# LagranTexPac: A Software Tool to Obtain the Dynamic Equations of Mechanical Systems

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Abstract—Modeling of a dynamical system is an important research area to obtain mathematical models of real systems. There are several approaches to get mathematical models, one of them is utilizing the Lagrangian of holonomic systems. This paper introduces LagranTexPac, a package Matlab-based framework aimed at obtaining, and simulate the equations of motion of a mechanical system. LagranTexPac uses the Lagrangian function of a system to automatically generate the system equations of motion for numerically simulate and saves all phase portraits. Motion constraints can be included in Lagrangian, and a summary report can be generated. Utility of this software tool includes any engineer area, in which is used a Lagrangian function to obtain the equations of motion of a dynamic system, e.g., aerospace, biomedical, robotics and mechanical engineering. Three application examples are included to illustrate the usefulness of this software tool.

*Index Terms*—Dynamical systems, contact dynamics, matlab®, code generation, LATEX report generation, simulation tool.

#### I. INTRODUCTION

Dynamical systems are studied in several areas of engineering, in which system modeling is necessary to study and analyze the system behaviors. Newton's laws can be used to model dynamical systems and the development procedure may be complex. An alternative approach used to model systems is the Lagrangian function, wherewith the Lagrangian equations are obtained. Last approach can be less complex than the first one

A few tool packages are available to obtain the Lagrangian equations in Matlab framework. Some years ago, there was no software package to help in the procedure to obtain the Lagrangian equations of dynamical systems. To derive equations of the dynamical system is a fatigued procedure.

In [1], [2] were shown Matlab® toolboxes that resolve Euler- Lagrange equation and provide ODE equations for numerically simulation of a system. Those toolboxes do not interact with LaTeX. In [3] was reported a package to generate some LaTeX features to use with Matlab<sup>®</sup>, in [4] was viewed a similar technique used to create a simple LaTeX files directly from Matlab<sup>®</sup> and in [5] was used a package to format the Matlab<sup>®</sup> code syntax inside a LATeX document.

Lagrangian is a mathematical function, which is a function of the generalized coordinates. These coordinates and their

time derivatives contain information about the dynamics of the system. The dynamical model of a generic system can be described as follows:

$$\dot{\mathbf{y}} = f\left(\mathbf{y}\right) \tag{1}$$

where  $y \in \mathbb{R}^n$  is a vector state and  $f: U \subseteq \mathbb{R}^n \to \mathbb{R}^n$  is a vector field and describes an autonomous system [6]-[8].

In [9]-[11] were introduced different approaches to determine differential equations of the system dynamics. Examples about robot manipulators can be found in [10], [12], in [7], [8] were shown examples about nonlinear dynamical systems, in [13], [14] were illustrated examples about triple-pendulum arm and constrain to the movement for a mechanical system, respectively.

In the triple-pendulum context, [13], [15] show the fatigued work to obtain differential equations of dynamic system, and can be seen some development mistakes like the forgotten term  $m_2 z_2^2$ , in [13] at the mass matrix. This was one of motivations to develop the this tool to avoid writing mistakes or forgetting mistakes, and another motivation was to provide a file with the differential equations to simulate in Matlab<sup>®</sup> framework.

An important feature of the tool is to implement equations of the system dynamics taking into account constraints to the motion, which is part of the contacts dynamics and this topic was not considered in [1], [2]. Studied system in [14] was considered as an application of this feature of tool. Another feature of tool is to produce a summarized report of all development steps to obtain the differential equations of the system dynamics, and the report is generated to be used in a LATEX framework. The command used to write a line in a LATEX file is reported in [4], in this tool it is created a function to write the whole report selecting the language between English, German, Spanish or Potuguese.

Therefore, the aim of Matlab<sup>®</sup> tool is to make possible a fast analysis and simulation of dynamic equations with or without constraints to the motion, avoiding written mistake and forgotten mistakes and to make possible checking intermediate steps.

This article is organized as follows: in Section II is shown a summary of the Euler-Lagrange's equation of motion; in Section III is briefly designed the development of the tool; in Section IV is shown three application examples and finally in Section V is shown conclusions.

#### II. EULER-LAGRANGE'S EQUATIONS OF MOTION.

Dynamical systems can be described by a set of simultaneous differential equations known as Lagrange's equations according to [6], [9] and the system dynamics can be defined

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as:

$$\frac{d}{dt}\left(\frac{\partial L(q_i, \dot{q}_i)}{\partial \dot{q}_i}\right) - \frac{\partial L(q_i, \dot{q}_i)}{\partial q_i} + \frac{\partial R(q_i, \dot{q}_i)}{\partial \dot{q}_i} = \tau_i; \forall i \ (2)$$

where *i* is the index for all generalized coordinates  $q_i$  and  $\tau_i$  are generalized forces; *R* is the Rayleigh dissipation function, and the Lagrangian function is defined as follows:

$$L = T - V \tag{3}$$

where T is the kinetic energy, and V is the potential for the free system, as well as for the constrained system [14]:

$$L = T - V + \Lambda \tag{4}$$

with

$$\Lambda = \sum_{i=1}^{P} \lambda_i \,\phi_i \tag{5}$$

for the system with constraints,  $\lambda_i$  is the Lagrange multiplicators and  $\varphi_i$  are the P constraint equations of the motion. In this context, when there are constraints motion the tool package analyze the free, the constraints and contact dynamics separately. The energy equations for a generic dynamical system are defined as follows:

$$2T = T_1 + T_2$$
  
 $T_1 = 2$  times rotation kinetic energy  
 $T_2 = 2$  times translation kinetic energy (6)  
 $V =$  All kinds of potential energy  
 $R =$  All kinds of dissipation

To avoid the division by two at kinetics energies was chosen this form of these energies; and did not have division on potential energies, and the gravitational term have to be included.

Applying Euler-Lagrange equation from Eq. (2) at Lagrangian function L from Eq. (3) results in a set of n second order differential equations [9], which can be rewritten in the matrix notation form [9], [10] and are given by:

$$M \cdot q + \underbrace{B(\underline{q} \ \underline{q})q}_{\underline{c}} + G(q) = \underline{\tau}$$
(7)

where  $M \in \mathbb{R}^{n \times n}$ ;  $\underline{C}, \underline{G}, \underline{\tau} \in \mathbb{R}^{n \times l}$ .

## III. STRUCTURE AND FLOW DIAGRAM OF TOOLBOX

Desired terms of kinetic energy must be entered as  $T(l), \forall l = \{1,2\}, e.g., 2T = m_1 \hat{\gamma}_l^2$  is needed to enter "T(1) = m\_1 \* dq(1)^2", with the variable, *num\_coord\_gen\_qq*, which previously define for the correct dimension of q(1) and *len\_T* for the correct dimension for T(l). Two main functions are performed by the package, which are:

1) Develops the dynamical system, the left side of Eq. (2).

2) Performs the numerical integrator, ode45 from Matlab<sup>®</sup>. Fig. 1 shows the flow diagram of the main program called

Base, which run all other functions. The Base is composed by severals sections, in the 'Preparations' section are defined parameters. In def\_coord were created the co- ordinate variables as were defined at Base, and def\_coord may not be edited. To create coordinates variables was used the variables *num\_coord\_gen\_qq* to define dimesions of  $q \in R^{num\_coord\_gen\_qq}$  - and num\_coord\_gen\_pp  $p \in R^{num\_coord\_gen\_pp}$ . In the script Energy are written the Lagrangian equations for energies 2T, V and R, separately for a better presentation of results. The script Barriers is used for the constraint equations. In the variable "strings\_latex" are stored all the strings for exhibited text. The parameter exec is used to split the flow to generate equations for numerical integration or to initiate the simulation. First of all, the equations must be generated and so exec = 0. After that, the Lagrangian equations are applied and an expression for the generalized forces is generated, and finally the result is separated in M matrix,  $\underline{C}$  and  $\underline{G}$  vectors which are described in Eq. (7).

In the case of a system with constraint, parameter *restrito* = 1, the same procedure is done considering the constraints. The variable "eq\_gen" is the core for the generation of the equations of the system dynamics. In this script, the equations are written as described in Eq. (1). Equations for the integrator are shown in the command window and stored on file and the "function a .mat dv = z\_Dinamica\_ode\_dy\_sel\_000 (t,y)" is created, where "000" has the variable sel value. This is exactly the name of the function called by the integrator at *eq\_int* script.

Fig. 2 shows the definition of *str\_latex*, which is useful and necessary to change some notations from Matlab<sup>®</sup> to LATEX e.g., "d2pp" to "\ddot p". Thus, the correct LATEX notation is applied and the report is easier to read. To do this procedure is used the *mega\_loc\_sub* script. The program inserts a pre-defined header and footer as comments, then it searches specific terms in *mega\_loc\_sub* based on an alphabet defined at *alfabeto*.

The *size\_alfabeto* must be correctly set and every term to be searched must be named as *alfabeto\_in\_000*, changing "000" sequentially and the new value must be *alfabeto\_out\_000*, always with 3 digits. Finally the *save\_file\_tex* function receives the. *tex* file, which name was constructed with *nome\_arquivo* as a base and adding some of the control variable to it, this file will be created at the path specified at *caminho* (path).

On the other side, the dynamical system is copied into dy, which is accessed by odeXY inside the  $eq_int$  script and this script is accessed from the Base when exec = 1 is defined. If the constrains have been set, the user may choose to integrate step by step with  $int_passos = 1$ , so the  $eq_int$  keeps verifying the collision.

The change at the free system behavior is updated as the user described at *restri*, this script must contain the first update of  $\lambda$ , part of this update is also given by *eq\_gen*, and should be complemented by the user with the equations of conservation of momentum (linear and/or angular) at the collision instant using the correct notation outside the *odeXY* script.

Finished integration, phase portrait graphs are printed and data is stored, which allows the user redo or create specific graphs, all figures are also stored and at the end the LATEX code for the stored figures are exhibited at the command window.

This code is also saved in a *.tex* which may be directly inserted at the report *.tex* file at the correct folder.



Fig. 2. Program data flow - .tex file creation with correct notation.

Although all automatic development and creation given by the program, the correct description of the constraints and how they interact with the free system, relationships and impact dynamics are not a concern of this package, and it should be correctly described by the project designer. A tutorial file describing the program and how to handle the generated code and report has also been developed.

#### IV. APPLICATION EXAMPLES

In this section are shown three application examples, which are based on references  $[1^{1}2]$ -[14]. Only the line before the figure and figure itself have been written by us. All development and equations have been directly inserted by our package as it is shown (therefore some equations are a little bit misaligned) writing only the "\input{file}" command.

## A. Example in Aerospace: Constrained Mass-Spring System

The constrained mass-spring system is shown at Fig. 3, it is detailed at [14] while investigation contacts dynamics of bodies. There is the object of mass  $m_1$  and the wall  $m_w$ , they are separated for the distance *d* while at the equilibrium points. We reproduce the same result from [14] with the interaction force  $\lambda$  from Lagrangian multipliers.

Fig. 4, Fig. 5 and Fig. 6 were created and inserted at the LATEX file automatically, LagranTexPac generates all combinations of coordinates of each system, coordinates q for free system and p for wall system. To avoid numerical errors, it is defined  $\varepsilon_1$ . If the distance between the mass and

the wall is smaller than  $\varepsilon_1$ , then the system is defined "constrained" and changes to the constrained equations. It returns if the distance becomes greater than  $\varepsilon_1$  or if the constrained force  $\lambda$  is smaller than  $\varepsilon_{\lambda 1}$ .

Fig. 4 shows the time dependent variables  $x_1$  to  $x_4$ , system oscillates until position  $x_1$  hits the wall and it starts to move as can be seen by the  $x_3$  and  $x_4$  appearing out of zero,  $x_3$  starts from zero of its own coordinates system and  $x_4$  starts with non-zero velocity after the collision momentum transfer. "*Free System*" line has value 1 when system is "free", no contact, and 0 when it does. The impact instant is also shown as a thin line just after "Free System" line reaches 0. While system is constrained, there is the contact force  $\lambda_1$ , this force begins at its highest value and decrease to 0 when the system is free once again. Fig. 5 and Fig. 6 are the phase portraits  $x_1$ ×  $x_2$  and  $x_3 \times x_4$  respectively, the impact instant is identified by the abrupt chance in both phase portraits.



Fig. 3. Constrained mass-spring system, with q1 = x1 and q1 = xw. source: [14].

System energies Kinetics:

$$2\mathcal{T} = \dot{q}_1^2 m_1 + \tag{8}$$

Potential:

$$V = \frac{k_1 q_1^2}{2}$$
(9)

Rayleigh dissipation:

$$\mathbf{R} = \mathbf{0} \tag{10}$$

## **Barriers system energies**

Kinetics:

$$2\mathcal{T}_{pp} = \dot{p}_1^2 m_w + \tag{11}$$

Potential:

$$V = \frac{k_w p_1^2}{2}$$
(12)

Rayleigh dissipation:

$$R = \frac{c_w \dot{p}_1^2}{2}$$
(13)

## Free system Euler-Lagrange

Euler-Lagrange term to T:

$$2tauT(1) = (2m_1)\ddot{q}_1 \tag{14}$$

Euler-Lagrange term to V:

$$tauV(1) = (-k_1)q_1$$
 (15)

Euler-Lagrange term to R:

$$tau\mathbf{R}(1) = 0 \tag{16}$$

Results for with kinetics, potential and dissipation energies:

$$tau(1) = m_1 \ddot{q}_1 + k_1 q_1 \tag{17}$$

Free system matrices

Matrix *M* for the system:

$$M(1,1) = m_1 \tag{18}$$

Vector  $\underline{C}$  for the system:

$$\mathbf{C}(1) = \mathbf{k}_1 \cdot \boldsymbol{q}_1 \tag{19}$$

Vector  $\underline{G}(\underline{q})$  for the system:

$$G(1) = (-g) \tag{20}$$

**Barriers system Euler-Lagrange** 

Euler-Lagrange term to T:

$$2tauT_{pp}(1) = (2m_w)\ddot{p}_1 \tag{21}$$

Euler-Lagrange term to V:

$$tauV_{pp}(1) = -k_w p_1 \tag{22}$$

Euler-Lagrange term to R:

$$tauR_{pp}(1) = c_w \dot{p}_1 \tag{23}$$

Results for  $\underline{\tau}$  with barriers kinetics, potential and dissipation energies:

$$tau_{pp}(1) = m_{w}\ddot{p}_{1} + c_{w}\dot{p}_{1} + k_{w}p_{1}$$
(24)

## **Barriers system matrices**

Matrix *M* for the barrier system:

$$M(1,1) = m_{w} \tag{25}$$

Vector <u>C</u> for the barrier system:

$$C(1) = c_w \dot{p}_1 + k_w p_1 \tag{26}$$

Vector  $\underline{G}(\underline{q})$  for the barrier system:

$$G(1) = (-g) \tag{27}$$

Additional Euler-Lagrange with constraints  $\Lambda$ Development to  $\tau_{\lambda}$ :

$$tau\lambda(1) = q_1 - p_1 - dist1$$
(28)

Development to  $\tau_{\Lambda}$  :

$$tau\Lambda(1) = \lambda_1 \tag{29}$$

Development to  $\tau_{pp_{\Lambda}}$ :

$$tau_{pp}\Lambda(1) = -\lambda_1 \tag{30}$$

Matrices with constraints  $\lambda$ To generalized coordinates <u>q</u>: Matrix *M* for the system:

$$M(1,1) = m_1 \tag{31}$$

Vector  $\underline{C}$  for the system:

$$C(1) = k_1 q_1 + \lambda_1 \tag{32}$$

Vector  $\underline{G(q)}$  for the system:

$$G(1) = (-g) \tag{33}$$

To barriers coordinates <u>p:</u> Matrix *M* for the barrier system:

$$M(1,1) = m_{w} \tag{34}$$

Vector <u>C</u> for the barrier system:

$$C(1) = c_w \dot{p}_1 + k_w p_1 - \lambda_1 \tag{35}$$

Vector  $\underline{G}(\underline{q})$  for the barrier system:

$$G(1) = (-g) \tag{36}$$

## Equations to $\lambda$ :

Definition of  $\underline{p}$  and derivatives as functions of  $\underline{q}$  and derivatives:

$$p_{1} = q_{1} - dist1$$

$$\dot{p}_{1} = \dot{q}_{1} \qquad (37)$$

$$\ddot{p}_{1} = \ddot{q}_{1}$$

Isolating  $\lambda_i$ :

$$\begin{aligned}
\lambda_{1} &= m_{w}\ddot{p}_{1} + c_{w}\dot{p}_{1} + k_{w}p_{1} \\
\lambda_{1} &= m_{w}\ddot{p}_{1} + c_{w}\dot{p}_{1} - k_{w}(dist1 - q_{1}) \\
\lambda_{1} &= m_{w}\ddot{p}_{1} + (c_{w}\dot{q}_{1} - k_{w}(dist1 - q_{1})) \\
\lambda_{1} &= m_{w}\ddot{q}_{1} + c_{w}\dot{q}_{1} + k_{w}q_{1} - dist1k_{w}
\end{aligned}$$
(38)

## Numerical simulation

Program with fixed integration step ode45.

Selected options:  $coef_restitui = 0$ , restrito = 1,  $int_passos = 1$ , tspan(end) = 3.500, dt = 0.010.

Parameters of this batch:  $m_1 = 10.00$ ;  $k_1 = 7.00$ ;  $m_w = 5.00$ ;  $c_w = 20.29$ ;  $k_w = 7.00$ ; dist1 = 0.005;  $\epsilon_1 = 0.0001$ ;  $\epsilon_{\lambda 1} = 0.0001$ .







Fig. 6. Phase portrait  $x3 \times x4$ .

B. Example in Biomedical: Triple Pendulum Arm



Fig. 7. Triple pendulum arm, with  $qi = \theta i$ ,  $\forall i = 1, 2, 3$ . source: [13].

Fig. 7 shows the triple pendulum arm system. In [13] treated about concepts in physiology and mechanics applied to rehabilitation studies. The system studied in [13] is an excellent application example how the extended equations may become complex and how the proposed package handles this problem writing the LATEX summarized report as clear as possible directly from Matlab<sup>®</sup> and this package tries to avoid writing mistakes and forgetting mistakes.

## System energies

Kinetics:

$$2T = L_{1}^{2} \dot{q}_{1}^{2} m_{2} + L_{1}^{2} \dot{q}_{1}^{2} m_{3} + L_{2}^{2} \dot{q}_{2}^{2} m_{3}$$

$$+ \dot{q}_{1}^{2} m_{1} z_{1}^{2} + \dot{q}_{2}^{2} m_{2} z_{2}^{2} + \dot{q}_{3}^{2} m_{3} z_{3}^{2}$$

$$+ 2L_{1} L_{2} \dot{q}_{1} \dot{q}_{2} m_{3} \left(\cos(q_{1}) \cos(q_{2}) + \sin(q_{1}) \sin(q_{2})\right) \qquad (39)$$

$$+ 2L_{1} \dot{q}_{1} \dot{q}_{2} m_{2} z_{2} \left(\cos(q_{1}) \cos(q_{2}) + \sin(q_{1}) \sin(q_{2})\right)$$

$$+ 2L_{1} \dot{q}_{1} \dot{q}_{3} m_{3} z_{3} \left(\cos(q_{1}) \cos(q_{3}) + \sin(q_{1}) \sin(q_{3})\right)$$

$$+ 2L_{2} \dot{q}_{2} \dot{q}_{3} m_{3} z_{3} \left(\cos(q_{2}) \cos(q_{3}) + \sin(q_{2}) \sin(q_{3})\right)$$

Potential:

$$V = g \begin{pmatrix} m_1 z_1 \cos(q_1) + m_2 z_2 \cos(q_2) + m_3 z_3 \cos(q_3) + \\ L_1 m_2 \cos(q_1) + L_1 m_3 \cos(q_1) + L_2 m_3 \cos(q_2) \end{pmatrix} (40)$$

Rayleigh dissipation:

$$\mathbf{R} = \mathbf{0} \tag{41}$$

## **Free system Euler-Lagrange** Euler-Lagrange term to T:

$$2tauT(1) = \left(2L_{1}^{2}m_{2} + 2L_{1}^{2}m_{3} + 2m_{1}z_{1}^{2}\right)q_{1}$$

$$+ \left(2L_{1}m_{2}z_{2}\left(\cos(q_{1})\cos(q_{2}) + \sin(q_{1})\sin(q_{2})\right)\right) + 2L_{1}L_{2}m_{3}\left(\cos(q_{1})\cos(q_{2}) + \sin(q_{1})\sin(q_{2})\right)\right) \ddot{q}_{2} + \left(-2L_{1}m_{2}z_{2}\left(\cos(q_{1})\sin(q_{2}) - \cos(q_{2})\sin(q_{1})\right)\right) - 2L_{1}L_{2}m_{3}\left(\cos(q_{1})\sin(q_{2}) - \cos(q_{2})\sin(q_{1})\right)\right) \dot{q}_{2}^{2} + \left(2L_{1}m_{3}z_{3}\left(\cos(q_{1})\cos(q_{3}) + \sin(q_{1})\sin(q_{3})\right)\right) \ddot{q}_{3} + \left(-2L_{1}m_{3}z_{3}\left(\cos(q_{1})\sin(q_{3}) - \cos(q_{3})\sin(q_{1})\right)\right) \dot{q}_{3}^{2}$$

$$(42)$$

$$2tauT(2) = \begin{pmatrix} 2L_{1}m_{2}z_{2} \left(\cos(q_{1})\cos(q_{2}) + \sin(q_{1})\sin(q_{2})\right) \\ +2L_{1}L_{2}m_{3} \left(\cos(q_{1})\cos(q_{2}) + \sin(q_{1})\sin(q_{2})\right) \end{pmatrix} \ddot{q}_{1} \\ + \begin{pmatrix} 2L_{1}m_{2}z_{2} \left(\cos(q_{1})\sin(q_{2}) - \cos(q_{2})\sin(q_{1})\right) \\ +2L_{1}L_{2}m_{3} \left(\cos(q_{1})\sin(q_{2}) - \cos(q_{2})\sin(q_{1})\right) \end{pmatrix} \dot{q}_{1}^{2} \\ + \left(2m_{3}L_{2}^{2} + 2m_{2}z_{2}^{2}\right) \ddot{q}_{2} \\ + \left(2L_{2}m_{3}z_{3} \left(\cos(q_{2})\cos(q_{3}) + \sin(q_{2})\sin(q_{3})\right)\right) \ddot{q}_{3} \\ + \left(-2L_{2}m_{3}z_{3} \left(\cos(q_{2})\sin(q_{3}) - \cos(q_{3})\sin(q_{2})\right)\right) \dot{q}_{3}^{2} \end{cases}$$

$$(43)$$

$$2tauT(3) = \left(2L_{1}m_{3}z_{3}\left(\cos\left(q_{1}\right)\cos\left(q_{3}\right) + \sin\left(q_{1}\right)\sin\left(q_{3}\right)\right)\right)\ddot{q}_{1} + \left(2L_{1}m_{3}z_{3}\left(\cos\left(q_{1}\right)\sin\left(q_{3}\right) - \cos\left(q_{3}\right)\sin\left(q_{1}\right)\right)\right)\dot{q}_{1}^{2} + \left(2L_{2}m_{3}z_{3}\left(\cos\left(q_{2}\right)\cos\left(q_{3}\right) + \sin\left(q_{2}\right)\sin\left(q_{3}\right)\right)\right)\ddot{q}_{2} + \left(2L_{2}m_{3}z_{3}\left(\cos\left(q_{2}\right)\sin\left(q_{3}\right) - \cos\left(q_{3}\right)\sin\left(q_{2}\right)\right)\right)\dot{q}_{2}^{2} + \left(2m_{3}z_{3}^{2}\right)\ddot{q}_{3}$$

$$(44)$$

Euler-Lagrange term to V:

$$tauV(1) = g\sin(q_1)(L_1m_2 + L_1m_3 + m_1z_1) \quad (45)$$

$$tauV(2) = g\sin(q_2)(L_2m_3 + m_2z_2)$$
(46)

$$tauV(3) = gm_3 z_3 \sin(q_3) \tag{47}$$

Euler-Lagrange term to R:

$$tauR(1) = 0 \tag{48}$$

$$tauR(2) = 0 \tag{49}$$

$$tauR(3) = 0 \tag{50}$$

Results for with kinetics, potential and dissipation energies:

$$tau(1) = \left(L_{1}^{2}m_{2} + L_{1}^{2}m_{3} + m_{1}z_{1}^{2}\right)\ddot{q}_{1}$$

$$+ \left(L_{1}L_{2}m_{3}\cos\left(q_{1} - q_{2}\right) + L_{1}m_{2}z_{2}\cos\left(q_{1} - q_{2}\right)\right)\ddot{q}_{2}$$

$$+ \left(L_{1}L_{2}m_{3}\sin\left(q_{1} - q_{2}\right) + L_{1}m_{2}z_{2}\sin\left(q_{1} - q_{2}\right)\right)\dot{q}_{2}^{2}$$

$$+ \left(L_{1}m_{3}z_{3}\cos\left(q_{1} - q_{3}\right)\right)\ddot{q}_{3}$$

$$+ \left(L_{1}m_{3}z_{3}\sin\left(q_{1} - q_{3}\right)\right)\dot{q}_{3}^{2}$$

$$- \left(L_{1}gm_{2}\sin\left(q_{1}\right) + L_{1}gm_{3}\sin\left(q_{1}\right) + gm_{1}z_{1}\sin\left(q_{1}\right)\right)$$
(51)

$$tau(2) = (L_{1}L_{2}m_{3}\cos(q_{1}-q_{2})+L_{1}m_{2}z_{2}\cos(q_{1}-q_{2}))\ddot{q}_{1} + (-L_{1}L_{2}m_{3}\sin(q_{1}-q_{2})-L_{1}m_{2}z_{2}\sin(q_{1}-q_{2}))\dot{q}_{1}^{2} + (m_{3}L_{2}^{2}+m_{2}z_{2}^{2})\ddot{q}_{2} + (L_{2}m_{3}z_{3}\cos(q_{2}-q_{3}))\ddot{q}_{3} + (L_{2}m_{3}z_{3}\sin(q_{2}-q_{3}))\dot{q}_{3}^{2} - (L_{2}gm_{3}\sin(q_{2})+gm_{2}z_{2}\sin(q_{2}))$$
(52)

$$tau(3) = (L_1 m_3 z_3 \cos(q_1 - q_3)) \ddot{q}_1 + (-L_1 m_3 z_3 \sin(q_1 - q_3)) \dot{q}_1^2 + (L_2 m_3 z_3 \cos(q_2 - q_3)) \ddot{q}_2 + (-L_2 m_3 z_3 \sin(q_2 - q_3)) \dot{q}_2^2 + (m_3 z_3^2) \ddot{q}_3 - g m_3 z_3 \sin(q_3)$$
(53)

## Free system matrices

Matrix *M* for the system:

$$M(1,1) = L_1^2 m_2 + L_1^2 m_3 + m_1 z_1^2$$
(54)

$$M(1,2) = L_1 \cos(q_1 - q_2) (L_2 m_3 + m_2 z_2)$$
 (55)

$$M(1,3) = L_1 m_3 z_3 \cos(q_1 - q_3)$$
(56)

$$M(2,1) = L_1 \cos(q_1 - q_2) (L_2 m_3 + m_2 z_2)$$
(57)

$$M(2,2) = m_3 L_2^2 + m_2 z_2^2$$
(58)

$$M(2,3) = L_2 m_3 z_3 \cos(q_2 - q_3)$$
(59)

$$M(3,1) = L_1 m_3 z_3 \cos(q_1 - q_3)$$
(60)

$$M(3,2) = L_2 m_3 z_3 \cos(q_2 - q_3)$$
(61)

$$M(3,3) = m_3 z_3^{\ 2} \tag{62}$$

Vector  $\underline{C}$  for the system:

$$C(1) = (L_1 L_2 m_3 \sin(q_1 - q_2) + L_1 m_2 z_2 \sin(q_1 - q_2)) \dot{q}_2^2 + (L_1 m_3 z_3 \sin(q_1 - q_3)) \dot{q}_3^2$$
(63)

$$C(2) = \left(-L_1 L_2 m_3 \sin(q_1 - q_2) - L_1 m_2 z_2 \sin(q_1 - q_2)\right) \dot{q}_1^2 + \left(L_2 m_3 z_3 \sin(q_2 - q_3)\right) \dot{q}_3^2$$
(64)

$$C(3) = (-L_1 m_3 z_3 \sin(q_1 - q_3)) \dot{q}_1^2 + (-L_2 m_3 z_3 \sin(q_2 - q_3)) \dot{q}_2^2$$
(65)

Vector  $\underline{G(q)}$  for the system:

$$G(1) = (-g)\sin(q_1)(L_1m_2 + L_1m_3 + m_1z_1) \quad (66)$$

$$G(2) = (-g)\sin(q_2)(L_2m_3 + m_2z_2)$$
(67)

$$G(3) = (-g)m_3 z_3 \sin(q_3)$$
(68)

## C. Example in Robotics: Polar Telescopic Robot Manipulator

Fig. 8 shows the polar telescopic robot manipulator system and is an example studied in [12], which presents the Lagrangian approach as a method to obtain the equations of motion of a robot.



Fig. 8. Polar Telescopic Robotic, with q1 = r,  $q2 = \theta$ . Source: [12].

#### 1) System energies

Kinetics:

$$2\mathbf{T} = \frac{\dot{q}_2^2 l_1^2 m_1}{3} + m_2 \left( \dot{q}_1^2 + \dot{q}_2^2 q_1^2 \right) + \frac{\dot{q}_2^2 l_2^2 m_2}{12}$$
(69)

Potential:

$$V = g \sin(q_2) \left( \frac{l_1 m_1}{2} + m_2 q_1 \right)$$
(70)

Rayleigh dissipation:

$$\mathbf{R} = \mathbf{0} \tag{71}$$

2) Free system Euler-Lagrange

Euler-Lagrange term to T:

$$2tauT(1) = (2m_2)\ddot{q}_1 + (-2m_2)q_1\dot{q}_2^2 \qquad (72)$$

$$2tauT(2) = (4m_2)\dot{q}_1q_1\dot{q}_2 + (2m_2)q_1^2\ddot{q}_2 + \left(\frac{2m_1l_1^2}{3} + \frac{m_2l_2^2}{6}\right)\ddot{q}_2$$
(73)

Euler-Lagrange term to V:

$$tauV(1) = -gm_2\sin(q_2) \tag{74}$$

$$tauV(2) = \left(-gm_2\cos(q_2)\right)q_1 - \frac{gl_1m_1\cos(q_2)}{2}$$
(75)

Euler-Lagrange term to R:

$$tauR(1) = 0 \tag{76}$$

$$tauR(2) = 0 \tag{77}$$

Results for with kinetics, potential and dissipation energies:

$$tau(1) = m_2 \ddot{q}_1 + \left(-m_2 q_1\right) \dot{q}_2^2 + gm_2 \sin\left(q_2\right)$$
(78)

$$tau(2) = \left(2m_2q_1\right)\dot{q}_1\dot{q}_2 + \left(\frac{m_1l_1^2}{3} + \frac{m_2l_2^2}{12} + m_2q_1^2\right)\ddot{q}_2 + \frac{g\cos(q_2)(l_1m_1 + 2m_2q_1)}{2}$$
(79)

#### 3) Free system matrices

Matrix *M* for the system:

$$M(1,1) = m_2 \tag{80}$$

$$M(1,2) = 0 (81)$$

$$M(2,1) = 0 (82)$$

$$M(2,2) = \frac{m_1 l_1^2}{3} + \frac{m_2 l_2^2}{12} + m_2 q_1^2$$
(83)

Vector  $\underline{C}$  for the system:

$$C(1) = (-m_2) q_1 \dot{q}_2^2 \tag{84}$$

$$C(2) = (2m_2)\dot{q}_1 q_1 \dot{q}_2 \tag{85}$$

Vector  $\underline{G}(\underline{q})$  for the system:

$$G(1) = (-g)\left(-m_2\sin\left(q_2\right)\right) \tag{86}$$

$$G(2) = (-g) \left( -\frac{\cos(q_2)(l_1m_1 + 2m_2q_1)}{2} \right)$$
(87)

## V. CONCLUSIONS

This paper has showed a software tool to obtain and simulate the equations of motion of holonomic mechanical systems. The tool provides a fast and precise development of the differential equations of holonomic dynamical systems and simulation of the system models of any dimension that can be modeled with Euler-Lagrange's equations [12]-[15]. This tool is particular useful for great order systems, which give many terms to differentiate and even have great dimension matrices

An important feature of this tool is to give an option to analyze all steps by looking at the summary report, otherwise, the user should manually write it in LATEX, which demands some effort, time and is susceptible to errors. Others features are: (i) Write the differential equation in a *.m* file for a numerical simulation and run it; (ii) Save all portrait phase graph and all time evolution of variables; (iii) For system with constraints, save and plot the contact force on time dependent graph; and finally (iv) Write a summarized report in LATEX.

This tool seems to apply in different engineering areas as shown in Section IV, as a start this tool proved useful in aerospace, biomedical and robotics and mechanical engineering. Reader is invited to access and test the tool. Files will be sent by authors' e-mails.

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