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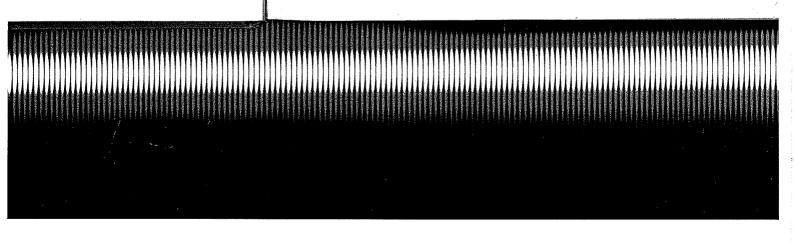
Solving Ordinary Differential Equations II

Stiff and Differential-Algebraic Problems

Second Revised Edition With 137 Figures



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VII.7 Computation of Multibody Mechanisms

Dynamics of multibody systems is of great importance in the fields of robotics, biomechanics, spacecraft control, road and rail vehicle design, and dynamics of machinery.

(W. Schiehlen 1990)

After having seen several different approaches for the numerical solution of constrained mechanical systems, we are interested in their efficiency when applied to a concrete situation. We consider two particular multibody mechanisms with constraints, one nonstiff and one stiff. General references for the computation of mechanical systems are Haug (1989) and Roberson & Schwertassek (1988).

Description of the Model

We first consider "Andrews' squeezer mechanism", which has become prominent through the work of Giles (1978) and Manning (1981), who promoted it as a test example for numerical codes; see also Ormrod & Andrews (1986). It consists of 7 rigid bodies connected by joints without friction in plane motion. It is represented in Fig. 7.1, which we have copied (with permission) from the book of Schiehlen (1990). The numerical constants, also taken from Schiehlen (1990), are displayed in Tables 7.1 and 7.2. The arrows in the right picture of Fig. 7.1 indicate the positions of the centres of gravity $C_1, \ldots, C_{\mathbf{f}}$. In Table 7.1 the spring coefficient of the spring connecting the point D with C is denoted by c_0 and the unstretched length is ℓ_0 . We suppose that the mechanism is driven by a motor, located at D, whose constant drive torque is given by mom = 0.033. The coordinate origin is the point D in Fig. 7.1 and the coordinates of the other fixed points A, B and C are given by

$$\begin{pmatrix} xa \\ ya \end{pmatrix} = \begin{pmatrix} -0.06934 \\ -0.00227 \end{pmatrix}, \begin{pmatrix} xb \\ yb \end{pmatrix} = \begin{pmatrix} -0.03635 \\ 0.03273 \end{pmatrix}, \begin{pmatrix} xc \\ yc \end{pmatrix} = \begin{pmatrix} 0.014 \\ 0.072 \end{pmatrix}. (7.1)$$

Pable 7.1. Geometrical parameters

d = 0.028	da = 0.0115	e = 0.02	
ea = 0.01421	f = 0.02	$f_2 = 0.01421$	
10:0	70:0	Ja 0.01441	
r = 0.007	ra = 0.00092	ss = 0.035	
sa = 0.01874	sb = 0.01043	sc = 0.018	
sd = 0.02	zt = 0.04	ta = 0.02308	
tb = 0.00916	u = 0.04	ua = 0.01228	
ub = 0.00449	$c_0 = 4530$	$\ell_0 = 0.07785$	

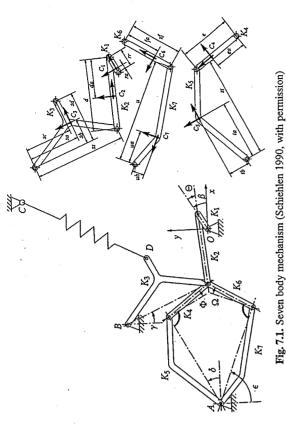


Table 7.2. Parameters of the 7 bodies

No.	masses m ₁ to m ₇	inertias I_1 to I_7
-	0.04325	$2.194 \cdot 10^{-6}$
7	0.00365	$4.410 \cdot 10^{-7}$
3	0.02373	5.255 · 10-6
4	0.00706	$5.667 \cdot 10^{-7}$
2	0.07050	1.169.10 ⁻⁵
9	0.00706	5.667 · 10 ⁻⁷
7	0.05498	$1.912 \cdot 10^{-5}$

In order to derive the equations of motion we use the angles (see Fig. 7.1)

$$q_1 = \beta, \quad q_2 = \Theta, \quad q_3 = \gamma, \quad q_4 = \Phi, \quad q_5 = \delta, \quad q_6 = \Omega, \quad q_7 = \varepsilon, \quad (7.2)$$

as position coordinates for the mechanical system. If (x_j, y_j) are the cartesian coordinates of the centre of gravity C_j (j = 1, ..., 7), the kinetic energy of the multibody system is

$$T = \sum_{j=1}^{7} m_j \frac{x_j^2 + y_j^2}{2} + \sum_{j=1}^{7} I_j \frac{\omega_j^2}{2}$$
 (7.3)

where ω_j is the total angle of rotation of the j th body and m_j , I_j are constants given in Table 7.2. The values of x_j, y_j , $\dot{x}_j^2 + \dot{y}_j^2$ and $\dot{\omega}_j$ can be obtained in terms

of (7.2) by simple geometry (see Fig. 7.1): 532

$$C_1: x_1 = ra \cdot \cos \beta$$

$$y_1 = ra \cdot \sin \beta$$

$$\dot{x}_1^2 + \dot{y}_1^2 = ra^2 \cdot \dot{\beta}^2$$

$$\dot{\omega}_1 = \dot{\beta}$$

$$C_2: x_2 = rr \cdot \cos \beta - da \cdot \cos(\beta + \Theta)$$

$$y_2 = rr \cdot \sin \beta - da \cdot \sin(\beta + \Theta)$$

$$\dot{x}_2^2 + \dot{y}_2^2 = (rr^2 - 2 \cdot da \cdot rr \cdot \cos \Theta + da^2) \cdot \dot{\beta}^2$$

$$+ 2 \cdot (-rr \cdot da \cdot \cos \Theta + da^2) \cdot \dot{\beta} \cdot \dot{\Theta} + da^2 \cdot \dot{\Theta}^2$$

$$\dot{\omega}_2 = \dot{\beta} + \dot{\Theta}$$

$$C_{3}: x_{3} = xb + sa \cdot \sin \gamma + sb \cdot \cos \gamma$$

$$y_{3} = yb - sa \cdot \cos \gamma + sb \cdot \sin \gamma$$

$$\dot{x}_{3}^{2} + \dot{y}_{2}^{2} = (sa^{2} + sb^{2}) \cdot \dot{\gamma}^{2}$$

$$\dot{\omega}_{3} = \dot{\gamma}$$

$$C_{4}: x_{4} = xa + zt \cdot \cos \delta + (e - ea) \cdot \sin(\Phi + \delta)$$

$$y_{4} = ya + zt \cdot \sin \delta - (e - ea) \cdot \cos(\Phi + \delta)$$

$$\dot{x}_{4}^{2} + \dot{y}_{4}^{2} = (e - ea)^{2} \cdot \dot{\Phi}^{2} + 2 \cdot ((e - ea)^{2} + zt \cdot (e - ea) \cdot \sin \Phi) \cdot \dot{\Phi} \cdot \dot{\delta}$$

$$+ (zt^{2} + 2 \cdot zt \cdot (e - ea) \cdot \sin \Phi + (e - ea)^{2}) \cdot \dot{\delta}^{2}$$

$$C_5: x_5 = xa + ta \cdot \cos \delta - tb \cdot \sin \delta$$

$$y_5 = ya + ta \cdot \sin \delta + tb \cdot \cos \delta$$

$$\dot{x}_5^2 + \dot{y}_5^2 = (ta^2 + tb^2) \cdot \dot{\delta}^2$$

$$\dot{\omega}_5 = \dot{\delta}$$

$$\begin{split} C_6: & \quad x_6 = xa + u \cdot \sin \varepsilon + (xf - fa) \cdot \cos(\Omega + \varepsilon) \\ & \quad y_6 = ya - u \cdot \cos \varepsilon + (xf - fa) \cdot \sin(\Omega + \varepsilon) \\ & \quad \dot{x}_6^2 + \dot{y}_6^2 = (xf - fa)^2 \cdot \dot{\Omega}^2 + 2 \cdot \left((xf - fa)^2 - u \cdot (xf - fa) \cdot \sin \Omega \right) \cdot \dot{\Omega} \cdot \dot{\varepsilon} \\ & \quad + \left((xf - fa)^2 - 2 \cdot u \cdot (xf - fa) \cdot \sin \Omega + u^2 \right) \cdot \dot{\varepsilon}^2 \\ & \quad \dot{\omega}_6 = \dot{\Omega} + \dot{\varepsilon} \end{split}$$

$$C_7: x_7 = xa + ua \cdot \sin \varepsilon - ub \cdot \cos \varepsilon$$

$$y_7 = ya - ua \cdot \cos \varepsilon - ub \cdot \sin \varepsilon$$

$$\dot{x}_7^2 + \dot{y}_7^2 = (ua^2 + ub^2) \cdot \dot{\varepsilon}^2$$

$$\dot{\omega}_7 = \dot{\varepsilon}$$

The potential energy of the system is due to the motor at the origin and to the spring

connecting the point D with C. By Hooke's law it is

$$U = -mom \cdot \beta + c_0 \frac{(\ell - \ell_0)^2}{2},$$
 (7.4)

where ℓ is the distance between D and C, namely

$$\ell = \sqrt{(xd - xc)^2 + (yd - yc)^2}$$

$$xd = xb + sc \cdot \sin \gamma + sd \cdot \cos \gamma$$

$$yd = yb - sc \cdot \cos \gamma + sd \cdot \sin \gamma.$$

three loops. The first loop connects O with B via K_1, K_2, K_3 ; the other two loops Finally, we have to formulate the algebraic constraints. The mechanism contains connect O with A, one via K_1, K_2, K_4, K_5 , the other via K_1, K_2, K_6, K_7 . For each loop we get two algebraic conditions:

$$rr \cdot \cos \beta - d \cdot \cos(\beta + \Theta) - ss \cdot \sin \gamma = xb$$

$$rr \cdot \sin \beta - d \cdot \sin(\beta + \Theta) + ss \cdot \cos \gamma = yb$$

$$rr \cdot \cos \beta - d \cdot \cos(\beta + \Theta) - e \cdot \sin(\Phi + \delta) - zt \cdot \cos \delta = xa$$

$$rr \cdot \sin \beta - d \cdot \sin(\beta + \Theta) + e \cdot \cos(\Phi + \delta) - zt \cdot \sin \delta = ya$$

$$rr \cdot \cos \beta - d \cdot \cos(\beta + \Theta) - zf \cdot \cos(\Omega + \varepsilon) - u \cdot \sin \varepsilon = xa$$

$$rr \cdot \sin \beta - d \cdot \sin(\beta + \Theta) - zf \cdot \sin(\Omega + \varepsilon) + u \cdot \cos \varepsilon = ya.$$

With the position coordinates q from (7.2) the equations (7.5) represent the constraint g(q)=0 where $g:\mathbb{R}^7\to\mathbb{R}^6$. Together with the kinetic energy T of (7.3) the potential energy U of (7.4) and $L=T-U-\lambda_1g_1-\ldots-\lambda_6g_6$ the equations of motion (1.46) are fully determined.

Fortran Subroutines

routines describing the differential-algebraic problem. The equations of motion are For the reader's convenience we include the essential parts of the FORTRAN sub-

$$M(q)\ddot{q} = f(q, \dot{q}) - G^{T}(q)$$

$$(7.6a)$$

$$0 = q(q)$$

$$(7.6b)$$

tion the variables $\mathbb{Q}(1), \ldots, \mathbb{Q}(7)$ correspond to $\beta, \ldots, \varepsilon$ (exactly as in (7.2)) and $\mathbb{QP}(1), \ldots, \mathbb{QP}(7)$ to their derivatives $\beta, \ldots, \varepsilon$. In all subroutines we have used where $q \in \mathbb{R}^7$ is the vector defined in (7.2) and $\lambda \in \mathbb{R}^6$. In the following descriphe abbreviations

SIBE = SIN
$$(q(1))$$
 COBE = COS $(q(1))$
SITH = SIN $(q(2))$ COTH = COS $(q(2))$
SIGA = SIN $(q(3))$ COGA = COS $(q(3))$

534

```
SIPH = SIN (q(4))

SIDE = SIN (q(5))

SIDE = SIN (q(5))

SIDE = SIN (q(6))

SIETH = SIN (q(7))

SIBETH = SIN (q(1)+q(5))

SIPHUE = SIN (q(4)+q(5))

SIPHUE = SIN (q(4)+q(5))

SIPHUE = SIN (q(6)+q(7))

SIPHUE = COS (q(4))

COBETH = COS (q(7))

COPHUE = COS (q(7))

COPHUE = COS (q(6))

COPHUE = COS (q(6))
```

The remaining parameters XA, YA,..., D, DA, E, EA,..., M1, I1, M2,... are those of (7.1) and Tables 7.1 and 7.2. They usually reside in a COMMON block. The elements of M(q) in (7.6) are given by

$$m_{ij} = \frac{\partial^2 L}{\partial \dot{q}_i \partial \dot{q}_j} = \frac{\partial^2 T}{\partial \dot{q}_i \partial \dot{q}_i}.$$

This matrix is symmetric and (due to the special arrangement of the coordinates) tridiagonal. The non-zero elements (on and below the diagonal) are

The *i*th component of the function f in (7.6) is defined by

$$f_i(q,\dot{q}) = \frac{\partial (T-U)}{\partial q_i} - \sum_{j=1}^7 \frac{\partial^2 (T-U)}{\partial \dot{q}_i \partial q_j} \cdot \dot{q}_j.$$

Written as FORTRAN statements we have

```
XD = SD*COGA + SC*SIGA + XB
YD = SD*SIGA - SC*COGA + YB
LANG = SQRT ((XD-XC)**2 + (YD-YC)**2)
FORCE = - CO * (LANG - LO)/LANG
FX = FORCE * (XD-XC)
FY = FORCE * (XD-XC)
FY = FORCE * (YD-YC)
F(1) = MOM - M2*DA*RR*THP* (THP+2*BEP)*SITH
F(2) = M2*DA*RR*BEP*2*SITH
F(2) = M2*CA*RR*BEP*2*SITH
F(3) = FX*GS*COGA - SD*SIGA) + FY*(SD*COGA + SC*SIGA)
F(4) = M4*ZT*(E-EA)*DEP*2*COPH
F(5) = - M4*ZT*(E-EA)*PHP*(PHP+2*DEP)*COPH
F(6) = - M6*U*(ZF-FA)*EPP*2*COM
F(7) = M6*U*(ZF-FA)*CMP*(OMP+2*EPP)*COOM
```

The algebraic constraints g(q) = 0 are given by the following six equations (see (7.5))

```
G(1) = RR*CDEE - D*COBETH - SS*SIGA - XB
G(2) = RR*SIBE - D*SIBETH + SS*COGA - YB
G(3) = RR*COBE - D*COBETH - ESSIPHDE - ZT*CODE - XA
G(4) = RR*SIBE - D*SIBETH + E*COPHDE - ZT*SIDE - YA
G(5) = RR*COBE - D*COBETH - ZF*COOMEP - U*SIEP - XA
G(6) = RR*SIBE - D*SIBETH - ZF*SIOMEP + U*COEP - YA
```

And here is the Jacobian matrix $G(q)=g_q(q)$. The non-zero entries of this 6×7 array are

```
      QQ(1,1) = - RR*SIBE + D*SIBETH
      QQ(4,2) = - D*COBETH

      QQ(1,2) = D*SIBETH
      QQ(4,4) = - E*SIPHDE

      QQ(2,1) = - SS*COGA
      QQ(6,1) = - RR*SIBE + D*SIBETH

      QQ(2,2) = - D*COBETH
      QQ(5,1) = - RR*SIBE + D*SIBETH

      QQ(2,3) = - SS*SIGA
      QQ(5,2) = D*SIBETH

      QQ(2,2) = - RR*SIBE + D*SIBETH
      QQ(5,2) = D*SIBETH

      QQ(3,2) = - RR*SIBE + D*SIBETH
      QQ(5,2) = D*SIBETH

      QQ(3,4) = - E*COPHDE
      QQ(6,1) = RR*COBE - D*COBETH

      QQ(3,4) = - E*COPHDE
      QQ(6,2) = - D*COBETH

      QQ(3,4) = - E*COPHDE
      QQ(6,2) = - D*COBETH

      QQ(3,4) = - E*COPHDE
      QQ(6,2) = - CF*COOMEP

      QQ(3,4) = - E*COPHDE
      QQ(6,2) = - CF*COOMEP
```

If we apply a numerical method to the index 1 formulation of the system, we also need the expression $g_{qq}(q)(\dot{q},\dot{q})$. It is given by

```
GQQ(1) = - RR*COBE*V(1)**2 + D*COBETH*(V(1)+V(2))**2 +

SS*SIGA*V(3)**2 +

GQQ(2) = - RR*SIBE*V(1)**2 + D*SIBETH*(V(1)+V(2))**2 -

* SS*COGA*V(3)**2 +

GQQ(3) = - RR*COBE*V(1)**2 + D*COBETH*(V(1)+V(2))**2 +

* E*SIPHDE*(V(4)+V(5))**2 + ZT*SUDE*V(5)**2 +

GQQ(4) = - RR*SIBE*V(1)**2 + D*SIBETH*(V(1)+V(2))**2 +

* E*COPHDE*(V(4)+V(5))**2 + ZT*SIDE*V(5)**2 +

* GQQ(5) = - RR*COBE*V(1)**2 + D*COBETH*(V(1)+V(2))**2 +

* ZF*COOMEF*(V(4)+V(7))**2 + U*SIDE*V(7)**2 +

* GQQ(6) = - RR*SIBE*V(1)**2 + D*SIBETH*(V(1)+V(2))**2 +

* ZF*COMEF*(V(6)+V(7))**2 - U*SIBETY(7)**2 +

* ZF*COMEF*(V(6)+V(7))**2 - U*COBE*V(7)**2 +

* ZF*SIBE*V(1)**2 + D*SIBETH*(V(1)+V(2))**2 +

* ZF*SIBE*V(1)**2 - U*COBE*V(7)**2 +

* ZF*SIBE*V(7)**2 - U*COBE*V(7)**2 +

* ZF*SIBE*V(7)**2
```

Computation of Consistent Initial Values

We first compute a solution of g(q)=0. Since g consists of 6 equations in 7 unknowns we can fix one of them arbitrarily, say $\Theta(0)=0$, and compute the remaining coordinates by Newton iterations. This gives

```
\begin{split} \beta(0) &= -0.0617138900142764496358948458001\\ \gamma(0) &= 0.455279819163070380255912382449\\ \Phi(0) &= 0.222668390165885884674473185609\\ \delta(0) &= 0.487364979543842550225598953530\\ \Omega(0) &= -0.222668390165885884674473185609\\ \varepsilon(0) &= 1.23054744454982119249735015568. \end{split}
```

536

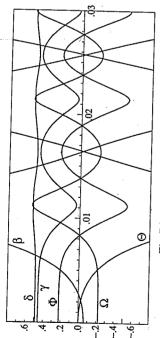


Fig. 7.2. Solution of 7 body mechanism

The condition G(q)q = 0 is satisfied if we put

$$\dot{\beta}(0) = \dot{\Theta}(0) = \dot{\gamma}(0) = \dot{\Phi}(0) = \dot{\delta}(0) = \dot{\Omega}(0) = \dot{\varepsilon}(0) = 0.$$
 (7.8)

The values of $\lambda(0)$ and $\ddot{q}(0)$ are then uniquely determined by (7.6a) and the twice differentiated constraint $0 = g_{qq}(q)(\dot{q},\dot{q}) + G(q)\ddot{q}$. We just have to solve a linear system with the matrix

$$\begin{pmatrix} M(q) & G^T(q) \\ G(q) & 0 \end{pmatrix}. \tag{7.9}$$

Observe that g_{qq} need not be evaluated, because $\dot{q}(0)=0$. Due to the choice $\Theta(0) = 0$ most components of $\lambda(0)$ and $\ddot{q}(0)$ vanish. Only the first two of these are different from zero and given by

$$\ddot{\beta}(0) = 14222.4439199541138705911625887$$

$$\ddot{\Theta}(0) = -10666.8329399655854029433719415$$

$$\lambda_1(0) = 98.5668703962410896057654982170$$
 (7.10)

The solution of this seven body mechanism is plotted (mod 2π) in Fig. 7.2 for $0 \le t \le 0.03$

 $\lambda_2(0) = -6.12268834425566265503114393122.$

Numerical Computations

We first transform (7.6) into a first order system by introducing the new variable $v=\dot{q}$. Our codes apply only to problems where the derivative is multiplied by a constant matrix. We therefore also consider $w=\ddot{q}$ as a variable so that (7.6a) becomes an algebraic relation. The various formulations of the problem, as discussed in Sect. VII.1, are now as follows:

VII.7 Computation of Multibody Mechanisms

Index 3 Formulation. With $v = \dot{q}$ and $w = \ddot{q}$ the system (7.6) can be written as

$$\begin{aligned} \dot{q} &= v & (7.11a) \\ \dot{v} &= w & (7.11b) \end{aligned}$$

$$0 = M(q)w - f(q, v) + G^{T}(q)\lambda$$
 (7.11c)

$$0 = g(q).$$
 (7.11d)

$$0 = g(q)$$
.

Index 2 Formulation. If we differentiate 0=g(q) once and replace (7.11d) by

0 = G(q)v,

(7.11e)

Index 1 Formulation. One more differentiation of (7.11e) yields

$$0 = g_{qq}(q)(v,v) + G(q)w,$$

(7.11f)

so that (7.11a,b,c,f) constitutes an index 1 problem.

We have applied several codes with many different tolerances between 10-2 and 10⁻¹⁰ to these formulations. The results are given in Fig. 7.3. We have plotted the computing time (on a SUN Spark 20 workstation) against the error of the (q, v)components at $x_{end} = 0.03$ (in double logarithmic scale).

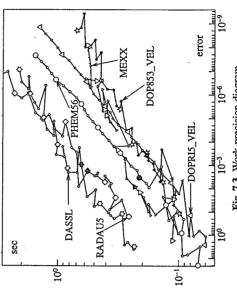


Fig. 7.3. Work-precision diagram

plicit methods such as DOPRI5 or DOP853 of Volume I. For this we have written a Explicit Runge-Kutta Methods. The index 1 formulation allows us to apply exfunction subroutine which solves in each call the linear system (7.11c,f) for w and λ and inserts the result into (7.11a,b). Since there is no stiffness in the obtained