OpenSim Moco Cheat Sheet for the Matlab Interface

MocoStudy

MocoProblem

Access the MocoProblem from the study.

```
problem = study.updProblem();
```

Set the model.

```
problem.setModel(Model('model_file.osim'));
```

Set variable bounds.
Set initial time to 0; final time between 0.5 and 1.5 s.

```
problem.setTimeBounds(MocoInitialBounds(0),
    MocoFinalBounds(0.5, 1.5));
```

The coordinate value must be between 0 and π over the phase, and its initial value is 0 and its final value is $\pi/2$.

```
problem.setStateInfo('/jointset/j0/q0/value',
    [0, pi], 0, pi/2);
```

The control for actuator '/tau0' must be within [-50, 50] over the phase.

```
problem.setControlInfo('/tau0', [-50, 50]);
```

Optimize static model properties.
Create parameter 'myparam' to optimize the mass of Body '/bodyset/b0' within [0.1, 0.5].

```
problem.addParameter(MocoParameter('myparam',
    '/bodyset/b0', 'mass', MocoBounds(0.1, 0.5)));
```

Add cost terms to the problem.

Control · ControlTracking · FinalTime
StateTracking · MarkerTracking · TranslationTracking
OrientationTracking · JointReaction

Minimize the sum of squared controls with weight 1.5.

```
problem.addCost(MocoControlCost('effort', 1.5));
```

Add path constraints to the problem.
Define time-dependent bounds for controls.

```
pathCon = MocoControlBoundConstraint();
problem.addPathConstraint(pathCon);
```

MocoSolver

Initialize the CasADi or Tropter solver.

```
solver = study.initCasADiSolver();
% alternative: solver = study.initTropterSolver();
```

Settings for Tropter and CasADi solvers.
Solve the problem on a grid of 51 mesh points.

```
solver.set_num_mesh_points(51);
```

Transcribe the optimal control problem with the Hermite-Simpson scheme (alternative: 'trapezoidal').

```
solver.set_transcription_scheme('hermite-simpson');
```

Loosen the convergence and constraint tolerances.

```
solver.set_convergence_tolerance(1e-3);
solver.set_constraint_tolerance(1e-3);
```

Stop optimization after 500 iterations.

```
solver.set_max_iterations(500);
```

By default, the Hessian is approximated from first derivatives. Set to 'exact' to use an exact Hessian.

```
solver.set_optim_hessian_approximation('exact');
```

Create a guess, randomize it, then set the guess.

```
guess = solver.createGuess(); guess.randomizeAdd();
solver.setGuess(guess);
```

Set the guess from a MocoTrajectory or MocoSolution file.

```
solver.setGuessFile('previous_solution.sto');
```

Settings for only CasADi solver.
By default, CasADi uses 'central' finite differences; 'forward' differences are faster but less accurate.

```
solver.set_finite_difference_scheme('forward');
```

Turn off parallel calculations.

```
solver.set_parallel(0);
```

Monitor solver progress by writing every 10th iterate to file.

```
solver.set_output_interval(10);
```

Solve the study and obtain a MocoSolution.

```
solution = study.solve();
```

Visualize the solution.

```
study.visualize(solution);
```

Compute outputs from the solution.

```
outputs = StdVectorString();
outputs.add('.*active_force_length_multiplier');
table = study.analyze(solution, outputs);
```

MocoTrajectory and MocoSolution

Create a MocoTrajectory.

```
traj = MocoTrajectory('MocoStudy_solution.sto');
traj = MocoTrajectory.createFromStatesControlsTables(
    states, controls);
```

Get time information.

```
traj.getNumTimes();
traj.getInitialTime(); trajectory.getFinalTime();
traj.getTimeMat();
```

Get names of variables.

```
traj.getStateNames(); traj.getControlNames();
traj.getMultiplierNames(); traj.getParameterNames();
```

Get the trajectory/value for a single variable by name.

```
traj.getStateMat(name); traj.getControlMat(name);
traj.getMultiplierMat(name); traj.getParameter(name);
```

Get the trajectories/values for all variables of a given type.

```
traj.getStatesTrajectoryMat();
traj.getControlsTrajectoryMat();
traj.getMultipliersTrajectoryMat();
traj.getParametersMat();
```

Change the number of times in the trajectory.

```
traj.resampleWithNumTimes(150);
```

Set variable values.

```
traj.setTime(times)
traj.setState(stateTraj); traj.setControl(controlTraj);
traj.setParameter(value);
traj.setStatesTrajectory(statesTraj);
traj.insertStatesTrajectory(subsetStates);
```

Randomize the variable values.

```
traj.randomizeAdd();
```

Export the trajectory.

```
traj.write('mocotrajectory.sto');
traj.exportToStatesTable()
traj.exportToStateTrajectory(mocoProblem)
```

Compare two trajectories.

```
traj.isNumericallyEqual(otherTraj);
traj.compareContinuousVariablesRMS(otherTraj);
traj.compareParametersRMS(otherTraj);
```

Functions on only MocoSolution.

```
solution.success(); solution.getStatus();
solution.getObjective(); solution.getNumIterations();
solution.getSolverDuration();
solution.unseal(); % Access a failed solution.
```

The Moco Optimal Control Problem

minimize	$\sum_j w_j J_j \left(t_0, t_f, y_0, y_f, x_0, x_f, \lambda_0, \lambda_f, p, \int_{t_0}^{t_f} s_{c,j}(t, y, x, \lambda, p)\, dt \right)$	`problem.addCost()`
subject to	$\dot{q} = u$	
	$M(q,p)\dot{u} + G(q,p)^T \lambda = f_{\text{app}}(t,y,x,p) - f_{\text{bias}}(q,u,p)$	
	$\dot{z}(t) = f_{\text{aux}}(t, y, x, \lambda, p)$	`problem.setModel()`
	$0 = \phi(q,p)$	
	$0 = \nu(q,u,p)$	
	$0 = \alpha(q,u,\dot{u},p)$	
	$g_L \le g(t,y,x,\lambda,p) \le g_U$	`problem.addPathConstraint()`
	$y_{0,L} \le y_0 \le y_{0,U}$ $\quad y_{f,L} \le y_f \le y_{f,U}$	`problem.setStateInfo()`
	$x_{0,L} \le x_0 \le x_{0,U}$ $\quad x_{f,L} \le x_f \le x_{f,U}$	`problem.setControlInfo()`
with respect to	$y \in [y_L, y_U]$ $\quad x \in [x_L, x_U]$	
	$t_0 \in [t_{0,L}, t_{0,U}]$ $\quad t_f \in [t_{f,L}, t_{f,U}]$	`problem.setTimeBounds()`
	$p \in [p_L, p_U]$	`problem.addParameter()`

- t: time
- $q(t)$: generalized coordinates
- $u(t)$: generalized speeds
- $z(t)$: auxiliary states (muscle fiber length and activation)
- $y(t) = (q(t), u(t), z(t))$
- $x(t)$: controls (including muscle activation)
- p: constant parameters
- λ: kinematic constraint multipliers
- w_j: weight for j-th cost
- J_j: the j-th cost
- $s_{c,j}$: integrand used in the j-th cost
- M: mass matrix
- f_{bias}: centripetal and coriolis forces
- G: kinematic constraint Jacobian
- f_{app}: applied forces (gravity, muscles, etc.)
- f_{aux}: auxiliary (muscle) dynamics
- ϕ: position-level (holonomic) kinematic constraints
- ν: velocity-level (non-holonomic) kinematic constraints
- α: acceleration-level kinematic constraints
- g: path constraints
- K: number of endpoint constraint goals
- subscript U: an upper bound
- subscript L: a lower bound