB(P1(J));
      AR(J)=DERIV(VR(J),T)+OR(P1(J))*VR(P2(J))
      -OR(P2(J))*VR(P1(J));
END;
/*
DO J=1 TO 3; LET(J="J");
LET(VB(J)=SQUISH(VB(J)); VR(J)=SQUISH(VR(J));
  AB(J)=SQUISH(AB(J)); AR(J)=SQUISH(AR(J));
END;
/*
/***** MOMENTS OF INERTIA *****/
/*
DO J=1 TO 3; LET(J="J");
DO K=1 TO 3; LET(K="K");
LET(IR(J,K)=0);
DO L=1 TO 3; LET(L="L");
LET(IR(J,K)=IR(J,K)+RHO(L)*CR(J,L)*CR(K,L));
END;
END;
/*
/***** INERTIA TORQUE MEASURE NUMBERS *****/
/*
DO J=1 TO 3; LET(J="J");
LET(TR(J)=U(P1(J))*U(P2(J))*(BETA(P1(J))-BETA(P2(J)))
      -UD(J)*BETA(J));
LET(TR(J)=0);
DO K=1 TO 3; LET(K="K");
LET(TR(J)=TR(J)+(IR(P1(J),K)*OR(K))*OR(P2(J))
      -(IR(P2(J),K)*OR(K))*OR(P1(J))
      -IR(J,K)*ALR(K));
END;
END;
/*
DO J=1 TO 3; LET(J="J");
LET(OB(J)=SQUISH(OB(J)); OR(J)=SQUISH(OR(J));
  TB(J)=SQUISH(TB(J)); TR(J)=SQUISH(TR(J));
ATOMIZE(ALR(J));
DO K=1 TO 3; LET(K="K");
ATOMIZE(IR(J,K); CR(J,K));
END;
END;
/*
/***** EQUATIONS OF MOTION *****/
/*
DO J=1 TO 7; LET(J="J");
LET(E(J)=0);
DO K=1 TO 3; LET(K="K");
LET(E(J)=E(J)+(F(K)-MB*AB(K))*DERIV(VB(K),U(J))
      -MR*AR(K)*DERIV(VR(K),U(J))
      +(T(K)+TB(K))*DERIV(OR(K),U(J))
      +TR(K)*DERIV(OR(K),U(J)));
END;
PRINT_OUT(E(J)); PUT SKIP(4);
SAVE(E(J));
END; PUT SKIP(4);
/*
DO J=1 TO 3; LET(J="J");
ATOMIZE(AB(J); AR(J); TB(J); TR(J));
END;
/*
DO J=1 TO 7; LET(J="J");
LET(XE(J)=EXPAND(E(J)));
DO K=1 TO 7; LET(K="K");
LET(X(J,K)=COEFF(XE(J),UD(K)));
LET(X(J,K)=EXPAND(REPLACE(X(J,K),CQ*CQ,1-SQ*SQ));
  X(J,K)=REPLACE(X(J,K),MR,0)+MR*COEFF(X(J,K),MR));
PRINT_OUT(X(J,K)); PUT SKIP(4);
PUNCH_F(X(J,K));
ATOMIZE(X(J,K));
END; PUT SKIP(4);

```

```

      ATOMIZE(XE(J));
      LET(Y(J)=-REPLACE(E(J),UD(1),0,UD(2),0,UD(3),0,
                        UD(4),0,UD(5),0,UD(6),0,UD(7),0));
      SAVE(Y(J));
      ATOMIZE(E(J));
      END;
/*
      DO J=1 TO 7; LET(J="J");
      PRINT_OUT(Y(J)); PUT SKIP(4);
      PUNCH_F(Y(J));
      ATOMIZE(Y(J));
      END;
/*
      END ROTOR1;

```

Appendix B: Listing of Punched Output

```

      X(1,1) = -BETA(1)-RHO(1) + MR*(CQ*B(3)*R(3)**2 + CQ*R(2)*B(2)**2 + S
      CQ*B(3)*R(2)**2-SQ*R(3)*B(2)**2-B(3)**2-R(2)**2-R(3)**2-B(2)**2)
      X(1,2) = MR*(-CQ*B(1)*R(2) + CQ*R(1)*R(2) + SQ*B(1)*R(3)-SQ*R(1)*R
      C(3) + B(1)*B(2)-R(1)*B(2))
      X(1,3) = MR*(-CQ*B(1)*R(3) + CQ*R(1)*R(3)-SQ*B(1)*R(2) + SQ*R(1)*R
      C(2) + B(3)*B(1)-B(3)*R(1))
      X(1,4) = -RHO(1) + MR*(CQ*B(3)*R(3) + CQ*R(2)*B(2) + SQ*B(3)*R(2)-
      CSQ*R(3)*R(2)-R(2)**2-R(3)**2)
      X(1,5) = 0
      X(1,6) = MR*(B(3)-CQ*R(3)-SQ*R(2))
      X(1,7) = MR*(-B(2) + CQ*R(2)-SQ*R(3))
      X(2,1) = MR*(-CQ*B(1)*R(2) + CQ*R(1)*R(2) + SQ*B(1)*R(3)-SQ*R(1)*R
      C(3) + B(1)*B(2)-R(1)*B(2))
      X(2,2) = -BETA(2)-RHO(2) + MR*(-CQ*SQ*R(2)*R(3)**2 + CQ*B(3)*R(3)**2
      C + SQ*B(3)*R(2)**2 + B(1)*R(1)**2-SQ**2*R(2)**2 + SQ**2*R(3)**2-B(3)
      C**2-B(1)**2-R(1)**2-R(3)**2)-RHO(3)*SQ**2 + RHO(2)*SQ**2
      X(2,3) = MR*(CQ*SQ*R(2)**2-CQ*SQ*R(3)**2-CQ*B(3)*R(2)-CQ*R(3)*B(2)
      C + SQ*B(3)*R(3)-SQ*R(2)*B(2) + B(3)*B(2) + R(2)*R(3)-SQ**2*R(2)*R(
      C3)**2 + RHO(3)*CQ*SQ-CQ*RHO(2)*SQ
      X(2,4) = MR*(-CQ*B(1)*R(2) + CQ*R(1)*R(2) + SQ*B(1)*R(3)-SQ*R(1)*R
      C(3))
      X(2,5) = MR*(-B(3) + CQ*R(3) + SQ*R(2))
      X(2,6) = 0
      X(2,7) = MR*(B(1)-R(1))
      X(3,1) = MR*(-CQ*B(1)*R(3) + CQ*R(1)*R(3)-SQ*B(1)*R(2) + SQ*R(1)*R
      C(2) + B(3)*B(1)-B(3)*R(1))
      X(3,2) = MR*(CQ*SQ*R(2)**2-CQ*SQ*R(3)**2-CQ*B(3)*R(2)-CQ*R(3)*B(2)
      C + SQ*B(3)*R(3)-SQ*R(2)*B(2) + B(3)*B(2) + R(2)*R(3)-SQ**2*R(2)*R(
      C3)**2 + RHO(3)*CQ*SQ-CQ*RHO(2)*SQ
      X(3,3) = -BETA(3)-RHO(3) + MR*(CQ*SQ*R(2)*R(3)**2 + CQ*R(2)*B(2)**2-
      CSQ*R(3)*R(2)**2 + B(1)*R(1)**2 + SQ**2*R(2)**2-SQ**2*R(3)**2-B(1)**2
      C-R(1)**2-R(2)**2-B(2)**2) + RHO(3)*SQ**2-RHO(2)*SQ**2
      X(3,4) = MR*(-CQ*B(1)*R(3) + CQ*R(1)*R(3)-SQ*B(1)*R(2) + SQ*R(1)*R
      C(2))
      X(3,5) = MR*(R(2)-CQ*R(2) + SQ*R(3))
      X(3,6) = MR*(-B(1) + R(1))
      X(3,7) = 0
      X(4,1) = -RHO(1) + MR*(CQ*B(3)*R(3) + CQ*R(2)*B(2) + SQ*B(3)*R(2)-
      CSQ*R(3)*R(2)-R(2)**2-R(3)**2)
      X(4,2) = MR*(-CQ*B(1)*R(2) + CQ*R(1)*R(2) + SQ*B(1)*R(3)-SQ*R(1)*R
      C(3))
      X(4,3) = MR*(-CQ*B(1)*R(3) + CQ*R(1)*R(3)-SQ*B(1)*R(2) + SQ*R(1)*R
      C(2))
      X(4,4) = -RHO(1) + MR*(-R(2)**2-R(3)**2)
      X(4,5) = 0
      X(4,6) = MR*(-CQ*R(3)-SQ*R(2))
      X(4,7) = MR*(CQ*R(2)-SQ*R(3))
      X(5,1) = 0
      X(5,2) = MR*(-B(3) + CQ*R(3) + SQ*R(2))
      X(5,3) = MR*(B(2)-CQ*R(2) + SQ*R(3))
      X(5,4) = 0
      X(5,5) = -MB-MR
      X(5,6) = 0
      X(5,7) = 0
      X(6,1) = MR*(B(3)-CQ*R(3)-SQ*R(2))
      X(6,2) = 0
      X(6,3) = MR*(-B(1) + R(1))
      X(6,4) = MR*(-CQ*B(3)-SQ*R(2))
      X(6,5) = 0
      X(6,6) = -MB-MR
      X(6,7) = 0
      X(7,1) = MR*(-B(2) + CQ*R(2)-SQ*R(3))

```

```

X(7,2) = MR*(B(1)-R(1))
X(7,3) = 0
X(7,4) = MR*(CQ*R(2)-SQ*R(3))
X(7,5) = 0
X(7,6) = 0
X(7,7) = -MB-MR

```

```

Y(1) = -T(1) + MR*(B(2)-CQ*R(2) + SQ*R(3))*(CQ*R(3)*U(4)**2 + SQ*R
C(2)*U(4)**2 + U(1)*(U(6) + CQ*R(3)*U(4) + SQ*R(2)*U(4)-U(1)*(B(3)-
CCQ*R(3)-SQ*R(2)) + U(3)*(B(1)-R(1))) + U(1)*(CQ*R(3)*U(4) + SQ*R(2
C)*U(4))-U(2)*(U(5) + U(2)*(B(3)-CQ*R(3)-SQ*R(2))-U(3)*(B(2)-CQ*R(2
C) + SQ*R(3))) + MR*(-B(3) + CQ*R(3) + SQ*R(2))*(CQ*R(2)*U(4)**2-S
CQ*R(3)*U(4)**2-U(1)*(U(7)-CQ*R(2)*U(4) + SQ*R(3)*U(4) + U(1)*(B(2)
C-CQ*R(2) + SQ*R(3))-U(2)*(B(1)-R(1))-U(1)*(-CQ*R(2)*U(4) + SQ*R(3
C)*U(4)) + U(3)*(U(5) + U(2)*(B(3)-CQ*R(3)-SQ*R(2))-U(3)*(B(2)-CQ*R
C(2) + SQ*R(3))) -U(2)*U(3)*(BETA(2)-BETA(3)) + U(2)*U(3)*(RHO(3)*C
CQ**2 + RHO(2)*SQ**2)-U(2)*U(3)*(RHO(3)*SQ**2 + CQ**2*RHO(2)) + U(2
C)**2*(-RHO(3)*CQ*SQ + CQ*RHO(2)*SQ)-U(3)**2*(-RHO(3)*CQ*SQ + CQ*RH
CO(2)*SQ)

```

```

Y(2) = -T(2) + MR*(B(3)-CQ*R(3)-SQ*R(2))*(U(2)*(U(7)-CQ*R(2)*U(4)
C+ SQ*R(3)*U(4) + U(1)*(B(2)-CQ*R(2) + SQ*R(3))-U(2)*(B(1)-R(1))) +
C U(2)*(-CQ*R(2)*U(4) + SQ*R(3)*U(4))-U(3)*(U(6) + CQ*R(3)*U(4) + S
CQ*R(2)*U(4)-U(1)*(B(3)-CQ*R(3)-SQ*R(2)) + U(3)*(B(1)-R(1))-U(3)*(
CCQ*R(3)*U(4) + SQ*R(2)*U(4))) + MR*(-B(1) + R(1))*(CQ*R(3)*U(4)**2
C + SQ*R(2)*U(4)**2 + U(1)*(U(6) + CQ*R(3)*U(4) + SQ*R(2)*U(4)-U(1)
C*(B(3)-CQ*R(3)-SQ*R(2)) + U(3)*(B(1)-R(1))) + U(1)*(CQ*R(3)*U(4) +
C SQ*R(2)*U(4))-U(2)*(U(5) + U(2)*(B(3)-CQ*R(3)-SQ*R(2))-U(3)*(B(2)
C-CQ*R(2) + SQ*R(3))) + RHO(1)*U(3)*(U(1) + U(4))-U(1)*U(3)*(BETA(
C3)-BETA(1))-U(2)*(U(1) + U(4))*(-RHO(3)*CQ*SQ + CQ*RHO(2)*SQ)-U(3)
C*(U(1) + U(4))*(RHO(3)*CQ**2 + RHO(2)*SQ**2) + (-RHO(3)*CQ*SQ + CQ
C*RHO(2)*SQ)*(U(1)*U(2)-U(2)*(U(1) + U(4))) + (-U(1)*U(3) + U(3)*(U
C(1) + U(4)))*(RHO(3)*SQ**2 + CQ**2*RHO(2))

```

```

Y(3) = -T(3) + MR*(B(1)-R(1))*(CQ*R(2)*U(4)**2-SQ*R(3)*U(4)**2-U(1
C)*(U(7)-CQ*R(2)*U(4) + SQ*R(3)*U(4) + U(1)*(B(2)-CQ*R(2) + SQ*R(3)
C)-U(2)*(B(1)-R(1))-U(1)*(-CQ*R(2)*U(4) + SQ*R(3)*U(4)) + U(3)*(U(
C5) + U(2)*(B(3)-CQ*R(3)-SQ*R(2))-U(3)*(B(2)-CQ*R(2) + SQ*R(3))) +
C MR*(-B(2) + CQ*R(2)-SQ*R(3))*(U(2)*(U(7)-CQ*R(2)*U(4) + SQ*R(3)*U
C(4) + U(1)*(B(2)-CQ*R(2) + SQ*R(3))-U(2)*(B(1)-R(1))) + U(2)*(-CQ*
CR(2)*U(4) + SQ*R(3)*U(4))-U(3)*(U(6) + CQ*R(3)*U(4) + SQ*R(2)*U(4)
C-U(1)*(B(3)-CQ*R(3)-SQ*R(2)) + U(3)*(B(1)-R(1))-U(3)*(CQ*R(3)*U(4)
C) + SQ*R(2)*U(4))-RHO(1)*U(2)*(U(1) + U(4))-U(1)*U(2)*(-BETA(2) +
C BETA(1)) + U(2)*(U(1) + U(4))*(RHO(3)*SQ**2 + CQ**2*RHO(2)) + U(3
C)*(U(1) + U(4))*(-RHO(3)*CQ*SQ + CQ*RHO(2)*SQ) + (-RHO(3)*CQ*SQ +
CCQ*RHO(2)*SQ)*(-U(1)*U(3) + U(3)*(U(1) + U(4))) + (RHO(3)*CQ**2 +
CRHO(2)*SQ**2)*(U(1)*U(2)-U(2)*(U(1) + U(4)))

```

```

Y(4) = MR*(-CQ*R(2) + SQ*R(3))*(CQ*R(3)*U(4)**2 + SQ*R(2)*U(4)**2
C+ U(1)*(U(6) + CQ*R(3)*U(4) + SQ*R(2)*U(4)-U(1)*(B(3)-CQ*R(3)-SQ*R
C(2)) + U(3)*(B(1)-R(1))) + U(1)*(CQ*R(3)*U(4) + SQ*R(2)*U(4))-U(2)
C*(U(5) + U(2)*(B(3)-CQ*R(3)-SQ*R(2))-U(3)*(B(2)-CQ*R(2) + SQ*R(3)
C)) + MR*(CQ*R(2)*U(4)**2-SQ*R(3)*U(4)**2-U(1)*(U(7)-CQ*R(2)*U(4) +
C SQ*R(3)*U(4) + U(1)*(B(2)-CQ*R(2) + SQ*R(3))-U(2)*(B(1)-R(1))-U(
C1)*(-CQ*R(2)*U(4) + SQ*R(3)*U(4)) + U(3)*(U(5) + U(2)*(B(3)-CQ*R(3)
C)-SQ*R(2))-U(3)*(B(2)-CQ*R(2) + SQ*R(3)))*(CQ*R(3) + SQ*R(2)) + U
C(2)*U(3)*(RHO(3)*CQ**2 + RHO(2)*SQ**2)-U(2)*U(3)*(RHO(3)*SQ**2 + C
CQ**2*RHO(2)) + U(2)**2*(-RHO(3)*CQ*SQ + CQ*RHO(2)*SQ)-U(3)**2*(-RH
CO(3)*CQ*SQ + CQ*RHO(2)*SQ)

```

```

Y(5) = -F(1) + MB*(-U(6)*U(3) + U(7)*U(2)) + MR*(U(2)*(U(7)-CQ*R(2
C)*U(4) + SQ*R(3)*U(4) + U(1)*(B(2)-CQ*R(2) + SQ*R(3))-U(2)*(B(1)-R
C(1))) + U(2)*(-CQ*R(2)*U(4) + SQ*R(3)*U(4))-U(3)*(U(6) + CQ*R(3)*U
C(4) + SQ*R(2)*U(4)-U(1)*(B(3)-CQ*R(3)-SQ*R(2)) + U(3)*(B(1)-R(1)))
C-U(3)*(CQ*R(3)*U(4) + SQ*R(2)*U(4)))

```

```

Y(6) = -F(2) + MB*(U(5)*U(3)-U(7)*U(1)) + MR*(CQ*R(2)*U(4)**2-SQ*R
C(3)*U(4)**2-U(1)*(U(7)-CQ*R(2)*U(4) + SQ*R(3)*U(4) + U(1)*(B(2)-CQ
C*R(2) + SQ*R(3))-U(2)*(B(1)-R(1))-U(1)*(-CQ*R(2)*U(4) + SQ*R(3)*U
C(4)) + U(3)*(U(5) + U(2)*(B(3)-CQ*R(3)-SQ*R(2))-U(3)*(B(2)-CQ*R(2)
C + SQ*R(3)))

```

```

Y(7) = -F(3) + MB*(-U(5)*U(2) + U(6)*U(1)) + MR*(CQ*R(3)*U(4)**2 +
C SQ*R(2)*U(4)**2 + U(1)*(U(6) + CQ*R(3)*U(4) + SQ*R(2)*U(4)-U(1)*(
CB(3)-CQ*R(3)-SQ*R(2)) + U(3)*(B(1)-R(1))) + U(1)*(CQ*R(3)*U(4) + S
CQ*R(2)*U(4))-U(2)*(U(5) + U(2)*(B(3)-CQ*R(3)-SQ*R(2))-U(3)*(B(2)-C
CQ*R(2) + SQ*R(3)))

```


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