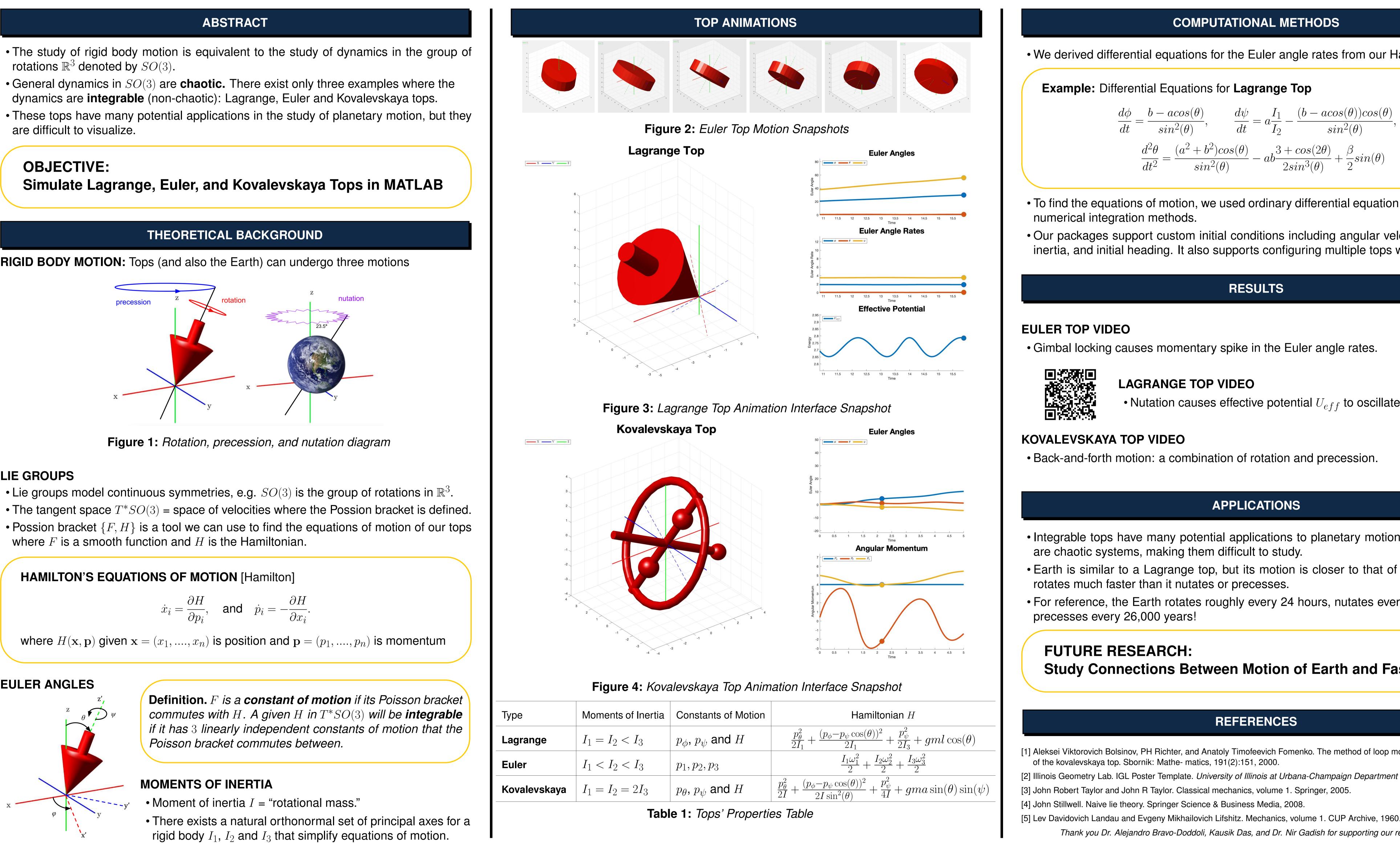


- rotations  $\mathbb{R}^3$  denoted by SO(3).
- General dynamics in SO(3) are **chaotic.** There exist only three examples where the dynamics are **integrable** (non-chaotic): Lagrange, Euler and Kovalevskaya tops.
- are difficult to visualize.

# **OBJECTIVE:** Simulate Lagrange, Euler, and Kovalevskaya Tops in MATLAB

**RIGID BODY MOTION:** Tops (and also the Earth) can undergo three motions



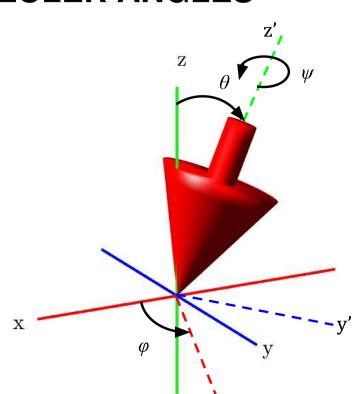
# LIE GROUPS

- Lie groups model continuous symmetries, e.g. SO(3) is the group of rotations in  $\mathbb{R}^3$ .
- where F is a smooth function and H is the Hamiltonian.

# HAMILTON'S EQUATIONS OF MOTION [Hamilton]

$$i = \frac{\partial H}{\partial p_i}$$
, and  $\dot{p}_i = -\frac{\partial H}{\partial x_i}$ .

# **EULER ANGLES**



# Let's Take It From the Top

# Ramon Diego, Haley Gipson, Huaidian Hou, and Anna Huang Mentors: Dr. Alejandro Bravo-Doddoli and Kausik Das

University of Michigan Laboratory of Geometry



# **COMPUTATIONAL METHODS**

• We derived differential equations for the Euler angle rates from our Hamiltonians

$\frac{d\psi}{dt}$ =	$=a\frac{I_1}{I_2}-$	$-\frac{(b-acos(\theta))cos(\theta)}{sin^2(\theta)},$
( heta)	$-ab\frac{3+}{2s}$	$rac{1}{\sin^3(\theta)} + rac{eta}{2} sin( heta)$

• To find the equations of motion, we used ordinary differential equation solvers and other

• Our packages support custom initial conditions including angular velocity, moments of inertia, and initial heading. It also supports configuring multiple tops within one file.

# RESULTS

• Nutation causes effective potential  $U_{eff}$  to oscillate.



# APPLICATIONS

• Integrable tops have many potential applications to planetary motion because planets

• Earth is similar to a Lagrange top, but its motion is closer to that of a fast top, i.e. it

• For reference, the Earth rotates roughly every 24 hours, nutates every 18.6 years, and

# Study Connections Between Motion of Earth and Fast Tops

# REFERENCES

[1] Aleksei Viktorovich Bolsinov, PH Richter, and Anatoly Timofeevich Fomenko. The method of loop molecules and the topology

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