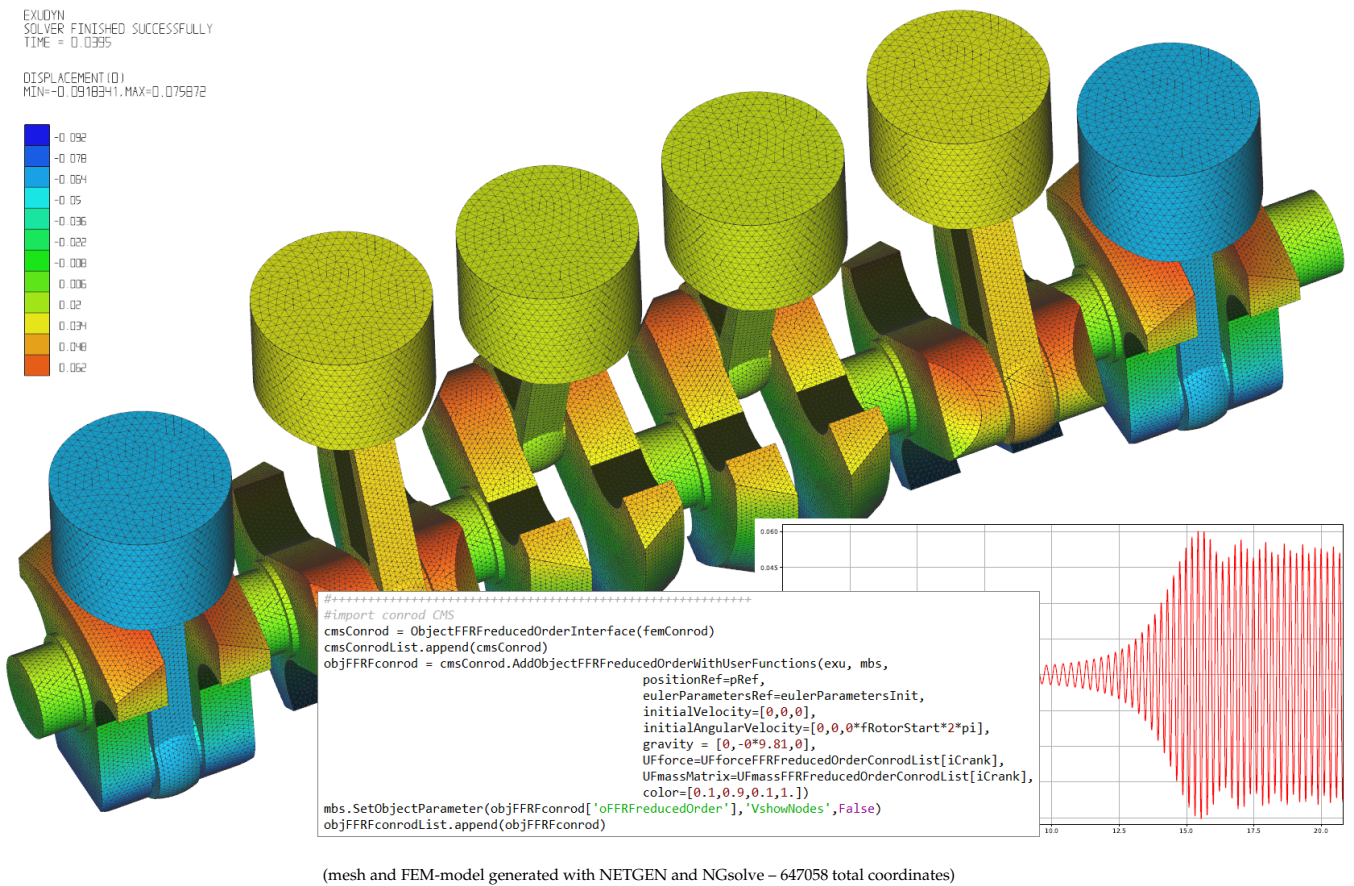


# Flexible Multibody Dynamics Systems with Python and C++

## EXUDYN USER DOCUMENTATION



EXUDYN version = 1.0.19

CHECK section 'CHANGES' for changes from previous versions!!!

University of Innsbruck, Department of Mechatronics, November 18, 2020,

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# Table of Contents

<b>1</b>	<b>Getting Started</b>	<b>1</b>
1.1	Getting started . . . . .	1
1.1.1	What is EXUDYN? . . . . .	1
1.1.2	Who is developing EXUDYN? . . . . .	2
1.1.3	How to install EXUDYN? . . . . .	2
1.1.4	Install with Windows MSI installer . . . . .	3
1.1.5	Install from Wheel . . . . .	3
1.1.6	Work without installation and editing <code>sys.path</code> . . . . .	4
1.1.7	Build and install EXUDYN under Windows 10? . . . . .	4
1.1.8	Build and install EXUDYN under UBUNTU? . . . . .	5
1.1.9	Uninstall EXUDYN . . . . .	6
1.1.10	How to install EXUDYN and using the C++ code (advanced)? . . . . .	6
1.1.11	Goals of EXUDYN . . . . .	6
1.1.12	Run a simple example in Spyder . . . . .	7
1.2	Trouble shooting and FAQ . . . . .	8
<b>2</b>	<b>Overview on EXUDYN</b>	<b>13</b>
2.1	Module structure . . . . .	13
2.1.1	Overview of modules . . . . .	13
2.1.2	Conventions: items, indices, coordinates . . . . .	14
2.2	Items: Nodes, Objects, Loads, Markers, Sensors, ... . . . .	16
2.2.1	Nodes . . . . .	16
2.2.2	Objects . . . . .	16
2.2.3	Markers . . . . .	17
2.2.4	Loads . . . . .	17
2.2.5	Sensors . . . . .	17
2.2.6	Reference coordinates and displacements . . . . .	18
2.3	Exudyn Basics . . . . .	18
2.3.1	Interaction with the EXUDYN module . . . . .	18
2.3.2	Simulation settings . . . . .	18
2.3.3	Visualization settings . . . . .	19
2.3.4	Solver . . . . .	20

2.3.5	Generating output and results . . . . .	20
2.3.6	Graphics pipeline . . . . .	23
2.3.7	Color and RGBA . . . . .	25
2.3.8	Generating animations . . . . .	25
2.4	C++ Code . . . . .	26
2.4.1	Focus of the C++ code . . . . .	26
2.4.2	C++ Code structure . . . . .	27
2.4.3	C++ Code: Modules . . . . .	28
2.4.4	Code style and conventions . . . . .	28
2.4.5	Notation conventions . . . . .	29
2.4.6	No-abbreviations-rule . . . . .	30
2.5	Changes . . . . .	30
<b>3</b>	<b>Tutorial</b>	<b>35</b>
<b>4</b>	<b>Python-C++ command interface</b>	<b>41</b>
4.1	EXUDYN . . . . .	42
4.2	SystemContainer . . . . .	43
4.3	MainSystem . . . . .	44
4.3.1	MainSystem: Node . . . . .	46
4.3.2	MainSystem: Object . . . . .	48
4.3.3	MainSystem: Marker . . . . .	49
4.3.4	MainSystem: Load . . . . .	50
4.3.5	MainSystem: Sensor . . . . .	51
4.4	SystemData . . . . .	52
4.4.1	SystemData: Access coordinates . . . . .	53
4.4.2	SystemData: Get object local-to-global (LTG) coordinate mappings . . . . .	55
4.5	Type definitions . . . . .	55
4.5.1	OutputVariableType . . . . .	55
4.5.2	ConfigurationType . . . . .	57
4.5.3	LinearSolverType . . . . .	57
<b>5</b>	<b>Python utility functions</b>	<b>59</b>
5.1	Module: basicUtilities . . . . .	59
5.2	Module: utilities . . . . .	61
5.3	Module: graphicsDataUtilities . . . . .	65
5.4	Module: rigidBodyUtilities . . . . .	68
5.5	Module: lieGroupBasics . . . . .	72
5.6	Module: FEM . . . . .	78
5.6.1	CLASS ObjectFFRFInterface (in module FEM) . . . . .	81
5.6.2	CLASS ObjectFFRFreducedOrderInterface (in module FEM) . . . . .	82
5.6.3	CLASS FEMinterface (in module FEM) . . . . .	84

5.7	Module: plot . . . . .	88
5.8	Module: processing . . . . .	89
<b>6</b>	<b>Objects, nodes, markers, loads and sensors reference manual</b>	<b>93</b>
6.1	Notation for item equations . . . . .	93
6.1.1	Reference and current coordinates . . . . .	95
6.1.2	Coordinate Systems . . . . .	95
6.2	Nodes . . . . .	96
6.2.1	NodePoint . . . . .	96
6.2.2	NodePoint2D . . . . .	98
6.2.3	NodeRigidBodyEP . . . . .	100
6.2.4	NodeRigidBodyRxyz . . . . .	102
6.2.5	NodeRigidBodyRotVecLG . . . . .	104
6.2.6	NodeRigidBody2D . . . . .	106
6.2.7	Node1D . . . . .	108
6.2.8	NodePoint2DSlope1 . . . . .	110
6.2.9	NodeGenericODE2 . . . . .	112
6.2.10	NodeGenericData . . . . .	113
6.2.11	NodePointGround . . . . .	114
6.3	Objects . . . . .	115
6.3.1	ObjectMassPoint . . . . .	115
6.3.2	ObjectMassPoint2D . . . . .	117
6.3.3	ObjectMass1D . . . . .	119
6.3.4	ObjectRotationalMass1D . . . . .	122
6.3.5	ObjectRigidBody . . . . .	125
6.3.6	ObjectRigidBody2D . . . . .	127
6.3.7	ObjectGenericODE2 . . . . .	130
6.3.8	ObjectFFRF . . . . .	133
6.3.9	ObjectFFRFReducedOrder . . . . .	137
6.3.10	ObjectANCFcable2D . . . . .	142
6.3.11	ObjectALEANCFcable2D . . . . .	144
6.3.12	ObjectGround . . . . .	147
6.3.13	ObjectConnectorSpringDamper . . . . .	148
6.3.14	ObjectConnectorCartesianSpringDamper . . . . .	151
6.3.15	ObjectConnectorRigidBodySpringDamper . . . . .	154
6.3.16	ObjectConnectorCoordinateSpringDamper . . . . .	157
6.3.17	ObjectConnectorDistance . . . . .	160
6.3.18	ObjectConnectorCoordinate . . . . .	162
6.3.19	ObjectConnectorCoordinateVector . . . . .	165
6.3.20	ObjectConnectorRollingDiscPenalty . . . . .	168
6.3.21	ObjectContactCoordinate . . . . .	172
6.3.22	ObjectContactCircleCable2D . . . . .	173

6.3.23	ObjectContactFrictionCircleCable2D . . . . .	175
6.3.24	ObjectJointGeneric . . . . .	177
6.3.25	ObjectJointSpherical . . . . .	182
6.3.26	ObjectJointRollingDisc . . . . .	184
6.3.27	ObjectJointRevolute2D . . . . .	187
6.3.28	ObjectJointPrismatic2D . . . . .	188
6.3.29	ObjectJointSliding2D . . . . .	190
6.3.30	ObjectJointALEMoving2D . . . . .	194
6.4	Markers . . . . .	198
6.4.1	MarkerBodyMass . . . . .	198
6.4.2	MarkerBodyPosition . . . . .	199
6.4.3	MarkerBodyRigid . . . . .	200
6.4.4	MarkerNodePosition . . . . .	201
6.4.5	MarkerNodeRigid . . . . .	202
6.4.6	MarkerNodeCoordinate . . . . .	203
6.4.7	MarkerNodeRotationCoordinate . . . . .	204
6.4.8	MarkerSuperElementPosition . . . . .	205
6.4.9	MarkerSuperElementRigid . . . . .	208
6.4.10	MarkerObjectODE2Coordinates . . . . .	211
6.4.11	MarkerBodyCable2DShape . . . . .	212
6.4.12	MarkerBodyCable2DCoordinates . . . . .	213
6.5	Loads . . . . .	214
6.5.1	LoadForceVector . . . . .	214
6.5.2	LoadTorqueVector . . . . .	215
6.5.3	LoadMassProportional . . . . .	216
6.5.4	LoadCoordinate . . . . .	217
6.6	Sensors . . . . .	218
6.6.1	SensorNode . . . . .	218
6.6.2	SensorObject . . . . .	219
6.6.3	SensorBody . . . . .	220
6.6.4	SensorSuperElement . . . . .	221
6.6.5	SensorLoad . . . . .	222
6.7	GraphicsData . . . . .	223
<b>7</b>	<b>EXUDYN Settings</b>	<b>225</b>
7.1	Simulation settings . . . . .	225
7.1.1	SolutionSettings . . . . .	225
7.1.2	NumericalDifferentiationSettings . . . . .	226
7.1.3	NewtonSettings . . . . .	227
7.1.4	GeneralizedAlphaSettings . . . . .	229
7.1.5	TimeIntegrationSettings . . . . .	230
7.1.6	StaticSolverSettings . . . . .	231

7.1.7	SimulationSettings . . . . .	232
7.2	Visualization settings . . . . .	233
7.2.1	VSettingsGeneral . . . . .	233
7.2.2	VSettingsWindow . . . . .	233
7.2.3	VSettingsOpenGL . . . . .	234
7.2.4	VSettingsContour . . . . .	236
7.2.5	VSettingsExportImages . . . . .	236
7.2.6	VSettingsNodes . . . . .	237
7.2.7	VSettingsBeams . . . . .	237
7.2.8	VSettingsBodies . . . . .	238
7.2.9	VSettingsConnectors . . . . .	238
7.2.10	VSettingsMarkers . . . . .	239
7.2.11	VSettingsLoads . . . . .	239
7.2.12	VSettingsSensors . . . . .	240
7.2.13	VisualizationSettings . . . . .	240
7.3	Solver substructures . . . . .	241
7.3.1	CSolverTimer . . . . .	241
7.3.2	SolverLocalData . . . . .	242
7.3.3	SolverIterationData . . . . .	243
7.3.4	SolverConvergenceData . . . . .	243
7.3.5	SolverOutputData . . . . .	244
7.3.6	MainSolverStatic . . . . .	245
7.3.7	MainSolverImplicitSecondOrder . . . . .	247
<b>8</b>	<b>3D Graphics Visualization</b>	<b>253</b>
8.1	Mouse input . . . . .	253
8.2	Keyboard input . . . . .	253
<b>9</b>	<b>System of equations</b>	<b>255</b>
9.1	LHS-RHS naming conventions in EXUDYN . . . . .	255
<b>10</b>	<b>Solver</b>	<b>257</b>
10.1	Jacobian computation . . . . .	257
10.2	Implicit trapezoidal rule solver . . . . .	258
10.3	Representation of coordinates and equations of motion . . . . .	258
10.4	Newmark method . . . . .	259
<b>11</b>	<b>License</b>	<b>263</b>



# Chapter 1

## Getting Started

The documentation for EXUDYN is split into this introductory section, including a quick start up, code structure and important hints, as well as a couple of sections containing references to the available Python interfaces to interact with EXUDYN and finally some information on theory (e.g., 'Solver').

EXUDYN is hosted on GitHub [2]:

- web: <https://github.com/jgerstmayr/EXUDYN/wiki>

For any comments, requests, issues, bug reports, send an email to:

- email: [reply.exudyn@gmail.com](mailto:reply.exudyn@gmail.com)

Thanks for your contribution!

### 1.1 Getting started

This section will show:

1. What is EXUDYN?
2. Who is developing EXUDYN?
3. How to install EXUDYN
4. How to link EXUDYN and Python
5. Goals of EXUDYN
6. Run a simple example in Spyder
7. FAQ – Frequently asked questions

#### 1.1.1 What is EXUDYN?

EXUDYN– (f)lEXible mUltibody **DY**Namics – **EX**tend yoUr **DY**Namics)

EXUDYN is a C++ based Python library for efficient simulation of flexible multibody dynamics systems. It is the follow up code of the previously developed multibody code HOTINT, which Johannes Gerstmayr started during his PhD-thesis. The open source code HOTINT reached limits of



further (efficient) development and it seemed impossible to continue from this code as it is outdated regarding programming techniques and the numerical formulation.

EXUDYN is designed to easily set up complex multibody models, consisting of rigid and flexible bodies with joints, loads and other components. It shall enable automatized model setup and parameter variations, which are often necessary for system design but also for analysis of technical problems. The broad usability of python allows to couple a multibody simulation with environments such as optimization, statistics, data analysis, machine learning and others.

The multibody formulation is mainly based on redundant coordinates. This means that computational objects (rigid bodies, flexible bodies, ...) are added as independent bodies to the system. Hereafter, connectors (e.g., springs or constraints) are used to interconnect the bodies. The connectors are using Markers on the bodies as interfaces, in order to transfer forces and displacements. For details on the interaction of nodes, objects, markers and loads see Section 2.2.

### 1.1.2 Who is developing EXUDYN?

EXUDYN is currently (11-2020) developed at the University of Innsbruck. In the first phase most of the core code is written by Johannes Gerstmayr, implementing ideas that followed out of the project HOTINT. 15 years of development led to a lot of lessons learned and after 20 years, a code must be re-designed.

Some specific codes regarding Pybind11 (by Jakob Wenzel, <https://github.com/pybind/pybind11>, thanks a lot!!!) interface and parallelization have been written by Stefan Holzinger, who also supported the upload to GitLab.

Important discussions with researchers from the community were important for the design and development of EXUDYN, where we like to mention Joachim Schöberl from TU-Vienna who boosted the design of the code with great concepts.

The cooperation and funding within the EU H2020-MSCA-ITN project 'Joint Training on Numerical Modelling of Highly Flexible Structures for Industrial Applications' contributes to the development of the code.

The following people have contributed to the examples:

- Stefan Holzinger, Michael Pieber, Joachim Schöberl, Manuel Schieferle, Martin Knapp, Lukas March, Dominik Sponring, David Wibmer, Andreas Zwölfer

– thanks a lot! –

### 1.1.3 How to install EXUDYN?

In order to run EXUDYN, you need an appropriate Python installation. We recommend to use

- Anaconda, 32bit, Python 3.6.5 or Anaconda 64bit, Python 3.6.5)<sup>1</sup>

---

<sup>1</sup>Anaconda 32/64bit with Python3.6 can be downloaded via the repository archive <https://repo.anaconda.com/archive/> choosing Anaconda3-5.2.0-Windows-x86.exe or Anaconda3-5.2.0-Windows-x86\_64.exe for 64bit.

- Spyder 3.2.8 with Python 3.6.5 32 bit (alternatively 64bit), which is included in the Anaconda installation<sup>2</sup>

If you plan to extend the C++ code, we recommend to use VS2017<sup>3</sup> to compile your code, which offers Python 3.7 compatibility. However, you should know that Python versions and the version of the module must be identical (e.g., Python 3.6 32 bit **both** in the EXUDYN module and in Spyder).

#### 1.1.4 Install with Windows MSI installer

The simplest way on Windows 10 (and maybe also Windows 7), which works well **if you installed only one python version** and if you installed Anaconda with the option '**Register Anaconda as my default Python 3.x**' or similar, then you can use the provided .msi installers in the main/dist directory:

- For the 64bits python 3.6 version, double click on (version may differ):  
exudyn-1.0.8.win-amd64-py3.6.msi
- Follow the instructions of the installer
- If python/ Anaconda is not found by the installer, provide the 'python directory' as the installation directory of Anaconda3, which usually is installed in:  
C:\ProgramData\Anaconda3

#### 1.1.5 Install from Wheel

The **standard way to install** the python package EXUDYN is to use the so-called 'wheels' (file ending .whl) provided at the directory wheels in the EXUDYN repository. First, open an Anaconda prompt:

- EITHER calling: START->Anaconda->... OR go to anaconda/Scripts folder and call activate.bat
- You can check your python version then, by running python<sup>4</sup>, the output reads like:

```
Python 3.6.5 [Anaconda, Inc.] (default, Mar 29 2018, 13:32:41) [MSC v.1900 64 bit
(AMD64)] on win32
...
```

- → type exit() to close python

**Go to the folder Exudyn\_git/main** (where setup.py lies) and choose the wheel in subdirectory main/dist according to your system (windows/UBUNTU), python version (3.6 or 3.7) and 32 or 64 bits.

For Windows this may read (version number 1.0.8 may be different):

- Python 3.6, 32bit: pip install dist\exudyn-1.0.8-cp36-cp36m-win32.whl

---

<sup>2</sup>It is important that Spyder, python and exudyn are **either** 32bit **or** 64bit. There will be a strange .DLL error, if you mix up 32/64bit. It is possible to install both, Anaconda 32bit and Anaconda 64bit – then you should follow the recommendations of paths as suggested by Anaconda installer

<sup>3</sup>previously, VS2019 was recommended: However, VS2019 has problems with the library 'Eigen' and therefore leads to erroneous results with the sparse solver. VS2017 can also be configured with Python 3.7 now.

<sup>4</sup>python3 under UBUNTU 18.04

- Python 3.6, 64bit: `pip install dist\exudyn-1.0.8-cp36-cp36m-win_amd64.whl`
- Python 3.7, 64bit: `pip install dist\exudyn-1.0.8-cp37-cp37m-win_amd64.whl`

For UBUNTU18.04 this may read (version number 1.0.8 may be different):

- Python 3.6, 64bit: `pip3 install dist\exudyn-1.0.8-cp36-cp36m-linux_x86_64.whl`

### 1.1.6 Work without installation and editing `sys.path`

The **uncommon and old way** ( $\rightarrow$  not recommended for EXUDYN versions  $\geq 1.0.0$ ) is to use Python's `sys` module to link to your `exudyn` (previously `WorkingRelease`) directory, for example:

---

```
import sys
sys.path.append('C:/DATA/cpp/EXUDYN_git/bin/EXUDYN32bitsPython36')
```

---

The folder `EXUDYN32bitsPython36` needs to be adapted to the location of the according EXUDYN package.

### 1.1.7 Build and install EXUDYN under Windows 10?

Note that there are a couple of pre-requisites, depending on your system and installed libraries. For Windows 10, the following steps proved to work:

- install your Anaconda distribution including Spyder
- close all Python programs (e.g. Spyder, Jupyter, ...)
- run an Anaconda prompt (may need to be run as administrator)
- if you cannot run Anaconda prompt directly, do:
  - open windows shell (`cmd.exe`) as administrator (`START`  $\rightarrow$  search for `cmd.exe`  $\rightarrow$  right click on app  $\rightarrow$  'run as administrator' if necessary)
  - go to your Scripts folder inside the Anaconda folder (e.g. `C:\ProgramData\Anaconda\Scripts`)
  - run 'activate.bat'
- go to 'main' of your cloned github folder of `exudyn`
- run: `python setup.py install`
- read the output; if there are errors, try to solve them by installing appropriate modules

You can also create your own wheels, doing the above steps to activate the according python version and then calling (requires installation of Microsoft Visual Studio; recommended: VS2017):

```
python setup.py bdist_wheel
```

This will add a wheel in the `dist` folder.

### 1.1.8 Build and install EXUDYN under UBUNTU?

Having a new UBUNTU 18.04 standard installation (e.g. using a VM virtual box environment), the following steps need to be done (python 3.6 is already installed on UBUNTU18.04, otherwise use `sudo apt install python3`)<sup>5</sup>:

First update ...

---

```
sudo apt-get update
```

---

Install necessary python libraries and pip3; matplotlib and scipy are not required for installation but used in EXUDYN examples:

---

```
sudo dpkg --configure -a
sudo apt install python3-pip
pip3 install numpy
pip3 install matplotlib
pip3 install scipy
```

---

Install pybind11 (needed for running the setup.py file derived from the pybind11 example):

---

```
pip3 install pybind11
```

---

If graphics is used (`#define USE_Glfw_GRAPHICS` in `BasicDefinitions.h`), you must install the according GLFW and OpenGL libs:

---

```
sudo apt-get install freeglut3 freeglut3-dev
sudo apt-get install mesa-common-dev
sudo apt-get install libglfw3 libglfw3-dev
sudo apt-get install libx11-dev xorg-dev libglew1.5 libglew1.5-dev libglu1-mesa libglu1-
    mesa-dev libgl1-mesa-glx libgl1-mesa-dev
```

---

With all of these libs, you can run the setup.py installer (go to `Exudyn_git/main` folder), which takes some minutes for compilation (the `-user` option is used to install in local user folder):

---

```
sudo python3 setup.py install --user
```

---

Congratulation! **Now, run a test example** (will also open an OpenGL window if successful):

```
python3 pythonDev/Examples/rigid3Dexample.py
```

You can also create a UBUNTU wheel which can be easily installed on the same machine (x64), same operating system (UBUNTU18.04) and with same python version (e.g., 3.6):

```
sudo pip3 install wheel
sudo python3 setup.py bdist_wheel
```

---

<sup>5</sup>see also the youtube video: [https://www.youtube.com/playlist?list=PLZduTa9mdcm0h5KVUqatD9GzVg\\_jt16fx](https://www.youtube.com/playlist?list=PLZduTa9mdcm0h5KVUqatD9GzVg_jt16fx)

### 1.1.9 Uninstall EXUDYN

To uninstall exudyn under Windows, run (may require admin rights):

```
pip uninstall exudyn
```

To uninstall under UBUNTU, run:

```
sudo pip3 uninstall exudyn
```

If you upgrade to a newer version, uninstall is usually not necessary!

### 1.1.10 How to install EXUDYN and using the C++ code (advanced)?

EXUDYN is still under intensive development of core modules. There are several ways to using the code, but you **cannot** install EXUDYN as compared to other executable programs and apps.

In order to make full usage of the C++ code and extending it, you can use:

- Windows / Microsoft Visual Studio 2017 and above:
  - get the files from git
  - put them into a local directory (recommended: C:/DATA/cpp/EXUDYN\_git)
  - start `main_sln.sln` with Visual Studio
  - compile the code and run `main/pythonDev/pytest.py` example code
  - adapt `pytest.py` for your applications
  - extend the C++ source code
  - link it to your own code
  - NOTE: on Linux systems, you mostly need to replace `/'` with `'\'`
- Linux, etc.: not fully supported yet; however, all external libraries are Linux-compatible and thus should run with minimum adaptation efforts.

### 1.1.11 Goals of EXUDYN

After the first development phase (planned in Q4/2021), it will

- be a small multibody library, which can be easily linked to other projects,
- allow to efficiently simulate small scale systems (compute 100000s time steps per second for systems with  $n_{DOF} < 10$ ),
- safe and widely accessible module for Python,
- allow to add user defined objects in C++,
- allow to add user defined solvers in Python).

Listing 1.1: My first example

```
import exudyn as exu                                #EXUDYN package including C++ core part
from exudyn.itemInterface import * #conversion of data to exudyn dictionaries

SC = exu.SystemContainer()                          #container of systems
mbs = SC.AddSystem()                               #add a new system to work with

nMP = mbs.AddNode(NodePoint2D(referenceCoordinates=[0,0]))
mbs.AddObject(ObjectMassPoint2D(physicsMass=10, nodeNumber=nMP))
mMP = mbs.AddMarker(MarkerNodePosition(nodeNumber = nMP))
mbs.AddLoad(Force(markerNumber = mMP, loadVector=[0.001,0,0]))

mbs.Assemble()                                     #assemble system and solve
simulationSettings = exu.SimulationSettings()
simulationSettings.timeIntegration.verboseMode=1 #provide some output
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', simulationSettings)
```

### 1.1.12 Run a simple example in Spyder

After performing the steps of the previous section, this section shows a simplistic model which helps you to check if EXUDYN runs on your computer.

In order to start, run the python interpreter Spyder. For the following example, either

- open Spyder and copy the example provided in Listing 1.1 into a new file, or
- open `myFirstExample.py` from your `EXUDYN32bitsPython366` directory

Hereafter, press the play button or F5 in Spyder.

If successful, the IPython Console of Spyder will print something like:

---

```
runfile('C:/DATA/cpp/EXUDYN_git/main/bin/EXUDYN32bitsPython36/myFirstExample.py',
        wdir='C:/DATA/cpp/EXUDYN_git/main/bin/EXUDYN32bitsPython36')
+++++
EXUDYN V1.0.1 solver: implicit second order time integration
STEP100, t = 1 sec, timeToGo = 0 sec, Nit/step = 1
solver finished after 0.0007824 seconds.
```

---

If you check your current directory (where `myFirstExample.py` lies), you will find a new file `coordinatesSolution.txt`, which contains the results of your computation (with default values for time integration). The beginning and end of the file should look like:

---

```
#Exudyn generalized alpha solver solution file
#simulation started=2019-11-14,20:35:12
```

---

<sup>6</sup>or any other directory according to your python version

```

#columns contain: time, ODE2 displacements, ODE2 velocities, ODE2 accelerations, AE
    coordinates, ODE2 velocities
#number of system coordinates [nODE2, nODE1, nAlgebraic, nData] = [2,0,0,0]
#number of written coordinates [nODE2, nVel2, nAcc2, nODE1, nVel1, nAlgebraic, nData] =
    [2,2,2,0,0,0,0]
#total columns exported (excl. time) = 6
#number of time steps (planned) = 100
#
0,0,0,0,0,0,0.0001,0
0.02,2e-08,0,2e-06,0,0.0001,0
0.03,4.5e-08,0,3e-06,0,0.0001,0
0.04,8e-08,0,4e-06,0,0.0001,0
0.05,1.25e-07,0,5e-06,0,0.0001,0

...

0.96,4.608e-05,0,9.6e-05,0,0.0001,0
0.97,4.7045e-05,0,9.7e-05,0,0.0001,0
0.98,4.802e-05,0,9.8e-05,0,0.0001,0
0.99,4.9005e-05,0,9.9e-05,0,0.0001,0
1,5e-05,0,0.0001,0,0.0001,0
#simulation finished=2019-11-14,20:35:12
#Solver Info: errorOccurred=0,converged=1,solutionDiverged=0,total time steps=100,total
    Newton iterations=100,total Newton jacobians=100

```

---

Within this file, the first column shows the simulation time and the following columns provide solution of coordinates, their derivatives and Lagrange multipliers on system level. As expected, the  $x$ -coordinate of the point mass has constant acceleration  $a = f/m = 0.001/10 = 0.0001$ , the velocity grows up to 0.0001 after 1 second and the point mass moves 0.00005 along the  $x$ -axis.

## 1.2 Trouble shooting and FAQ

### Known issues:

1. Sometimes the `exudyn` module cannot be loaded into Python. There are several reasons and workarounds:

- You mixed up 32 and 64 bits (see below) version
- You are using an `exudyn` version for Python  $x_1.y_1$  (e.g., 3.6. $z_1$ ) different from the Python  $x_2.y_2$  version in your Anaconda (e.g., 3.7. $z_2$ ); note that  $x_1 = x_2$  and  $y_1 = y_2$  must be obeyed while  $z_1$  and  $z_2$  may be different
- `ModuleNotFoundError: No module named 'exudynCPP'`:  
A known reason is that your CPU<sup>7</sup> does not support AVX2, while `exudyn` is compiled with the AVX2 option;

---

<sup>7</sup>modern Intel Core-i3, Core-i5 and Core-i7 processors as well as AMD processors, especially Zen and Zen-2 architectures should have no problems with AVX; however, low-cost Celeron and Pentium processors **will not support AVX**, e.g., Intel Celeron G3900, Intel core 2 quad q6600, Intel Pentium Gold G5400T; check the system settings of your computer to find out the processor type; typical CPU manufacturer pages or Wikipedia provide information on this

→ **workaround** to solve the AVX problem: use the Python 3.6 32bits version, which is compiled without AVX2; you can also compile for your specific Python version without AVX if you adjust the `setup.py` file in the main folder.

The `ModuleNotFoundError` may also happen if something went wrong during installation (paths, problems with Anaconda, ..) → very often a new installation of Anaconda and EXUDYN helps.

## 2. When importing EXUDYN in python (windows) I get the error (or similar):

---

```
Traceback (most recent call last):
  File "C:\DATA\cpp\EXUDYN_git\main\pythonDev\pytest.py", line 18, in <module>
    import exudyn as exu
ImportError: DLL load failed: %1 is no valid Win32 application.
```

---

→ probably this is a 32/64bit problem. Your Python installation and EXUDYN need to be **BOTH** either 64bit OR 32bit (Check in your python help; `exudyn` in `WorkingRelease64` is the 64bit version, in `WorkingRelease` it is the 32bit version) and the Python installation and EXUDYN need to have **BOTH** the same version and 1st subversion number (e.g., 3.6.5 should be compatible with 3.6.2).

## 3. I do not understand the python errors – how can I find the reason of the error or crash?

- First, you should read all error messages and warnings: from the very first to the last message. Very often, there is a definite line number which shows the error. Note, that if you are executing a string (or module) as a python code, the line numbers refer to the local line number inside the script or module.
- If everything fails, try to execute only part of the code to find out where the first error occurs. By omitting parts of the code, you should find the according source of the error.
- If you think, it is a bug: send an email with a representative code snippet, version, etc. to [reply.exudyn@gmail.com](mailto:reply.exudyn@gmail.com)

## 4. Spyder console hangs up, does not show error messages, ...:

- very often a new start of Spyder helps; most times, it is sufficient to restart the kernel or to just press the 'x' in your IPython console, which closes the current session and restarts the kernel (this is much faster than restarting Spyder);
- restarting the IPython console also brings back all error messages

## List of Frequently asked questions:

### 1. Where do I find the '.exe' file?

- EXUDYN is only available via the python interface as `exudyn.pyd` library, which is located in folder: `main/bin/WorkingRelease`. This means that you need to run python (best: Spyder) and import the EXUDYN module.

### 2. I get the error message 'check potential mixing of different (object, node, marker, ...) indices', what does it mean?

- probably you used wrong item indices, see beginning of Section 4.



- E.g., an object number `oNum = mbs.AddObject(...)` is used at a place where a `NodeIndex` is expected: `mbs.AddObject(MassPoint(nodeNumber=oNum, ...))`
- Usually, this is an ERROR in your code, it does not make sense to mix up these indices!
- In the exceptional case, that you want to convert numbers, see beginning of Section 4.

### 3. Why does type auto completion does not work for mbs (Main system)?

- UPDATE 2020-06-01: with Spyder 4, using Python 3.7, type auto completion works much better, but may find too many completions.
- most python environments (e.g., with Spyder 3) only have information up to the first sub-structure, e.g., `SC=exu.SystemContainer()` provides full access to `SC` in the type completion, but `mbs=SC.AddSystem()` is at the second sub-structure of the module and is not accessible. WORKAROUND: type `mbs=MainSystem()` **before** the `mbs=SC.AddSystem()` command and the interpreter will know what type `mbs` is. This also works for settings, e.g., simulation settings 'Newton'.

### 4. How to add graphics?

- Graphics (lines, text, 3D triangular / STL mesh) can be added to all `BodyGraphicsData` items in objects. Graphics objects which are fixed with the background can be attached to a `ObjectGround` object. Moving objects must be attached to the `BodyGraphicsData` of a moving body. Other moving bodies can be realized, e.g., by adding a `ObjectGround` and changing its reference with time.

### 5. What is the difference between `MarkerBodyPosition` and `MarkerBodyRigid`?

- Position markers (and nodes) do not have information on the orientation (rotation). For that reason, there is a difference between position based and rigid-body based markers. In case of a rigid body attached to ground with a `SpringDamper`, you can use both, `MarkerBodyPosition` or `MarkerBodyRigid`, markers. For a prismatic joint, you will need a `MarkerBodyRigid`.

### 6. I get an error in `SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', simulationSettings)` but no further information – how can I solve it?

- Typical time integration errors may look like:  

```
File "C:/DATA/cpp/EXUDYN_git/main/pythonDev/...<file name>", line XXX, in <module>
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', simulationSettings)
SystemError: <built-in method TimeIntegrationSolve of PyCapsule object at 0x0CC63590> returned
a result with an error set
```
- The prechecks, which are performed to enable a crash-free simulation are insufficient for your model.
- As a first try, restart the IPython console in order to get all error messages, which may be blocked due to a previous run of EXUDYN.
- Very likely, you are using python user functions inside EXUDYN: They lead to an internal python error, which is not caught by EXUDYN. However, you can just check all your user functions, if they will run without EXUDYN. E.g., a load user function `UFloat(t,load)`, which tries to access `load[4]` will fail internally.

- Use the `print(...)` command in python at many places to find a possible error in user functions (e.g., put `print("Start user function XYZ")` at the beginning of every user function.
- It is also possible, that you are using inconsistent data, which leads to the crash. In that case, you should try to change your model: omit parts and find out which part is causing your error
- see also *I do not understand the python errors – how can I find the cause?*

#### 7. Why can't I get the focus of the simulation window on startup (render window hidden)?

- Starting EXUDYN out of Spyder might not bring the simulation window to front, because of specific settings in Spyder(version 3.2.8), e.g., Tools→Preferences→Editor→Advanced settings: uncheck 'Maintain focus in the Editor after running cells or selections'; Alternatively, set `SC.visualizationSettings.window.alwaysOnTop=True` **before** starting the renderer with `exu.StartRenderer()`



# Chapter 2

## Overview on EXUDYN

### 2.1 Module structure

This section will show:

- Overview of modules
- Conventions: dimension of nodes, objects and vectors
- Coordinates: reference coordinates and displacements
- Nodes, Objects, Markers and Loads

#### 2.1.1 Overview of modules

Currently, the module structure is simple:

- Python parts:
  - `itemInterface`: contains the interface, which transfers python classes (e.g., of a `NodePoint`) to dictionaries that can be understood by the C++ module
  - `exudynUtilities`: contains helper classes in Python, which allows simpler working with EXUDYN
- C++ parts, see Figs. 2.1 and 2.2:
  - `exudyn`<sup>1</sup>: on this level, there are just very few functions: `SystemContainer()`, `StartRenderer()`, `StopRenderer()`
  - `SystemContainer`: contains the systems (most important), solvers (static, dynamics, ...), visualization settings
  - `mbs`: system created with `mbs = SC.AddSystem()`, this structure contains everything that defines a solvable multibody system; a large set of nodes, objects, markers, loads can be added to the system, see Section 6;

---

<sup>1</sup>For versions < 1.0.0: there is a second module, called `exudynFast`, which deactivates all range-, index- or memory allocation checks at the gain of higher speed (probably 30 percent in regular cases and up to 100 percent in the 64 bit version). This module is included by `import exudynFast as exu` and can be used same as `exudyn`. To check the version, just type `exu.__doc__` and you will see a note on 'exudynFast' in the `exudynFast` module.

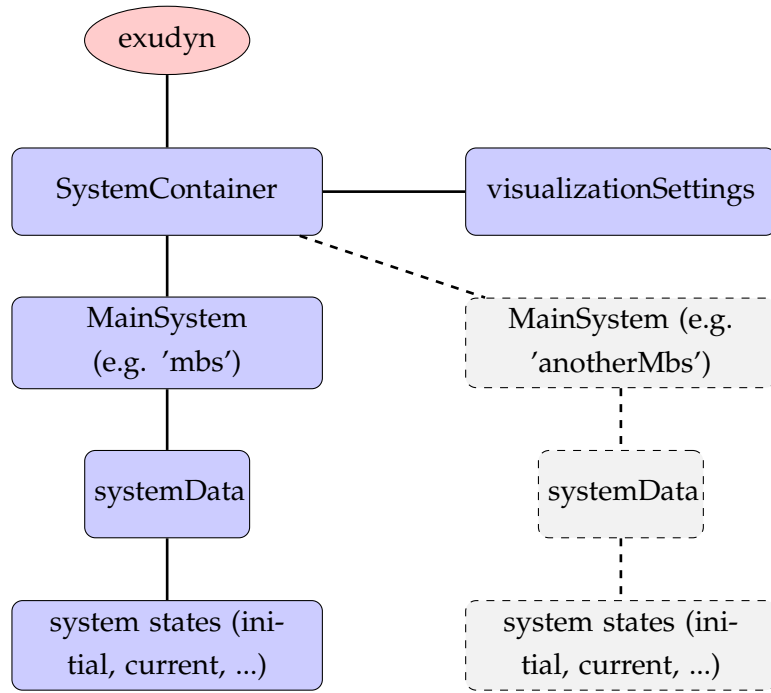


Figure 2.1: Overview of exudyn module.

- `mbs.systemData`: contains the initial, current, visualization, ... states of the system and holds the items, see Fig. 2.2

### 2.1.2 Conventions: items, indices, coordinates

In this documentation, we will use the term **item** to identify nodes, objects, markers and loads:

$$\text{item} \in \{\text{node}, \text{object}, \text{marker}, \text{load}\} \quad (2.1)$$

#### Indices: arrays and vector starting with 0:

As known from Python, all **indices** of arrays, vectors, etc. are starting with 0. This means that the first component of the vector  $v=[1, 2, 3]$  is accessed with  $v[0]$  in Python (and also in the C++ part of EXUDYN). The range is usually defined as  $\text{range}(0, 3)$ , in which '3' marks the index after the last valid component of an array or vector.

#### Dimensionality of objects and vectors:

As a convention, quantities in EXUDYN are 3D, such as nodes, objects, markers, loads, measured quantities, etc. For that reason, we denote planar nodes, objects, etc. with the suffix '2D', but 3D objects do not get this suffix.

Output and input to objects, markers, loads, etc. is usually given by 3D vectors (or matrices), such as (local) position, force, torque, rotation, etc. However, initial and reference values for nodes depend on their dimensionality. As an example, consider a `NodePoint2D`:

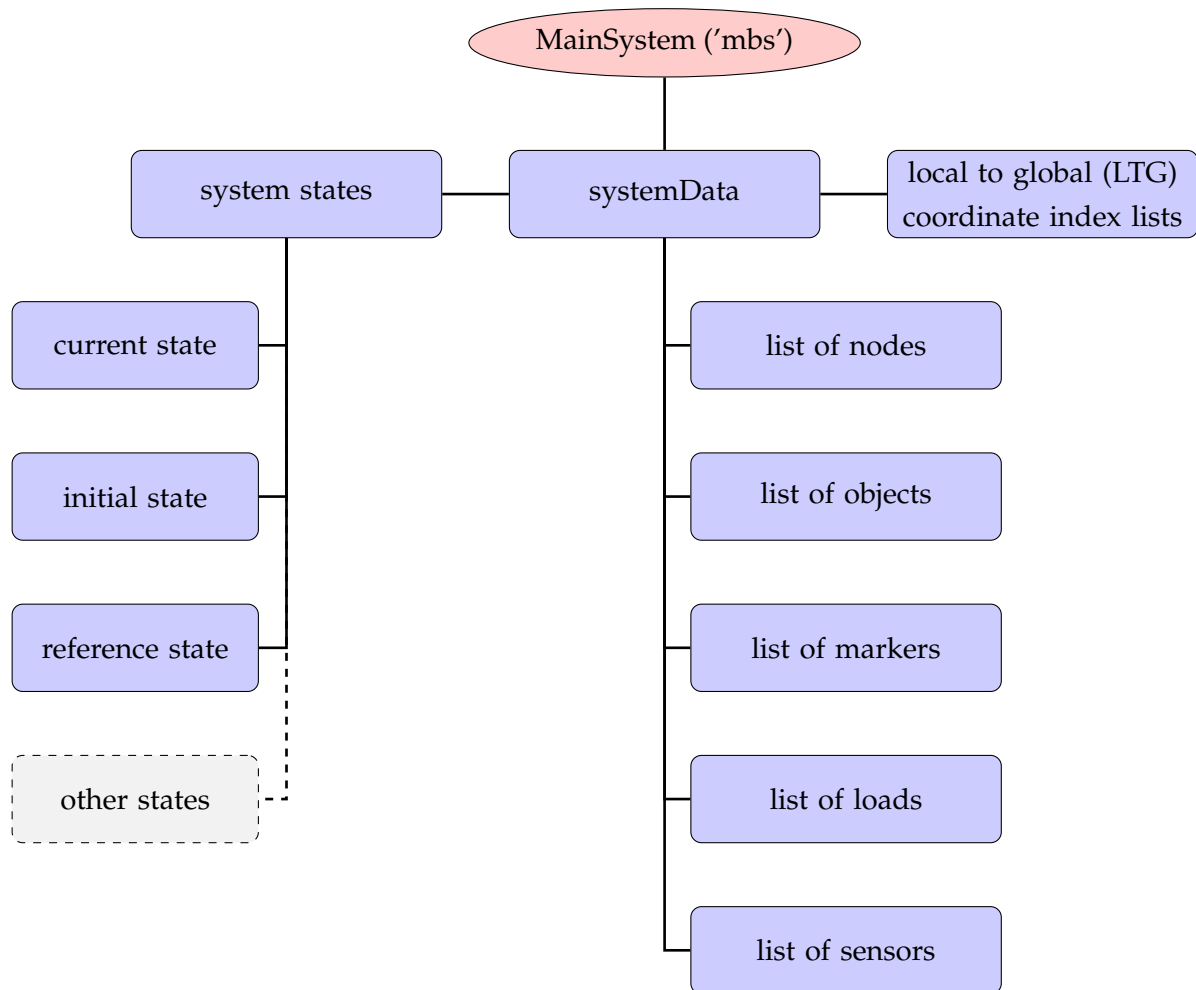


Figure 2.2: Overview of systemData, which connects items and states. Note that access to items is provided via functions in `system`.

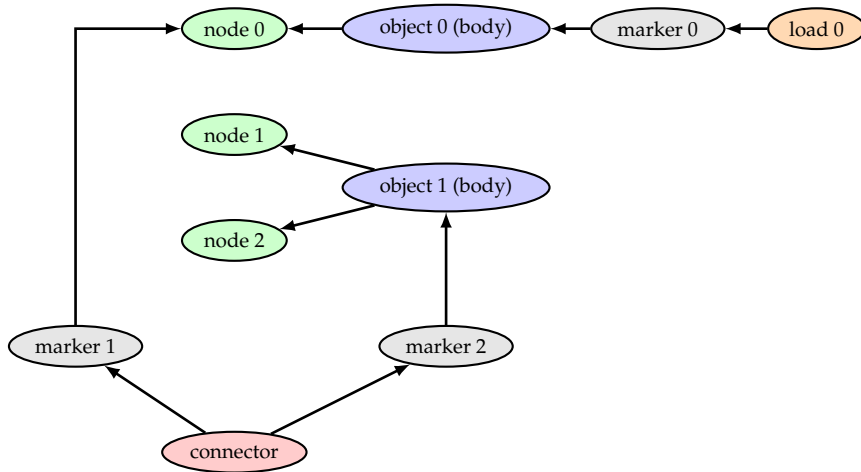


Figure 2.3: Typical interaction of items in a multibody system. Note that both, bodies and connectors/constraints are (computational) objects. The arrows indicate, that, e.g., object 1 has node 1 and node 2 (indices) and that marker 0 is attached to object 0, while load 0 uses marker 0 to apply the load. Sensors could additionally be attached to certain items.

- `referenceCoordinates` is a 2D vector (but could be any dimension in general nodes)
- measuring the current position of `NodePoint2D` gives a 3D vector
- when attaching a `MarkerNodePosition` and a `LoadForceVector`, the force will be still a 3D vector

Furthermore, the local position in 2D objects is provided by a 3D vector. Usually, the dimensionality is given in the reference manual. User errors in the dimensionality will be usually detected either by the python interface (i.e., at the time the item is created) or by the system-preprocessor

## 2.2 Items: Nodes, Objects, Loads, Markers, Sensors, ...

In this section, the most important part of EXUDYN are provided. An overview of the interaction of the items is given in Fig. 2.3

### 2.2.1 Nodes

Nodes provide the coordinates (and the degrees of freedom) to the system. They have no mass, stiffness or whatsoever assigned. Without nodes, the system has no unknown coordinates. Adding a node provides (for the system unknown) coordinates. In addition we also need equations for every nodal coordinate – otherwise the system cannot be computed (NOTE: this is currently not checked by the preprocessor).

### 2.2.2 Objects

Objects are ‘computational objects’ and they provide equations to your system. Objects additionally often provide derivatives and have measurable quantities (e.g. displacement) and they provide access, which can be used to apply, e.g., forces.

Objects can be a:

- general object (e.g. a controller, user defined object, ...; no example yet)
- body: has a mass or mass distribution; markers can be placed on bodies; loads can be applied; constraints can be attached via markers; bodies can be:
  - ground object: has no nodes
  - simple body: has one node (e.g. mass point, rigid body)
  - finite element and more complicated body (e.g. FFRF-object): has more than one node
- connector: uses markers to connect nodes and/or bodies; adds additional terms to system equations either based on stiffness/damping or with constraints (and Lagrange multipliers). Possible connectors:
  - algebraic constraint (e.g. constrain two coordinates:  $q_1 = q_2$ )
  - classical joint
  - spring-damper or penalty constraint

### 2.2.3 Markers

Markers are interfaces between objects/nodes and constraints/loads. A constraint (which is also an object) or load cannot act directly on a node or object without a marker. As a benefit, the constraint or load does not need to know whether it is applied, e.g., to a node or to a local position of a body.

Typical situations are:

- Node – Marker – Load
- Node – Marker – Constraint (object)
- Body(object) – Marker – Load
- Body1 – Marker1 – Joint(object) – Marker2 – Body2

### 2.2.4 Loads

Loads are used to apply forces and torques to the system. The load values are static values. However, you can use Python functionality to modify loads either by linearly increasing them during static computation or by using the 'preStepPyExecute' structure in order to modify loads in every integration step depending on time or on measured quantities (thus, creating a controller).

### 2.2.5 Sensors

Sensors are only used to measure output variables (values) in order to simpler generate the requested output quantities. They have a very weak influence on the system, because they are only evaluated after certain solver steps as requested by the user.



## 2.2.6 Reference coordinates and displacements

Nodes usually have separated reference and initial quantities. Here, `referenceCoordinates` are the coordinates for which the system is defined upon creation. Reference coordinates are needed, e.g., for definition of joints and for the reference configuration of finite elements. In many cases it marks the undeformed configuration (e.g., with finite elements), but not, e.g., for `ObjectConnectorSpringDamper`, which has its own reference length.

Initial displacement (or rotation) values are provided separately, in order to start a system from a configuration different from the reference configuration. As an example, the initial configuration of a `NodePoint` is given by `referenceCoordinates + initialCoordinates`, while the initial state of a dynamic system additionally needs `initialVelocities`.

## 2.3 Exudyn Basics

This section will show:

- Interaction with the EXUDYN module
- Simulation settings
- Visualization settings
- Generating output and results
- Graphics pipeline
- Generating animations

### 2.3.1 Interaction with the EXUDYN module

It is important that the EXUDYN module is basically a state machine, where you create items on the C++ side using the Python interface. This helps you to easily set up models using many other Python modules (numpy, sympy, matplotlib, ...) while the computation will be performed in the end on the C++ side in a very efficient manner.

#### Where do objects live?

Whenever a system container is created with `SC = exu.SystemContainer()`, the structure `SC` lives in C++ and will be modified via the python interface. Usually, the system container will hold at least one system, usually called `mbs`. Commands such as `mbs.AddNode(...)` add objects to the system `mbs`. The system will be prepared for simulation by `mbs.Assemble()` and can be solved (e.g., using `SC.TimeIntegrationSolve(...)`) and evaluated hereafter using the results files. Using `mbs.Reset()` will clear the system and allows to set up a new system. Items can be modified (`ModifyObject(...)`) after first initialization, even during simulation.

### 2.3.2 Simulation settings

The simulation settings consists of a couple of substructures, e.g., for `solutionSettings`, `staticSolver`, `timeIntegration` as well as a couple of general options – for details see Sections 7.1.1 – 7.1.7.

Simulation settings are needed for every solver. They contain solver-specific parameters (e.g., the way how load steps are applied), information on how solution files are written, and very specific control parameters, e.g., for the Newton solver.

The simulation settings structure is created with

---

```
simulationSettings = exu.SimulationSettings()
```

---

Hereafter, values of the structure can be modified, e.g.,

---

```
#10 seconds of simulation time:
simulationSettings.timeIntegration.endTime = 10

#1000 steps for time integration:
simulationSettings.timeIntegration.numberOfSteps = 1000

#assigns a new tolerance for Newton's method:
simulationSettings.timeIntegration.newton.relativeTolerance = 1e-9

#write some output while the solver is active (SLOWER):
simulationSettings.timeIntegration.verboseMode = 2

#write solution every 0.1 seconds:
simulationSettings.solutionSettings.solutionWritePeriod = 0.1

#use sparse matrix storage and solver (package Eigen):
simulationSettings.linearSolverType = exu.LinearSolverType.EigenSparse
```

---

### 2.3.3 Visualization settings

Visualization settings are used for user interaction with the model. E.g., the nodes, markers, loads, etc., can be visualized for every model. There are default values, e.g., for the size of nodes, which may be inappropriate for your model. Therefore, you can adjust those parameters. In some cases, huge models require simpler graphics representation, in order not to slow down performance – e.g., the number of faces to represent a cylinder should be small if there are 10000s of cylinders drawn. Even computation performance can be slowed down, if visualization takes lots of CPU power. However, visualization is performed in a separate thread, which usually does not influence the computation exhaustively. Details on visualization settings and its substructures are provided in Sections 7.2.1 – 7.2.13.

The visualization settings structure can be accessed in the system container SC (access per reference, no copying!), accessing every value or structure directly, e.g.,

---

```
SC.visualizationSettings.nodes.defaultSize = 0.001          #draw nodes very small

#change openGL parameters; current values can be obtained from SC.GetRenderState()
#change zoom factor:
SC.visualizationSettings.openGL.initialZoom = 0.2

#set the center point of the scene (can be attached to moving object):
SC.visualizationSettings.openGL.initialCenterPoint = [0.192, -0.0039, -0.075]
```

---

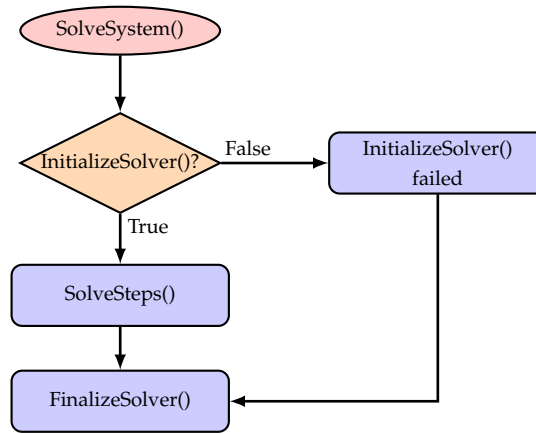


Figure 2.4: Basic solver flow chart for SolveSystem(). This flow chart is the same for static solver and for time integration.

```

#turn of auto-fit:
SC.visualizationSettings.general.autoFitScene = False

#change smoothness of a cylinder:
SC.visualizationSettings.general.cylinderTiling = 100

#make round objects flat:
SC.visualizationSettings.openGL.shadeModelSmooth = False

#turn on coloured plot, using y-component of displacements:
SC.visualizationSettings.contour.outputVariable = exu.OutputVariableType.
    Displacement
SC.visualizationSettings.contour.outputVariableComponent = 1 #0=x, 1=y, 2=z

```

---

### 2.3.4 Solver

Both in the static as well as in the dynamic case, the solver runs in a loop to solve a nonlinear system of (differential and/or algebraic) equations over a given time or load interval. For the time integration (dynamic solver), Fig. 2.4 shows the basic loops for the solution process. The inner loops are shown in Fig. 2.6 and Fig. 2.7. The static solver behaves very similar, while no velocities or accelerations need to be solved and time is replaced by load steps.

Settings for the solver substructures, like timer, output, iterations, etc. are described in Sections 7.3.1 – 7.3.5. The description of interfaces for solvers starts in Section 7.3.6.

### 2.3.5 Generating output and results

The solvers provide a number of options in solutionSettings to generate a solution file. As a default, exporting solution to the solution file is activated with a writing period of 0.01 seconds.

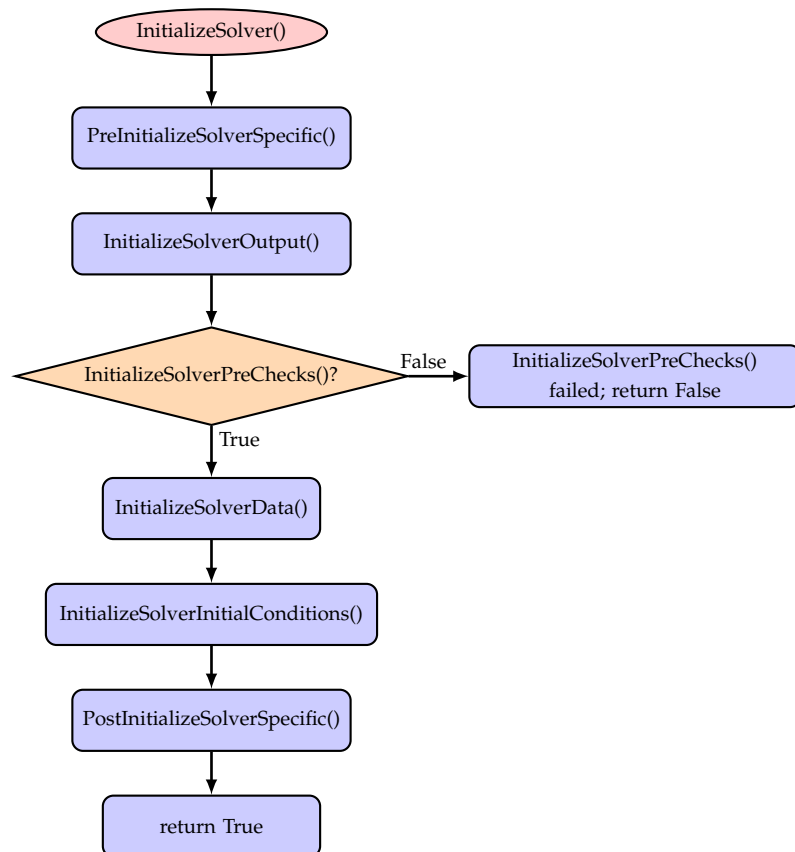


Figure 2.5: Basic solver flow chart for function `InitializeSolver()`.

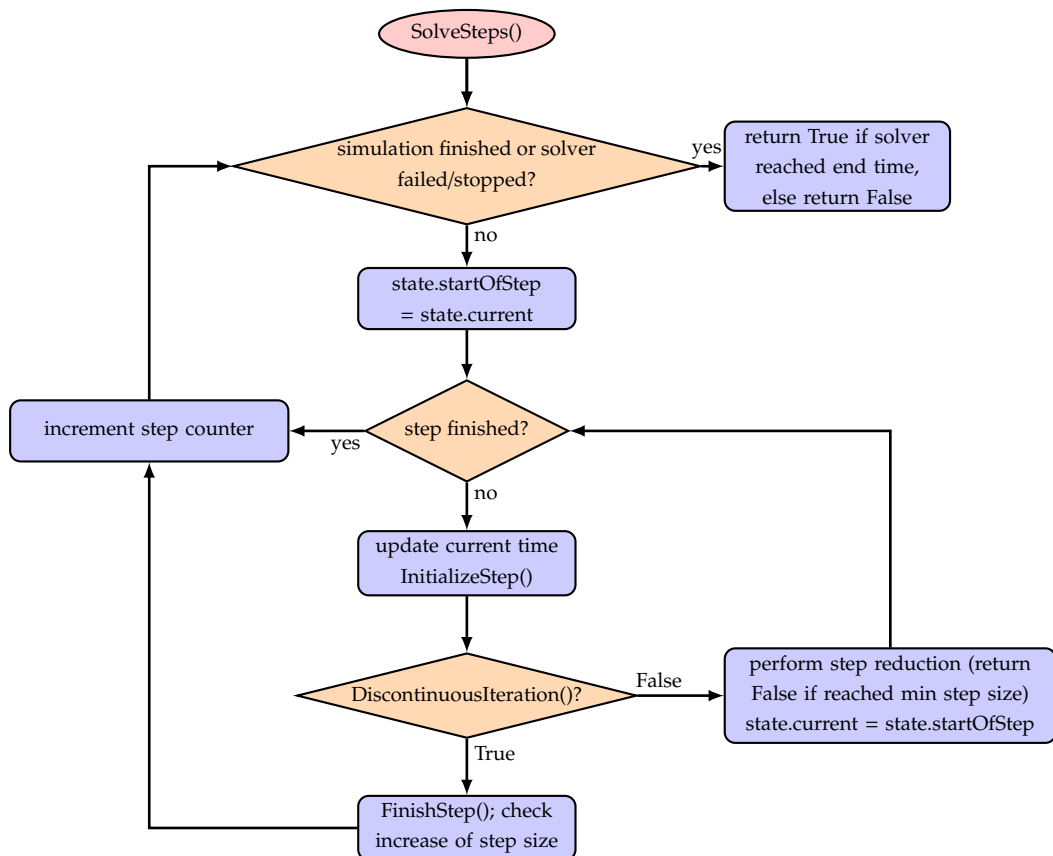


Figure 2.6: Solver flow chart for SolveSteps(), which is the inner loop of the solver.

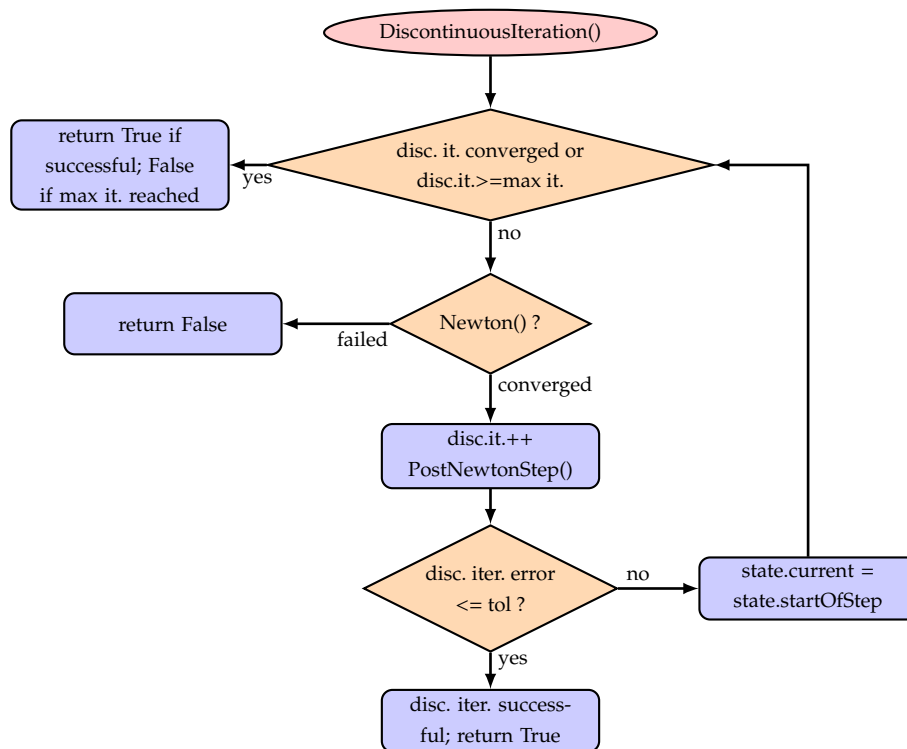


Figure 2.7: Solver flow chart for `DiscontinuousIteration()`, which is run for every solved step inside the static/dynamic solvers. If the `DiscontinuousIteration()` returns False, `SolveSteps()` will try to reduce the step size.

Typical output settings are:

---

```

#create a new simulationSettings structure:
simulationSettings = exu.SimulationSettings()

#activate writing to solution file:
simulationSettings.solutionSettings.writeSolutionToFile = True
#write results every 1ms:
simulationSettings.solutionSettings.solutionWritePeriod = 0.001

#assign new filename to solution file
simulationSettings.solutionSettings.coordinatesSolutionFileName= "myOutput.txt"

#do not export certain coordinates:
simulationSettings.solutionSettings.exportDataCoordinates = False
  
```

---

### 2.3.6 Graphics pipeline

The user cannot interact with the visualization part for now. There are basically two loops during simulation, which feed the graphics pipeline. The solver runs a loop:

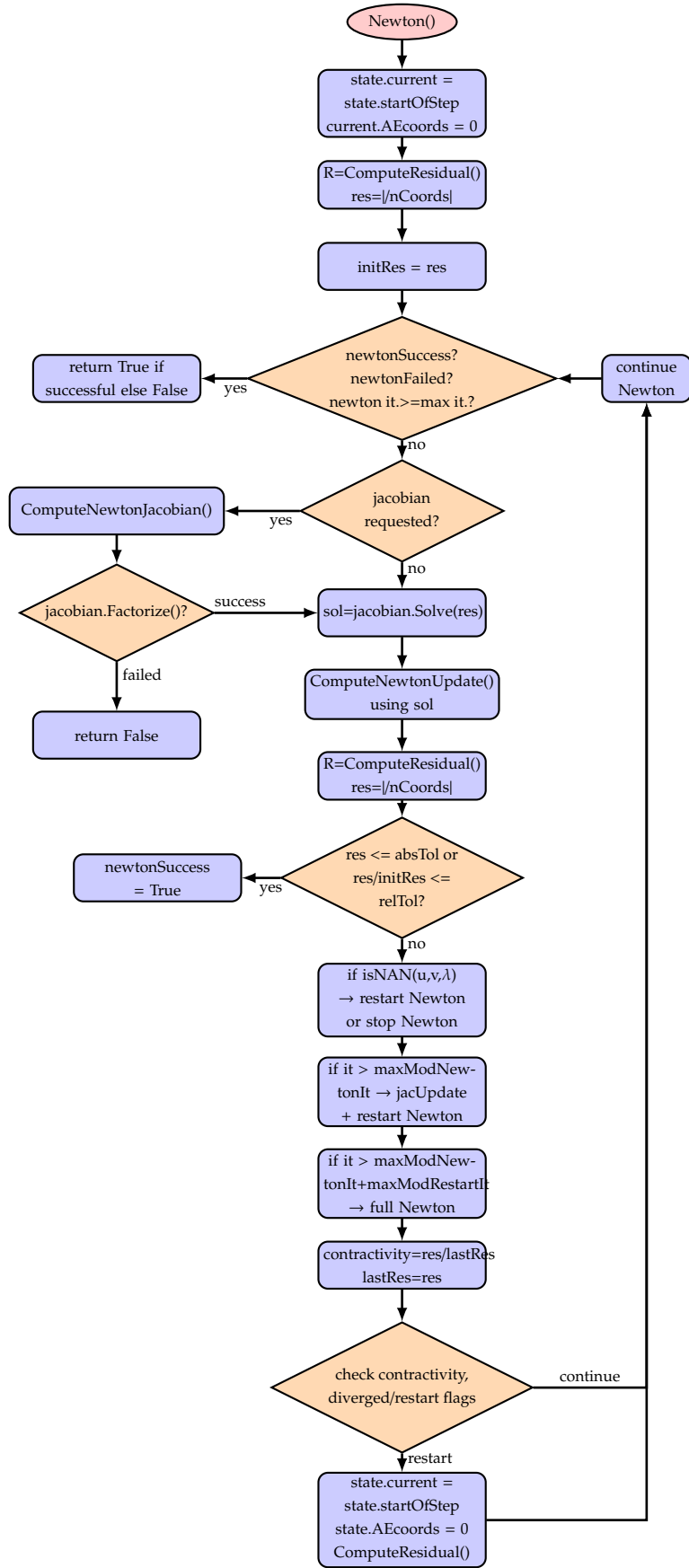


Figure 2.8: Solver flow chart for Newton(), which is run inside the DiscontinuousIteration(). The shown case is valid for newtonResidualMode = 0.

- compute new step
- finish computation step; results are in current state
- copy current state to visualization state (thread safe)
- signal graphics pipeline that new visualization data is available

The OpenGL graphics thread runs the following loop:

- render OpenGL scene with a given graphicsData structure (containing lines, faces, text, ...)
- go idle for some milliseconds
- check if OpenGL rendering needs an update (e.g. due to user interaction)
  - if update is needed, the visualization of all items is updated – stored in a graphicsData structure)
- check if new visualization data is available and the time since last update is larger than a prescribed value, the graphicsData structure is updated with the new visualization state

### 2.3.7 Color and RGBA

Many functions and objects include color information. In order to allow transparency, all colors contain a list of 4 RGBA values, all values being in the range [0..1]:

- red (R) channel
- green (G) channel
- blue (B) channel
- alpha (A) value, representing transparency (A=0: fully transparent, A=1: solid)

E.g., red color with no transparency is obtained by the color=[1,0,0,1]. Color predefinitions are found in `exudynGraphicsDataUtilities.py`, e.g., `color4red` or `color4steelblue` as well a list of 10 colors `color4list`, which is convenient to be used in a loop creating objects.

### 2.3.8 Generating animations

In many dynamics simulations, it is very helpful to create animations in order to better understand the motion of bodies. Specifically, the animation can be used to visualize the model much slower or faster than the model is computed.

Animations are created based on a series of images (frames, snapshots) taken during simulation. It is important, that the current view is used to record these images – this means that the view should not be changed during the recording of images. To turn on recording of images during solving, set the following flag to a positive value

- `simulationSettings.solutionSettings.recordImagesInterval = 0.01`

which means, that after every 0.01 seconds of simulation time, an image of the current view is taken and stored in the directory and filename (without filename ending) specified by

- `SC.visualizationSettings.exportImages.saveImageFileName = "myFolder/frame"`



By default, a consecutive numbering is generated for the image, e.g., 'frame0000.tga, frame0001.tga,...'. Note that '.tga' files contain raw image data and therefore can become very large.

To create animation files, an external tool FFMPEG is used to efficiently convert a series of images into an animation. In windows, simple DOS batch files can do the job to convert frames given in the local directory to animations, e.g.:

---

```
echo off
REM 2019-12-23, Johannes Gerstmayr
REM helper file for EXUDYN to convert all frame00000.tga, frame00001.tga, ... files to a
    video
REM for higher quality use crf option (standard: -crf 23, range: 0-51, lower crf value
    means higher quality)

IF EXIST animation.mp4 (
    echo "animation.mp4 already exists! rename the file"
) ELSE (
    "C:\Program Files (x86)\FFMPEG\bin\ffmpeg.exe" -r 25 -start_number 0 -i frame%%05d.
        tga -c:v libx264 -vf "fps=25,format=yuv420p" animation.mp4
)

```

---

After the video has been created, you should delete the single images:

---

```
REM 2019-12-23, Johannes Gerstmayr
REM helper file for EXUDYN
REM delete all .tga images of current directory

del *.tga

```

---

## 2.4 C++ Code

This section covers some information on the C++ code. For more information see the Open source code and use doxygen.

Exudyn was developed for the efficient simulation of flexible multi-body systems. Exudyn was designed for rapid implementation and testing of new formulations and algorithms in multibody systems, whereby these algorithms can be easily implemented in efficient C++ code. The code is applied to industry-related research projects and applications.

### 2.4.1 Focus of the C++ code

**Four principles:**

1. developer-friendly
2. error minimization

3. efficiency
4. user-friendliness

The focus is therefore on:

- A developer-friendly basic structure regarding the C++ class library and the possibility to add new components.
- The basic libraries are slim, but extensively tested; only the necessary components are available
- Complete unit tests are added to new program parts during development; for more complex processes, tests are available in Python
- In order to implement the sometimes difficult formulations and algorithms without errors, error avoidance is always prioritized.
- To generate efficient code, classes for parallelization (vectorization and multithreading) are provided. We live the principle that parallelization takes place on multi-core processors with a central main memory, and thus an increase in efficiency through parallelization is only possible with small systems, as long as the program runs largely in the cache of the processor cores. Vectorization is tailored to SIMD commands as they have Intel processors, but could also be extended to GPGPUs in the future.
- The user interface (Python) provides a 1:1 image of the system and the processes running in it, which can be controlled with the extensive possibilities of Python.

#### 2.4.2 C++ Code structure

The functionality of the code is based on systems (MainSystem/CSystem) representing the multibody system or similar physical systems to be simulated. Parts of the core structure of Exudyn are:

- CSystem / MainSystem: a multibody system which consists of nodes, objects, markers, loads, etc.
- SystemContainer: holds a set of systems; connects to visualization (container)
- node: used to hold coordinates (unknowns)
- (computational) object: leads to equations, using nodes
- marker: defines a consistent interface to objects (bodies) and nodes; write access ('AccessFunction') – provides jacobian and read access ('OutputVariable')
- load: acts on an object or node via a marker
- computational objects: efficient objects for computation = bodies, connectors, connectors, loads, nodes, ...
- visualization objects: interface between computational objects and 3D graphics
- main (manager) objects: do all tasks (e.g. interface to visualization objects, GUI, python, ...) which are not needed during computation
- static solver, kinematic solver, time integration
- python interface via pybind11; items are accessed with a dictionary interface; system structures and settings read/written by direct access to the structure (e.g. SimulationSettings, Visualization-Settings)
- interfaces to linear solvers; future: optimizer, eigenvalue solver, ... (mostly external or in python)

### 2.4.3 C++ Code: Modules

The following internal modules are used, which are represented by directories in `main/src`:

- Autogenerated: item (nodes, objects, markers and loads) classes split into main (management, python connection), visualization and computation
- Graphics: a general data structure for 2D and 3D graphical objects and a tiny openGL visualization; linkage to GLFW
- Linalg: Linear algebra with vectors and matrices; separate classes for small vectors (SlimVector), large vectors (Vector and ResizableVector), vectors without copying data (LinkedDataVector), and vectors with constant size (ConstVector)
- Main: mainly contains SystemContainer, System and ObjectFactory
- Objects: contains the implementation part of the autogenerated items
- Pymodules: manually created libraries for linkage to python via pybind; remaining linking to python is located in autogenerated folder
- pythonGenerator: contains python files for automatic generation of C++ interfaces and python interfaces of items;
- Solver: contains all solvers for solving a CSystem
- System: contains core item files (e.g., MainNode, CNode, MainObject, CObject, ...)
- Tests: files for testing of internal linalg (vector/matrix), data structure libraries (array, etc.) and functions
- Utilities: array structures for administrative/managing tasks (indices of objects ... bodies, forces, connectors, ...); basic classes with templates and definitions

The following main external libraries are linked to Exudyn:

- LEST: for testing of internal functions (e.g. linalg)
- GLFW: 3D graphics with openGL; cross-platform capabilities
- Eigen: linear algebra for large matrices, linear solvers, sparse matrices and link to special solvers
- pybind11: linking of C++ to python

### 2.4.4 Code style and conventions

This section provides general coding rules and conventions, partly applicable to the C++ and python parts of the code. Many rules follow common conventions (e.g., google code style, but not always – see notation):

- write simple code (no complicated structures or uncommon coding)
- write readable code (e.g., variables and functions with names that represent the content or functionality; AVOID abbreviations)
- put a header in every file, according to Doxygen format
- put a comment to every (global) function, member function, data member, template parameter

- ALWAYS USE curly brackets for single statements in 'if', 'for', etc.; example: `if (i<n) {i += 1;}`
- use Doxygen-style comments (use `/// Qt style and '@ date' with '@' instead of 'for commands)`
- use Doxygen (with preceeding '@') 'test' for tests, 'todo' for todos and 'bug' for bugs
- USE 4-spaces-tab
- use C++11 standards when appropriate, but not exhaustively
- ONE class ONE file rule (except for some collectors of single implementation functions)
- add complete unit test to every function (every file has link to LEST library)
- avoid large classes (>30 member functions; > 15 data members)
- split up god classes (>60 member functions)
- mark changed code with your name and date
- REPLACE tabs by spaces: Extras->Options->C/C++->Tabstopps: tab stop size = 4 (=standard)  
+ KEEP SPACES=YES

### 2.4.5 Notation conventions

The following notation conventions are applied (**no exceptions!**):

- use lowerCamelCase for names of variables (including class member variables), consts, c-define variables, ...; EXCEPTION: for algorithms following formulas, e.g.,  $f = M * q_t t + K * q$ , GBar, ...
- use UpperCamelCase for functions, classes, structs, ...
- Special cases for CamelCase: write 'ODEsystem', BUT: 'ODE1Equations'
- '[...]Init' ... in arguments, for initialization of variables; e.g. 'valueInit' for initialization of member variable 'value'
- use American English throughout: Visualization, etc.
- for (abbreviations) in capital letters, e.g. ODE, use a lower case letter afterwards:
- do not use consecutive capitalized words, e.g. DO NOT WRITE 'ODEAE'
- for functions use `ODEComputeCoords()`, for variables avoid 'ODE' at beginning: use `nODE` or write `odeCoords`
- do not use '\_' within variable or function names; exception: derivatives
- use name which exactly describes the function/variable: 'numberOfItems' instead of 'size' or 'l'
- examples for variable names: `secondOrderSize`, `massMatrix`, `mThetaTheta`
- examples for function/class names: `SecondOrderSize`, `EvaluateMassMatrix`, `Position(const Vector3D& localPosition)`
- use the `Get/Set...()` convention if data is retrieved from a class (`Get`) or something is set in a class (`Set`); Use `const T& Get()/T& Get` if direct access to variables is needed; Use `Get/Set` for pybind11
- example `Get/Set`: `Real* GetDataPointer()`, `Vector::SetAll(Real)`, `GetTransposed()`, `SetRotationalParan`, `SetColor(...)`, ...
- use 'Real' instead of double or float: for compatibility, also for AVX with SP/DP
- use 'Index' for array/vector size and index instead of `size_t` or `int`
- item: object, node, marker, load: anything handled within the computational/visualization systems

### 2.4.6 No-abbreviations-rule

The code uses a **minimum set of abbreviations**; however, the following abbreviation rules are used throughout: In general: DO NOT ABBREVIATE function, class or variable names: `GetDataPointer()` instead of `GetPtr()`; exception: `cnt`, `i`, `j`, `k`, `x` or `v` in cases where it is really clear (5-line member functions).

Exceptions to the NO-ABBREVIATIONS-RULE:

- ODE ... ordinary differential equations;
- ODE2 ... marks parts related to second order differential equations (SOS2, EvalF2 in HOTINT)
- ODE1 ... marks parts related to first order differential equations (ES, EvalF in HOTINT)
- AE ... algebraic equations (IS, EvalG in HOTINT); write 'Aecoordinates' for 'algebraicEquation-sCoordinates'
- 'C[...]' ... Computational, e.g. for `ComputationalNode` ==> use 'CNode'
- min, max ... minimum and maximum
- write time derivatives with underscore: `_t`, `_tt`; example: `Position_t`, `Position_tt`, ...
- write space-wise derivatives with underscore: `_x`, `_xx`, `_y`, ...
- if a scalar, write coordinate derivative with underscore: `_q`, `_v` (derivative w.r.t. velocity coordinates)
- for components, elements or entries of vectors, arrays, matrices: use 'item' throughout
- '[...].Init' ... in arguments, for initialization of variables; e.g. 'valueInit' for initialization of member variable 'value'

## 2.5 Changes

The following list covers changes in the python interface and functionality:

- **Version 1.0.17 → Version 1.0.18**
  - **removed** AVX compilation flag for Python 3.6 32bits version due to incompatibility on older Celeron processors
  - added acceleration sensor functionality to most objects and nodes
  - added Lie group utilities (see `exudyn.lieGroup`)
  - added processing utilities for parameter variation and optimization (see `exudyn.processing`)
- **Version 1.0.12 → Version 1.0.13**

corrected LHS (left-hand-side) and RHS (right-hand-side) terminology (issue: 330), see Section 9.1:

  - objects, connectors, etc., use LHS conventions: all terms (mass, stiffness, elastic forces, damping) are computed at LHS of equation
  - forces are written at the RHS
  - system quantities are always written on RHS:  $m\ddot{q} = f_{RHS}$

- **Version 1.0.8 → Version 1.0.9**

change from Index in `mbs.AddNode(...)`, `mbs.AddObject`, ... to special 'item indices' (issue: 333):

- before: `mbs.AddNode(...)` → Index; **now**: `mbs.AddNode(...)` → NodeIndex
- before: `mbs.AddObject(...)` → Index; **now**: `mbs.AddObject(...)` → ObjectIndex
- before: `mbs.AddMarker(...)` → Index; **now**: `mbs.AddMarker(...)` → MarkerIndex
- before: `mbs.AddLoad(...)` → Index; **now**: `mbs.AddLoad(...)` → LoadIndex
- before: `mbs.AddSensor(...)` → Index; **now**: `mbs.AddSensor(...)` → SensorIndex
- Functions previously requiring an itemNumber have been changed to the according itemIndex, e.g., `mbs.SetNodeParameter(nodeNumber=..., ...)` now requires a nodeNumber of type NodeIndex in order to avoid mistakes due to wrong types of indices.
- for further details and specific usage, see beginning of Section 4!

finally removed functions `mbs.CallNodeFunction(...)` and `mbs.CallObjectFunction(...)` (issue: 288)

removed functions `mbs.GetNodeByName(...)`, `GetObjectByName(...)`, etc. (issue: 445)

- **Version 1.0.6 → Version 1.0.7**

autocreate directories (issue: 431):

- directories (folders) will be created for given paths
- this applies, e.g., to sensor's fileName or simulation settings coordinatesSolutionFileName
- previously, a non-existing directory led to an exception

- **Version 0.1.368 → Version 1.0.0**

**Major changes** in the python interface, as the utilities moved into the exudyn package:

- `from itemInterface import *` → `from exudyn.itemInterface import *`
- `from exudynUtilities import *` → `from exudyn.utilities import *`
- `from exudynBasicUtilities import *` → `from exudyn.basicUtilities import *`
- `from exudynFEM import *` → `from exudyn.FEM import *`
- `from exudynGraphicsDataUtilities import *` → `from exudyn.graphicsDataUtilities import *`
- `from exudynGUI import *` → `from exudyn.GUI import *`
- `from exudynLieGroupIntegration import *` → `from exudyn.lieGroupIntegration import *`
- `from exudynRigidBodyUtilities import *` → `from exudyn.rigidBodyUtilities import *`
- `from exudynRobotics import *` → `from exudyn.robotics import *`

- **Version 0.1.360 → Version 0.1.361**

Changes in the python interface:

- `simulationSettings.timeIntegration.preStepPyExecute` and `simulationSettings.staticSolver.preStepPyExecute` are deprecated, DON'T USE any more

– Use `mbs.SetPreStepUserFunction(...)` instead!

- **Version 0.1.352 → Version 0.1.353**

Changes in the renderer screen:

- Keys '0' and 'KEYPAD 0' → not available any more (set default rotation x/y)
- Use keys CTRL+'1', SHIFT+CTRL+'1', CTRL+'2', ... → keys for new standard views!

- **Version 0.1.288 → Version 0.1.289**

Changes in the python interface (**ESSENTIAL!**):

- Added time 't' as additional first argument in user functions: `ObjectCoordinateSpringD`  
`ObjectConnectorCoordinateSpringDamper`, `ObjectConnectorCartesianSpringDamper`

- **Version 0.1.287 → Version 0.1.288**

Changes in the python interface (**ESSENTIAL!**):

- changed the name of **initialDisplacements** to **initialCoordinates** in all Nodes for consistency reasons with rotation parameters!

- **Version 0.1.282 → Version 0.1.284**

Changes in the python interface:

- all `bodyFixed` parameters in `MarkerRigidBody`, which were inactive so far, have been eliminated

- **Version 0.1.260 → Version 0.1.263**

Changes in the python interface:

- `mbs.systemData.GetCurrentTime()` → `mbs.systemData.GetTime()`
- `mbs.systemData.GetVisualizationTime()` → `mbs.systemData.GetTime(configurationType=exu.C`

- **Version 0.1.244 → Version 0.1.245**

Changes in the implementation / solver (LEADS TO DIFFERENT RESULTS):

- **Solvers updated:** static solver and time integration have been updated; old solvers are still available with the 'OldSolver' extension

Changes in the python interface (new functions / interface to call the old solvers):

- `SC.SolveStaticOldSolver(...)`
- `SC.TimeIntegrationSolve(mbs, 'GeneralizedAlphaOldSolver', simulationSettings)`

- **Version 0.1.243 → Version 0.1.244**

Changes in the python interface:

- `simulationSettings.staticSolver.pauseAfterEachStep`  
→ `simulationSettings.pauseAfterEachStep` (merged with `timeIntegration.pauseAfterEachStep`)

- **Version 0.1.238 → Version 0.1.240**

Changes in the implementation / solver (LEADS TO DIFFERENT RESULTS):

- **generalizedAlpha**: corrected initialization of algorithmic acceleration for discontinuous iteration
- **time integration**: corrected time  $t$  for evaluation of RHS from beginning to end of time step (improves accuracy for time-dependent loads significantly)

Changes in the python interface:

- `simulationSettings.timeIntegration.pauseAfterEachStep`  
→ `simulationSettings.pauseAfterEachStep`
- ADDED: `simulationSettings.timeIntegration.verboseModeFile`
- ADDED: `simulationSettings.staticSolver.verboseModeFile`





## Chapter 3

# Tutorial

This section will show:

- A basic tutorial for a 1D mass and spring-damper with initial displacements, shortest possible model with practically no special settings
- A more advanced 2D rigid-body model (*coming soon*)
- Links to examples section

The python source code of this section can be found in the file:

```
main/pythonDev/Examples/springDamperTutorial.py
```

This tutorial will set up a mass point and a spring damper, dynamically compute the solution and evaluate the reference solution.

We import the exudyn library and the interface for all nodes, objects, markers, loads and sensors:

---

```
import exudyn as exu
from exudyn.itemInterface import *
import numpy as np #for postprocessing
```

---

Next, we need a SystemContainer, which contains all computable systems and add a new system. Per default, you always should name your system 'mbs' (multibody system), in order to copy/paste code parts from other examples, tutorials and other projects:

---

```
SC = exu.SystemContainer()
mbs = SC.AddSystem()
```

---

In order to check, which version you are using, you can printout the current EXUDYN version. This version is in line with the issue tracker and marks the number of open/closed issues added to EXUDYN:

---

```
print('EXUDYN version='+exu.__version__)
```

---

Using the powerful Python language, we can define some variables for our problem, which will also be used for the analytical solution:

---

```

L=0.5           #reference position of mass
mass = 1.6      #mass in kg
spring = 4000   #stiffness of spring-damper in N/m
damper = 8      #damping constant in N/(m/s)
f =80          #force on mass

```

---

For the simple spring-mass-damper system, we need initial displacements and velocities:

---

```

u0=-0.08       #initial displacement
v0=1           #initial velocity
x0=f/spring     #static displacement
print('resonance frequency = '+str(np.sqrt(spring/mass)))
print('static displacement = '+str(x0))

```

---

We first need to add nodes, which provide the coordinates (and the degrees of freedom) to the system. The following line adds a 3D node for 3D mass point<sup>1</sup>:

---

```

n1=mbs.AddNode(Point(referenceCoordinates = [L,0,0],
                    initialCoordinates = [u0,0,0],
                    initialVelocities = [v0,0,0]))

```

---

Here, `Point` (`=NodePoint`) is a Python class, which takes a number of arguments defined in the reference manual. The arguments here are `referenceCoordinates`, which are the coordinates for which the system is defined. The initial configuration is given by `referenceCoordinates` + `initialCoordinates`, while the initial state additionally gets `initialVelocities`. The command `mbs.AddNode(...)` returns a `NodeIndex` `n1`, which basically contains an integer, which can only be used as node number. This node number will be used later on to use the node in the object or in the marker.

While `Point` adds 3 unknown coordinates to the system, which need to be solved, we also can add ground nodes, which can be used similar to nodes, but they do not have unknown coordinates – and therefore also have no initial displacements or velocities. The advantage of ground nodes (and ground bodies) is that no constraints are needed to fix these nodes. Such a ground node is added via:

---

```

nGround=mbs.AddNode(NodePointGround(referenceCoordinates = [0,0,0]))

```

---

In the next step, we add an object<sup>2</sup>, which provides equations for coordinates. The `MassPoint` needs at least a mass (kg) and a node number to which the mass point is attached. Additionally, graphical objects could be attached:

---

```

massPoint = mbs.AddObject(MassPoint(physicsMass = mass, nodeNumber = n1))

```

---

In order to apply constraints and loads, we need markers. These markers are used as local positions (and frames), where we can attach a constraint later on. In this example, we work on the coordinate

---

<sup>1</sup>Note: `Point` is an abbreviation for `NodePoint`, defined in `itemInterface.py`.

<sup>2</sup>For the moment, we just need to know that objects either depend on one or more nodes, which are usually bodies and finite elements, or they can be connectors, which connect (the coordinates of) objects via markers, see Section 2.1.

level, both for forces as well as for constraints. Markers are attached to the according ground and regular node number, additionally using a coordinate number (0 ... first coordinate):

---

```
groundMarker=mbs.AddMarker(MarkerNodeCoordinate(nodeNumber= nGround ,
                                                    coordinate = 0))

#marker for springDamper for first (x-)coordinate:
nodeMarker = mbs.AddMarker(MarkerNodeCoordinate(nodeNumber= n1,
                                                    coordinate = 0))
```

---

This means that loads can be applied to the first coordinate of node n1 via marker with number nodeMarker, which is in fact of type MarkerIndex.

Now we add a spring-damper to the markers with numbers groundMarker and the nodeMarker, providing stiffness and damping parameters:

---

```
mbs.AddObject(CoordinateSpringDamper(markerNumbers = [groundMarker, nodeMarker],
                                       stiffness = spring,
                                       damping = damper))
```

---

A load is added to marker nodeMarker, with a scalar load with value f:

---

```
nLoad = mbs.AddLoad(LoadCoordinate(markerNumber = nodeMarker ,
                                     load = f))
```

---

Finally, a sensor is added to the coordinate constraint object nC, requesting the outputvariable Force:

---

```
mbs.AddSensor(SensorObject(objectNumber=nC, fileName='groundForce.txt',
                           outputVariableType=exu.OutputVariableType.Force))
```

---

Note that sensors can be attached, e.g., to nodes, bodies, objects (constraints) or loads. As our system is fully set, we can print the overall information and assemble the system to make it ready for simulation:

---

```
print(mbs)
mbs.Assemble()
```

---

We will use time integration and therefore define a number of steps (fixed step size; must be provided) and the total time span for the simulation:

---

```
steps = 1000 #number of steps to show solution
tEnd = 1     #end time of simulation
```

---

All settings for simulation, see according reference section, can be provided in a structure given from exu.SimulationSettings(). Note that this structure will contain all default values, and only non-default values need to be provided:

---

```
simulationSettings = exu.SimulationSettings()
simulationSettings.solutionSettings.solutionWritePeriod = 5e-3 #output interval
general
```

---

---

```
simulationSettings.solutionSettings.sensorsWritePeriod = 5e-3 #output interval of
    sensors
simulationSettings.timeIntegration.numberOfSteps = steps
simulationSettings.timeIntegration.endTime = tEnd
```

---

We are using a generalized alpha solver, where numerical damping is needed for index 3 constraints. As we have only spring-dampers, we can set the spectral radius to 1, meaning no numerical damping:

---

```
simulationSettings.timeIntegration.generalizedAlpha.spectralRadius = 1
```

---

In order to visualize the results online, a renderer can be started. As our computation will be very fast, it is a good idea to wait for the user to press SPACE, before starting the simulation (uncomment second line):

---

```
exu.StartRenderer() #start graphics visualization
#mbs.WaitForUserToContinue() #wait for pressing SPACE bar to continue
```

---

As the simulation is still very fast, we will not see the motion of our node. Using e.g. steps=10000000 in the lines above allows you online visualize the resulting oscillations.

Finally, we start the solver, by telling which system to be solved, solver type and the simulation settings:

---

```
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', simulationSettings)
```

---

After simulation, our renderer needs to be stopped (otherwise it would stay in background and prohibit further simulations). Sometimes you would like to wait until closing the render window, using WaitForRenderEngineStopFlag():

---

```
#SC.WaitForRenderEngineStopFlag()#wait for pressing 'Q' to quit
exu.StopRenderer() #safely close rendering window!
```

---

There are several ways to evaluate results, see the reference pages. In the following we take the final value of node n1 and read its 3D position vector:

---

```
#evaluate final (=current) output values
u = mbs.GetNodeOutput(n1, exu.OutputVariableType.Position)
print('displacement=',u)
```

---

The following code generates a reference (exact) solution for our example:

---

```
import matplotlib.pyplot as plt
import matplotlib.ticker as ticker

omega0 = np.sqrt(spring/mass) #eigen frequency of undamped system
dRel = damper/(2*np.sqrt(spring*mass)) #dimensionless damping
omega = omega0*np.sqrt(1-dRel**2) #eigen freq of damped system
C1 = u0-x0 #static solution needs to be considered!
C2 = (v0+omega0*dRel*C1) / omega #C1, C2 are coeffs for solution
```

---

```

refSol = np.zeros((steps+1,2))
for i in range(0,steps+1):
    t = tEnd*i/steps
    refSol[i,0] = t
    refSol[i,1] = np.exp(-omega0*dRel*t)*(C1*np.cos(omega*t)+C2*np.sin(omega*t))+x0

plt.plot(refSol[:,0], refSol[:,1], 'r-', label='displacement (m); exact solution')

```

---

Now we can load our results from the default solution file `coordinatesSolution.txt`, which is in the same directory as your python tutorial file. For convenient reading the file containing commented lines, we use a numpy feature and finally plot the displacement of coordinate 0 or our mass point<sup>3</sup>:

```

data = np.loadtxt('coordinatesSolution.txt', comments='#', delimiter=',')
plt.plot(data[:,0], data[:,1], 'b-', label='displacement (m); numerical solution')

```

---

The sensor result can be loaded in the same way. The sensor output format contains time in the first column and sensor values in the remaining columns. The number of columns depends on the sensor and the output quantity (scalar, vector, ...):

```

data = np.loadtxt('groundForce.txt', comments='#', delimiter=',')
plt.plot(data[:,0], data[:,1]*1e-3, 'g-', label='force (kN)')

```

---

In order to get a nice plot within Spyder, the following options can be used<sup>4</sup>:

```

ax=plt.gca() # get current axes
ax.grid(True, 'major', 'both')
ax.xaxis.set_major_locator(ticker.MaxNLocator(10))
ax.yaxis.set_major_locator(ticker.MaxNLocator(10))
plt.legend() #show labels as legend
plt.tight_layout()
plt.show()

```

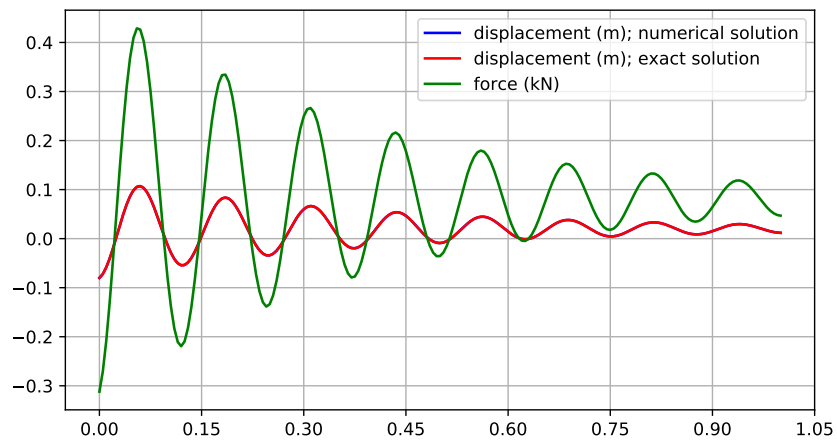
---

The matplotlib output should look like this:

---

<sup>3</sup>`data[:,0]` contains the simulation time, `data[:,1]` contains displacement of (global) coordinate 0, `data[:,2]` contains displacement of (global) coordinate 1, ...)

<sup>4</sup>note, in some environments you need finally the command `plt.show()`



Further examples can be found in your copy of exudyn:

`main/pythonDev/Examples`

`main/pythonDev/TestModels`

## Chapter 4

# Python-C++ command interface

This section lists the basic interface functions which can be used to set up a EXUDYN model in Python.

To import the module, just include the EXUDYN module in Python (for compatibility with examples and other users, we recommend to use the 'exu' abbreviation throughout)<sup>1</sup>:

```
import exudyn as exu
```

In addition, you may work with a convenient interface for your items, therefore also always include the line

```
from exudyn.itemInterface import *
```

The exudyn module will usually hold one SystemContainer, which is a class that is initialized by assigning a system container to a variable, usually denoted as 'SC':

```
SC = exu.SystemContainer()
```

Furthermore, there are a couple of commands available directly in the EXUDYN module, given in the following subsections. Regarding the **(basic) module access**, functions are related to the 'exudyn = exu' module, see these examples:

---

```
import exudyn as exu
from exudyn.itemInterface import *
SC = exu.SystemContainer()
exu.InfoStat() #prints some statistics if available
exu.Go() #creates a systemcontainer and main system
nInvalid = exu.InvalidIndex() #the invalid index, depends on architecture and version
```

---

Understanding the usage of functions for python object 'SystemContainer' provided by EXUDYN, the following examples might help:

---

<sup>1</sup>note that there is a second module, called exudynFast, which does deactivates all range-, index- or memory allocation checks at the gain of higher speed (probably 30 percent in regular cases). To check which version you have, just type `exu.__doc__` and you will see a note on 'exudynFast' in the exudynFast module.



---

```

import exudyn as exu
from exudyn.itemInterface import *
SC = exu.SystemContainer()
mbs = SC.AddSystem()
nSys = SC.NumberOfSystems()
print(nSys)
SC.Reset()

```

---

Many functions will work with node numbers ('NodeIndex'), object numbers ('ObjectIndex'), marker numbers ('MarkerIndex') and others. These numbers are special objects, which have been introduced in order to avoid mixing up, e.g., node and object numbers. For example, the command `mbs.AddNode(...)` returns a `NodeIndex`. For these indices, the following rules apply:

- `mbs.Add[Node|Object|...](...)` returns a specific `NodeIndex`, `ObjectIndex`, ...
- You can create any item index, e.g., using `ni = NodeIndex(42)` or `oi = ObjectIndex(42)`
- You can convert any item index, e.g., `NodeIndex ni` into an integer number using `int(ni)`
- Still, you can use integers as initialization for item numbers, e.g.,  
`mbs.AddObject(MassPoint(nodeNumber=13, ...))`  
However, it must be a pure integer type.
- You can also print item indices, e.g., `print(ni)` as it converts to string by default
- If you are unsure about the type of an index, use `ni.GetTypeString()` to show the index type

## 4.1 EXUDYN

These are the access functions to the EXUDYN module.

function/structure name	description
<code>Go()</code>	Creates a <code>SystemContainer</code> <code>SC</code> and a main system <code>mbs</code>
<code>InfoStat()</code>	Print some global (debug) information: linear algebra, memory allocation, threads, computational efficiency, etc.
<code>StartRenderer(verbose = false)</code>	Start OpenGL rendering engine (in separate thread); use <code>verbose=True</code> to output information during OpenGL window creation
<code>StopRenderer()</code>	Stop OpenGL rendering engine
<code>SetOutputPrecision(numberOfDigits)</code>	Set the precision (integer) for floating point numbers written to console (reset when simulation is started!)
<code>SetLinalgOutputFormatPython(flagPythonFormat)</code>	true: use python format for output of vectors and matrices; false: use matlab format
<code>InvalidIndex()</code>	This function provides the invalid index, which depends on the kind of 32-bit, 64-bit signed or unsigned integer; e.g. node index or item index in list
<code>SetWriteToConsole(flag)</code>	set flag to write (true) or not write to console; default = true

SetWriteToFile(filename, flagWriteToFile = true, flagAppend = false)	set flag to write (true) or not write to console; default value of flagWriteToFile = false; flagAppend appends output to file, if set true; in order to finalize the file, write <code>exu.SetWriteToFile("", False)</code> to close the output file <b>EXAMPLE:</b> <code>exu.SetWriteToConsole(False) #no output to console</code> <code>exu.SetWriteToFile(filename='testOutput.log', flagWriteToFile=True, flagAppend=False)</code> <code>exu.Print('print this to file')</code> <code>exu.SetWriteToFile("", False) #terminate writing to file which closes the file</code>
SetPrintDelayMilliseconds(delayMilliseconds)	add some delay (in milliseconds) to printing to console, in order to let Spyder process the output; default = 0
Print()	this allows printing via exudyn with similar syntax as in python <code>print(args)</code> except for keyword arguments: <code>print('test=',42)</code> ; allows to redirect all output to file given by <code>SetWriteToFile(...)</code> ; does not output in case that <code>SetWriteToConsole</code> is set to false
variables	this dictionary may be used by the user to store exudyn-wide data in order to avoid global python variables; usage: <code>exu.variables["myvar"] = 42</code>
sys	this dictionary is used by the system, e.g. for testsuite or solvers to store exudyn-wide data in order to avoid global python variables

## 4.2 SystemContainer

The SystemContainer is the top level of structures in EXUDYN. The container holds all systems, solvers and all other data structures for computation. Currently, only one container shall be used. In future, multiple containers might be usable at the same time.

Example:

```
import exudyn as exu
SC = exu.SystemContainer()
mbs = SC.AddSystem()
```

function/structure name	description
AddSystem()	add a new computational system
Reset()	delete all systems and reset SystemContainer (including graphics)
NumberOfSystems()	obtain number of systems available in system container
GetSystem(systemNumber)	obtain systems with index from system container
WaitForRenderEngineStopFlag()	Wait for user to stop render engine (Press 'Q' or Escape-key)

RenderEngineZoomAll()	Send zoom all signal, which will perform zoom all at next redraw request
GetRenderState()	Get dictionary with current render state (openGL zoom, modelview, etc.) <b>EXAMPLE:</b> <code>SC = exu.SystemContainer() renderState = SC.GetRenderState() print(renderState['zoom'])</code>
SetRenderState(renderState)	Set current render state (openGL zoom, modelview, etc.) with given dictionary; usually, this dictionary has been obtained with GetRenderState <b>EXAMPLE:</b> <code>SC = exu.SystemContainer() SC.GetRenderState(renderState)</code>
RedrawAndSaveImage()	Redraw openGL scene and save image (command waits until process is finished)
TimeIntegrationSolve(mainSystem, solverName, simulationSettings)	Call time integration solver for given system with solverName ('RungeKutta1'...explicit solver, 'GeneralizedAlpha'...implicit solver); use simulationSettings to individually configure the solver <b>EXAMPLE:</b> <code>simSettings = exu.SimulationSettings() simSettings.timeIntegration.numberOfSteps = 1000 simSettings.timeIntegration.endTime = 2 simSettings.timeIntegration.verboseMode = 1 SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', simSettings)</code>
StaticSolve(mainSystem, simulationSettings)	Call solver to compute a static solution of the system, considering acceleration and velocity coordinates to be zero (initial velocities may be considered by certain objects) <b>EXAMPLE:</b> <code>simSettings = exu.SimulationSettings() simSettings.staticSolver.newton.relativeTolerance = 1e-6 SC.StaticSolve(mbs, simSettings)</code>
visualizationSettings	this structure is read/writeable and contains visualization settings, which are immediately applied to the rendering window. <b>EXAMPLE:</b> <code>SC = exu.SystemContainer() SC.visualizationSettings.autoFitScene=False</code>

## 4.3 MainSystem

This is the structure which defines a (multibody) system. In C++, there is a MainSystem (links to python) and a System (computational part). For that reason, the name is MainSystem on the python

side, but it is often just called 'system'. It can be created, visualized and computed. Use the following functions for system manipulation.

Usage:

```
import exudyn as exu
SC = exu.SystemContainer()
mbs = SC.AddSystem()
```

function/structure name	description
Assemble()	assemble items (nodes, bodies, markers, loads, ...); Calls CheckSystemIntegrity(...), AssembleCoordinates(), AssembleLTGLists(), and AssembleInitializeSystemCoordinates()
AssembleCoordinates()	assemble coordinates: assign computational coordinates to nodes and constraints (algebraic variables)
AssembleLTGLists()	build local-to-global (ltg) coordinate lists for objects (used to build global ODE2RHS, MassMatrix, etc. vectors and matrices)
AssembleInitializeSystemCoordinates()	initialize all system-wide coordinates based on initial values given in nodes
Reset()	reset all lists of items (nodes, bodies, markers, loads, ...) and temporary vectors; deallocate memory
WaitForUserToContinue()	interrupt further computation until user input -> 'pause' function
SendRedrawSignal()	this function is used to send a signal to the renderer that the scene shall be redrawn because the visualization state has been updated
GetRenderEngineStopFlag()	get the current stop simulation flag; true=user wants to stop simulation
SetRenderEngineStopFlag()	set the current stop simulation flag; set to false, in order to continue a previously user-interrupted simulation
SetPreStepUserFunction()	Sets a user function PreStepUserFunction(mbs, t) executed at beginning of every computation step; in normal case return True; return False to stop simulation after current step <b>EXAMPLE:</b> <pre>def PreStepUserFunction(mbs, t):     print(mbs.systemData.NumberOfNodes())     if(t&gt;1):         return False     return True mbs.SetPreStepUserFunction(PreStepUserFunction)</pre>
__repr__()	return the representation of the system, which can be, e.g., printed <b>EXAMPLE:</b> <pre>print(mbs)</pre>

systemIsConsistent	this flag is used by solvers to decide, whether the system is in a solvable state; this flag is set to false as long as Assemble() has not been called; any modification to the system, such as Add...(), Modify...(), etc. will set the flag to false again; this flag can be modified (set to true), if a change of e.g. an object (change of stiffness) or load (change of force) keeps the system consistent, but would normally lead to systemIsConsistent=False
interactiveMode	set this flag to true in order to invoke a Assemble() command in every system modification, e.g. AddNode, AddObject, ModifyNode, ...; this helps that the system can be visualized in interactive mode.
variables	this dictionary may be used by the user to store model-specific data, in order to avoid global python variables in complex models; mbs.variables["myvar"] = 42
sys	this dictionary is used by exudyn python libraries, e.g., solvers, to avoid global python variables
solverSignalJacobianUpdate	this flag is used by solvers to decide, whether the jacobian should be updated; at beginning of simulation and after jacobian computation, this flag is set automatically to False; use this flag to indicate system changes, e.g. during time integration
systemData	Access to SystemData structure; enables access to number of nodes, objects, ... and to (current, initial, reference, ...) state variables (ODE2, AE, Data,...)

#### 4.3.1 MainSystem: Node

This section provides functions for adding, reading and modifying nodes. Nodes are used to define coordinates (unknowns to the static system and degrees of freedom if constraints are not present). Nodes can provide various types of coordinates for second/first order differential equations (ODE2/ODE1), algebraic equations (AE) and for data (history) variables – which are not providing unknowns in the nonlinear solver but will be solved in an additional nonlinear iteration for e.g. contact, friction or plasticity.

function/structure name	description
-------------------------	-------------

AddNode(pyObject)	<p>add a node with nodeDefinition from Python node class; returns (global) node index (type NodeIndex) of newly added node; use int(nodeIndex) to convert to int, if needed (but not recommended in order not to mix up index types of nodes, objects, markers, ...)</p> <p><b>EXAMPLE:</b></p> <pre> item = Rigid2D( referenceCoordinates= [1,0.5,0], initialVelocities= [10,0,0]) mbs.AddNode(item) nodeDict = {'nodeType': 'Point', 'referenceCoordinates': [1.0, 0.0, 0.0], 'initialCoordinates': [0.0, 2.0, 0.0], 'name': 'example node'} mbs.AddNode(nodeDict) </pre>
GetNodeNumber(nodeName)	<p>get node's number by name (string)</p> <p><b>EXAMPLE:</b></p> <pre>n = mbs.GetNodeNumber('example node')</pre>
GetNode(nodeNumber)	<p>get node's dictionary by node number (type NodeIndex)</p> <p><b>EXAMPLE:</b></p> <pre>nodeDict = mbs.GetNode(0)</pre>
ModifyNode(nodeNumber, nodeDict)	<p>modify node's dictionary by node number (type NodeIndex)</p> <p><b>EXAMPLE:</b></p> <pre>mbs.ModifyNode(nodeNumber, nodeDict)</pre>
GetNodeDefaults(typeName)	<p>get node's default values for a certain nodeType as (dictionary)</p> <p><b>EXAMPLE:</b></p> <pre>nodeType = 'Point' nodeDict = mbs.GetNodeDefaults(nodeType)</pre>
GetNodeOutput(nodeNumber, variableType, configuration = ConfigurationType.Current)	<p>get the output of the node specified with the OutputVariableType; default configuration = 'current'; output may be scalar or array (e.g. displacement vector)</p> <p><b>EXAMPLE:</b></p> <pre>mbs.GetNodeOutput(nodeNumber=0, variableType='exu.OutputVariable.Displacement')</pre>
GetNodeODE2Index(nodeNumber)	<p>get index in the global ODE2 coordinate vector for the first node coordinate of the specified node</p> <p><b>EXAMPLE:</b></p> <pre>mbs.GetNodeODE2Index(nodeNumber=0)</pre>
GetNodeParameter(nodeNumber, parameterName)	<p>get nodes's parameter from node number (type NodeIndex) and parameterName; parameter names can be found for the specific items in the reference manual</p>
SetNodeParameter(nodeNumber, parameterName, value)	<p>set parameter 'parameterName' of node with node number (type NodeIndex) to value; parameter names can be found for the specific items in the reference manual</p>

### 4.3.2 MainSystem: Object

This section provides functions for adding, reading and modifying objects, which can be bodies (mass point, rigid body, finite element, ...), connectors (spring-damper or joint) or general objects. Objects provided terms to the residual of equations resulting from every coordinate given by the nodes. Single-noded objects (e.g. mass point) provides exactly residual terms for its nodal coordinates. Connectors constrain or penalize two markers, which can be, e.g., position, rigid or coordinate markers. Thus, the dependence of objects is either on the coordinates of the marker-objects/nodes or on nodes which the objects possess themselves.

function/structure name	description
AddObject(pyObject)	add an object with objectDefinition from Python object class; returns (global) object number (type ObjectIndex) of newly added object <b>EXAMPLE:</b> <pre>item = MassPoint(name='heavy object', nodeNumber=0, physicsMass=100) mbs.AddObject(item) objectDict = {'objectType': 'MassPoint', 'physicsMass': 10, 'nodeNumber': 0, 'name': 'example object'} mbs.AddObject(objectDict)</pre>
GetObjectNumber(objectName)	get object's number by name (string) <b>EXAMPLE:</b> <pre>n = mbs.GetObjectNumber('heavy object')</pre>
GetObject(objectNumber)	get object's dictionary by object number (type ObjectIndex) <b>EXAMPLE:</b> <pre>objectDict = mbs.GetObject(0)</pre>
ModifyObject(objectNumber, objectDict)	modify object's dictionary by object number (type ObjectIndex) <b>EXAMPLE:</b> <pre>mbs.ModifyObject(objectNumber, objectDict)</pre>
GetObjectDefaults(typeName)	get object's default values for a certain objectType as (dictionary) <b>EXAMPLE:</b> <pre>objectType = 'MassPoint' objectDict = mbs.GetObjectDefaults(objectType)</pre>
GetObjectOutput(objectNumber, variableType)	get object's current output variable from object number (type ObjectIndex) and OutputVariableType; can only be computed for exu.ConfigurationType.Current configuration!

GetObjectOutputBody(objectNumber, variableType, localPosition, configuration = ConfigurationType.Current)	get body's output variable from object number (type ObjectIndex) and OutputVariableType <b>EXAMPLE:</b> u = mbs.GetObjectOutputBody(objectNumber = 1, variableType = exu.OutputVariableType.Position, localPosition=[1,0,0], configuration = exu.ConfigurationType.Initial)
GetObjectOutputSuperElement(objectNumber, variableType, meshNodeNumber, configuration = ConfigurationType.Current)	get output variable from mesh node number of object with type SuperElement (GenericODE2, FFRF, FFRFReduced - CMS) with specific OutputVariableType; the meshNodeNumber is the object's local node number, not the global node number! <b>EXAMPLE:</b> u = mbs.GetObjectOutputSuperElement(objectNumber = 1, variableType = exu.OutputVariableType.Position, meshNodeNumber = 12, configuration = exu.ConfigurationType.Initial)
GetObjectParameter(objectNumber, parameterName)	get objects's parameter from object number (type ObjectIndex) and parameterName; parameter names can be found for the specific items in the reference manual
SetObjectParameter(objectNumber, parameterName, value)	set parameter 'parameterName' of object with object number (type ObjectIndex) to value; parameter names can be found for the specific items in the reference manual

### 4.3.3 MainSystem: Marker

This section provides functions for adding, reading and modifying markers. Markers define how to measure primal kinematical quantities on objects or nodes (e.g., position, orientation or coordinates themselves), and how to act on the quantities which are dual to the kinematical quantities (e.g., force, torque and generalized forces). Markers provide unique interfaces for loads, sensors and constraints in order to address these quantities independently of the structure of the object or node (e.g., rigid or flexible body).

function/structure name	description
AddMarker(pyObject)	add a marker with markerDefinition from Python marker class; returns (global) marker number (type MarkerIndex) of newly added marker <b>EXAMPLE:</b> item = MarkerNodePosition(name='my marker', nodeNumber=1) mbs.AddMarker(item) markerDict = {'markerType': 'NodePosition', 'nodeNumber': 0, 'name': 'position0'} mbs.AddMarker(markerDict)



GetMarkerNumber(markerName)	get marker's number by name (string) <b>EXAMPLE:</b> <code>n = mbs.GetMarkerNumber('my marker')</code>
GetMarker(markerNumber)	get marker's dictionary by index <b>EXAMPLE:</b> <code>markerDict = mbs.GetMarker(0)</code>
ModifyMarker(markerNumber, markerDict)	modify marker's dictionary by index <b>EXAMPLE:</b> <code>mbs.ModifyMarker(markerNumber, markerDict)</code>
GetMarkerDefaults(typeName)	get marker's default values for a certain markerType as (dictionary) <b>EXAMPLE:</b> <code>markerType = 'NodePosition'</code> <code>markerDict = mbs.GetMarkerDefaults(markerType)</code>
GetMarkerParameter(markerNumber, parameterName)	get markers's parameter from markerNumber and parameterName; parameter names can be found for the specific items in the reference manual
SetMarkerParameter(markerNumber, parameterName, value)	set parameter 'parameterName' of marker with markerNumber to value; parameter names can be found for the specific items in the reference manual

#### 4.3.4 MainSystem: Load

This section provides functions for adding, reading and modifying operating loads. Loads are used to act on the quantities which are dual to the primal kinematic quantities, such as displacement and rotation. Loads represent, e.g., forces, torques or generalized forces.

function/structure name	description
AddLoad(pyObject)	add a load with loadDefinition from Python load class; returns (global) load number (type LoadIndex) of newly added load <b>EXAMPLE:</b> <code>item = mbs.AddLoad(LoadForceVector(loadVector=[1,0,0], markerNumber=0, name='heavy load'))</code> <code>mbs.AddLoad(item)</code> <code>loadDict = {'loadType': 'ForceVector', 'markerNumber': 0, 'loadVector': [1.0, 0.0, 0.0], 'name': 'heavy load'}</code> <code>mbs.AddLoad(loadDict)</code>
GetLoadNumber(loadName)	get load's number by name (string) <b>EXAMPLE:</b> <code>n = mbs.GetLoadNumber('heavy load')</code>
GetLoad(loadNumber)	get load's dictionary by index <b>EXAMPLE:</b> <code>loadDict = mbs.GetLoad(0)</code>

ModifyLoad(loadNumber, loadDict)	modify load's dictionary by index <b>EXAMPLE:</b> <code>mbs.ModifyLoad(loadNumber, loadDict)</code>
GetLoadDefaults(typeName)	get load's default values for a certain loadType as (dictionary) <b>EXAMPLE:</b> <code>loadType = 'ForceVector'</code> <code>loadDict = mbs.GetLoadDefaults(loadType)</code>
GetLoadValues(loadNumber)	Get current load values, specifically if user-defined loads are used; can be scalar or vector-valued return value
GetLoadParameter(loadNumber, parameterName)	get loads's parameter from loadNumber and parameterName; parameter names can be found for the specific items in the reference manual
SetLoadParameter(loadNumber, parameterName, value)	set parameter 'parameterName' of load with loadNumber to value; parameter names can be found for the specific items in the reference manual

#### 4.3.5 MainSystem: Sensor

This section provides functions for adding, reading and modifying operating sensors. Sensors are used to measure information in nodes, objects, markers, and loads for output in a file.

function/structure name	description
AddSensor(pyObject)	add a sensor with sensor definition from Python sensor class; returns (global) sensor number (type SensorIndex) of newly added sensor <b>EXAMPLE:</b> <code>item = mbs.AddSensor(SensorNode(sensorType=exu.SensorType.Node, nodeNumber=0, name='test sensor'))</code> <code>mbs.AddSensor(item)</code> <code>sensorDict = {'sensorType': 'Node', 'nodeNumber': 0, 'fileName': 'sensor.txt', 'name': 'test sensor'}</code> <code>mbs.AddSensor(sensorDict)</code>
GetSensorNumber(sensorName)	get sensor's number by name (string) <b>EXAMPLE:</b> <code>n = mbs.GetSensorNumber('test sensor')</code>
GetSensor(sensorNumber)	get sensor's dictionary by index <b>EXAMPLE:</b> <code>sensorDict = mbs.GetSensor(0)</code>
ModifySensor(sensorNumber, sensorDict)	modify sensor's dictionary by index <b>EXAMPLE:</b> <code>mbs.ModifySensor(sensorNumber, sensorDict)</code>

GetSensorDefaults(typeName)	get sensor's default values for a certain sensorType as (dictionary) <b>EXAMPLE:</b> <code>sensorType = 'Node'</code> <code>sensorDict = mbs.GetSensorDefaults(sensorType)</code>
GetSensorValues(sensorNumber, configuration = ConfigurationType.Current)	get sensors's values for configuration; can be a scalar or vector-valued return value!
GetSensorParameter(sensorNumber, parameterName)	get sensors's parameter from sensorNumber and parameterName; parameter names can be found for the specific items in the reference manual
SetSensorParameter(sensorNumber, parameterName, value)	set parameter 'parameterName' of sensor with sensorNumber to value; parameter names can be found for the specific items in the reference manual

## 4.4 SystemData

This is the data structure of a system which contains Objects (bodies/constraints/...), Nodes, Markers and Loads. The SystemData structure allows advanced access to this data, which HAS TO BE USED WITH CARE, as unexpected results and system crash might happen.

Usage:

#obtain current ODE2 system vector (e.g. after static simulation finished):

`u = mbs.systemData.GetODE2Coordinates()`

#set initial ODE2 vector for next simulation:

`mbs.systemData.SetODE2Coordinates(coordinates=u, configurationType=exu.ConfigurationType.Initial)`

function/structure name	description
NumberOfLoads()	return number of loads in system <b>EXAMPLE:</b> <code>print(mbs.systemData.NumberOfLoads())</code>
NumberOfMarkers()	return number of markers in system <b>EXAMPLE:</b> <code>print(mbs.systemData.NumberOfMarkers())</code>
NumberOfNodes()	return number of nodes in system <b>EXAMPLE:</b> <code>print(mbs.systemData.NumberOfNodes())</code>
NumberOfObjects()	return number of objects in system <b>EXAMPLE:</b> <code>print(mbs.systemData.NumberOfObjects())</code>
NumberOfSensors()	return number of sensors in system <b>EXAMPLE:</b> <code>print(mbs.systemData.NumberOfSensors())</code>
GetTime(configurationType exu.ConfigurationType.Current)	= get configuration dependent time. <b>EXAMPLE:</b> <code>mbs.systemData.GetTime(exu.ConfigurationType.Initial)</code>

SetTime(newTime, configurationType exu.ConfigurationType.Current)	=	set configuration dependent time; use this access with care, e.g. in user-defined solvers. <b>EXAMPLE:</b> <code>mbs.systemData.SetTime(10., exu.ConfigurationType.Initial)</code>
GetCurrentTime()		DEPRICATED; get current (simulation) time; time is updated in time integration solvers and in static solver; use this function e.g. during simulation to define time-dependent loads <b>EXAMPLE:</b> <code>mbs.systemData.GetCurrentTime()</code>
SetVisualizationTime()		DEPRICATED; set time for render window (visualization) <b>EXAMPLE:</b> <code>mbs.systemData.SetVisualizationTime(1.3)</code>
Info()		print detailed system information for every item; for short information use print(mbs) <b>EXAMPLE:</b> <code>mbs.systemData.Info()</code>

#### 4.4.1 SystemData: Access coordinates

This section provides access functions to global coordinate vectors. Assigning invalid values or using wrong vector size might lead to system crash and unexpected results.

function/structure name		description
GetODE2Coordinates(configuration exu.ConfigurationType.Current)	=	get ODE2 system coordinates (displacements) for given configuration (default: exu.Configuration.Current) <b>EXAMPLE:</b> <code>uCurrent = mbs.systemData.GetODE2Coordinates()</code>
SetODE2Coordinates(coordinates, configuration exu.ConfigurationType.Current)	=	set ODE2 system coordinates (displacements) for given configuration (default: exu.Configuration.Current); invalid vector size may lead to system crash! <b>EXAMPLE:</b> <code>mbs.systemData.SetODE2Coordinates(uCurrent)</code>
GetODE2Coordinates_t(configuration exu.ConfigurationType.Current)	=	get ODE2 system coordinates (velocities) for given configuration (default: exu.Configuration.Current) <b>EXAMPLE:</b> <code>vCurrent = mbs.systemData.GetODE2Coordinates_t()</code>
SetODE2Coordinates_t(coordinates, configuration exu.ConfigurationType.Current)	=	set ODE2 system coordinates (velocities) for given configuration (default: exu.Configuration.Current); invalid vector size may lead to system crash! <b>EXAMPLE:</b> <code>mbs.systemData.SetODE2Coordinates_t(vCurrent)</code>
GetODE2Coordinates_tt(configuration exu.ConfigurationType.Current)	=	get ODE2 system coordinates (accelerations) for given configuration (default: exu.Configuration.Current) <b>EXAMPLE:</b> <code>vCurrent = mbs.systemData.GetODE2Coordinates_tt()</code>

SetODE2Coordinates_tt(coordinates, configuration = exu.ConfigurationType.Current)	=	set ODE2 system coordinates (accelerations) for given configuration (default: exu.Configuration.Current); invalid vector size may lead to system crash! <b>EXAMPLE:</b> <code>mbs.systemData.SetODE2Coordinates_tt(aCurrent)</code>
GetODE1Coordinates(configuration = exu.ConfigurationType.Current)	=	get ODE1 system coordinates (displacements) for given configuration (default: exu.Configuration.Current) <b>EXAMPLE:</b> <code>qCurrent = mbs.systemData.GetODE1Coordinates()</code>
SetODE1Coordinates(coordinates, configuration = exu.ConfigurationType.Current)	=	set ODE1 system coordinates (displacements) for given configuration (default: exu.Configuration.Current); invalid vector size may lead to system crash! <b>EXAMPLE:</b> <code>mbs.systemData.SetODE1Coordinates(qCurrent)</code>
GetAECordinates(configuration = exu.ConfigurationType.Current)	=	get algebraic equations (AE) system coordinates for given configuration (default: exu.Configuration.Current) <b>EXAMPLE:</b> <code>lambdaCurrent = mbs.systemData.GetAECordinates()</code>
SetAECordinates(coordinates, configuration = exu.ConfigurationType.Current)	=	set algebraic equations (AE) system coordinates for given configuration (default: exu.Configuration.Current); invalid vector size may lead to system crash! <b>EXAMPLE:</b> <code>mbs.systemData.SetAECordinates(lambdaCurrent)</code>
GetDataCoordinates(configuration = exu.ConfigurationType.Current)	=	get system data coordinates for given configuration (default: exu.Configuration.Current) <b>EXAMPLE:</b> <code>dataCurrent = mbs.systemData.GetDataCoordinates()</code>
SetDataCoordinates(coordinates, configuration = exu.ConfigurationType.Current)	=	set system data coordinates for given configuration (default: exu.Configuration.Current); invalid vector size may lead to system crash! <b>EXAMPLE:</b> <code>mbs.systemData.SetDataCoordinates(dataCurrent)</code>
GetSystemState(configuration = exu.ConfigurationType.Current)	=	get system state for given configuration (default: exu.Configuration.Current); state vectors do not include the non-state derivatives ODE1_t and ODE2_tt and the time; function is copying data - not highly efficient; format of pyList: [ODE2Coords, ODE2Coords_t, ODE1Coords, AEcoords, dataCoords] <b>EXAMPLE:</b> <code>sysStateList = mbs.systemData.GetSystemState()</code>
SetSystemState(systemStateList, configuration = exu.ConfigurationType.Current)	=	set system data coordinates for given configuration (default: exu.Configuration.Current); invalid list of vectors / vector size may lead to system crash; write access to state vectors (but not the non-state derivatives ODE1_t and ODE2_tt and the time); function is copying data - not highly efficient; format of pyList: [ODE2Coords, ODE2Coords_t, ODE1Coords, AEcoords, dataCoords] <b>EXAMPLE:</b> <code>mbs.systemData.SetDataCoordinates(sysStateList, configuration = exu.ConfigurationType.Initial)</code>

## 4.4.2 SystemData: Get object local-to-global (LTG) coordinate mappings

This section provides access functions the LTG-lists for every object (body, constraint, ...) in the system.

function/structure name	description
GetObjectLTGODE2(objectNumber)	get local-to-global coordinate mapping (list of global coordinate indices) for ODE2 coordinates; only available after Assemble() <b>EXAMPLE:</b> <code>ltgObject4 = mbs.systemData.GetObjectLTGODE2(4)</code>
GetObjectLTGODE1(objectNumber)	get local-to-global coordinate mapping (list of global coordinate indices) for ODE1 coordinates; only available after Assemble() <b>EXAMPLE:</b> <code>ltgObject4 = mbs.systemData.GetObjectLTGODE1(4)</code>
GetObjectLTGAE(objectNumber)	get local-to-global coordinate mapping (list of global coordinate indices) for algebraic equations (AE) coordinates; only available after Assemble() <b>EXAMPLE:</b> <code>ltgObject4 = mbs.systemData.GetObjectLTGODE2(4)</code>
GetObjectLTGData(objectNumber)	get local-to-global coordinate mapping (list of global coordinate indices) for data coordinates; only available after Assemble() <b>EXAMPLE:</b> <code>ltgObject4 = mbs.systemData.GetObjectLTGData(4)</code>

## 4.5 Type definitions

This section defines a couple of structures, which are used to select, e.g., a configuration type or a variable type. In the background, these types are integer numbers, but for safety, the types should be used as type variables.

Conversion to integer is possible:

```
x = int(exu.OutputVariableType.Displacement)
```

and also conversion from integer:

```
varType = exu.OutputVariableType(8)
```

### 4.5.1 OutputVariableType

This section shows the OutputVariableType structure, which is used for selecting output values, e.g. for GetObjectOutput(...) or for selecting variables for contour plot.

Available output variables and the interpretation of the output variable can be found at the object definitions. The OutputVariableType does not provide information about the size of the output variable, which can be either scalar or a list (vector). For vector output quantities, the contour plot option offers an additional parameter for selection of the component of the OutputVariableType.

function/structure name	description
_None	no value; used, e.g., to select no output variable in contour plot
Distance	e.g., measure distance in spring damper connector
Position	measure 3D position, e.g., of node or body
Displacement	measure displacement; usually difference between current position and reference position
Velocity	measure (translational) velocity of node or object
Acceleration	measure (translational) acceleration of node or object
RotationMatrix	measure rotation matrix of rigid body node or object
AngularVelocity	measure angular velocity of node or object
AngularVelocityLocal	measure local (body-fixed) angular velocity of node or object
AngularAcceleration	measure angular acceleration of node or object
Rotation	measure, e.g., scalar rotation of 2D body, Euler angles of a 3D object or rotation within a joint
Coordinates	measure the coordinates of a node or object; coordinates usually just contain displacements, but not the position values
Coordinates_t	measure the time derivative of coordinates (= velocity coordinates) of a node or object
Coordinates_tt	measure the second time derivative of coordinates (= acceleration coordinates) of a node or object
SlidingCoordinate	measure sliding coordinate in sliding joint
Director1	measure a director (e.g. of a rigid body frame), or a slope vector in local 1 or x-direction
Director2	measure a director (e.g. of a rigid body frame), or a slope vector in local 2 or y-direction
Director3	measure a director (e.g. of a rigid body frame), or a slope vector in local 3 or z-direction
Force	measure force, e.g., in joint or beam (resultant force)
Torque	measure torque, e.g., in joint or beam (resultant couple/-moment)
Strain	measure strain, e.g., axial strain in beam
Stress	measure stress, e.g., axial stress in beam
Curvature	measure curvature; may be scalar or vectorial: twist and curvature
DisplacementLocal	measure local displacement, e.g. in local joint coordinates
VelocityLocal	measure local (translational) velocity, , e.g. in local joint coordinates
ForceLocal	measure local force, e.g., in joint or beam (resultant force)
TorqueLocal	measure local torque, e.g., in joint or beam (resultant couple/moment)
ConstraintEquation	evaluates constraint equation (=current deviation or drift of constraint equation)
EndOfEnumList	this marks the end of the list, usually not important to the user

### 4.5.2 ConfigurationType

This section shows the ConfigurationType structure, which is used for selecting a configuration for reading or writing information to the module. Specifically, the ConfigurationType.Current configuration is usually used at the end of a solution process, to obtain result values, or the ConfigurationType.Initial is used to set initial values for a solution process.

function/structure name	description
_None	no configuration; usually not valid, but may be used, e.g., if no configurationType is required
Initial	initial configuration prior to static or dynamic solver; is computed during mbs.Assemble() or AssembleInitializeSystemCoordinates()
Current	current configuration during and at the end of the computation of a step (static or dynamic)
Reference	configuration used to define deformable bodies (reference configuration for finite elements) or joints (configuration for which some joints are defined)
StartOfStep	during computation, this refers to the solution at the start of the step = end of last step, to which the solver falls back if convergence fails
Visualization	this is a state completely de-coupled from computation, used for visualization
EndOfEnumList	this marks the end of the list, usually not important to the user

### 4.5.3 LinearSolverType

This section shows the LinearSolverType structure, which is used for selecting output values, e.g. for GetObjectOutput(...) or for selecting variables for contour plot.

function/structure name	description
_None	no value; used, e.g., if no solver is selected
EXUdense	use dense matrices and according solvers for densely populated matrices (usually the CPU time grows cubically with the number of unknowns)
EigenSparse	use sparse matrices and according solvers; additional overhead for very small systems; specifically, memory allocation is performed during a factorization process





# Chapter 5

## Python utility functions

This chapter describes in every subsection the functions and classes of the utility modules. These modules help to create multibody systems with the EXUDYN core module. Functions are implemented in Python and can be easily changed, extended and also verified by the user. They are much slower than the functions available in the C++ core. Some matrix computations with larger matrices implemented in numpy and scipy, however, are parallelized and therefore very efficient.

Functions have been implemented, if not otherwise noted, by Johannes Gerstmayr.

### 5.1 Module: basicUtilities

def **DiagonalMatrix**(*rowsColumns*, *value*=1)

- **function description:** create a diagonal or identity matrix; used for interface.py, avoiding the need for numpy
- **input:**
  - rowsColumns*: provides the number of rows and columns
  - value*: initialization value for diagonal terms
- **output:** list of lists representing a matrix

---

def **NormL2**(*vector*)

- **function description:** compute L2 norm for vectors without switching to numpy or math module
- **input:** vector as list or in numpy format
- **output:** L2-norm of vector

---

def **VSum**(*vector*)

- **function description:** compute sum of all values of vector
- **input:** vector as list or in numpy format
- **output:** sum of all components of vector

---

def **VAdd**(*v0*, *v1*)

- **function description:** add two vectors instead using numpy
- **input:** vectors *v0* and *v1* as list or in numpy format
- **output:** component-wise sum of *v0* and *v1*

---

def **VSub**(*v0*, *v1*)

- **function description:** subtract two vectors instead using numpy: result = *v0-v1*
- **input:** vectors *v0* and *v1* as list or in numpy format
- **output:** component-wise difference of *v0* and *v1*

---

def **VMult**(*v0*, *v1*)

- **function description:** scalar multiplication of two vectors instead using numpy: result = *v0'\*v1*
- **input:** vectors *v0* and *v1* as list or in numpy format
- **output:** sum of all component wise products: *c0[0]\*v1[0] + v0[1]\*v1[0] + ...*

---

def **ScalarMult**(*scalar*, *v*)

- **function description:** multiplication vectors with scalar: result = *s\*v*
- **input:** value *scalar* and vector *v* as list or in numpy format
- **output:** scalar multiplication of all components of *v*: [*scalar\*v[0]*, *scalar\*v[1]*, ...]

---

def **Normalize**(*v*)

- **function description:** take a 3D vector and return a normalized 3D vector (L2Norm=1)
- **input:** vector *v* as list or in numpy format
- **output:** vector *v* multiplied with scalar such that L2-norm of vector is 1

---

def **Vec2Tilde**(*v*)

- **function description:** apply tilde operator (skew) to 3D-vector and return skew matrix
- **input:** 3D vector *v* as list or in numpy format

- **output:** matrix as list of lists containing the skew-symmetric matrix computed from *v*: 
$$\begin{bmatrix} 0 & -v[2] & v[1] \\ v[2] & 0 & -v[0] \\ -v[1] & v[0] & 0 \end{bmatrix}$$

---

def **Tilde2Vec**(*m*)

- **function description:** take skew symmetric matrix and return vector (inverse of Skew(...))
- **input:** list of lists containing a skew-symmetric matrix (3x3)
- **output:** list containing the vector *v* (inverse function of Vec2Tilde(...))

## 5.2 Module: utilities

def **PlotLineCode**(*index*)

- **function description:** helper functions for matplotlib, returns a list of 28 line codes to be used in plot, e.g. 'r-' for red solid line
- **input:** index in range(0:28)
- **output:** a color and line style code for matplotlib plot

---

def **FillInSubMatrix**(*subMatrix*, *destinationMatrix*, *destRow*, *destColumn*)

- **function description:** fill submatrix into given destinationMatrix; all matrices must be numpy arrays
- **input:**
  - subMatrix*: input matrix, which is filled into destinationMatrix
  - destinationMatrix*: the subMatrix is entered here
  - destRow*: row destination of subMatrix
  - destColumn*: column destination of subMatrix
- **output:** destinationMatrix is changed after function call
- **notes:** may be erased in future!

---

def **SweepSin**(*t*, *t1*, *f0*, *f1*)

- **function description:** compute sin sweep at given time *t*
- **input:**
  - t*: evaluate of sweep at time *t*
  - t1*: end time of sweep frequency range
  - f0*: start of frequency interval [*f0*,*f1*] in Hz
  - f1*: end of frequency interval [*f0*,*f1*] in Hz
- **output:** evaluation of sin sweep (in range -1..+1)

---

def **SweepCos**(*t*, *t1*, *f0*, *f1*)

- **function description:** compute cos sweep at given time  $t$
- **input:**
  - $t$ : evaluate of sweep at time  $t$
  - $t1$ : end time of sweep frequency range
  - $f0$ : start of frequency interval  $[f0, f1]$  in Hz
  - $f1$ : end of frequency interval  $[f0, f1]$  in Hz
- **output:** evaluation of cos sweep (in range -1..+1)

---

def **FrequencySweep**( $t, t1, f0, f1$ )

- **function description:** frequency according to given sweep functions SweepSin, SweepCos
- **input:**
  - $t$ : evaluate of frequency at time  $t$
  - $t1$ : end time of sweep frequency range
  - $f0$ : start of frequency interval  $[f0, f1]$  in Hz
  - $f1$ : end of frequency interval  $[f0, f1]$  in Hz
- **output:** frequency in Hz

---

def **RoundMatrix**( $matrix, threshold=1e-14$ )

- **function description:** set all entries in matrix to zero which are smaller than given treshold; operates directly on matrix
- **input:** matrix as np.array, treshold as positive value
- **output:** changes matrix

---

def **ComputeSkewMatrix**( $v$ )

- **function description:** compute  $(3 \times 3 \times n)$  skew matrix from  $(3 \times n)$  vector; used for ObjectFFRF and CMS implementation
- **input:** a vector  $v$  in np.array format, containing  $3 \times n$  components
- **output:**  $(3 \times 3 \times n)$  skew matrix in np.array format

---

def **CheckInputVector**( $vector, length=-1$ )

- **function description:** check if input is list or array with according length; if  $length == -1$ , the length is not checked; raises Exception if the check fails
- **input:**
  - $vector$ : a vector in np.array or list format

*length*: desired length of vector; if *length*=-1, it is ignored

- **output**: None

---

def **CheckInputIndexArray**(*indexArray*, *length*=-1)

- **function description**: check if input is list or array with according length and positive indices; if *length*==1, the length is not checked; raises Exception if the check fails
- **input**:
  - indexArray*: a vector in np.array or list format
  - length*: desired length of vector; if *length*=-1, it is ignored
- **output**: None

---

def **LoadSolutionFile**(*fileName*)

- **function description**: read coordinates solution file (exported during static or dynamic simulation with option `exu.SimulationSettings().solutionSettings.coordinatesSolutionFileName='...'`) into dictionary:
- **input**: *fileName*: string containing directory and filename of stored coordinatesSolutionFile
- **output**: dictionary with 'data': the matrix of stored solution vectors, 'columnsExported': a list with binary values showing the exported columns [nODE2, nVel2, nAcc2, nODE1, nVel1, nAlgebraic, nData], 'nColumns': the number of data columns and 'nRows': the number of data rows

---

def **SetSolutionState**(*exu*, *mbs*, *solution*, *row*, *configuration*)

- **function description**: load selected row of solution dictionary (previously loaded with LoadSolutionFile) into specific state

---

def **SetVisualizationState**(*exu*, *mbs*, *solution*, *row*)

- **function description**: load selected row of solution dictionary into visualization state and redraw
- **input**:
  - exu*: the exudyn library
  - mbs*: the system, where the state is applied to
  - solution*: solution dictionary previously loaded with LoadSolutionFile
  - row*: the according row of the solution file which is visualized
- **output**: renders the scene in mbs and changes the visualization state in mbs

---

def **AnimateSolution**(*exu*, *SC*, *mbs*, *solution*, *rowIncrement*=1, *timeout*=0.04, *createImages*=False, *runLoop*=False)

- **function description**: consecutively load the rows of a solution file and visualize the result

- **input:**
  - exu*: the exudyn library
  - SC*: the system container, where the mbs lives in
  - mbs*: the system used for animation
  - solution*: solution dictionary previously loaded with LoadSolutionFile; will be played from first to last row
  - rowIncrement*: can be set larger than 1 in order to skip solution frames: e.g. *rowIncrement*=10 visualizes every 10th row (frame)
  - timeout*: in seconds is used between frames in order to limit the speed of animation; e.g. use *timeout*=0.04 to achieve approximately 25 frames per second
  - createImages*: creates consecutively images from the animation, which can be converted into an animation
  - runLoop*: if True, the animation is played in a loop until 'q' is pressed in render window
- **output:** renders the scene in mbs and changes the visualization state in mbs continuously

---

```
def DrawSystemGraph(mbs, showLoads=True, showSensors=True, useItemNames=False, useItemTypes=False)
```

- **function description:** helper function which draws system graph of a MainSystem (mbs); several options let adjust the appearance of the graph
- **input:**
  - showLoads*: toggle appearance of loads in mbs
  - showSensors*: toggle appearance of sensors in mbs
  - useItemNames*: if True, object names are shown instead of basic object types (Node, Load, ...)
  - useItemTypes*: if True, object type names (ObjectMassPoint, ...) are shown instead of basic object types (Node, Load, ...)
- **output:** nothing

---

```
def GenerateStraightLineANCFcable2D(mbs, positionOfNode0, positionOfNode1, numberOfElements, cableTemplate, massProportionalLoad=[0,0,0], fixedConstraintsNode0=[0,0,0,0], fixedConstraintsNode1=[0,0,0,0], vALE=0, ConstrainAleCoordinate=True)
```

- **function description:** generate cable elements along straight line with certain discretization
- **input:**
  - mbs*: the system where ANCF cables are added
  - positionOfNode0*: 3D position (list or np.array) for starting point of line
  - positionOfNode1*: 3D position (list or np.array) for end point of line
  - numberOfElements*: for discretization of line
  - cableTemplate*: a ObjectANCFcable2D object, containing the desired cable properties; cable length and node numbers are set automatically
  - massProportionalLoad*: a 3D list or np.array, containing the gravity vector or zero

*fixedConstraintsNode0*: a list of 4 binary values, indicating the coordinate constraints on the first node (x,y-position and x,y-slope)

*fixedConstraintsNode1*: a list of 4 binary values, indicating the coordinate constraints on the last node (x,y-position and x,y-slope)

*vALE*: used for ObjectALEANCFcable2D objects

- **output**: returns a list [cableNodeList, cableObjectList, loadList, cableNodePositionList, cableCoordinateConstraintList]

---

def **GenerateSlidingJoint**(*mbs, cableObjectList, markerBodyPositionOfSlidingBody, localMarkerIndexOfStartCable=0, slidingCoordinateStartPosition=0*)

- **function description**: generate a sliding joint from a list of cables, marker to a sliding body, etc.

---

def **GenerateAleSlidingJoint**(*mbs, cableObjectList, markerBodyPositionOfSlidingBody, AleNode, localMarkerIndexOfStartCable=0, AleSlidingOffset=0, activeConnector=True, penaltyStiffness=0*)

- **function description**: generate an ALE sliding joint from a list of cables, marker to a sliding body, etc.

## 5.3 Module: graphicsDataUtilities

def **GraphicsDataRectangle**(*xMin, yMin, xMax, yMax, color=[0.,0.,0.,1.]*)

- **function description**: generate graphics data for 2D rectangle
- **input**: minimal and maximal cartesian coordinates in (x/y) plane; color provided as list of 4 RGBA values
- **output**: graphicsData dictionary, to be used in visualization of EXUDYN objects

---

def **GraphicsDataOrthoCubeLines**(*xMin, yMin, zMin, xMax, yMax, zMax, color=[0.,0.,0.,1.]*)

- **function description**: generate graphics data for orthogonal cube drawn with lines
- **input**: minimal and maximal cartesian coordinates for orthogonal cube; color provided as list of 4 RGBA values
- **output**: graphicsData dictionary, to be used in visualization of EXUDYN objects

---

def **GraphicsDataOrthoCube**(*xMin, yMin, zMin, xMax, yMax, zMax, color=[0.,0.,0.,1.]*)

- **function description**: generate graphics data for orthogonal 3D cube with min and max dimensions
- **input**: minimal and maximal cartesian coordinates for orthogonal cube; color provided as list of 4 RGBA values
- **output**: graphicsData dictionary, to be used in visualization of EXUDYN objects



---

def **GraphicsDataOrthoCubePoint**(*centerPoint*, *size*, *color*=[0.,0.,0.,1.])

- **function description:** generate graphics data for orthogonal 3D cube with center point and size
- **input:** center point and size of cube (as 3D list or np.array); color provided as list of 4 RGBA values
- **output:** graphicsData dictionary, to be used in visualization of EXUDYN objects

---

def **GraphicsDataCube**(*pList*, *color*=[0.,0.,0.,1.], *faces*=[1,1,1,1,1,1])

- **function description:** generate graphics data for general cube with endpoints, according to given vertex definition
- **input:**
  - pList*: is a list of points [[x0,y0,z0],[x1,y1,z1],...]
  - color*: provided as list of 4 RGBA values
  - faces*: includes the list of six binary values (0/1), denoting active faces (value=1); set index to zero to hide face
- **output:** graphicsData dictionary, to be used in visualization of EXUDYN objects

---

def **SwitchTripletOrder**(*vector*)

- **function description:** switch order of three items in a list; mostly used for reverting normals in triangles
- **input:** 3D vector as list or as np.array
- **output:** interchanged 2nd and 3rd component of list

---

def **GraphicsDataSphere**(*point*, *radius*, *color*=[0.,0.,0.,1.], *nTiles*=8)

- **function description:** generate graphics data for a sphere with point p and radius
- **input:**
  - point*: center of sphere (3D list or np.array)
  - radius*: positive value
  - color*: provided as list of 4 RGBA values
  - nTiles*: used to determine resolution of sphere  $\geq 3$ ; use larger values for finer resolution
- **output:** graphicsData dictionary, to be used in visualization of EXUDYN objects

---

def **GraphicsDataCylinder**(*pAxis*, *vAxis*, *radius*, *color*=[0.,0.,0.,1.], *nTiles*=16, *angleRange*=[0,2\*np.pi], *lastFace*=True, *cutPlain*=True)

- **function description:** generate graphics data for a cylinder with given axis, radius and color; nFaces gives the number of tiles (minimum=3)
- **input:**

*pAxis*: axis point of one face of cylinder (3D list or np.array)  
*vAxis*: vector representing the cylinder's axis (3D list or np.array)  
*radius*: positive value representing radius of cylinder  
*color*: provided as list of 4 RGBA values  
*nTiles*: used to determine resolution of cylinder  $\geq 3$ ; use larger values for finer resolution  
*angleRange*: given in rad, to draw only part of cylinder (halfcylinder, etc.); for full range use  $[0, 2 * \pi]$   
*lastFace*: if *angleRange*  $\neq [0, 2 * \pi]$ , then the faces of the open cylinder are shown with *lastFace* = True  
*cutPlain*: only used for *angleRange*  $\neq [0, 2 * \pi]$ ; if True, a plane is cut through the part of the cylinder; if False, the cylinder becomes a cake shape ...

- **output**: graphicsData dictionary, to be used in visualization of EXUDYN objects

---

```
def GraphicsDataRigidLink(p0, p1, axis0=[0,0,0], axis1=[0,0,0], radius=[0.1,0.1], thickness=0.05, width=[0.05,0.05],
color=[0.,0.,0.,1.], nTiles=16)
```

- **function description**: generate graphics data for a planar Link between the two joint positions, having two axes
- **input**:
  - p0*: joint0 center position
  - p1*: joint1 center position
  - axis0*: direction of rotation axis at *p0*, if drawn as a cylinder; [0,0,0] otherwise
  - axis1*: direction of rotation axis of *p1*, if drawn as a cylinder; [0,0,0] otherwise
  - radius*: list of two radii [*radius0*, *radius1*], being the two radii of the joints drawn by a cylinder or sphere
  - width*: list of two widths [*width0*, *width1*], being the two widths of the joints drawn by a cylinder; ignored for sphere
  - thickness*: the thickness of the link (shaft) between the two joint positions; thickness in z-direction or diameter (cylinder)
  - color*: provided as list of 4 RGBA values
  - nTiles*: used to determine resolution of cylinder  $\geq 3$ ; use larger values for finer resolution
- **output**: graphicsData dictionary, to be used in visualization of EXUDYN objects

---

```
def GraphicsDataFromSTLfileTxt(fileName, color=[0.,0.,0.,1.], verbose=False)
```

- **function description**: generate graphics data from STL file (text format!) and use color for visualization
- **input**:
  - fileName*: string containing directory and filename of STL-file (in text / SCII format) to load
  - color*: provided as list of 4 RGBA values
  - verbose*: if True, useful information is provided during reading
- **output**: interchanged 2nd and 3rd component of list

## 5.4 Module: rigidBodyUtilities

def **ComputeOrthonormalBasis**(*vector0*)

- **function description:** compute orthogonal basis vectors (normal1, normal2) for given vector0 (non-unique solution!); if vector0 == [0,0,0], then any normal basis is returned

---

def **GramSchmidt**(*vector0*, *vector1*)

- **function description:** compute Gram-Schmidt projection of given 3D vector 1 on vector 0 and return normalized triad (vector0, vector1, vector0 x vector1)

---

def **Skew**(*vector*)

- **function description:** compute skew symmetric 3x3-matrix from 3x1- or 1x3-vector

---

def **Skew2Vec**(*skew*)

- **function description:**  
convert skew symmetric matrix m to vector  
*def Skew2Vec(m):*

---

def **ComputeSkewMatrix**(*v*)

- **function description:** compute (3 x 3\*n) skew matrix from (3\*n) vector

---

def **EulerParameters2G**(*eulerParameters*)

- **function description:** convert Euler parameters (ep) to G-matrix ( $=\partial\omega/\partial\mathbf{p}_i$ )
- **input:** vector of 4 eulerParameters as list or np.array
- **output:** 3x4 matrix G as np.array

---

def **EulerParameters2GLocal**(*eulerParameters*)

- **function description:** convert Euler parameters (ep) to local G-matrix ( $=\partial^b\omega/\partial\mathbf{p}_i$ )
- **input:** vector of 4 eulerParameters as list or np.array
- **output:** 3x4 matrix G as np.array

---

def **EulerParameters2RotationMatrix**(*eulerParameters*)

- **function description:** compute rotation matrix from eulerParameters

- **input:** vector of 4 eulerParameters as list or np.array
- **output:** 3x3 rotation matrix as np.array

---

def **RotationMatrix2EulerParameters**(*rotationMatrix*)

- **function description:** compute Euler parameters from given rotation matrix
- **input:** 3x3 rotation matrix as list of lists or as np.array
- **output:** vector of 4 eulerParameters as np.array

---

def **AngularVelocity2EulerParameters\_t**(*angularVelocity*, *eulerParameters*)

- **function description:**  
compute time derivative of Euler parameters from (global) angular velocity vector  
note that for Euler parameters  $\mathbf{p}$ , we have  $\boldsymbol{\omega} = \mathbf{G}\mathbf{p}_t \Rightarrow \mathbf{G}^T \boldsymbol{\omega} = \mathbf{G}^T \cdot \mathbf{G} \cdot \mathbf{p}_t \Rightarrow \mathbf{G}^T \mathbf{G} = 4(\mathbf{I}_{4 \times 4} - \mathbf{p} \cdot \mathbf{p}^T) \mathbf{p}_t = 4(\mathbf{I}_{4 \times 4}) \mathbf{p}_t$
- **input:**  
*angularVelocity*: 3D vector of angular velocity in global frame, as lists or as np.array  
*eulerParameters*: vector of 4 eulerParameters as np.array or list
- **output:** vector of time derivatives of 4 eulerParameters as np.array

---

def **RotationVector2RotationMatrix**(*rotationVector*)

- **function description:** rotation matrix from rotation vector, see appendix B in [?]
- **input:** 3D rotation vector as list or np.array
- **output:** 3x3 rotation matrix as np.array

---

def **RotationMatrix2RotationVector**(*rotationMatrix*)

- **function description:** compute rotation vector from rotation matrix
- **input:** 3x3 rotation matrix as list of lists or as np.array
- **output:** vector of 3 components of rotation vector as np.array

---

def **ComputeRotationAxisFromRotationVector**(*rotationVector*)

- **function description:** compute rotation axis from given rotation vector
- **input:** 3D rotation vector as np.array
- **output:** 3D vector as np.array representing the rotation axis

---

def **RotXYZ2RotationMatrix**(*rot*)

- **function description:** compute rotation matrix from consecutive xyz rotations (Tait-Bryan angles);  $A=A_x*A_y*A_z$ ;  $rot=[rotX, rotY, rotZ]$
- **input:** 3D vector of Tait-Bryan rotation parameters  $[X,Y,Z]$  in radiant
- **output:** 3x3 rotation matrix as np.array

---

def **RotationMatrix2RotXYZ**(*rotationMatrix*)

- **function description:** convert rotation matrix to xyz Euler angles (Tait-Bryan angles);  $A=A_x*A_y*A_z$ ;
- **input:** 3x3 rotation matrix as list of lists or np.array
- **output:** vector of Tait-Bryan rotation parameters  $[X,Y,Z]$  (in radiant) as np.array

---

def **AngularVelocity2RotXYZ\_t**(*angularVelocity, rotation*)

- **function description:** compute time derivatives of angles RotXYZ from (global) angular velocity vector and given rotation
- **input:**
  - angularVelocity*: global angular velocity vector as list or np.array
  - rotation*: 3D vector of Tait-Bryan rotation parameters  $[X,Y,Z]$  in radiant
- **output:** time derivative of vector of Tait-Bryan rotation parameters  $[X,Y,Z]$  (in radiant) as np.array

---

def **RotXYZ2EulerParameters**(*alpha*)

- **function description:** compute four Euler parameters from given RotXYZ angles, see [?]
- **input:** alpha: 3D vector as np.array containing RotXYZ angles
- **output:**
  - 4D vector as np.array containing four Euler parameters
  - entry zero of output represent the scalar part of Euler parameters

---

def **RotationMatrixX**(*angleRad*)

- **function description:** compute rotation matrix w.r.t. X-axis (first axis)
- **input:** angle around X-axis in radiant
- **output:** 3x3 rotation matrix as np.array

---

def **RotationMatrixY**(*angleRad*)

- **function description:** compute rotation matrix w.r.t. Y-axis (second axis)
- **input:** angle around Y-axis in radiant
- **output:** 3x3 rotation matrix as np.array

---

def **RotationMatrixZ**(*angleRad*)

- **function description:** compute rotation matrix w.r.t. Z-axis (third axis)
- **input:** angle around Z-axis in radiant
- **output:** 3x3 rotation matrix as np.array

---

def **HomogeneousTransformation**(*A, r*)

- **function description:** compute homogeneous transformation matrix from rotation matrix A and translation vector r

---

def **HTtranslate**(*r*)

- **function description:** homogeneous transformation for translation with vector r

---

def **HT0**()

- **function description:** identity homogeneous transformation:

---

def **HTrotateX**(*angle*)

- **function description:** homogeneous transformation for rotation around axis X (first axis)

---

def **HTrotateY**(*angle*)

- **function description:** homogeneous transformation for rotation around axis X (first axis)

---

def **HTrotateZ**(*angle*)

- **function description:** homogeneous transformation for rotation around axis X (first axis)

---

def **HT2translation**(*T*)

- **function description:** return translation part of homogeneous transformation

---

def **HT2rotationMatrix**(T)

- **function description:** return rotation matrix of homogeneous transformation

---

def **InverseHT**(T)

- **function description:** return inverse homogeneous transformation such that  $\text{inv}(T)^*T = \text{np.eye}(4)$

---

def **AddRigidBody**(mainSys, inertia, nodeType, position=[0,0,0], velocity=[0,0,0], rotationMatrix=[], rotationParameters=[], angularVelocity=[0,0,0], gravity=[0,0,0], graphicsDataList=[])

- **function description:**
  - adds a node (with str(exu.NodeType. ...)) and body for a given rigid body
  - either the initial rotation is given by the rotationMatrix (while rotationParameters=[]) or by rotationParameters (while rotationMatrix=[]) (non empty)
  - position ... initial position, etc.
  - all quantities (esp. velocity and angular velocity) are given in global coordinates!
  - returns node number and body number
  - adds gravity force, i.e.,  $m \cdot \text{gravity}$

## 5.5 Module: lieGroupBasics

def **Sinc**(x)

- **function description:** compute the cardinal sine function in radians
- **input:** scalar float or int value
- **output:** float value in radians

---

def **Cot**(x)

- **function description:** compute the cotangent function  $\cot(x)=1/\tan(x)$  in radians
- **input:** scalar float or int value
- **output:** float value in radians

---

def **R3xSO3Matrix2RotationMatrix**(G)

- **function description:** computes 3x3 rotation matrix from 7x7 R3xSO(3) matrix, see [?]
- **input:** G: 7x7 matrix as np.array

- **output:** 3x3 rotation matrix as np.array

---

def **R3xSO3Matrix2Translation**(G)

- **function description:** computes translation part of R3xSO(3) matrix, see [?]
- **input:** G: 7x7 matrix as np.array
- **output:** 3D vector as np.array containg translational part of R3xSO(3)

---

def **R3xSO3Matrix**( $x$ ,  $R$ )

- **function description:** builds 7x7 matrix as element of the Lie group R3xSO(3), see [?]
- **input:**
  - $x$ : 3D vector as np.array representing the translation part corresponding to R3
  - $R$ : 3x3 rotation matrix as np.array
- **output:** 7x7 matrix as np.array

---

def **ExpSO3**( $\Omega$ )

- **function description:** compute the matrix exponential map on the Lie group SO(3), see [?]
- **input:** 3D rotation vector as np.array
- **output:** 3x3 matrix as np.array

---

def **ExpS3**( $\Omega$ )

- **function description:** compute the quaternion exponential map on the Lie group S(3), see [?, ?]
- **input:** 3D rotation vector as np.array
- **output:**
  - 4D vector as np.array containing four Euler parameters
  - entry zero of output represent the scalar part of Euler parameters

---

def **LogSO3**( $R$ )

- **function description:** compute the matrix logarithmic map on the Lie group SO(3), see [?, ?]
- **input:** 3x3 rotation matrix as np.array
- **output:** 3x3 skew symmetric matrix as np.array

---

def **TEpSO3**( $\Omega$ )



- **function description:** compute the tangent operator corresponding to  $\text{ExpSO3}$ , see [?]
- **input:** 3D rotation vector as `np.array`
- **output:** 3x3 matrix as `np.array`

---

def **TExpSO3Inv**(*Omega*)

- **function description:**
  - compute the inverse of the tangent operator  $\text{TExpSO3}$ , see [?]
  - this function was improved, see `coordinateMaps.pdf` by Stefan Holzinger
- **input:** 3D rotation vector as `np.array`
- **output:** 3x3 matrix as `np.array`

---

def **ExpSE3**(*x*)

- **function description:** compute the matrix exponential map on the Lie group  $\text{SE}(3)$ , see [?]
- **input:** 6D incremental motion vector as `np.array`
- **output:** 4x4 homogeneous transformation matrix as `np.array`

---

def **LogSE3**(*H*)

- **function description:** compute the matrix logarithm on the Lie group  $\text{SE}(3)$ , see [?]
- **input:** 4x4 homogeneous transformation matrix as `np.array`
- **output:** 4x4 skew symmetric matrix as `np.array`

---

def **TExpSE3**(*x*)

- **function description:** compute the tangent operator corresponding to  $\text{ExpSE3}$ , see [?]
- **input:** 6D incremental motion vector as `np.array`
- **output:** 6x6 matrix as `np.array`

---

def **TExpSE3Inv**(*x*)

- **function description:** compute the inverse of tangent operator  $\text{TExpSE3}$ , see [?]
- **input:** 6D incremental motion vector as `np.array`
- **output:** 6x6 matrix as `np.array`

---

def **ExpR3xSO3**(*x*)

- **function description:** compute the matrix exponential map on the Lie group  $R3 \times SO(3)$ , see [?]
- **input:** 6D incremental motion vector as `np.array`
- **output:** 7x7 matrix as `np.array`

---

def **TExpR3xSO3**(*x*)

- **function description:** compute the tangent operator corresponding to `ExpR3xSO3`, see [?]
- **input:** 6D incremental motion vector as `np.array`
- **output:** 6x6 matrix as `np.array`

---

def **TExpR3xSO3Inv**(*x*)

- **function description:** compute the inverse of tangent operator `TExpR3xSO3`
- **input:** 6D incremental motion vector as `np.array`
- **output:** 6x6 matrix as `np.array`

---

def **CompositionRuleDirectProductR3AndS3**(*q0, incrementalMotionVector*)

- **function description:** compute composition operation for pairs in the Lie group  $R3 \times S3$
- **input:**
  - q0*: 7D vector as `np.array` containing position coordinates and Euler parameters
  - incrementalMotionVector*: 6D incremental motion vector as `np.array`
- **output:** 7D vector as `np.array` containing composed position coordinates and composed Euler parameters

---

def **CompositionRuleSemiDirectProductR3AndS3**(*q0, incrementalMotionVector*)

- **function description:** compute composition operation for pairs in the Lie group  $R3$  semiTimes  $S3$  (corresponds to  $SE(3)$ )
- **input:**
  - q0*: 7D vector as `np.array` containing position coordinates and Euler parameters
  - incrementalMotionVector*: 6D incremental motion vector as `np.array`
- **output:** 7D vector as `np.array` containing composed position coordinates and composed Euler parameters

---

def **CompositionRuleDirectProductR3AndR3RotVec**(*q0, incrementalMotionVector*)

- **function description:**
  - compute composition operation for pairs in the group obtained from the direct product of  $R3$  and  $R3$ , see [?]

the rotation vector is used as rotation parametrizations

this composition operation can be used in formulations which represent the translational velocities in the global (inertial) frame

– **input:**

$q0$ : 6D vector as np.array containing position coordinates and rotation vector

*incrementalMotionVector*: 6D incremental motion vector as np.array

– **output:** 7D vector as np.array containing composed position coordinates and composed rotation vector

---

def **CompositionRuleSemiDirectProductR3AndR3RotVec**( $q0$ , *incrementalMotionVector*)

– **function description:**

compute composition operation for pairs in the group obtained from the direct product of R3 and R3.

the rotation vector is used as rotation parametrizations

this composition operation can be used in formulations which represent the translational velocities in the local (body-attached) frame

– **input:**

$q0$ : 6D vector as np.array containing position coordinates and rotation vector

*incrementalMotionVector*: 6D incremental motion vector as np.array

– **output:** 6D vector as np.array containing composed position coordinates and composed rotation vector

---

def **CompositionRuleDirectProductR3AndR3RotXYZAngles**( $q0$ , *incrementalMotionVector*)

– **function description:**

compute composition operation for pairs in the group obtained from the direct product of R3 and R3.

Cardan-Tait/Bryan (CTB) angles are used as rotation parametrizations

this composition operation can be used in formulations which represent the translational velocities in the global (inertial) frame

– **input:**

$q0$ : 6D vector as np.array containing position coordinates and Cardan-Tait/Bryan angles

*incrementalMotionVector*: 6D incremental motion vector as np.array

– **output:** 6D vector as np.array containing composed position coordinates and composed Cardan-Tait/Bryan angles

---

def **CompositionRuleSemiDirectProductR3AndR3RotXYZAngles**( $q0$ , *incrementalMotionVector*)

– **function description:**

compute composition operation for pairs in the group obtained from the direct product of R3 and R3.

Cardan-Tait/Bryan (CTB) angles are used as rotation parametrizations

this composition operation can be used in formulations which represent the translational velocities in the local (body-attached) frame

– **input:**

$q0$ : 6D vector as np.array containing position coordinates and Cardan-Tait/Bryan angles

*incrementalMotionVector*: 6D incremental motion vector as np.array

– **output:** 6D vector as np.array containing composed position coordinates and composed Cardan-Tait/Bryan angles

---

def **CompositionRuleForEulerParameters**( $q, p$ )

– **function description:**

compute composition operation for Euler parameters (unit quaternions)

this composition operation is quaternion multiplication, see [?]

– **input:**

$q$ : 4D vector as np.array containing Euler parameters

$p$ : 4D vector as np.array containing Euler parameters

– **output:** 4D vector as np.array containing composed (multiplied) Euler parameters

---

def **CompositionRuleForRotationVectors**( $v0, \Omega$ )

– **function description:** compute composition operation for rotation vectors  $v0$  and  $\Omega$ , see [?]

– **input:**

$v0$ : 3D rotation vector as np.array

$\Omega$ : 3D (incremental) rotation vector as np.array

– **output:** 3D vector as np.array containing composed rotation vector  $v$

---

def **CompositionRuleRotXYZAnglesRotationVector**( $\alpha0, \Omega$ )

– **function description:** compute composition operation for RotXYZ angles, see [?]

– **input:**

$\alpha0$ : 3D vector as np.array containing RotXYZ angles

$\Omega$ : 3D vector as np.array containing the (incremental) rotation vector

– **output:** 3D vector as np.array containing composed RotXYZ angles

## 5.6 Module: FEM

def **CompressedRowSparseToDenseMatrix**(*sparseData*)

- **function description:** convert zero-based sparse matrix data to dense numpy matrix
- **input:** *sparseData*: format (per row): [row, column, value] ==> converted into dense format
- **output:** a dense matrix as np.array

---

def **MapSparseMatrixIndices**(*matrix*, *sorting*)

- **function description:** resort a sparse matrix (internal CSR format) with given sorting for rows and columns; changes matrix directly! used for ANSYS matrix import

---

def **VectorDiadicUnitMatrix3D**(*v*)

- **function description:** compute diadic product of vector *v* and a 3D unit matrix =  $\text{diadic}(v, I_{3 \times 3})$ ; used for ObjectFFRF and CMS implementation

---

def **CyclicCompareReversed**(*list1*, *list2*)

- **function description:** compare cyclic two lists, reverse second list; return True, if any cyclic shifted lists are same, False otherwise

---

def **AddEntryToCompressedRowSparseArray**(*sparseData*, *row*, *column*, *value*)

- **function description:**  
add entry to compressedRowSparse matrix, avoiding duplicates  
value is either added to existing entry (avoid duplicates) or a new entry is appended

---

def **CSRtoRowsAndColumns**(*sparseMatrixCSR*)

- **function description:** compute rows and columns of a compressed sparse matrix and return as tuple: (rows, columns)

---

def **CSRtoScipySparseCSR**(*sparseMatrixCSR*)

- **function description:** convert internal compressed CSR to scipy.sparse csr matrix

---

def **ScipySparseCSRtoCSR**(*scipyCSR*)

- **function description:** convert scipy.sparse csr matrix to internal compressed CSR

---

def **ResortIndicesOfCSRmatrix**(*mXXXXYYZZ*, *numberOfRows*)

– **function description:**

resort indices of given CSR matrix in XXXYYYZZZ format to XYZXYZXYZ format; *numberOfRows* must be equal to columns  
needed for import from NGSolve

---

def **ConvertHexToTrigs**(*nodeNumbers*)

– **function description:**

convert list of Hex8/C3D8 element with 8 nodes in *nodeNumbers* into triangle-List  
also works for Hex20 elements, but does only take the corner nodes!

---

def **ConvertDenseToCompressedRowMatrix**(*denseMatrix*)

– **function description:** convert numpy.array dense matrix to (internal) compressed row format

---

def **ReadMatrixFromAnsysMMF**(*fileName*, *verbose*=False)

– **function description:**

This function reads either the mass or stiffness matrix from an Ansys Matrix Market Format (MMF). The corresponding matrix can either be exported as dense matrix or sparse matrix.

– **input:** *fileName* of MMF file

– **output:** internal compressed row sparse matrix (as (nrows x 3) numpy array)

– **author:**

Stefan Holzinger

*Note:*

A MMF file can be created in Ansys by placing the following APDL code inside the solution tree in Ansys Workbench:

!!

! APDL code that exports sparse stiffness and mass matrix in MMF format. If

! the dense matrix is needed, replace \*SMAT with \*DMAT in the following

! APDL code.

! Export the stiffness matrix in MMF format

\*SMAT,MatKD,D,IMPORT,FULL,file.full,STIFF

\*EXPORT,MatKD,MMF,fileNameStiffnessMatrix,,,

! Export the mass matrix in MMF format

+++++

read sorting vector for ANSYS mass and stiffness matrices and return sorting vector as np.array  
the file contains sorting for nodes and applies this sorting to the DOF (assuming 3 DOF per node!)  
the resulting sorted vector is already converted to 0-based indices

+++++

---

```
def ReadElementsFromAnsysTxt(fileName, verbose=False)
```

- **function description:** This function reads the nodal coordinates exported from Ansys.
- **input:** fileName (file name ending must be .txt!)
- **output:** element connectivity as numpy array
- **author:** Stefan Holzinger
- **notes:**

The elements can be exported from Ansys by creating a named selection of the body whos mesh should to exported by choosing its geometry. Next, create a second named selction by using a worksheet. Add the named selection that was created first into the worksheet of the second named selection. Inside the worksheet, choose 'convert' and convert the first created named selection to 'mesh element' (Netzelement in german) and click on generate to create the second named selection. Next, right click on the second named selection tha was created and choose 'export' and save the elements as .txt file.

+++++

### 5.6.1 CLASS ObjectFFRFInterface (in module FEM)

**class description:** compute terms necessary for ObjectFFRF class used internally in FEMInterface to compute ObjectFFRF object this class holds all data for ObjectFFRF user functions

```
def __init__(self, femInterface)
```

- **classFunction:**
  - initialize ObjectFFRFInterface with FEMInterface class
  - initializes the ObjectFFRFInterface with nodes, modes, surface description and systemmatrices from FEMInterface
  - data is then transfered to mbs object with classFunction AddObjectFFRF(...)

---

```
def AddObjectFFRF(self, exu, mbs, positionRef=[0,0,0], eulerParametersRef=[1,0,0,0], initialVelocity=[0,0,0], initialAngularVelocity=[0,0,0], gravity=[0,0,0], constrainRigidBodyMotion=True, massProportionalDamping=0, stiffnessProportionalDamping=0, color=[0.1,0.9,0.1,1.])
```

- **classFunction:** add according nodes, objects and constraints for FFRF object to MainSystem mbs
- **input:**
  - exu: the exudyn module
  - mbs: a MainSystem object
  - positionRef: reference position of created ObjectFFRF (set in rigid body node underlying to ObjectFFRF)



*eulerParametersRef*: reference euler parameters of created ObjectFFRF (set in rigid body node underlying to ObjectFFRF)

*initialVelocity*: initial velocity of created ObjectFFRF (set in rigid body node underlying to ObjectFFRF)

*initialAngularVelocity*: initial angular velocity of created ObjectFFRF (set in rigid body node underlying to ObjectFFRF)

*gravity*: set [0,0,0] if no gravity shall be applied, or to the gravity vector otherwise

*constrainRigidBodyMotion*: set True in order to add constraint (Tisserand frame) in order to suppress rigid motion of mesh nodes

*color*: provided as list of 4 RGBA values

add object to mbs as well as according nodes

---

**def Ufforce**(*self, exu, mbs, t, q, q\_t*)

- **classFunction**: optional forceUserFunction for ObjectFFRF (per default, this user function is ignored)

---

**def UfmassGenericODE2**(*self, exu, mbs, t, q, q\_t*)

- **classFunction**: optional massMatrixUserFunction for ObjectFFRF (per default, this user function is ignored)

## 5.6.2 CLASS ObjectFFRFreducedOrderInterface (in module FEM)

**class description**: compute terms necessary for ObjectFFRFreducedOrder class used internally in FEMinterface to compute ObjectFFRFreducedOrder dictionary this class holds all data for ObjectFFRFreducedOrder user functions

**def \_\_init\_\_**(*self, femInterface, roundMassMatrix=1e-13, roundStiffNessMatrix=1e-13*)

- **classFunction**:
  - initialize ObjectFFRFreducedOrderInterface with FEMinterface class
  - initializes the ObjectFFRFreducedOrderInterface with nodes, modes, surface description and reduced system matrices from FEMinterface
  - data is then transfered to mbs object with classFunction AddObjectFFRFreducedOrderWithUserFunctions(...)
- **input**:
  - femInterface*: must provide nodes, surfaceTriangles, modeBasis, massMatrix, stiffness
  - roundMassMatrix*: use this value to set entries of reduced mass matrix to zero which are below the treshold
  - roundStiffNessMatrix*: use this value to set entries of reduced stiffness matrix to zero which are below the treshold

---

```
def AddObjectFFRFReducedOrderWithUserFunctions(self, exu, mbs, positionRef=[0,0,0], eulerParametersRef=[1,0,0,0],
initialVelocity=[0,0,0], initialAngularVelocity=[0,0,0], gravity=[0,0,0], UForce=0, UMassMatrix=0, massProportionalDamping=0, stiffnessProportionalDamping=0, color=[0.1,0.9,0.1,1.])
```

- **classFunction**: add according nodes, objects and constraints for ObjectFFRFReducedOrder object to Main-System mbs

- **input**:

*exu*: the exudyn module

*mbs*: a MainSystem object

*positionRef*: reference position of created ObjectFFRFReducedOrder (set in rigid body node underlying to ObjectFFRFReducedOrder)

*eulerParametersRef*: reference euler parameters of created ObjectFFRFReducedOrder (set in rigid body node underlying to ObjectFFRFReducedOrder)

*initialVelocity*: initial velocity of created ObjectFFRFReducedOrder (set in rigid body node underlying to ObjectFFRFReducedOrder)

*initialAngularVelocity*: initial angular velocity of created ObjectFFRFReducedOrder (set in rigid body node underlying to ObjectFFRFReducedOrder)

*gravity*: set [0,0,0] if no gravity shall be applied, or to the gravity vector otherwise

*UForce*: provide a user function, which computes the quadratic velocity vector and applied forces; usually this function reads like:

```
def UForceFFRFReducedOrder(t, qReduced, qReduced_t):
    return cms.UForceFFRFReducedOrder(exu, mbs, t, qReduced, qReduced_t)
```

*UMassMatrix*: provide a user function, which computes the quadratic velocity vector and applied forces; usually this function reads like:

```
def UMassFFRFReducedOrder(t, qReduced, qReduced_t):
    return cms.UMassFFRFReducedOrder(exu, mbs, t, qReduced, qReduced_t)
```

*massProportionalDamping*: Rayleigh damping factor for mass proportional damping, added to floating frame/modal coordinates only

*stiffnessProportionalDamping*: Rayleigh damping factor for stiffness proportional damping, added to floating frame/modal coordinates only

*color*: provided as list of 4 RGBA values

---

```
def UMassFFRFReducedOrder(self, exu, mbs, t, qReduced, qReduced_t)
```

- **classFunction**: CMS mass matrix user function; qReduced and qReduced\_t contain the coordinates of the rigid body node and the modal coordinates in one vector!

---

```
def UForceFFRFReducedOrder(self, exu, mbs, t, qReduced, qReduced_t)
```

- **classFunction**: CMS force matrix user function; qReduced and qReduced\_t contain the coordinates of the rigid body node and the modal coordinates in one vector!

### 5.6.3 CLASS FEMinterface (in module FEM)

**class description:** general interface to different FEM / mesh imports and export to EXUDYN functions use this class to import meshes from different meshing or FEM programs (NETGEN/NGsolve, ABAQUS, ANSYS, ..) and store it in a unique format do mesh operations, compute eigenmodes and reduced basis, etc. load/store the data efficiently with LoadFromFile(...), SaveToFile(...) if import functions are slow export to EXUDYN objects

**def \_\_init\_\_(self)**

- **classFunction:** initialize all data of the FEMinterface by, e.g., `fem = FEMinterface()`

---

**def SaveToFile(self, fileName)**

- **classFunction:** save all data (nodes, elements, ...) to a data filename; this function is much faster than the text-based import functions
- **input:** use filename without ending ==> ".npy" will be added

---

**def LoadFromFile(self, fileName)**

- **classFunction:**  
load all data (nodes, elements, ...) from a data filename previously stored with SaveToFile(...).  
this function is much faster than the text-based import functions
- **input:** use filename without ending ==> ".npy" will be added

---

**def ImportFromAbaqusInputFile(self, fileName, typeName='Part', name='Part-1', verbose=False)**

- **classFunction:**  
import nodes and elements from Abaqus input file and create surface elements  
node numbers in elements are converted from 1-based indices to python's 0-based indices  
only works for Hex8, Hex20, Tet4 and Tet10 (C3D4, C3D8, C3D10, C3D20) elements  
return node numbers as numpy array

---

**def ReadMassMatrixFromAbaqus(self, fileName, type='SparseRowColumnValue')**

- **classFunction:**  
*read mass matrix from compressed row text format (exported from Abaqus); in order to export system matrices, write the following lines in your Abaqus input file:*  
  
\*STEP  
\*MATRIX GENERATE, STIFFNESS, MASS  
\*MATRIX OUTPUT, STIFFNESS, MASS, FORMAT=COORDINATE  
\*End Step

---

```
def ReadStiffnessMatrixFromAbaqus(self, fileName, type='SparseRowColumnValue')
```

- **classFunction**: read stiffness matrix from compressed row text format (exported from Abaqus)

---

```
def ImportMeshFromNGsolve(self, mesh, density, youngsModulus, poissosRatio, verbose=False)
```

- **classFunction**: import mesh from NETGEN/NGsolve and setup mechanical problem

- **input**:

*mesh*: a previously created ngs.mesh (NGsolve mesh, see examples)

*youngsModulus*: Young's modulus used for mechanical model

*poissosRatio*: Poisson's ratio used for mechanical model

*density*: density used for mechanical model

*verbose*: set True to print out some status information

- **notes**:

The interface to NETGEN/NGsolve has been created together with Joachim SchÄ¶berl, main developer of NETGEN/NGsolve; Thank's a lot!

download NGsolve at: <https://ngsolve.org/>

NGsolve needs Python 3.7 (64bit) ==> use according EXUDYN version!

note that node/element indices in the NGsolve mesh are 1-based and need to be converted to 0-base!

---

```
def GetMassMatrix(self, sparse=True)
```

- **classFunction**: get sparse mass matrix in according format

---

```
def GetStiffnessMatrix(self, sparse=True)
```

- **classFunction**: get sparse stiffness matrix in according format

---

```
def NumberOfNodes(self)
```

- **classFunction**: get total number of nodes

---

```
def GetNodePositionsAsArray(self)
```

- **classFunction**: get node points as array; only possible, if there exists only one type of Position nodes

---

```
def NumberOfCoordinates(self)
```

- **classFunction**: get number of total nodal coordinates

---

**def GetNodeAtPoint**(*self, point, tolerance=1e-5, raiseException=True*)

- **classFunction**:
  - get node number for node at given point, e.g.  $p=[0.1,0.5,-0.2]$ , using a tolerance (+/-) if coordinates are available only with reduced accuracy
  - if not found, it returns an invalid index

---

**def GetNodesInPlane**(*self, point, normal, tolerance=1e-5*)

- **classFunction**:
  - get node numbers in plane defined by point  $p$  and (normalized) normal vector  $n$  using a tolerance for the distance to the plane
  - if not found, it returns an empty list

---

**def GetNodesOnCircle**(*self, point, normal, r, tolerance=1e-5*)

- **classFunction**:
  - get node numbers on a circle, by point  $p$ , (normalized) normal vector  $n$  (which is the axis of the circle) and radius  $r$
  - using a tolerance for the distance to the plane
  - if not found, it returns an empty list

---

**def GetSurfaceTriangles**(*self*)

- **classFunction**: return surface trigs as node number list (for drawing in EXUDYN)

---

**def VolumeToSurfaceElements**(*self, verbose=False*)

- **classFunction**:
  - generate surface elements from volume elements
  - stores the surface in `self.surface`
  - only works for one element list and one type ('Hex8') of elements

---

**def GetGyroscopicMatrix**(*self, rotationAxis=2, sparse=True*)

- **classFunction**: get gyroscopic matrix in according format;  $rotationAxis=[0,1,2] = [x,y,z]$

---

**def ScaleMassMatrix**(*self, factor*)

- **classFunction**: scale (=multiply) mass matrix with factor

---

**def ScaleStiffnessMatrix**(*self, factor*)

- **classFunction**: scale (=multiply) stiffness matrix with factor

---

**def AddElasticSupportAtNode**(*self, nodeNumber, springStiffness=[1e8,1e8,1e8]*)

- **classFunction**:  
modify stiffness matrix to add elastic support (joint, etc.) to a node; nodeNumber zero based (as everywhere in the code...)  
springStiffness must have length according to the node size

---

**def AddNodeMass**(*self, nodeNumber, addedMass*)

- **classFunction**: modify mass matrix by adding a mass to a certain node, modifying directly the mass matrix

---

**def ComputeEigenmodes**(*self, nModes, excludeRigidBodyModes=0, useSparseSolver=True*)

- **classFunction**:  
compute nModes smallest eigenvalues and eigenmodes from mass and stiffnessMatrix  
store mode vector in modeBasis, but exclude a number of 'excludeRigidBodyModes' rigid body modes from modeBasis  
if excludeRigidBodyModes > 0, then the computed modes is nModes + excludeRigidBodyModes, from which excludeRigidBodyModes smallest eigenvalues are excluded

---

**def GetEigenFrequenciesHz**(*self*)

- **classFunction**: return list of eigenvalues in Hz of previously computed eigenmodes

---

**def ComputeCampbellDiagram**(*self, terminalFrequency, nEigenfrequencies=10, frequencySteps=25, rotationAxis=2, plotDiagram=False, verbose=False*)

- **classFunction**:  
compute Campbell diagram for given mechanical system  
create a first order system  $Ax + Bx = 0$  with  $x = [q, \dot{q}]'$  and compute eigenvalues  
takes mass M, stiffness K and gyroscopic matrix G from FEMinterface  
currently only uses dense matrices, so it is limited to approx. 5000 unknowns!

– **input:**

*terminalFrequency*: frequency in Hz, up to which the campbell diagram is computed

*nEigenfrequencies*: gives the number of computed eigenfrequencies(modes), in addition to the rigid body mode 0

*frequencySteps*: gives the number of increments (gives frequencySteps+1 total points in campbell diagram)

*rotationAxis*: [0,1,2] = [x,y,z] provides rotation axis

*plotDiagram*: if True, plots diagram for nEigenfrequencies before terminating

*verbose*: if True, shows progress of computation

– **output:**

[listFrequencies, campbellFrequencies]

*listFrequencies*: list of computed frequencies

*campbellFrequencies*: array of campbell frequencies per eigenfrequency of system

---

def **CheckConsistency**(self)

- **classFunction**: perform some consistency checks

---

def **ReadMassMatrixFromAnsys**(self, fileName, dofMappingVectorFile, sparse=True, verbose=False)

- **classFunction**: read mass matrix from CSV format (exported from Ansys)

---

def **ReadStiffnessMatrixFromAnsys**(self, fileName, dofMappingVectorFile, sparse=True, verbose=False)

- **classFunction**: read stiffness matrix from CSV format (exported from Ansys)

---

def **ReadNodalCoordinatesFromAnsys**(self, fileName, verbose=False)

- **classFunction**: read nodal coordinates (exported from Ansys as .txt-File)

---

def **ReadElementsFromAnsys**(self, fileName, verbose=False)

- **classFunction**: read elements (exported from Ansys as .txt-File)

## 5.7 Module: plot

def **PlotSensor**(mbs, sensorNumbers, components=0, \*\*kwargs)

- **function description**: helper for matplotlib in order to easily visualize sensor output

- **input:**

*mbs*: must be a valid MainSystem (mbs)

*sensorNumbers*: consists of one or a list of sensor numbers (type SensorIndex) as returned by the mbs function AddSensor(...)

*components*: consists of one or a list of components according to the component of the sensor to be plotted;

*\*kwargs*: additional options, e.g.:

xLabel -> string for text at x-axis

yLabel -> string for text at y-axis

– **output**: plots the sensor data

– **example**:

```
s0=mbs.AddSensor(SensorNode(nodeNumber=0))
```

```
s1=mbs.AddSensor(SensorNode(nodeNumber=1))
```

```
Plot(mbs, s0, 0)
```

```
Plot(mbs, sensorNumbers=[s0,s1], components=[0,2], xlabel='time in seconds')
```

## 5.8 Module: processing

```
def ProcessParameterList(parameterFunction, parameterList, addComputationIndex, useMultiProcessing, **kwargs)
```

– **function description**: processes parameterFunction for given parameters in parameterList, see ParameterVariation

– **input**:

*parameterFunction*: function, which takes the form parameterFunction(parameterDict) and which returns any values that can be stored in a list (e.g., a floating point number)

*parameterList*: list of parameter sets (as dictionaries) which are fed into the parameter variation, e.g., ['mass': 10, 'mass':20, ...]

*addComputationIndex*: if True, key 'computationIndex' is added to every parameterDict in the call to parameterFunction(), which allows to generate independent output files for every parameter, etc.

*useMultiProcessing*: if True, the multiprocessing lib is used for parallelized computation; WARNING: be aware that the function does not check if your function runs independently; DO NOT use GRAPHICS and DO NOT write to same output files, etc.!

*numberOfThreads*: default: same as number of cpus (threads); used for multiprocessing lib;

– **output**: returns values containing the results according to parameterList

– **notes**: options are passed from ParameterVariation

---

```
def ParameterVariation(parameterFunction, parameters, **kwargs)
```

– **function description**:

calls successively the function parameterFunction(parameterDict) with variation of parameters in given range; parameterDict is a dictionary, containing the current values of parameters,



e.g., `parameterDict={'mass':13, 'stiffness':12000}` to be computed and returns a value or a list of values which is then stored for each parameter

– **input:**

*parameterFunction*: function, which takes the form `parameterFunction(parameterDict)` and which returns any values that can be stored in a list (e.g., a floating point number)

*parameters*: given as a dictionary, consist of name and triple of (begin, end and steps) same as in `np.linspace(...)`, e.g. `'mass':(10,50,10)` for a mass varied from 10 to 50, using 10 steps

*useLogSpace*: (optional) if True, the parameters are varied at a logarithmic scale, e.g., `[1, 10, 100]` instead linear `[1, 50.5, 100]`

*debugMode*: if True, additional print out is done

*addComputationIndex*: if True, key `'computationIndex'` is added to every `parameterDict` in the call to `parameterFunction()`, which allows to generate independent output files for every parameter, etc.

*useMultiProcessing*: if True, the multiprocessing lib is used for parallelized computation; WARNING: be aware that the function does not check if your function runs independently; DO NOT use GRAPHICS and DO NOT write to same output files, etc.!

*numberOfThreads*: default: same as number of cpus (threads); used for multiprocessing lib;

– **output:**

returns `[parameterList, values]`, containing, e.g., `parameterList={'mass':[1,1,1,2,2,2,3,3,3], 'stiffness':[4,5,6,4,5,6,4,5,6]}` and the result values of the parameter variation according to the `parameterList`, `values=[7,8,9,3,4,5,6,7,8]` (depends on solution of problem ..., can also contain tuples, etc.)

– **example:** `ParameterVariation(parameters={'mass':(1,10,10), 'stiffness':(1000,10000,10)}, parameterFunction=Test, useMultiProcessing=True)`

---

def **GeneticOptimization**(*objectiveFunction*, *parameters*, *\*\*kwargs*)

– **function description:** compute minimum of given `objectiveFunction`

– **input:**

*objectiveFunction*: function, which takes the form `parameterFunction(parameterDict)` and which returns a value or list (or numpy array) which reflects the size of the objective to be minimized

*parameters*: given as a dictionary, consist of name and tuple of (begin, end), e.g. `'mass':(10,50)`

*initialPopulationSize*: number of random initial individuals (default: 100)

*numberOfGenerations*: number of generations (default: 10)

*numberOfChildren*: number childrens of surviving population (default: 8)

*useGenCrossing*: if True, the children are generated from parents by gen-crossover (default: True)

*survivingIndividuals*: number of surviving individuals after children are born (default: 8)

*rangeReductionFactor*: (not implemented yet!) reduction of mutation range relative to ranges of last step (default: 0.7)

*randomizerInitialization*: (not implemented yet!) initialize randomizer at beginning in order to get reproducible results (Default: True)

*distanceFactor*: (not implemented yet!) children only survive at a certain distance of the current range (Default: 0.1)

*debugMode*: if True, additional print out is done

*addComputationIndex*: if True, key 'computationIndex' is added to every parameterDict in the call to parameterFunction(), which allows to generate independent output files for every parameter, etc.

*useMultiProcessing*: if True, the multiprocessing lib is used for parallelized computation; WARNING: be aware that the function does not check if your function runs independently; DO NOT use GRAPHICS and DO NOT write to same output files, etc.!

*numberOfThreads*: default: same as number of cpus (threads); used for multiprocessing lib;

– **output:**

returns [optimumParameter, optimumValue, parametersAll, valuesAll], containing the optimum parameter set 'optimumParameter', optimum value 'optimumValue', the whole list of parameters parametersAll with according objective values 'valuesAll'

values=[7,8,9 ,3,4,5, 6,7,8] (depends on solution of problem ..., can also contain tuples, etc.)

– **notes:** This function is still under development and shows an experimental state!

– **example:** GeneticOptimization(objectiveFunction = fOpt, parameters={'mass':(1,10), 'stiffness':(1000,10000)})

---

def **PlotOptimizationResults**(parameterList, valueList)

– **function description:** visualize results of optimization

– **input:**

*parameterList*: taken from output parametersAll of GeneticOptimization, containing a list of parameter set dictionaries

*valueList*: taken from output valuesAll of GeneticOptimization; containing a list of floats that result from the objective function

– **output:** creates a figure for every parameter in parameterList



## Chapter 6

# Objects, nodes, markers, loads and sensors reference manual

This chapter includes the reference manual for all objects (bodies/constraints), nodes, markers, loads and sensors.

### 6.1 Notation for item equations

The following subscripts are used to define configurations of a quantity, e.g.,  $\mathbf{x}$ :

- $\mathbf{x}_{\text{config}}$  ...  $\mathbf{x}$  in any configuration
- $\mathbf{x}_{\text{ref}}$  ...  $\mathbf{x}$  in reference configuration, e.g., reference coordinates:  $\mathbf{c}_{\text{ref}}$
- $\mathbf{x}_{\text{ini}}$  ...  $\mathbf{x}$  in initial configuration, e.g., initial displacements:  $\mathbf{u}_{\text{ini}}$
- $\mathbf{x}_{\text{cur}}$  ...  $\mathbf{x}$  in current configuration
- $\mathbf{x}_{\text{vis}}$  ...  $\mathbf{x}$  in visualization configuration
- $\mathbf{x}_{\text{start of step}}$  ...  $\mathbf{x}$  in start of step configuration

As written in the introduction, the coordinates can have the types:

- ODE2 ... second order differential equations coordinates
- ODE1 ... first order differential equations coordinates; CURRENTLY NOT AVAILABLE
- AE ... algebraic equations coordinates
- Data ... data coordinates (history variables)

Time is usually defined as 'time' or  $t$ . The cross product or vector product ' $\times$ ' is often replaced by the skew symmetric matrix using the tilde ' $\sim$ ' symbol,

$$\mathbf{a} \times \mathbf{b} = \tilde{\mathbf{a}} \mathbf{b} = -\tilde{\mathbf{b}} \mathbf{a} \quad (6.1)$$

The following table contains the common notation:

python name (or description)	symbol	description
displacement coordinates	$\mathbf{q} = [q_0, \dots, q_n]^T$	vector of $n$ displacement based coordinates in any configuration
rotation coordinates	$\boldsymbol{\psi} = [\psi_0, \dots, \psi_n]^T$	vector of $\eta$ <b>rotation based coordinates</b> in any configuration; this coordinates are added to reference rotation parameters to provide the current rotation parameters

algebraic coordinates	$\mathbf{z} = [z_0, \dots, z_m]^T$	vector of $m$ algebraic coordinates if not Lagrange multipliers in any configuration
Lagrange multipliers	$\boldsymbol{\lambda} = [\lambda_0, \dots, \lambda_m]^T$	vector of $m$ Lagrange multipliers (=algebraic coordinates) in any configuration
data coordinates	$\mathbf{x} = [x_0, \dots, x_l]^T$	vector of $l$ data coordinates in any configuration
<b>python name: OutputVariable</b>	<b>symbol</b>	<b>description</b>
Coordinate	$\mathbf{c} = [c_0, \dots, c_n]^T$	coordinate vector with $n$ generalized coordinates $c_i$ in any configuration
Coordinate_t	$\dot{\mathbf{c}} = [\dot{c}_0, \dots, \dot{c}_n]^T$	time derivative of coordinate vector
Displacement	${}^0\mathbf{u} = [u_0, u_1, u_2]^T$	global displacement vector with 3 displacement coordinates $u_i$ in any configuration; in 1D or 2D objects, some of these coordinates may be zero
Rotation	$\boldsymbol{\theta} = [\theta_0, \dots, \theta_n]^T$	vector of <b>rotation parameters</b> (e.g., Euler parameters, Tait Bryan angles, ...) with $n$ coordinates $\theta_i$ in any configuration
RotationMatrix	${}^{0b}\mathbf{A} = \begin{bmatrix} A_{00} & A_{01} & A_{02} \\ A_{10} & A_{11} & A_{12} \\ A_{20} & A_{21} & A_{22} \end{bmatrix}$	a 3D rotation matrix, which transforms local (e.g., body $b$ ) to global coordinates (0): ${}^0\mathbf{x} = {}^{0b}\mathbf{A} {}^b\mathbf{x}$
Position	${}^0\mathbf{p} = [p_0, p_1, p_2]^T$	global position vector with 3 position coordinates $p_i$ in any configuration
Velocity	${}^0\mathbf{v} = {}^0\dot{\mathbf{u}} = [v_0, v_1, v_2]^T$	global velocity vector with 3 displacement coordinates $v_i$ in any configuration
AngularVelocity	${}^0\boldsymbol{\omega} = [\omega_0, \dots, \omega_2]^T$	global angular velocity vector with 3 coordinates $\omega_i$ in any configuration
Acceleration	${}^0\mathbf{a} = {}^0\ddot{\mathbf{u}} = [a_0, a_1, a_2]^T$	global acceleration vector with 3 displacement coordinates $a_i$ in any configuration
AngularAcceleration	${}^0\boldsymbol{\alpha} = {}^0\dot{\boldsymbol{\omega}} = [\alpha_0, \dots, \alpha_2]^T$	global angular acceleration vector with 3 coordinates $\alpha_i$ in any configuration
VelocityLocal	${}^b\mathbf{v} = [v_0, v_1, v_2]^T$	local (body-fixed) velocity vector with 3 displacement coordinates $v_i$ in any configuration
AngularVelocityLocal	${}^b\boldsymbol{\omega} = [\omega_0, \dots, \omega_2]^T$	local (body-fixed) angular velocity vector with 3 coordinates $\omega_i$ in any configuration
Force	${}^0\mathbf{f} = [f_0, \dots, f_2]^T$	vector of 3 force components in global coordinates
Torque	${}^0\boldsymbol{\tau} = [\tau_0, \dots, \tau_2]^T$	vector of 3 torque components in global coordinates
<b>python name: input to nodes, markers, etc.</b>	<b>symbol</b>	<b>description</b>
referenceCoordinates	$\mathbf{c}_{\text{ref}} = [c_0, \dots, c_n]_{\text{ref}}^T = [c_{\text{Ref},0}, \dots, c_{\text{Ref},n}]_{\text{ref}}^T$	$n$ coordinates of reference configuration (can usually be set at initialization of nodes)
initialCoordinates	$\mathbf{c}_{\text{ini}}$	initial coordinates with generalized or mixed displacement/rotation quantities (can usually be set at initialization of nodes)

localPosition	${}^b\mathbf{p} = [{}^b p_0, {}^b p_1, {}^b p_2]^T$	local (body-fixed) position vector with 3 position coordinates $p_i$ in any configuration; used for local position of markers, sensors, etc.
---------------	---	--

### 6.1.1 Reference and current coordinates

An important fact on the coordinates is upon the splitting of quantities (e.g. position, rotation parameters, etc.) into reference and current (initial/visualization/...) coordinates. For the current position of a point node we have, e.g.,

$$\mathbf{p}_{\text{cur}} = \mathbf{p}_{\text{ref}} + \mathbf{u}_{\text{cur}} \quad (6.2)$$

The same holds, e.g., for rotation parameters,

$$\boldsymbol{\theta}_{\text{cur}} = \boldsymbol{\theta}_{\text{ref}} + \boldsymbol{\psi}_{\text{cur}} \quad (6.3)$$

### 6.1.2 Coordinate Systems

The left indices provide information about the coordinate system, e.g.,

$${}^0\mathbf{u} \quad (6.4)$$

is the displacement vector in the global (inertial) coordinate system 0, while

$${}^{m1}\mathbf{u} \quad (6.5)$$

represents the displacement vector in marker 1 ( $m1$ ) coordinates. Typical coordinate systems:

- ${}^0\mathbf{u}$  ... global coordinates
- ${}^b\mathbf{u}$  ... body-fixed, local coordinates
- ${}^{m0}\mathbf{u}$  ... local coordinates of (the body or node of) marker  $m0$
- ${}^{m1}\mathbf{u}$  ... local coordinates of (the body or node of) marker  $m1$

To transform the local coordinates  ${}^{m0}\mathbf{u}$  of marker 0 into global coordinates  ${}^0\mathbf{x}$ , we use

$${}^0\mathbf{u} = {}^{0,m0}\mathbf{A} {}^{m0}\mathbf{u} \quad (6.6)$$

in which  ${}^{0,m0}\mathbf{A}$  is the transformation matrix of (the body or node of) the underlying marker 0.

## 6.2 Nodes

### 6.2.1 NodePoint

A 3D point node for point masses or solid finite elements which has 3 displacement degrees of freedom for second order differential equations (ODE2).

**Additional information for NodePoint:**

- The Node has the following types = Position
- **Short name** for Python = **Point**
- **Short name** for Python (visualization object) = **VPoint**

The item **NodePoint** with type = 'Point' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector3D	3	[0.,0.,0.]	reference coordinates of node, e.g. ref. coordinates for finite elements; global position of node without displacement
initialCoordinates	Vector3D	3	[0.,0.,0.]	initial displacement coordinate
initialVelocities	Vector3D	3	[0.,0.,0.]	initial velocity coordinate
visualization	VNodePoint			parameters for visualization of item

The item VNodePoint has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size (diameter, dimensions of underlying cube, etc.) for item; size == -1.f means that default size is used
color	Float4	4	[-1.,-1.,-1.,-1.]	Default RGBA color for nodes; 4th value is alpha-transparency; R=-1.f means, that default color is used

**Detailed information on NodePoint:**

input parameter	symbol	description see tables above
referenceCoordinates	$\mathbf{q}_{\text{ref}} = [q_0, q_1, q_2]_{\text{ref}}^T = \mathbf{p}_{\text{ref}} = [r_0, r_1, r_2]^T$	
initialCoordinates	$\mathbf{q}_{\text{ini}} = [q_0, q_1, q_2]_{\text{ini}}^T = \mathbf{u}_{\text{ini}} = [u_0, u_1, u_2]_{\text{ini}}^T$	
initialVelocities	$\dot{\mathbf{q}}_{\text{ini}} = \mathbf{v}_{\text{ini}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2]_{\text{ini}}^T$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Position	$\mathbf{p}_{\text{config}} = [p_0, p_1, p_2]_{\text{config}}^T = \mathbf{u}_{\text{config}} + \mathbf{p}_{\text{ref}}$	global 3D position vector of node; $\mathbf{u}_{\text{ref}} = 0$

Displacement	$\mathbf{u}_{\text{config}} = [q_0, q_1, q_2]^T_{\text{config}}$	global 3D displacement vector of node
Velocity	$\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2]^T_{\text{config}}$	global 3D velocity vector of node
Acceleration	$\mathbf{a}_{\text{config}} = \ddot{\mathbf{q}}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, \ddot{q}_2]^T_{\text{config}}$	global 3D acceleration vector of node
Coordinates	$\mathbf{c}_{\text{config}} = \mathbf{u}_{\text{config}} = [q_0, q_1, q_2]^T_{\text{config}}$	coordinate vector of node
Coordinates_t	$\dot{\mathbf{c}}_{\text{config}} = \mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2]^T_{\text{config}}$	velocity coordinates vector of node
Coordinates_tt	$\ddot{\mathbf{c}}_{\text{config}} = \mathbf{a}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, \ddot{q}_2]^T_{\text{config}}$	acceleration coordinates vector of node

**Description of Item:** The node provides  $n_c = 3$  displacement coordinates. Equations of motion need to be provided by an according object (e.g., MassPoint, finite elements, ...). Usually, the nodal coordinates are provided in the global frame. However, the coordinate system is defined by the object (e.g. MassPoint uses global coordinates, but floating frame of reference objects use local frames). Note that for this very simple node, coordinates are identical to the nodal displacements, same for time derivatives. This is not the case, e.g. for nodes with orientation.

**Example** for NodePoint: see ObjectMassPoint



## 6.2.2 NodePoint2D

A 2D point node for point masses or solid finite elements which has 2 displacement degrees of freedom for second order differential equations.

**Additional information for NodePoint2D:**

- The Node has the following types = **Position2D**, **Position**
- **Short name** for Python = **Point2D**
- **Short name** for Python (visualization object) = **VPoint2D**

The item **NodePoint2D** with type = 'Point2D' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector2D	2	[0.,0.]	reference coordinates of node ==> e.g. ref. coordinates for finite elements; global position of node without displacement
initialCoordinates	Vector2D	2	[0.,0.]	initial displacement coordinate
initialVelocities	Vector2D	2	[0.,0.]	initial velocity coordinate
visualization	VNodePoint2D			parameters for visualization of item

The item **VNodePoint2D** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size (diameter, dimensions of underlying cube, etc.) for item; size == -1.f means that default size is used
color	Float4	4	[-1.,-1.,-1.,-1.]	Default RGBA color for nodes; 4th value is alpha-transparency; R=-1.f means, that default color is used

**Detailed information on NodePoint2D:**

input parameter	symbol	description see tables above
referenceCoordinates	$\mathbf{q}_{\text{ref}} = [q_0, q_1]^T_{\text{ref}} = \mathbf{p}_{\text{ref}} = [r_0, r_1]^T$	
initialCoordinates	$\mathbf{q}_{\text{ini}} = [q_0, q_1]^T_{\text{ini}} = [u_0, u_1]^T_{\text{ini}}$	
initialVelocities	$\dot{\mathbf{q}}_{\text{ini}} = \mathbf{v}_{\text{ini}} = [\dot{q}_0, \dot{q}_1]^T_{\text{ini}}$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Position	$\mathbf{p}_{\text{config}} = [p_0, p_1, 0]^T_{\text{config}} = \mathbf{u}_{\text{config}} + \mathbf{p}_{\text{ref}}$	global 3D position vector of node; $\mathbf{u}_{\text{ref}} = 0$
Displacement	$\mathbf{u}_{\text{config}} = [q_0, q_1, 0]^T_{\text{config}}$	global 3D displacement vector of node
Velocity	$\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, 0]^T_{\text{config}}$	global 3D velocity vector of node
Acceleration	$\mathbf{a}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, 0]^T_{\text{config}}$	global 3D acceleration vector of node
Coordinates	$\mathbf{c}_{\text{config}} = [q_0, q_1]^T_{\text{config}}$	coordinate vector of node

Coordinates_t	$\dot{\mathbf{c}}_{\text{config}} = [\dot{q}_0, \dot{q}_1]_{\text{config}}^T$	velocity coordinates vector of node
Coordinates_tt	$\ddot{\mathbf{c}}_{\text{config}} = \mathbf{a}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1]_{\text{config}}^T$	acceleration coordinates vector of node

**Description of Item:**

The node provides  $n_c = 2$  displacement coordinates. Equations of motion need to be provided by an according object (e.g., MassPoint2D). Coordinates are identical to the nodal displacements, except for the third coordinate  $u_2$ , which is zero, because  $q_2$  does not exist.

Note that for this very simple node, coordinates are identical to the nodal displacements, same for time derivatives. This is not the case, e.g. for nodes with orientation.

**Example** for NodePoint: see ObjectMassPoint2D

### 6.2.3 NodeRigidBodyEP

A 3D rigid body node based on Euler parameters for rigid bodies or beams; the node has 3 displacement coordinates (displacements of center of mass - COM: ux,uy,uz) and four rotation coordinates (Euler parameters = quaternions).

#### Additional information for NodeRigidBodyEP:

- The Node has the following types = Position, Orientation, RigidBody, RotationEulerParameters
- **Short name** for Python = **RigidEP**
- **Short name** for Python (visualization object) = **VRigidEP**

The item **NodeRigidBodyEP** with type = 'RigidBodyEP' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector7D	7	[0.,0.,0., 0.,0.,0.,0.]	reference coordinates (3 position coordinates and 4 Euler parameters) of node ==> e.g. ref. coordinates for finite elements or reference position of rigid body (e.g. for definition of joints)
initialCoordinates	Vector7D	7	[0.,0.,0., 0.,0.,0.,0.]	initial displacement coordinates and 4 Euler parameters relative to reference coordinates
initialVelocities	Vector7D	7	[0.,0.,0., 0.,0.,0.,0.]	initial velocity coordinates: time derivatives of initial displacements and Euler parameters
visualization	VNodeRigidBodyEP			parameters for visualization of item

The item VNodeRigidBodyEP has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size (diameter, dimensions of underlying cube, etc.) for item; size == -1.f means that default size is used
color	Float4	4	[-1.,-1.,-1.,-1.]	Default RGBA color for nodes; 4th value is alpha-transparency; R=-1.f means, that default color is used

#### Detailed information on NodeRigidBodyEP:

input parameter	symbol	description see tables above
referenceCoordinates	$\mathbf{q}_{\text{ref}} = [q_0, q_1, q_2, \psi_0, \psi_1, \psi_2, \psi_3]^T = [\mathbf{p}_{\text{ref}}^T, \boldsymbol{\psi}_{\text{ref}}^T]^T$	
initialCoordinates	$\mathbf{q}_{\text{ini}} = [q_0, q_1, q_2, \psi_0, \psi_1, \psi_2, \psi_3]^T = [\mathbf{u}_{\text{ini}}^T, \boldsymbol{\psi}_{\text{ini}}^T]^T$	
initialVelocities	$\dot{\mathbf{q}}_{\text{ini}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2, \dot{\psi}_0, \dot{\psi}_1, \dot{\psi}_2, \dot{\psi}_3]^T = [\dot{\mathbf{u}}_{\text{ini}}^T, \dot{\boldsymbol{\psi}}_{\text{ini}}^T]^T$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{\text{config}} = {}^0[p_0, p_1, p_2]^T_{\text{config}} = {}^0\mathbf{u}_{\text{config}} + {}^0\mathbf{p}_{\text{ref}}$	global 3D position vector of node; $\mathbf{u}_{\text{ref}} = 0$
Displacement	${}^0\mathbf{u}_{\text{config}} = [q_0, q_1, q_2]^T_{\text{config}}$	global 3D displacement vector of node
Velocity	${}^0\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2]^T_{\text{config}}$	global 3D velocity vector of node
Acceleration	${}^0\mathbf{a}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, \ddot{q}_2]^T_{\text{config}}$	global 3D acceleration vector of node
Coordinates	$\mathbf{c}_{\text{config}} = [q_0, q_1, q_2, \psi_0, \psi_1, \psi_2, \psi_3]^T_{\text{config}}$	coordinate vector of node, having 3 displacement coordinates and 4 Euler parameters
Coordinates_t	$\dot{\mathbf{c}}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2, \dot{\psi}_0, \dot{\psi}_1, \dot{\psi}_2, \dot{\psi}_3]^T_{\text{config}}$	velocity coordinates vector of node
Coordinates_tt	$\ddot{\mathbf{c}}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, \ddot{q}_2, \ddot{\psi}_0, \ddot{\psi}_1, \ddot{\psi}_2, \ddot{\psi}_3]^T_{\text{config}}$	acceleration coordinates vector of node
RotationMatrix	$[A_{00}, A_{01}, A_{02}, A_{10}, \dots, A_{21}, A_{22}]^T_{\text{config}}$	vector with 9 components of the rotation matrix ${}^{0b}\mathbf{A}_{\text{config}}$ in row-major format, in any configuration; the rotation matrix transforms local ( <i>b</i> ) to global (0) coordinates
Rotation	$[\varphi_0, \varphi_1, \varphi_2]^T_{\text{config}}$	vector with 3 components of the Euler angles in xyz-sequence ( ${}^{0b}\mathbf{A}_{\text{config}} =: \mathbf{A}_0(\varphi_0) \cdot \mathbf{A}_1(\varphi_1) \cdot \mathbf{A}_2(\varphi_2)$ ), recomputed from rotation matrix
AngularVelocity	${}^0\boldsymbol{\omega}_{\text{config}} = {}^0[\omega_0, \omega_1, \omega_2]^T_{\text{config}}$	global 3D angular velocity vector of node
AngularVelocityLocal	${}^b\boldsymbol{\omega}_{\text{config}} = {}^b[\omega_0, \omega_1, \omega_2]^T_{\text{config}}$	local (body-fixed) 3D angular velocity vector of node
AngularAcceleration	${}^0\boldsymbol{\alpha}_{\text{config}} = {}^0[\alpha_0, \alpha_1, \alpha_2]^T_{\text{config}}$	global 3D angular acceleration vector of node

**Description of Item:** All coordinates  $\mathbf{c}_{\text{config}}$  lead to second order differential equations, but there is one additional constraint equation for the quaternions. The additional constraint equation, which needs to be provided by the object, reads

$$1 - \sum_{i=0}^3 \theta_i^2 = 0. \quad (6.7)$$

The rotation matrix  ${}^{0b}\mathbf{A}_{\text{config}}$  transforms local (body-fixed) 3D positions  ${}^b\mathbf{p} = {}^b[p_0, p_1, p_2]^T$  to global 3D positions,

$${}^0\mathbf{p}_{\text{config}} = {}^{0b}\mathbf{A}_{\text{config}} {}^b\mathbf{p} \quad (6.8)$$

Note that the Euler parameters  $\boldsymbol{\theta}_{\text{cur}}$  are computed as sum of current coordinates plus reference coordinates,

$$\boldsymbol{\theta}_{\text{cur}} = \boldsymbol{\psi}_{\text{cur}} + \boldsymbol{\psi}_{\text{ref}}. \quad (6.9)$$

The rotation matrix is defined as function of the rotation parameters  $\boldsymbol{\theta} = [\theta_0, \theta_1, \theta_2, \theta_3]^T$

$${}^{0b}\mathbf{A} = \begin{bmatrix} -2\theta_3^2 - 2\theta_2^2 + 1 & -2\theta_3\theta_0 + 2\theta_2\theta_1 & 2 * \theta_3\theta_1 + 2 * \theta_2\theta_0 \\ 2\theta_3\theta_0 + 2\theta_2\theta_1 & -2\theta_3^2 - 2\theta_1^2 + 1 & 2\theta_3\theta_2 - 2\theta_1\theta_0 \\ -2\theta_2\theta_0 + 2\theta_3\theta_1 & 2\theta_3\theta_2 + 2\theta_1\theta_0 & -2\theta_2^2 - 2\theta_1^2 + 1 \end{bmatrix} \quad (6.10)$$

The derivatives of the angular velocity vectors w.r.t. the rotation velocity coordinates  $\dot{\boldsymbol{\theta}} = [\dot{\theta}_0, \dot{\theta}_1, \dot{\theta}_2, \dot{\theta}_3]^T$  lead to the **G** matrices, as used in the equations of motion for rigid bodies,

$${}^0\boldsymbol{\omega} = {}^0\mathbf{G} \dot{\boldsymbol{\theta}}, \quad (6.11)$$

$${}^b\boldsymbol{\omega} = {}^b\mathbf{G} \dot{\boldsymbol{\theta}}. \quad (6.12)$$

## 6.2.4 NodeRigidBodyRxyz

A 3D rigid body node based on Euler / Tait-Bryan angles for rigid bodies or beams; all coordinates lead to second order differential equations; NOTE that this node has a singularity if the second rotation parameter reaches  $\psi_1 = (2k - 1)\pi/2$ , with  $k \in \mathbb{N}$  or  $-k \in \mathbb{N}$ .

**Additional information for NodeRigidBodyRxyz:**

- The Node has the following types = Position, Orientation, RigidBody, RotationRxyz
- **Short name** for Python = **RigidRxyz**
- **Short name** for Python (visualization object) = **VRigidRxyz**

The item **NodeRigidBodyRxyz** with type = 'RigidBodyRxyz' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector6D	6	[0.,0.,0., 0.,0.,0.]	reference coordinates (3 position and 3 xyz Euler angles) of node ==> e.g. ref. coordinates for finite elements or reference position of rigid body (e.g. for definition of joints)
initialCoordinates	Vector6D	6	[0.,0.,0., 0.,0.,0.]	initial displacement coordinates: ux,uy,uz and 3 Euler angles (xyz) relative to reference coordinates
initialVelocities	Vector6D	6	[0.,0.,0., 0.,0.,0.]	initial velocity coordinate: time derivatives of ux,uy,uz and of 3 Euler angles (xyz)
visualization	VNodeRigidBodyRxyz			parameters for visualization of item

The item VNodeRigidBodyRxyz has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size (diameter, dimensions of underlying cube, etc.) for item; size == -1.f means that default size is used
color	Float4	4	[-1.,-1.,-1.,-1.]	Default RGBA color for nodes; 4th value is alpha-transparency; R=-1.f means, that default color is used

**Detailed information on NodeRigidBodyRxyz:**

input parameter	symbol	description see tables above
referenceCoordinates	$\mathbf{q}_{\text{ref}} = [q_0, q_1, q_2, \psi_0, \psi_1, \psi_2]_{\text{ref}}^T = [\mathbf{p}_{\text{ref}}^T, \boldsymbol{\psi}_{\text{ref}}^T]^T$	
initialCoordinates	$\mathbf{q}_{\text{ini}} = [q_0, q_1, q_2, \psi_0, \psi_1, \psi_2]_{\text{ini}}^T = [\mathbf{u}_{\text{ini}}^T, \boldsymbol{\psi}_{\text{ini}}^T]^T$	
initialVelocities	$\dot{\mathbf{q}}_{\text{ini}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2, \dot{\psi}_0, \dot{\psi}_1, \dot{\psi}_2]_{\text{ini}}^T = [\dot{\mathbf{u}}_{\text{ini}}^T, \dot{\boldsymbol{\psi}}_{\text{ini}}^T]^T$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{\text{config}} = {}^0[p_0, p_1, p_2]^T_{\text{config}} = {}^0\mathbf{u}_{\text{config}} + {}^0\mathbf{p}_{\text{ref}}$	global 3D position vector of node; $\mathbf{u}_{\text{ref}} = 0$
Displacement	${}^0\mathbf{u}_{\text{config}} = [q_0, q_1, q_2]^T_{\text{config}}$	global 3D displacement vector of node
Velocity	${}^0\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2]^T_{\text{config}}$	global 3D velocity vector of node
Acceleration	${}^0\mathbf{a}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, \ddot{q}_2]^T_{\text{config}}$	global 3D acceleration vector of node
Coordinates	$\mathbf{c}_{\text{config}} = [q_0, q_1, q_2, \psi_0, \psi_1, \psi_2]^T_{\text{config}}$	coordinate vector of node, having 3 displacement coordinates and 3 Euler angles
Coordinates_t	$\dot{\mathbf{c}}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2, \dot{\psi}_0, \dot{\psi}_1, \dot{\psi}_2]^T_{\text{config}}$	velocity coordinates vector of node
Coordinates_tt	$\ddot{\mathbf{c}}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, \ddot{q}_2, \ddot{\psi}_0, \ddot{\psi}_1, \ddot{\psi}_2]^T_{\text{config}}$	acceleration coordinates vector of node
RotationMatrix	$[A_{00}, A_{01}, A_{02}, A_{10}, \dots, A_{21}, A_{22}]^T_{\text{config}}$	vector with 9 components of the rotation matrix ${}^{0b}\mathbf{A}_{\text{config}}$ in row-major format, in any configuration; the rotation matrix transforms local ( $b$ ) to global ( $0$ ) coordinates
Rotation	$[\varphi_0, \varphi_1, \varphi_2]^T_{\text{config}} = [\psi_0, \psi_1, \psi_2]^T_{\text{ref}} + [\psi_0, \psi_1, \psi_2]^T_{\text{config}}$	vector with 3 components of the Euler / Tait-Bryan angles in xyz-sequence ( ${}^{0b}\mathbf{A}_{\text{config}} =: \mathbf{A}_0(\varphi_0) \cdot \mathbf{A}_1(\varphi_1) \cdot \mathbf{A}_2(\varphi_2)$ ), recomputed from rotation matrix
AngularVelocity	${}^0\boldsymbol{\omega}_{\text{config}} = {}^0[\omega_0, \omega_1, \omega_2]^T_{\text{config}}$	global 3D angular velocity vector of node
AngularVelocityLocal	${}^b\boldsymbol{\omega}_{\text{config}} = {}^b[\omega_0, \omega_1, \omega_2]^T_{\text{config}}$	local (body-fixed) 3D angular velocity vector of node
AngularAcceleration	${}^0\boldsymbol{\alpha}_{\text{config}} = {}^0[\alpha_0, \alpha_1, \alpha_2]^T_{\text{config}}$	global 3D angular acceleration vector of node

**Description of Item:** The node has 3 displacement coordinates  $[q_0, q_1, q_2]^T$  and 3 rotation coordinates  $[\psi_0, \psi_1, \psi_2]^T$  for consecutive rotations around the 0, 1 and 2-axis ( $x$ ,  $y$  and  $z$ ). All coordinates  $\mathbf{c}_{\text{config}}$  lead to second order differential equations. The rotation matrix  ${}^{0b}\mathbf{A}_{\text{config}}$  transforms local (body-fixed) 3D positions  ${}^b\mathbf{p} = {}^b[p_0, p_1, p_2]^T$  to global 3D positions,

$${}^0\mathbf{p}_{\text{config}} = {}^{0b}\mathbf{A}_{\text{config}} {}^b\mathbf{p} \quad (6.13)$$

Note that the Euler angles  $\boldsymbol{\theta}_{\text{cur}}$  are computed as sum of current coordinates plus reference coordinates,

$$\boldsymbol{\theta}_{\text{cur}} = \boldsymbol{\psi}_{\text{cur}} + \boldsymbol{\psi}_{\text{ref}}. \quad (6.14)$$

The rotation matrix is defined as function of the rotation parameters  $\boldsymbol{\theta} = [\theta_0, \theta_1, \theta_2]^T$

$${}^{0b}\mathbf{A} = \mathbf{A}_0(\theta_0)\mathbf{A}_1(\theta_1)\mathbf{A}_2(\theta_2) \quad (6.15)$$

The derivatives of the angular velocity vectors w.r.t. the rotation velocity coordinates  $\dot{\boldsymbol{\theta}} = [\dot{\theta}_0, \dot{\theta}_1, \dot{\theta}_2]^T$  lead to the  $\mathbf{G}$  matrices, as used in the equations of motion for rigid bodies,

$${}^0\boldsymbol{\omega} = {}^0\mathbf{G} \dot{\boldsymbol{\theta}}, \quad (6.16)$$

$${}^b\boldsymbol{\omega} = {}^b\mathbf{G} \dot{\boldsymbol{\theta}}. \quad (6.17)$$

## 6.2.5 NodeRigidBodyRotVecLG

A 3D rigid body node based on rotation vector and Lie group methods for rigid bodies or beams; the node has 3 displacement coordinates and three rotation coordinates.

**Additional information for NodeRigidBodyRotVecLG:**

- The Node has the following types = Position, Orientation, RigidBody, RotationRotationVector, RotationLieGroup
- **Short name** for Python = **RigidRotVecLG**
- **Short name** for Python (visualization object) = **VRigidRotVecLG**

The item **NodeRigidBodyRotVecLG** with type = 'RigidBodyRotVecLG' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector6D	3	[0.,0.,0., 0.,0.,0.]	reference coordinates (position and rotation vector $\mathbf{v}$ ) of node ==> e.g. ref. coordinates for finite elements or reference position of rigid body (e.g. for definition of joints)
initialCoordinates	Vector6D	3	[0.,0.,0., 0.,0.,0.]	initial displacement coordinates $\mathbf{u}$ and rotation vector $\mathbf{v}$ relative to reference coordinates
initialVelocities	Vector6D	3	[0.,0.,0., 0.,0.,0.]	initial velocity coordinate: time derivatives of displacement and angular velocity vector
visualization	VNodeRigidBodyRotVecLG			parameters for visualization of item

The item VNodeRigidBodyRotVecLG has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size (diameter, dimensions of underlying cube, etc.) for item; size == -1.f means that default size is used
color	Float4	4	[-1.,-1.,-1.,-1.]	Default RGBA color for nodes; 4th value is alpha-transparency; R=-1.f means, that default color is used

**Detailed information on NodeRigidBodyRotVecLG:**

input parameter	symbol	description see tables above
referenceCoordinates	$\mathbf{q}_{\text{ref}} = [q_0, q_1, q_2, v_0, v_1, v_2]^T_{\text{ref}} = [\mathbf{p}_{\text{ref}}^T, \mathbf{v}_{\text{ref}}^T]^T$	
initialCoordinates	$\mathbf{q}_{\text{ini}} = [q_0, q_1, q_2, v_0, v_1, v_2]^T_{\text{ini}} = [\mathbf{u}_{\text{ini}}^T, \mathbf{v}_{\text{ini}}^T]^T$	
initialVelocities	$\dot{\mathbf{q}}_{\text{ini}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2, \dot{v}_0, \dot{v}_1, \dot{v}_2]^T_{\text{ini}} = [\dot{\mathbf{u}}_{\text{ini}}^T, \dot{\mathbf{v}}_{\text{ini}}^T]^T$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{\text{config}} = {}^0[p_0, p_1, p_2]^T_{\text{config}} = {}^0\mathbf{u}_{\text{config}} + {}^0\mathbf{p}_{\text{ref}}$	global 3D position vector of node; $\mathbf{u}_{\text{ref}} = 0$
Displacement	${}^0\mathbf{u}_{\text{config}} = [q_0, q_1, q_2]^T_{\text{config}}$	global 3D displacement vector of node
Velocity	${}^0\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2]^T_{\text{config}}$	global 3D velocity vector of node
Coordinates	$\mathbf{c}_{\text{config}} = [q_0, q_1, q_2, v_0, v_1, v_2]^T_{\text{config}}$	coordinate vector of node, having 3 displacement coordinates and 3 Euler angles
Coordinates_t	$\dot{\mathbf{c}}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2, \dot{v}_0, \dot{v}_1, \dot{v}_2]^T_{\text{config}}$	velocity coordinates vector of node
RotationMatrix	$[A_{00}, A_{01}, A_{02}, A_{10}, \dots, A_{21}, A_{22}]^T_{\text{config}}$	vector with 9 components of the rotation matrix ${}^{0b}\mathbf{A}_{\text{config}}$ in row-major format, in any configuration; the rotation matrix transforms local ( <i>b</i> ) to global ( <i>0</i> ) coordinates
Rotation	$[\varphi_0, \varphi_1, \varphi_2]^T_{\text{config}}$	vector with 3 components of the Euler / Tait-Bryan angles in xyz-sequence ( ${}^{0b}\mathbf{A}_{\text{config}} =: \mathbf{A}_0(\varphi_0) \cdot \mathbf{A}_1(\varphi_1) \cdot \mathbf{A}_2(\varphi_2)$ ), recomputed from rotation matrix
AngularVelocity	${}^0\boldsymbol{\omega}_{\text{config}} = {}^0[\omega_0, \omega_1, \omega_2]^T_{\text{config}}$	global 3D angular velocity vector of node
AngularVelocityLocal	${}^b\boldsymbol{\omega}_{\text{config}} = {}^b[\omega_0, \omega_1, \omega_2]^T_{\text{config}}$	local (body-fixed) 3D angular velocity vector of node

**Description of Item:** For a detailed description on the rigid body dynamics formulation using this node, see Holzinger and Gerstmayr [4].

The node has 3 displacement coordinates  $[q_0, q_1, q_2]^T$  and three rotation coordinates, which is the rotation vector

$$\boldsymbol{\nu} = \boldsymbol{\varphi}\mathbf{n} = \boldsymbol{\nu}_{\text{config}} + \boldsymbol{\nu}_{\text{ref}}, \quad (6.18)$$

with the rotation angle  $\varphi$  and the rotation axis  $\mathbf{n}$ . All coordinates  $\mathbf{c}_{\text{config}}$  lead to second order differential equations, however the rotation vector cannot be used as a conventional parameterization. It must be computed within a nonlinear update, using appropriate Lie group methods.

The rotation matrix  ${}^{0b}\mathbf{A}_{\text{config}}$  transforms local (body-fixed) 3D positions  ${}^b\mathbf{p} = {}^b[p_0, p_1, p_2]^T$  to global 3D positions,

$${}^0\mathbf{p}_{\text{config}} = {}^{0b}\mathbf{A}_{\text{config}} {}^b\mathbf{p} \quad (6.19)$$

A Lie group integrator must be used with this node, which is why the is used, the rotation parameter velocities are identical to the local angular velocity  ${}^b\boldsymbol{\omega}$  and thus the matrix  ${}^b\mathbf{G}$  becomes the identity matrix.



## 6.2.6 NodeRigidBody2D

A 2D rigid body node for rigid bodies or beams; the node has 2 displacement degrees of freedom and one rotation coordinate (rotation around z-axis:  $\psi$ ). All coordinates are ODE2, used for second order differential equations.

### Additional information for NodeRigidBody2D:

- The Node has the following types = Position2D, Orientation2D, Position, Orientation, RigidBody
- **Short name** for Python = **Rigid2D**
- **Short name** for Python (visualization object) = **VRigid2D**

The item **NodeRigidBody2D** with type = 'RigidBody2D' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector3D	3	[0.,0.,0.]	reference coordinates (x-pos,y-pos and rotation) of node ==> e.g. ref. coordinates for finite elements; global position of node without displacement
initialCoordinates	Vector3D	3	[0.,0.,0.]	initial displacement coordinates and angle (relative to reference coordinates)
initialVelocities	Vector3D	3	[0.,0.,0.]	initial velocity coordinates
visualization	VNodeRigidBody2D			parameters for visualization of item

The item VNodeRigidBody2D has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size (diameter, dimensions of underlying cube, etc.) for item; size == -1.f means that default size is used
color	Float4	4	[-1.,-1.,-1.,-1.]	Default RGBA color for nodes; 4th value is alpha-transparency; R=-1.f means, that default color is used

### Detailed information on NodeRigidBody2D:

input parameter	symbol	description see tables above
referenceCoordinates	$\mathbf{q}_{\text{ref}} = [q_0, q_1, \psi_0]_{\text{ref}}^T$	
initialCoordinates	$\mathbf{q}_{\text{ini}} = [q_0, q_1, \psi_0]_{\text{ini}}^T$	
initialVelocities	$\dot{\mathbf{q}}_{\text{ini}} = [\dot{q}_0, \dot{q}_1, \dot{\psi}_0]_{\text{ini}}^T = [v_0, v_1, \omega_2]_{\text{ini}}^T$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{\text{config}} = {}^0[p_0, p_1, 0]_{\text{config}}^T = {}^0\mathbf{u}_{\text{config}} + {}^0\mathbf{p}_{\text{ref}}$	global 3D position vector of node; $\mathbf{u}_{\text{ref}} = 0$

Displacement	${}^0\mathbf{u}_{\text{config}} = [q_0, q_1, 0]_{\text{config}}^T$	global 3D displacement vector of node
Velocity	${}^0\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, 0]_{\text{config}}^T$	global 3D velocity vector of node
Acceleration	${}^0\mathbf{a}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, 0]_{\text{config}}^T$	global 3D acceleration vector of node
AngularVelocity	${}^0\boldsymbol{\omega}_{\text{config}} = [0, 0, \dot{\psi}_0]_{\text{config}}^T$	global 3D angular velocity vector of node
Coordinates	$\mathbf{c}_{\text{config}} = [q_0, q_1, \psi_0]_{\text{config}}^T$	coordinate vector of node, having 2 displacement coordinates and 1 angle
Coordinates_t	$\dot{\mathbf{c}}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{\psi}_0]_{\text{config}}^T$	velocity coordinates vector of node
Coordinates_tt	$\ddot{\mathbf{c}}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, \ddot{\psi}_0]_{\text{config}}^T$	acceleration coordinates vector of node
RotationMatrix	$[A_{00}, A_{01}, A_{02}, A_{10}, \dots, A_{21}, A_{22}]_{\text{config}}^T$	vector with 9 components of the rotation matrix ${}^{0b}\mathbf{A}_{\text{config}}$ in row-major format, in any configuration; the rotation matrix transforms local ( $b$ ) to global ( $0$ ) coordinates
Rotation	$[\theta_0]_{\text{config}}^T = [\psi_0]_{\text{ref}}^T + [\psi_0]_{\text{config}}^T$	vector with 1 angle around out of plane axis
AngularVelocityLocal	${}^b\boldsymbol{\omega}_{\text{config}} = [0, 0, \dot{\psi}_0]_{\text{config}}^T$	local (body-fixed) 3D angular velocity vector of node
AngularAcceleration	${}^0\boldsymbol{\alpha}_{\text{config}} = [0, 0, \ddot{\psi}_0]_{\text{config}}^T$	global 3D angular acceleration vector of node

**Description of Item:** The node provides 2 displacement coordinates (displacement of center of mass, COM,  $(q_0, q_1)$ ) and 1 rotation parameter ( $\theta_0$ ). According equations need to be provided by an according object (e.g., RigidBody2D). Using the rotation parameter  $\theta_{0\text{config}} = \psi_{0\text{ref}} + \psi_{0\text{config}}$ , the rotation matrix is defined as

$${}^{0b}\mathbf{A}_{\text{config}} = \begin{bmatrix} \cos(\theta_0) & -\sin(\theta_0) & 0 \\ \sin(\theta_0) & \cos(\theta_0) & 0 \\ 0 & 0 & 1 \end{bmatrix}_{\text{config}} \quad (6.20)$$

**Example** for NodeRigidBody2D: see ObjectRigidBody2D

### 6.2.7 Node1D

A node with one ODE2 coordinate for one dimensional (1D) problems; use e.g. for scalar dynamic equations (Mass1D) and mass-spring-damper mechanisms, representing either translational or rotational degrees of freedom: in most cases, Node1D is equivalent to NodeGenericODE2 using one coordinate, however, it offers a transformation to 3D translational or rotational motion and allows to couple this node to 2D or 3D bodies.

#### Additional information for Node1D:

- The Node has the following types = GenericODE2

The item **Node1D** with type = '1D' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector		[0.]	reference coordinate of node (in vector form)
initialCoordinates	Vector		[0.]	initial displacement coordinate (in vector form)
initialVelocities	Vector		[0.]	initial velocity coordinate (in vector form)
visualization	VNode1D			parameters for visualization of item

The item VNode1D has the following parameters:

Name	type	size	default value	description
show	bool		False	set true, if item is shown in visualization and false if it is not shown; The node1D is represented as reference position and displacement along the global x-axis, which must not agree with the representation in the object using the Node1D

#### Detailed information on Node1D:

input parameter	symbol	description see tables above
referenceCoordinates	$[q_0]_{\text{ref}}^T$	
initialCoordinates	$[q_0]_{\text{ini}}^T$	
initialVelocities	$[\dot{q}_0]_{\text{ini}}^T$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Coordinates	$\mathbf{q}_{\text{config}} = [q_0]_{\text{config}}^T$	ODE2 coordinate of node (in vector form)
Coordinates_t	$\dot{\mathbf{q}}_{\text{config}} = [\dot{q}_0]_{\text{config}}^T$	ODE2 velocity coordinate of node (in vector form)
Coordinates_tt	$\ddot{\mathbf{q}}_{\text{config}} = [\ddot{q}_0]_{\text{config}}^T$	ODE2 acceleration coordinate of node (in vector form)

**Description of Item:** The current position/rotation coordinate of the 1D node is computed from

$$p_0 = q_{0\text{ref}} + q_{0\text{cur}} \quad (6.21)$$

The coordinate leads to one second order differential equation. The graphical representation and the (internal) position of the node is

$$p_{\text{config}} = \begin{bmatrix} p_{0\text{config}} \\ 0 \\ 0 \end{bmatrix} \quad (6.22)$$

The (internal) velocity vector is  $[p_{0\text{config}}, 0, 0]^T$ .

### 6.2.8 NodePoint2DSlope1

A 2D point/slope vector node for planar Bernoulli-Euler ANCF (absolute nodal coordinate formulation) beam elements; the node has 4 displacement degrees of freedom (2 for displacement of point node and 2 for the slope vector 'slopex'); all coordinates lead to second order differential equations; the slope vector defines the directional derivative w.r.t the local axial (x) coordinate, denoted as (); in straight configuration aligned at the global x-axis, the slope vector reads  $\mathbf{r}' = [r'_x \ r'_y]^T = [1 \ 0]^T$ .

#### Additional information for NodePoint2DSlope1:

- The Node has the following types = Position2D, Orientation2D, Point2DSlope1, Position, Orientation
- **Short name** for Python = **Point2DS1**
- **Short name** for Python (visualization object) = **VPoint2DS1**

The item **NodePoint2DSlope1** with type = 'Point2DSlope1' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector4D	4	[0.,0.,1.,0.]	reference coordinates (x-pos,y-pos; x-slopex, y-slopex) of node; global position of node without displacement
initialCoordinates	Vector4D	4	[0.,0.,0.,0.]	initial displacement coordinates: ux, uy and x/y 'displacements' of slopex
initialVelocities	Vector4D	4	[0.,0.,0.,0.]	initial velocity coordinates
visualization	VNodePoint2DSlope1			parameters for visualization of item

The item **VNodePoint2DSlope1** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size (diameter, dimensions of underlying cube, etc.) for item; size == -1.f means that default size is used
color	Float4	4	[-1.,-1.,-1.,-1.]	Default RGBA color for nodes; 4th value is alpha-transparency; R=-1.f means, that default color is used

**Detailed information on NodePoint2DSlope1:** The following output parameters are available as **Output-VariableType** in sensors and other functions:

output parameter	symbol	description
Position		global 3D position vector of node (=displacement+reference position)
Displacement		global 3D displacement vector of node
Velocity		global 3D velocity vector of node
Coordinates		coordinates vector of node (2 displacement coordinates + 2 slope vector coordinates)

Coordinates_t		velocity coordinates vector of node (derivative of the 2 displacement coordinates + 2 slope vector coordinates)
Coordinates_tt		acceleration coordinates vector of node (derivative of the 2 displacement coordinates + 2 slope vector coordinates)

### 6.2.9 NodeGenericODE2

A node containing a number of ODE2 variables; use e.g. for scalar dynamic equations (Mass1D) or for the ALECable element.

**Additional information for NodeGenericODE2:**

- The Node has the following types = GenericODE2

The item **NodeGenericODE2** with type = 'GenericODE2' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector		[]	generic reference coordinates of node; must be consistent with numberOfODE2Coordinates
initialCoordinates	Vector		[]	initial displacement coordinates; must be consistent with numberOfODE2Coordinates
initialCoordinates_t	Vector		[]	initial velocity coordinates; must be consistent with numberOfODE2Coordinates
numberOfODE2Coordinates	Index		0	number of generic ODE2 coordinates
visualization	VNodeGenericODE2			parameters for visualization of item

The item **VNodeGenericODE2** has the following parameters:

Name	type	size	default value	description
show	bool		False	set true, if item is shown in visualization and false if it is not shown

**Detailed information on NodeGenericODE2:**

input parameter	symbol	description see tables above
referenceCoordinates	$\mathbf{q}_{\text{ref}} = [q_0, \dots, q_{n_c}]_{\text{ref}}^T$	
initialCoordinates	$\mathbf{q}_{\text{ini}} = [q_0, \dots, q_{n_c}]_{\text{ini}}^T$	
initialCoordinates_t	$\dot{\mathbf{q}}_{\text{ini}} = [\dot{q}_0, \dots, \dot{q}_{n_c}]_{\text{ini}}^T$	
numberOfODE2Coordinates	$n_c$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Coordinates	$\mathbf{q}_{\text{config}} = [q_0, \dots, q_{n_c}]_{\text{config}}^T$	coordinates vector of node
Coordinates_t	$\dot{\mathbf{q}}_{\text{config}} = [\dot{q}_0, \dots, \dot{q}_{n_c}]_{\text{config}}^T$	velocity coordinates vector of node
Coordinates_tt	$\ddot{\mathbf{q}}_{\text{config}} = [\ddot{q}_0, \dots, \ddot{q}_{n_c}]_{\text{config}}^T$	acceleration coordinates vector of node

### 6.2.10 NodeGenericData

A node containing a number of data (history) variables; use e.g. for contact (active set), friction or plasticity (history variable).

**Additional information for NodeGenericData:**

- The Node has the following types = GenericData

The item **NodeGenericData** with type = 'GenericData' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
initialCoordinates	Vector		[]	initial data coordinates
numberOfDataCoordinates	Index		0	number of generic data coordinates (history variables)
visualization	VNodeGenericData			parameters for visualization of item

The item VNodeGenericData has the following parameters:

Name	type	size	default value	description
show	bool		False	set true, if item is shown in visualization and false if it is not shown

**Detailed information on NodeGenericData:**

input parameter	symbol	description see tables above
initialCoordinates	$\mathbf{x}_{\text{ini}} = [x_0, \dots, x_{n_c}]_{\text{ini}}^T$	
numberOfDataCoordinates	$n_c$	

**The following output parameters are available as OutputVariableType in sensors and other functions:**

output parameter	symbol	description
Coordinates	$\mathbf{x}_{\text{config}} = [x_0, \dots, x_{n_c}]_{\text{config}}^T$	data coordinates (history variables) vector of node



### 6.2.11 NodePointGround

A 3D point node fixed to ground. The node can be used as NodePoint, but it does not generate coordinates. Applied or reaction forces do not have any effect.

**Additional information for NodePointGround:**

- The Node has the following types = Ground, Position2D, Position, GenericODE2
- **Short name** for Python = **PointGround**
- **Short name** for Python (visualization object) = **VPointGround**

The item **NodePointGround** with type = 'PointGround' has the following parameters:

Name	type	size	default value	description
name	String		"	node's unique name
referenceCoordinates	Vector3D	3	[0.,0.,0.]	reference coordinates of node ==> e.g. ref. coordinates for finite elements; global position of node without displacement
visualization	VNodePointGround			parameters for visualization of item

The item VNodePointGround has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size (diameter, dimensions of underlying cube, etc.) for item; size == -1.f means that default size is used
color	Float4	4	[-1.,-1.,-1.,-1.]	Default RGBA color for nodes; 4th value is alpha-transparency; R=-1.f means, that default color is used

**Detailed information on NodePointGround:**

input parameter	symbol	description see tables above
referenceCoordinates	$\mathbf{q}_{\text{ref}} = [q_0, q_1, q_2]_{\text{ref}}^T = \mathbf{p}_{\text{ref}} = [r_0, r_1, r_2]^T$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	$\mathbf{p}_{\text{config}} = [p_0, p_1, p_2]_{\text{config}}^T = \mathbf{p}_{\text{ref}}$	global 3D position vector of node (=reference position)
Displacement	$\mathbf{u}_{\text{config}} = [0, 0, 0]_{\text{config}}^T$	zero 3D vector
Velocity	$\mathbf{v}_{\text{config}} = [0, 0, 0]_{\text{config}}^T$	zero 3D vector
Coordinates	$\mathbf{c}_{\text{config}} = []$	vector of length zero
Coordinates_t	$\dot{\mathbf{c}}_{\text{config}} = []$	vector of length zero

## 6.3 Objects

### 6.3.1 ObjectMassPoint

A 3D mass point which is attached to a position-based node, usually NodePoint.

**Additional information for ObjectMassPoint:**

- The Object has the following types = Body, SingleNoded
- Requested node type = Position
- **Short name** for Python = **MassPoint**
- **Short name** for Python (visualization object) = **VMassPoint**

The item **ObjectMassPoint** with type = 'MassPoint' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
physicsMass	UReal		0.	mass [SI:kg] of mass point
nodeNumber	NodeIndex		MAXINT	node number (type NodeIndex) for mass point
visualization	VObjectMassPoint			parameters for visualization of item

The item VObjectMassPoint has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
graphicsData	BodyGraphicsData			Structure contains data for body visualization; data is defined in special list / dictionary structure

**Detailed information on ObjectMassPoint:**

input parameter	symbol	description see tables above
physicsMass	$m$	
nodeNumber	$n0$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{\text{config}}({}^b\mathbf{p}) = {}^0\mathbf{u}_{\text{config}} + {}^0\mathbf{p}_{\text{ref}} + {}^{0b}\mathbf{A} {}^b\mathbf{p}$	global position vector of translated local position; local (body) coordinate system = global coordinate system
Displacement	${}^0\mathbf{u}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2]_{\text{config}}^T$	global displacement vector of mass point
Velocity	${}^0\mathbf{v}_{\text{config}} = {}^0\dot{\mathbf{u}}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, \ddot{q}_2]_{\text{config}}^T$	global velocity vector of mass point
Acceleration	${}^0\mathbf{a}_{\text{config}} = {}^0\ddot{\mathbf{u}}_{\text{config}} = [\ddot{\ddot{q}}_0, \ddot{\ddot{q}}_1, \ddot{\ddot{q}}_2]_{\text{config}}^T$	global acceleration vector of mass point

**Description of Item:**

intermediate variables	symbol	description
node position	${}^0\mathbf{p}_{\text{config}} = {}^0\mathbf{p}(n_0)_{\text{config}}$	position of mass point which is provided by node $n_0$ in any configuration
node displacement	${}^0\mathbf{u}_{\text{config}} = [q_0, q_1, q_2]_{\text{config}}^T = {}^0\mathbf{u}(n_0)_{\text{config}}$	displacement of mass point which is provided by node $n_0$ in any configuration
node velocity	${}^0\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, \dot{q}_2]_{\text{config}}^T = {}^0\mathbf{v}(n_0)_{\text{config}}$	velocity of mass point which is provided by node $n_0$ in any configuration
transformation matrix	${}^{0b}\mathbf{A} = \mathbf{I}^{3 \times 3}$	transformation of local body ( $b$ ) coordinates to global (0) coordinates; this is the constant unit matrix, because local = global coordinates for the mass point
residual forces	${}^0\mathbf{f} = [f_0, f_1, f_2]^T$	residual of all forces on mass point
applied forces	${}^0\mathbf{f}_a = [f_0, f_1, f_2]^T$	applied forces (loads, connectors, joint reaction forces, ...)

Equations of motion:

$$\begin{bmatrix} m & 0 & 0 \\ 0 & m & 0 \\ 0 & 0 & m \end{bmatrix} \begin{bmatrix} \ddot{q}_0 \\ \ddot{q}_1 \\ \ddot{q}_2 \end{bmatrix} = \begin{bmatrix} f_0 \\ f_1 \\ f_2 \end{bmatrix}. \quad (6.23)$$

For example, a LoadCoordinate on coordinate 1 of the node would add a term in  $f_1$  on the RHS.

Position-based markers can measure position  $\mathbf{p}_{\text{config}}$ . The **position jacobian**

$$\mathbf{J}_{\text{pos}} = \partial \mathbf{p}_{\text{cur}} / \partial \mathbf{c}_{\text{cur}} = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 0 \end{bmatrix} \quad (6.24)$$

transforms the action of global applied forces  ${}^0\mathbf{f}_a$  of position-based markers on the coordinates  $\mathbf{c}$

$$\mathbf{Q} = \mathbf{J}_{\text{pos}} {}^0\mathbf{f}_a. \quad (6.25)$$

Example for ObjectMassPoint:

---

```

node = mbs.AddNode(NodePoint(referenceCoordinates = [1,1,0],
                             initialCoordinates=[0.5,0,0],
                             initialVelocities=[0.5,0,0]))
mbs.AddObject(MassPoint(nodeNumber = node, physicsMass=1))

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result
testError = mbs.GetNodeOutput(node, exu.OutputVariableType.Position)[0] - 2
#final x-coordinate of position shall be 2

```

---

### 6.3.2 ObjectMassPoint2D

A 2D mass point which is attached to a position-based 2D node.

**Additional information for ObjectMassPoint2D:**

- The Object has the following types = Body, SingleNoded
- Requested node type = Position2D + Position
- **Short name** for Python = **MassPoint2D**
- **Short name** for Python (visualization object) = **VMassPoint2D**

The item **ObjectMassPoint2D** with type = 'MassPoint2D' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
physicsMass	UReal		0.	mass [SI:kg] of mass point
nodeNumber	NodeIndex		MAXINT	node number (type NodeIndex) for mass point
visualization	VObjectMassPoint2D			parameters for visualization of item

The item **VObjectMassPoint2D** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
graphicsData	BodyGraphicsData			Structure contains data for body visualization; data is defined in special list / dictionary structure

**Detailed information on ObjectMassPoint2D:**

input parameter	symbol	description see tables above
physicsMass	$m$	
nodeNumber	$n0$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{\text{config}}({}^b\mathbf{p}) = {}^0\mathbf{u}_{\text{config}} + {}^0\mathbf{p}_{\text{ref}} + {}^{0b}\mathbf{A} {}^b\mathbf{p}$	global position vector of translated local position; local (body) coordinate system = global coordinate system
Displacement	${}^0\mathbf{u}_{\text{config}} = [q_0, q_1, 0]^T_{\text{config}}$	global displacement vector of mass point
Velocity	${}^0\mathbf{v}_{\text{config}} = {}^0\dot{\mathbf{u}}_{\text{config}} = [\dot{q}_0, \dot{q}_1, 0]^T_{\text{config}}$	global velocity vector of mass point
Acceleration	${}^0\mathbf{a}_{\text{config}} = {}^0\ddot{\mathbf{u}}_{\text{config}} = [\ddot{q}_0, \ddot{q}_1, 0]^T_{\text{config}}$	global acceleration vector of mass point

**Description of Item:**

intermediate variables	symbol	description
------------------------	--------	-------------

node position	${}^0\mathbf{p}_{\text{config}} = {}^0\mathbf{p}(n_0)_{\text{config}}$	position of mass point which is provided by node $n_0$ in any configuration
node displacement	${}^0\mathbf{u}_{\text{config}} = [\dot{q}_0, \dot{q}_1, 0]_{\text{config}}^T = {}^0\mathbf{u}(n_0)_{\text{config}}$	displacement of mass point which is provided by node $n_0$ in any configuration
node velocity	${}^0\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, 0]_{\text{config}}^T = {}^0\mathbf{v}(n_0)_{\text{config}}$	velocity of mass point which is provided by node $n_0$ in any configuration
transformation matrix	${}^{0b}\mathbf{A} = \mathbf{I}^{3 \times 3}$	transformation of local body ( $b$ ) coordinates to global (0) coordinates; this is the constant unit matrix, because local = global coordinates for the mass point
residual forces	${}^0\mathbf{f} = [f_0, f_1]^T$	residual of all forces on mass point
applied forces	${}^0\mathbf{f}_a = [f_0, f_1, f_2]^T$	applied forces (loads, connectors, joint reaction forces, ...)

**Equations of motion:**

$$\begin{bmatrix} m & 0 \\ 0 & m \end{bmatrix} \begin{bmatrix} \ddot{q}_0 \\ \ddot{q}_1 \end{bmatrix} = \begin{bmatrix} f_0 \\ f_1 \end{bmatrix}. \quad (6.26)$$

For example, a LoadCoordinate on coordinate 1 of the node would add a term in  $f_1$  on the RHS.

Position-based markers can measure position  $\mathbf{p}_{\text{config}}$ . The **position jacobian**

$$\mathbf{J}_{\text{pos}} = \partial \mathbf{p}_{\text{cur}} / \partial \mathbf{c}_{\text{cur}} = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \end{bmatrix} \quad (6.27)$$

transforms the action of global applied forces  ${}^0\mathbf{f}_a$  of position-based markers on the coordinates  $\mathbf{c}$

$$\mathbf{Q} = \mathbf{J}_{\text{pos}} {}^0\mathbf{f}_a. \quad (6.28)$$

**Example for ObjectMassPoint2D:**

---

```

node = mbs.AddNode(NodePoint2D(referenceCoordinates = [1,1],
                                initialCoordinates=[0.5,0],
                                initialVelocities=[0.5,0]))
mbs.AddObject(MassPoint2D(nodeNumber = node, physicsMass=1))

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result
testError = mbs.GetNodeOutput(node, exu.OutputVariableType.Position)[0] - 2
#final x-coordinate of position shall be 2

```

---

### 6.3.3 ObjectMass1D

A 1D (translational) mass which is attached to Node1D. Note, that the mass does not need to have the interpretation as a translational mass.

**Additional information for ObjectMass1D:**

- The Object has the following types = Body, SingleNoded
- Requested node type = GenericCODE2
- **Short name** for Python = **Mass1D**
- **Short name** for Python (visualization object) = **VMass1D**

The item **ObjectMass1D** with type = 'Mass1D' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
physicsMass	UReal		0.	mass [SI:kg] of mass
nodeNumber	NodeIndex		MAXINT	node number (type NodeIndex) for Node1D
referencePosition	Vector3D	3	[0.,0.,0.]	a reference position, used to transform the 1D coordinate to a position
referenceRotation	Matrix3D		[[1,0,0], [0,1,0], [0,0,1]]	the constant body rotation matrix, which transforms body-fixed (b) to global (0) coordinates
visualization	VObjectMass1D			parameters for visualization of item

The item VObjectMass1D has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
graphicsData	BodyGraphicsData			Structure contains data for body visualization; data is defined in special list / dictionary structure

**Detailed information on ObjectMass1D:**

input parameter	symbol	description see tables above
physicsMass	$m$	
nodeNumber	$n0$	
referencePosition	${}^0\mathbf{p}_0$	
referenceRotation	${}^{0b}\mathbf{A}_0 \in \mathbb{R}^{3 \times 3}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{\text{config}}$	global position vector; for interpretation see intermediate variables
Displacement	${}^0\mathbf{u}_{\text{config}}$	global displacement vector; for interpretation see intermediate variables

Velocity	${}^0\mathbf{v}_{\text{config}}$	global velocity vector; for interpretation see intermediate variables
RotationMatrix	${}^{0b}\mathbf{A}$	vector with 9 components of the rotation matrix (row-major format)
Rotation		vector with 3 components of the Euler angles in xyz-sequence ( $R=R_x*R_y*R_z$ ), recomputed from rotation matrix ${}^{0b}\mathbf{A}$
AngularVelocity	${}^0\boldsymbol{\omega}_{\text{config}}$	angular velocity of body
AngularVelocityLocal	${}^b\boldsymbol{\omega}_{\text{config}}$	local (body-fixed) 3D velocity vector of node

#### Description of Item:

intermediate variables	symbol	description
position coordinate	$p_{0\text{config}} = c_{0\text{config}} + c_{0\text{ref}}$	position coordinate of node (nodal coordinate $c_0$ ) in any configuration
displacement coordinate	$u_{0\text{config}} = c_{0\text{config}}$	displacement coordinate of mass node in any configuration
velocity coordinate	$u_{0\text{config}}$	velocity coordinate of mass node in any configuration
Position	${}^0\mathbf{p}_{\text{config}} = {}^0\mathbf{p}_0 + {}^{0b}\mathbf{A}_0 \begin{bmatrix} p_0 \\ 0 \\ 0 \end{bmatrix}_{\text{config}}$	(translational) position of mass object in any configuration
Displacement	${}^0\mathbf{u}_{\text{config}} = {}^{0b}\mathbf{A}_0 \begin{bmatrix} q_0 \\ 0 \\ 0 \end{bmatrix}_{\text{config}}$	(translational) displacement of mass object in any configuration
Velocity	${}^0\mathbf{v}_{\text{config}} = {}^{0b}\mathbf{A}_0 \begin{bmatrix} \dot{q}_0 \\ 0 \\ 0 \end{bmatrix}_{\text{config}}$	(translational) velocity of mass object in any configuration
residual force	$f$	residual of all forces on mass object
applied force	${}^0\mathbf{f}_a = [f_0, f_1, f_2]^T$	3D applied force (loads, connectors, joint reaction forces, ...)
applied torque	${}^0\boldsymbol{\tau}_a = [\tau_0, \tau_1, \tau_2]^T$	3D applied torque (loads, connectors, joint reaction forces, ...)

A rigid body marker (e.g., MarkerBodyRigid) may be attached to this object and forces/torques can be applied. However, torques will have no effect and forces will only have effect in 'direction' of the coordinate.

#### Equations of motion:

$$m \cdot \ddot{q}_0 = f. \quad (6.29)$$

Note that  $f$  is computed from all connectors and loads upon the object. E.g., a 3D force vector  ${}^0\mathbf{f}_a$  is transformed to  $f$  as

$$f = {}^b[1, 0, 0] {}^{b0}\mathbf{A}_0 {}^0\mathbf{f}_a \quad (6.30)$$

Thus, the **position jacobian** reads

$$\mathbf{J}_{pos} = \partial \mathbf{p}_{\text{cur}} / \partial q_{0\text{cur}} = {}^b[1, 0, 0] {}^{b0}\mathbf{A}_0 \quad (6.31)$$

**Example** for ObjectMass1D:

---

```
node = mbs.AddNode(Node1D(referenceCoordinates = [1],
                           initialCoordinates=[0.5],
                           initialVelocities=[0.5]))
mass = mbs.AddObject(Mass1D(nodeNumber = node, physicsMass=1))

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result, get current mass position at local position [0,0,0]
testError = mbs.GetObjectOutputBody(mass, exu.OutputVariableType.Position, [0,0,0])
            [0] - 2
#final x-coordinate of position shall be 2
```

---



### 6.3.4 ObjectRotationalMass1D

A 1D rotational inertia (mass) which is attached to Node1D.

**Additional information for ObjectRotationalMass1D:**

- The Object has the following types = Body, SingleNoded
- Requested node type = GenericODE2
- **Short name** for Python = **Rotor1D**
- **Short name** for Python (visualization object) = **VRotor1D**

The item **ObjectRotationalMass1D** with type = 'RotationalMass1D' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
physicsInertia	UReal		0.	inertia components [SI:kgm <sup>2</sup> ] of rotor / rotational mass
nodeNumber	NodeIndex		MAXINT	node number (type NodeIndex) of Node1D, providing rotation coordinate $\psi_0 = c_0$
referencePosition	Vector3D	3	[0.,0.,0.]	a constant reference position, used to assign joint constraints accordingly and for drawing
referenceRotation	Matrix3D		[[1,0,0], [0,1,0], [0,0,1]]	an intermediate rotation matrix, which transforms the 1D coordinate into 3D, see description
visualization	VObjectRotationalMass1D			parameters for visualization of item

The item VObjectRotationalMass1D has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
graphicsData	BodyGraphicsData			Structure contains data for body visualization; data is defined in special list / dictionary structure

**Detailed information on ObjectRotationalMass1D:**

input parameter	symbol	description see tables above
physicsInertia	$J$	
nodeNumber	$n0$	
referencePosition	${}^0\mathbf{p}_0$	
referenceRotation	${}^{0i}\mathbf{A}_0 \in \mathbb{R}^{3 \times 3}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{\text{config}}$	global position vector; for interpretation see intermediate variables

Displacement	${}^0\mathbf{u}_{\text{config}}$	global displacement vector; for interpretation see intermediate variables
Velocity	${}^0\mathbf{v}_{\text{config}}$	global velocity vector; for interpretation see intermediate variables
RotationMatrix	${}^{0b}\mathbf{A}$	vector with 9 components of the rotation matrix (row-major format)
Rotation		vector with 3 components of the Euler angles in xyz-sequence ( $R=R_x*R_y*R_z$ ), recomputed from rotation matrix ${}^{0b}\mathbf{A}$
AngularVelocity	${}^0\boldsymbol{\omega}_{\text{config}}$	angular velocity of body
AngularVelocityLocal	${}^b\boldsymbol{\omega}_{\text{config}}$	local (body-fixed) 3D velocity vector of node

#### Description of Item:

intermediate variables	symbol	description
position coordinate	$\theta_{0\text{config}} = c_{0\text{config}} + c_{0\text{ref}}$	total rotation coordinate of node (e.g., Node1D) in any configuration (nodal coordinate $c_0$ )
displacement coordinate	$\psi_{0\text{config}} = c_{0\text{config}}$	change of rotation coordinate of mass node (e.g., Node1D) in any configuration (nodal coordinate $c_0$ )
velocity coordinate	$\dot{\psi}_{0\text{config}}$	rotation velocity coordinate of mass node (e.g., Node1D) in any configuration
Position	${}^0\mathbf{p}_{\text{config}} = {}^0\mathbf{p}_0$	constant (translational) position of mass object in any configuration
Displacement	${}^0\mathbf{u}_{\text{config}} = [0, 0, 0]^T$	(translational) displacement of mass object in any configuration
Velocity	${}^0\mathbf{v}_{\text{config}} = [0, 0, 0]^T$	(translational) velocity of mass object in any configuration
AngularVelocity	${}^0\boldsymbol{\omega}_{\text{config}} = {}^{0i}\mathbf{A}_0 \begin{bmatrix} 0 \\ 0 \\ \dot{\psi}_0 \end{bmatrix}^T$	
AngularVelocityLocal	${}^b\boldsymbol{\omega}_{\text{config}} = \begin{bmatrix} 0 \\ 0 \\ \dot{\psi}_0 \end{bmatrix}^T$	
RotationMatrix	${}^{0b}\mathbf{A} = {}^{0i}\mathbf{A}_0 \begin{bmatrix} \cos(\theta_0) & -\sin(\theta_0) & 0 \\ \sin(\theta_0) & \cos(\theta_0) & 0 \\ 0 & 0 & 1 \end{bmatrix}$	transformation of local body ( $b$ ) coordinates to global ( $0$ ) coordinates
residual force	$\boldsymbol{\tau}$	residual of all forces on mass object
applied force	${}^0\mathbf{f}_a = [f_0, f_1, f_2]^T$	3D applied force (loads, connectors, joint reaction forces, ...)
applied torque	${}^0\boldsymbol{\tau}_a = [\tau_0, \tau_1, \tau_2]^T$	3D applied torque (loads, connectors, joint reaction forces, ...)

A rigid body marker (e.g., MarkerBodyRigid) may be attached to this object and forces/torques can be applied. However, forces will have no effect and torques will only have effect in 'direction' of the coordinate.

**Equations of motion:**

$$J \cdot \ddot{\psi}_0 = \tau. \quad (6.32)$$

Note that  $\tau$  is computed from all connectors and loads upon the object. E.g., a 3D torque vector  ${}^0\tau_a$  is transformed to  $\tau$  as

$$\tau = {}^b[0, 0, 1] {}^{b0}\mathbf{A}_0 {}^0\tau_a \quad (6.33)$$

Thus, the **rotation jacobian** reads

$$\mathbf{J}_{rot} = \partial \omega_{cur} / \partial \dot{q}_{0,cur} = {}^b[0, 0, 1] {}^{b0}\mathbf{A}_0 \quad (6.34)$$

**Example** for ObjectRotationalMass1D:

---

```

node = mbs.AddNode(Node1D(referenceCoordinates = [1], #\psi_0ref
                           initialCoordinates=[0.5],   #\psi_0ini
                           initialVelocities=[0.5]))    #\psi_t0ini
rotor = mbs.AddObject(Rotor1D(nodeNumber = node, physicsInertia=1))

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result
#check result, get current rotor z-rotation at local position [0,0,0]
testError = mbs.GetObjectOutputBody(rotor, exu.OutputVariableType.Rotation, [0,0,0])
- 2
#final z-angle of rotor shall be 2

```

---

### 6.3.5 ObjectRigidBody

A 3D rigid body which is attached to a 3D rigid body node. Equations of motion with the displacements  $[u_x \ u_y \ u_z]^T$  of the center of mass and the rotation parameters (Euler parameters)  $\mathbf{q}$ , the mass  $m$ , inertia  $\mathbf{J} = [J_{xx}, J_{xy}, J_{xz}, J_{yx}, J_{yy}, J_{yz}, J_{zx}, J_{zy}, J_{zz}]$  and the residual of all forces and moments  $[R_x \ R_y \ R_z \ R_{q0} \ R_{q1} \ R_{q2} \ R_{q3}]^T$  are given as ...; REMARK: Use the class RigidBodyInertia and AddRigidBody(...) of exudynRigidBodyUtilities.py to handle inertia, COM and mass.

#### Additional information for ObjectRigidBody:

- The Object has the following types = Body, SingleNoded
- Requested node type = Position + Orientation + RigidBody
- **Short name** for Python = **RigidBody**
- **Short name** for Python (visualization object) = **VRigidBody**

The item **ObjectRigidBody** with type = 'RigidBody' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
physicsMass	UReal		0.	mass [SI:kg] of rigid body
physicsInertia	Vector6D		[0.,0.,0., 0.,0.,0.]	inertia components [SI:kgm <sup>2</sup> ]: $[J_{xx}, J_{yy}, J_{zz}, J_{yz}, J_{xz}, J_{xy}]$ of rigid body w.r.t. to the reference point of the body, NOT w.r.t. to center of mass; use the class RigidBodyInertia and AddRigidBody(...) of exudynRigidBodyUtilities.py to handle inertia, COM and mass
physicsCenterOfMass	Vector3D	3	[0.,0.,0.]	local position of center of mass (COM); if the vector of the COM is [0,0,0], the computation will not consider additional terms for the COM and it is faster
nodeNumber	NodeIndex		MAXINT	node number (type NodeIndex) for rigid body node
visualization	VObjectRigidBody			parameters for visualization of item

The item VObjectRigidBody has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
graphicsDataUserFunction	PyFunctionGraphicsData		0	A python function which returns a body-GraphicsData object, which is a list of graphics data in a dictionary computed by the user function; the graphics elements need to be defined in the local body coordinates and are transformed by mbs to global coordinates
graphicsData	BodyGraphicsData			Structure contains data for body visualization; data is defined in special list / dictionary structure

### Detailed information on ObjectRigidBody:

input parameter	symbol	description see tables above
physicsMass	$m$	
physicsInertia	$J$	
physicsCenterOfMass	${}^b\mathbf{p}_{COM}$	
nodeNumber	$n0$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{config}({}^b\mathbf{p}) = {}^0\mathbf{u}_{config} + {}^0\mathbf{p}_{ref} + {}^{0b}\mathbf{A} {}^b\mathbf{p}$	global position vector of body-fixed point given by local position vector
Displacement	${}^0\mathbf{u}_{config} + {}^{0b}\mathbf{A} {}^b\mathbf{p}$	global displacement vector of body-fixed point given by local position vector
Velocity	${}^0\mathbf{v}_{config}({}^b\mathbf{p}) = {}^0\dot{\mathbf{u}}_{config} + {}^{0b}\mathbf{A} ({}^b\boldsymbol{\omega} \times {}^b\mathbf{p}_{config})$	global velocity vector of body-fixed point given by local position vector
RotationMatrix		vector with 9 components of the rotation matrix (row-major format)
Rotation		vector with 3 components of the Euler angles in xyz-sequence (R=Rx*Ry*Rz), recomputed from rotation matrix
AngularVelocity	${}^0\boldsymbol{\omega}_{config}$	angular velocity of body
AngularVelocityLocal	${}^b\boldsymbol{\omega}_{config}$	local (body-fixed) 3D velocity vector of node
Acceleration	${}^0\mathbf{a}_{config}({}^b\mathbf{p}) = {}^0\ddot{\mathbf{u}} + {}^0\boldsymbol{\alpha} \times ({}^{0b}\mathbf{A} {}^b\mathbf{p}) + {}^0\boldsymbol{\omega} \times ({}^0\boldsymbol{\omega} \times ({}^{0b}\mathbf{A} {}^b\mathbf{p}))$	global acceleration vector of body-fixed point given by local position vector
AngularAcceleration	${}^0\boldsymbol{\alpha}_{config}$	angular acceleration vector of body

**Description of Item:** equations coming soon!

### 6.3.6 ObjectRigidBody2D

A 2D rigid body which is attached to a rigid body 2D node. The body obtains coordinates, position, velocity, etc. from the underlying 2D node

**Additional information for ObjectRigidBody2D:**

- The Object has the following types = Body, SingleNoded
- Requested node type = Position2D + Orientation2D + Position + Orientation
- **Short name** for Python = **RigidBody2D**
- **Short name** for Python (visualization object) = **VRigidBody2D**

The item **ObjectRigidBody2D** with type = 'RigidBody2D' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
physicsMass	UReal		0.	mass [SI:kg] of rigid body
physicsInertia	UReal		0.	inertia [SI:kgm <sup>2</sup> ] of rigid body w.r.t. center of mass
nodeNumber	NodeIndex		MAXINT	node number (type NodeIndex) for 2D rigid body node
visualization	VObjectRigidBody2D			parameters for visualization of item

The item VObjectRigidBody2D has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
graphicsDataUserFunction	PyFunctionGraphicsData		0	A python function which returns a body-GraphicsData object, which is a list of graphics data in a dictionary computed by the user function; the graphics elements need to be defined in the local body coordinates and are transformed by mbs to global coordinates
graphicsData	BodyGraphicsData			Structure contains data for body visualization; data is defined in special list / dictionary structure

**Detailed information on ObjectRigidBody2D:**

input parameter	symbol	description see tables above
physicsMass	$m$	
physicsInertia	$J$	
nodeNumber	$n_0$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
------------------	--------	-------------

Position	${}^0\mathbf{p}_{\text{config}}({}^b\mathbf{p}) = {}^0\mathbf{u}_{\text{config}} + {}^0\mathbf{p}_{\text{ref}} + {}^{0b}\mathbf{A} {}^b\mathbf{p}$	global position vector of body-fixed point given by local position vector
Displacement	${}^0\mathbf{u}_{\text{config}} + {}^{0b}\mathbf{A} {}^b\mathbf{p}$	global displacement vector of body-fixed point given by local position vector
Velocity	${}^0\mathbf{v}_{\text{config}}({}^b\mathbf{p}) = {}^0\dot{\mathbf{u}}_{\text{config}} + {}^{0b}\mathbf{A} ({}^b\boldsymbol{\omega} \times {}^b\mathbf{p}_{\text{config}})$	global velocity vector of body-fixed point given by local position vector
Rotation	$\theta_{0\text{config}}$	scalar rotation angle of body
AngularVelocity	${}^0\boldsymbol{\omega}_{\text{config}}$	angular velocity vector of body
RotationMatrix	$\text{vec}({}^{0b}\mathbf{A}) = [A_{00}, A_{01}, A_{02}, A_{10}, \dots, A_{21}, A_{22}]^T_{\text{config}}$	rotation matrix in vector form (stored in row-major order)
Acceleration	${}^0\mathbf{a}_{\text{config}}({}^b\mathbf{p}) = {}^0\ddot{\mathbf{u}} + {}^0\boldsymbol{\alpha} \times ({}^{0b}\mathbf{A} {}^b\mathbf{p}) + {}^0\boldsymbol{\omega} \times ({}^0\boldsymbol{\omega} \times ({}^{0b}\mathbf{A} {}^b\mathbf{p}))$	global acceleration vector of body-fixed point given by local position vector
AngularAcceleration	${}^0\boldsymbol{\alpha}_{\text{config}}$	angular acceleration vector of body

### Description of Item:

intermediate variables	symbol	description
COM position	${}^0\mathbf{p}_{\text{config}} = {}^0\mathbf{p}(n_0)_{\text{config}}$	position of center of mass (COM) which is provided by node $n_0$ in any configuration
COM displacement	${}^0\mathbf{u}_{\text{config}} = [\dot{q}_0, \dot{q}_1, 0]^T_{\text{config}} = {}^0\mathbf{u}(n_0)_{\text{config}}$	displacement of center of mass which is provided by node $n_0$ in any configuration
COM velocity	${}^0\mathbf{v}_{\text{config}} = [\dot{q}_0, \dot{q}_1, 0]^T_{\text{config}} = {}^0\mathbf{v}(n_0)_{\text{config}}$	velocity of center of mass which is provided by node $n_0$ in any configuration
body rotation	${}^0\theta_{0\text{config}} = \theta_0(n_0)_{\text{config}} = \psi_0(n_0)_{\text{ref}} + \psi_0(n_0)_{\text{config}}$	rotation of body as provided by node $n_0$ in any configuration
body rotation matrix	${}^{0b}\mathbf{A}_{\text{config}} = {}^{0b}\mathbf{A}(n_0)_{\text{config}}$	rotation matrix which transforms local to global coordinates as given by node
local position	${}^b\mathbf{p} = [{}^bp_0, {}^bp_1, 0]^T$	local position as used by markers or sensors
body angular velocity	${}^0\boldsymbol{\omega}_{\text{config}} = {}^0[\omega_0(n_0), 0, 0]^T_{\text{config}}$	rotation of body as provided by node $n_0$ in any configuration
(generalized) coordinates	$\mathbf{c}_{\text{config}} = [q_0, q_1, \psi_0]^T$	generalized coordinates of body (= coordinates of node)
generalized forces	${}^0\mathbf{f} = [f_0, f_1, \tau_2]^T$	generalized forces applied to body
applied forces	${}^0\mathbf{f}_a = [f_0, f_1, 0]^T$	applied forces (loads, connectors, joint reaction forces, ...)
applied torques	${}^0\boldsymbol{\tau}_a = [0, 0, \tau_2]^T$	applied torques (loads, connectors, joint reaction forces, ...)

### Equations of motion:

$$\begin{bmatrix} m & 0 & 0 \\ 0 & m & 0 \\ 0 & 0 & J \end{bmatrix} \begin{bmatrix} \ddot{q}_0 \\ \ddot{q}_1 \\ \ddot{\psi}_0 \end{bmatrix} = \begin{bmatrix} f_0 \\ f_1 \\ \tau_2 \end{bmatrix} = \mathbf{f}. \quad (6.35)$$

For example, a LoadCoordinate on coordinate 2 of the node would add a torque  $\tau_2$  on the RHS.

Position-based markers can measure position  $\mathbf{p}_{\text{config}}$  depending on the local position  ${}^b\mathbf{p}$ . The **position jacobian** depends on the local position  ${}^b\mathbf{p}$  and is defined as,

$$\mathbf{J}_{\text{pos}} = \partial\mathbf{p}_{\text{cur}}/\partial\mathbf{c}_{\text{cur}} = \begin{bmatrix} 1 & 0 & -\sin(\theta) {}^bp_0 - \cos(\theta) {}^bp_1 \\ 0 & 1 & \cos(\theta) {}^bp_0 - \sin(\theta) {}^bp_1 \\ 0 & 0 & 0 \end{bmatrix} \quad (6.36)$$

which transforms the action of global forces  ${}^0\mathbf{f}$  of position-based markers on the coordinates  $\mathbf{c}$ ,

$$\mathbf{Q} = \mathbf{J}_{pos} {}^0\mathbf{f}_a \quad (6.37)$$

The **rotation jacobian**

$$\mathbf{J}_{rot} = \partial \mathbf{p}_{cur} / \partial \mathbf{c}_{cur} = \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 1 \end{bmatrix} \quad (6.38)$$

transforms the action of global torques  ${}^0\boldsymbol{\tau}$  of orientation-based markers on the coordinates  $\mathbf{c}$ ,

$$\mathbf{Q} = \mathbf{J}_{rot} {}^0\boldsymbol{\tau}_a \quad (6.39)$$

**Example** for ObjectRigidBody2D:

---

```
node = mbs.AddNode(NodeRigidBody2D(referenceCoordinates = [1,1,0.25*np.pi],
                                initialCoordinates=[0.5,0,0],
                                initialVelocities=[0.5,0,0.75*np.pi]))
mbs.AddObject(RigidBody2D(nodeNumber = node, physicsMass=1, physicsInertia=2))

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result
testError = mbs.GetNodeOutput(node, exu.OutputVariableType.Position)[0] - 2
testError+= mbs.GetNodeOutput(node, exu.OutputVariableType.Coordinates)[2] - 0.75*np.
    pi
#final x-coordinate of position shall be 2, angle theta shall be np.pi
```

---



### 6.3.7 ObjectGenericODE2

A system of  $n$  second order ordinary differential equations (ODE2), having a mass matrix, damping/gyroscopic matrix, stiffness matrix and generalized forces. It can combine generic nodes, or node points. User functions can be used to compute mass matrix and generalized forces depending on given coordinates. NOTE that all matrices, vectors, etc. must have the same dimensions  $n$  or  $(n \times n)$ , or they must be empty  $(0 \times 0)$ , except for the mass matrix which always needs to have dimensions  $(n \times n)$ .

#### Additional information for ObjectGenericODE2:

- The Object has the following types = Body, MultiNoded, SuperElement
- Requested node type: read detailed information of item

The item **ObjectGenericODE2** with type = 'GenericODE2' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
nodeNumbers	ArrayNodeIndex		[]	node numbers which provide the coordinates for the object (consecutively as provided in this list)
massMatrix	NumpyMatrix		Matrix[]	mass matrix of object in python numpy format
stiffnessMatrix	NumpyMatrix		Matrix[]	stiffness matrix of object in python numpy format
dampingMatrix	NumpyMatrix		Matrix[]	damping matrix of object in python numpy format
forceVector	NumpyVector		[]	generalized force vector added to RHS
forceUserFunction	PyFunctionVectorScalar2Vector		0	A python user function which computes the generalized user force vector for the ODE2 equations; The function takes the time, coordinates $q$ (without reference values) and coordinate velocities $\dot{q}$ ; Example for python function with numpy stiffness matrix $K$ : <code>def f(t, q, q_t): return np.dot(K, q)</code>
massMatrixUserFunction	PyFunctionMatrixScalar2Vector		0	A python user function which computes the mass matrix instead of the constant mass matrix; The function takes the time, coordinates $q$ (without reference values) and coordinate velocities $\dot{q}$ ; Example (academic) for python function with numpy stiffness matrix $M$ : <code>def f(t, q, q_t): return (q[0]+1)*M</code>
coordinateIndexPerNode	ArrayIndex		[]	this list contains the local coordinate index for every node, which is needed, e.g., for markers; the list is generated automatically every time parameters have been changed
visualization	VObjectGenericODE2			parameters for visualization of item

The item **VObjectGenericODE2** has the following parameters:

Name	type	size	default value	description
------	------	------	---------------	-------------

show	bool		True	set true, if item is shown in visualization and false if it is not shown
color	Float4	4	[-1.,-1.,-1.,-1.]	RGBA color for object; 4th value is alpha-transparency; R=-1.f means, that default color is used
triangleMesh	NumpyMatrixI		MatrixI[]	a matrix, containg node number triples in every row, referring to the node numbers of the GenericODE2 object; the mesh uses the nodes to visualize the underlying object; contour plot colors are still computed in the local frame!
showNodes	bool		False	set true, nodes are drawn uniquely via the mesh, eventually using the floating reference frame, even in the visualization of the node is show=False; node numbers are shown with indicator 'NF'
graphicsDataUserFunction	PyFunctionGraphicsData		0	A python function which returns a body-GraphicsData object, which is a list of graphics data in a dictionary computed by the user function; the graphics data is draw in global coordinates; it can be used to implement user element visualization, e.g., beam elements or simple mechanical systems; note that this user function may significantly slow down visualization

#### Detailed information on ObjectGenericODE2:

input parameter	symbol	description see tables above
nodeNumbers	$\mathbf{n}_n = [n_0, \dots, n_n]^T$	
massMatrix	$\mathbf{M} \in \mathbb{R}^{n \times n}$	
stiffnessMatrix	$\mathbf{K} \in \mathbb{R}^{n \times n}$	
dampingMatrix	$\mathbf{D} \in \mathbb{R}^{n \times n}$	
forceVector	$\mathbf{f} \in \mathbb{R}^n$	
forceUserFunction	$\mathbf{f}_{user} \in \mathbb{R}^n$	
massMatrixUserFunction	$\mathbf{M}_{user} \in \mathbb{R}^{n \times n}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Coordinates		all ODE2 coordinates
Coordinates_t		all ODE2 velocity coordinates
Force		generalized forces for all coordinates (residual of all forces except mass*accleration; corresponds to ComputeODE2LHS)

**Description of Item:** An object with node numbers  $[n_0, \dots, n_n]$  and according numbers of nodal coordinates  $[n_{c_0}, \dots, n_{c_n}]$ , the total number of equations (=coordinates) of the object is

$$n = \sum_i n_{c_i}. \quad (6.40)$$

Equations of motion:

$$\mathbf{M}\ddot{\mathbf{q}} + \mathbf{D}\dot{\mathbf{q}} + \mathbf{K}\mathbf{q} = \mathbf{f} + \mathbf{f}_{user}(t, \mathbf{q}, \dot{\mathbf{q}}) \quad (6.41)$$

Note that the user function  $\mathbf{f}_{user}(t, \mathbf{q}, \dot{\mathbf{q}})$  may be empty (=0).

In case that a user mass matrix is specified, Eq. (6.41) is replaced with

$$\mathbf{M}_{user}(t, \mathbf{q}, \dot{\mathbf{q}})\ddot{\mathbf{q}} + \mathbf{D}\dot{\mathbf{q}} + \mathbf{K}\mathbf{q} = \mathbf{f} + \mathbf{f}_{user}(t, \mathbf{q}, \dot{\mathbf{q}}) \quad (6.42)$$

CoordinateLoads are integrated for each ODE2 coordinate on the RHS of the latter equation. **Example** for ObjectGenericODE2:

---

```
#set up a mechanical system with two nodes; it has the structure: |~~M0~~M1
nMass0 = mbs.AddNode(NodePoint(referenceCoordinates=[0,0,0]))
nMass1 = mbs.AddNode(NodePoint(referenceCoordinates=[1,0,0]))

mass = 0.5 * np.eye(3)      #mass of nodes
stif = 5000 * np.eye(3)     #stiffness of nodes
damp = 50 * np.eye(3)       #damping of nodes
Z = 0. * np.eye(3)          #matrix with zeros
#build mass, stiffness and damping matrices (:
M = np.block([[mass,      0.*np.eye(3)],
              [0.*np.eye(3), mass      ] ])
K = np.block([[2*stif, -stif],
              [-stif,  stif] ])
D = np.block([[2*damp, -damp],
              [-damp,  damp] ])

oGenericODE2 = mbs.AddObject(ObjectGenericODE2(nodeNumbers=[nMass0,nMass1],
                                                massMatrix=M,
                                                stiffnessMatrix=K,
                                                dampingMatrix=D))

mNode1 = mbs.AddMarker(MarkerNodePosition(nodeNumber=nMass1))
mbs.AddLoad(Force(markerNumber = mNode1, loadVector = [10, 0, 0])) #static solution
                        =10*(1/5000+1/5000)=0.0004

#assemble and solve system for default parameters
mbs.Assemble()

sims=exu.SimulationSettings()
sims.timeIntegration.generalizedAlpha.spectralRadius=1
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', sims)

#check result at default integration time
testError = mbs.GetNodeOutput(nMass1, exu.OutputVariableType.Position)[0] -
            (1.0039999999354785)
```

---

### 6.3.8 ObjectFFRF

This object is used to represent equations modelled by the floating frame of reference formulation (FFRF). It contains a RigidBodyNode (always node 0) and a list of other nodes representing the finite element nodes used in the FFRF. Note that temporary matrices and vectors are subject of change in future.

**Additional information for ObjectFFRF:**

- The Object has the following types = Body, MultiNoded, SuperElement
- Requested node type: read detailed information of item

The item **ObjectFFRF** with type = 'FFRF' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
nodeNumbers	ArrayNodeIndex		[]	node numbers which provide the coordinates for the object (consecutively as provided in this list); the $(n_f + 1)$ nodes represent the nodes of the FE mesh (except for node 0); the global nodal position needs to be reconstructed from the rigid-body motion of the reference frame
massMatrixFF	PyMatrixContainer		PyMatrixContainer[]	body-fixed and ONLY flexible coordinates part of mass matrix of object given in python numpy format (sparse (CSR) or dense, converted to sparse matrix); internally data is stored in triplet format
stiffnessMatrixFF	PyMatrixContainer		PyMatrixContainer[]	body-fixed and ONLY flexible coordinates part of stiffness matrix of object in python numpy format (sparse (CSR) or dense, converted to sparse matrix); internally data is stored in triplet format
dampingMatrixFF	PyMatrixContainer		PyMatrixContainer[]	body-fixed and ONLY flexible coordinates part of damping matrix of object in python numpy format (sparse (CSR) or dense, converted to sparse matrix); internally data is stored in triplet format
forceVector	NumpyVector		[]	generalized, global force vector added to RHS; the rigid body part $\mathbf{f}_r$ is directly applied to rigid body coordinates while the flexible part $\mathbf{f}_{ff}$ is transformed from global to local coordinates
forceUserFunction	PyFunctionVectorScalar2Vector		0	A python user function which computes the generalized user force vector for the ODE2 equations; The function takes the time, coordinates $\mathbf{q}$ (without reference values) and coordinate velocities $\mathbf{q}_t$ ; Example for python function with numpy matrix $\mathbf{K}$ : def f(t, q, q_t): return np.dot(K, q)

massMatrixUserFunction	PyFunctionMatrix	Scalar2Vector	0	A python user function which computes the TOTAL mass matrix (including reference node) and adds the local constant mass matrix; this function takes the time, coordinates $q$ (without reference values) and coordinate velocities $\dot{q}_t$ ; Example (academic) for python function with numpy matrix $M$ : def f(t, q, $\dot{q}_t$ ): return ( $q[0]+1$ )* $M$
computeFFRFterms	bool		True	flag decides whether the standard FFRF terms are computed; use this flag for user-defined definition of FFRF terms in mass matrix and quadratic velocity vector
coordinateIndexPerNode	ArrayIndex		[]	this list contains the local coordinate index for every node, which is needed, e.g., for markers; the list is generated automatically every time parameters have been changed
objectIsInitialized	bool		False	flag used to correctly initialize all FFRF matrices; as soon as this flag is set false, FFRF matrices and terms are recomputed
physicsMass	UReal		0.	total mass [SI:kg] of FFRF object, auto-computed from mass matrix $M$
physicsInertia	Matrix3D		[[1,0,0], [0,1,0], [0,0,1]]	inertia tensor [SI:kgm <sup>2</sup> ] of rigid body w.r.t. to the reference point of the body, auto-computed from the mass matrix $M_{ff}$
physicsCenterOfMass	Vector3D	3	[0.,0.,0.]	local position of center of mass (COM); auto-computed from mass matrix $M$
PHItM	NumpyMatrix		Matrix[]	projector matrix; may be removed in future
referencePositions	NumpyVector		[]	vector containing the reference positions of all flexible nodes
tempVector	NumpyVector		[]	temporary vector
tempCoordinates	NumpyVector		[]	temporary vector containing coordinates
tempCoordinates_t	NumpyVector		[]	temporary vector containing velocity coordinates
tempRefPosSkew	NumpyMatrix		Matrix[]	temporary matrix with skew symmetric local (deformed) node positions
tempVelSkew	NumpyMatrix		Matrix[]	temporary matrix with skew symmetric local node velocities
visualization	VObjectFFRF			parameters for visualization of item

The item VObjectFFRF has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown; use visualizationSettings.bodies.deformationScaleFactor to draw scaled (local) deformations; the reference frame node is shown with additional letters RF

color	Float4	4	[-1.,-1.,-1.,-1.]	RGBA color for object; 4th value is alpha-transparency; R=-1.f means, that default color is used
triangleMesh	NumpyMatrixI		MatrixI[]	a matrix, containing node number triples in every row, referring to the node numbers of the GenericODE2 object; the mesh uses the nodes to visualize the underlying object; contour plot colors are still computed in the local frame!
showNodes	bool		False	set true, nodes are drawn uniquely via the mesh, eventually using the floating reference frame, even in the visualization of the node is show=False; node numbers are shown with indicator 'NF'

#### Detailed information on ObjectFFRF:

input parameter	symbol	description see tables above
nodeNumbers	$\mathbf{n}_n = [n_0, \dots, n_{n_f}]^T$	
massMatrixFF	$\mathbf{M}_{ff} \in \mathbb{R}^{n_{cf} \times n_{cf}}$	
stiffnessMatrixFF	$\mathbf{K}_{ff} \in \mathbb{R}^{n_{cf} \times n_{cf}}$	
dampingMatrixFF	$\mathbf{D}_{ff} \in \mathbb{R}^{n_{cf} \times n_{cf}}$	
forceVector	$\mathbf{f} \in \mathbb{R}^{n_c}$	
forceUserFunction	$\mathbf{f}_{user} \in \mathbb{R}^{n_c}$	
massMatrixUserFunction	$\mathbf{M}_{user} \in \mathbb{R}^{n_c \times n_c}$	
physicsMass	$m$	
physicsInertia	$\mathbf{J}_r \in \mathbb{R}^{3 \times 3}$	
physicsCenterOfMass	${}^b \mathbf{p}_{COM}$	
PHItTM	$\Phi_t^T \in \mathbb{R}^{n_{cf} \times 3}$	
referencePositions	$\mathbf{x}_f \in \mathbb{R}^{n_f}$	
tempVector	$\mathbf{v}_{temp} \in \mathbb{R}^{n_f}$	
tempCoordinates	$\mathbf{c}_{temp} \in \mathbb{R}^{n_f}$	
tempCoordinates_t	$\dot{\mathbf{c}}_{temp} \in \mathbb{R}^{n_f}$	
tempRefPosSkew	$\tilde{\mathbf{p}}_f \in \mathbb{R}^{n_{cf} \times 3}$	
tempVelSkew	$\dot{\tilde{\mathbf{c}}}_f \in \mathbb{R}^{n_{cf} \times 3}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Coordinates		all ODE2 coordinates
Coordinates_t		all ODE2 velocity coordinates
Force		generalized forces for all coordinates (residual of all forces except mass*acceleration; corresponds to ComputeODE2LHS)

#### Description of Item:

intermediate variables	symbol	description
object coordinates	$\mathbf{c} = [\mathbf{c}_r, \mathbf{q}_f]^T$	object coordinates

rigid body coordinates	$\mathbf{c}_r = [q_0, q_1, q_2, \psi_0, \psi_1, \psi_2, \psi_3]^T$	rigid body coordinates in case of Euler parameters
rotation coordinates	$\boldsymbol{\theta}_{\text{cur}} = [\psi_0, \psi_1, \psi_2, \psi_3]_{\text{ref}}^T + [\psi_0, \psi_1, \psi_2, \psi_3]_{\text{cur}}^T$	rigid body coordinates in case of Euler parameters
flexible coordinates	$\mathbf{q}_f$	flexible, body-fixed coordinates
flexible coordinates transformation matrix	${}^{0b}\mathbf{A}_{bd} = \text{diag}([{}^{0b}\mathbf{A}, \dots, {}^{0b}\mathbf{A}])$	block diagonal transformation matrix, which transforms all flexible coordinates from local to global coordinates

Consider an object with  $n = 1 + n_f$  nodes,  $n_f$  being the 'flexible' nodes. It has node numbers  $[n_0, \dots, n_{n_f}]$  and according numbers of nodal coordinates  $[n_{c_0}, \dots, n_{c_n}]$ . This gives  $n_c$  total nodal coordinates,

$$n_c = \sum_{i=0}^{n_f} n_{c_i}. \quad (6.43)$$

whereof the flexible coordinates are

$$n_{c_f} = \sum_{i=1}^{n_f} n_{c_i}. \quad (6.44)$$

The total number of equations (=coordinates) of the object is  $n_c$ . The first node  $n_0$  represents the rigid body motion of the underlying reference frame with  $n_{c_r} = n_{c_0}$  coordinates (e.g.,  $n_{c_r} = 6$  coordinates for Euler angles and  $n_{c_r} = 7$  coordinates in case of Euler parameters).

**Equations of motion**, in case that `computeFFRFterms = True`:

$$\left( \mathbf{M}_{\text{user}}(t, \mathbf{c}, \dot{\mathbf{c}}) + \begin{bmatrix} \mathbf{M}_{tt} & \mathbf{M}_{tr} & \mathbf{M}_{tf} \\ & \mathbf{M}_{rr} & \mathbf{M}_{rf} \\ \text{sym.} & & \mathbf{M}_{ff} \end{bmatrix} \right) \dot{\mathbf{c}} + \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & \mathbf{D}_{ff} \end{bmatrix} \dot{\mathbf{c}} + \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & \mathbf{K}_{ff} \end{bmatrix} \mathbf{c} = \mathbf{f}_Q(\mathbf{c}, \dot{\mathbf{c}}) + \begin{bmatrix} \mathbf{f}_r \\ {}^{0b}\mathbf{A}_{bd}^T \mathbf{f}_{ff} \end{bmatrix} + \mathbf{f}_{\text{user}}(t, \mathbf{c}, \dot{\mathbf{c}}) \quad (6.45)$$

In case that `computeFFRFterms = False`, the mass terms  $\mathbf{M}_{tt} \dots \mathbf{M}_{ff}$  are zero (not computed) and the quadratic velocity vector  $\mathbf{f}_Q = \mathbf{0}$ . Note that the user functions  $\mathbf{f}_{\text{user}}(t, \mathbf{c}, \dot{\mathbf{c}})$  and  $\mathbf{M}_{\text{user}}(t, \mathbf{c}, \dot{\mathbf{c}})$  may be empty (=0). The detailed equations of motion for this element can be found in [8].

CoordinateLoads are integrated for each ODE2 coordinate on the RHS of the latter equation.

If the rigid body node is using Euler parameters  $\boldsymbol{\theta} = [\theta_0, \theta_1, \theta_2, \theta_3]^T$ , an **additional constraint** (constraint nr. 0) is added automatically for the Euler parameter norm, reading

$$1 - \sum_{i=0}^3 \theta_i^2 = 0. \quad (6.46)$$

In order to suppress the rigid body motion of the mesh nodes, you should apply a `ObjectConnectorCoordinateVector` object with the following constraint equations which impose constraints of a so-called Tisserand frame, giving 3 constraints for the position of the center of mass

$$\Phi_t^T \mathbf{M} \mathbf{c}_f = 0 \quad (6.47)$$

and 3 constraints for the rotation,

$$\tilde{\mathbf{x}}_f^T \mathbf{M} \mathbf{c}_f = 0 \quad (6.48)$$

### 6.3.9 ObjectFFRFReducedOrder

This object is used to represent modally reduced flexible bodies using the floating frame of reference formulation (FFRF) and the component mode synthesis. It contains a RigidBodyNode (always node 0) and a NodeGenericODE2 representing the modal coordinates.

#### Additional information for ObjectFFRFReducedOrder:

- The Object has the following types = Body, MultiNoded, SuperElement
- Requested node type: read detailed information of item
- **Short name** for Python = CMSObject
- **Short name** for Python (visualization object) = VCMSObject

The item **ObjectFFRFReducedOrder** with type = 'FFRFReducedOrder' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
nodeNumbers	ArrayNodeIndex		[]	node numbers of rigid body node and NodeGenericODE2 for modal coordinates; the global nodal position needs to be reconstructed from the rigid-body motion of the reference frame, the modal coordinates and the mode basis
massMatrixReduced	PyMatrixContainer		PyMatrixContainer[]	body-fixed and ONLY flexible coordinates part of reduced mass matrix; provided as MatrixContainer(sparse/dense matrix)
stiffnessMatrixReduced	PyMatrixContainer		PyMatrixContainer[]	body-fixed and ONLY flexible coordinates part of reduced stiffness matrix; provided as MatrixContainer(sparse/dense matrix)
dampingMatrixReduced	PyMatrixContainer		PyMatrixContainer[]	body-fixed and ONLY flexible coordinates part of reduced damping matrix; provided as MatrixContainer(sparse/dense matrix)
forceUserFunction	PyFunctionVector	Scalar2Vector	0	A python user function which computes the generalized user force vector for the ODE2 equations; The function takes the time, coordinates q (without reference values) and coordinate velocities q_t; Example for python function with numpy matrix K: def f(t, q, q_t): return np.dot(K, q)
massMatrixUserFunction	PyFunctionMatrix	Scalar2Vector	0	A python user function which computes the TOTAL mass matrix (including reference node) and adds the local constant mass matrix; this function takes the time, coordinates q (without reference values) and coordinate velocities q_t; Example (academic) for python function with numpy matrix M: def f(t, q, q_t): return (q[0]+1)*M
computeFFRFterms	bool		True	flag decides whether the standard FFRF/CMS terms are computed; use this flag for user-defined definition of FFRF terms in mass matrix and quadratic velocity vector



modeBasis	NumpyMatrix		Matrix[]	mode basis, which transforms reduced coordinates to (full) nodal coordinates, written as a single vector $[u_{x,n_0}, u_{y,n_0}, u_{z,n_0}, \dots, u_{x,n_n}, u_{y,n_n}, u_{z,n_n}]^T$
outputVariableModeBasis	NumpyMatrix		Matrix[]	mode basis, which transforms reduced coordinates to output variables per mode; $s_{OV}$ is the size of the output variable, e.g., 6 for stress modes ( $S_{xx}, \dots, S_{xy}$ )
outputVariableTypeModeBasis	OutputVariableType		OutputVariableType	there must be the output variable type of the outputVariableModeBasis, e.g. exu.OutputVariableType.Stress
referencePositions	NumpyVector		[]	vector containing the reference positions of all flexible nodes, needed for graphics
physicsMass	UReal		0.	total mass [SI:kg] of FFRF object, auto-computed from mass matrix <b>M</b>
physicsInertia	Matrix3D		[[1,0,0], [0,1,0], [0,0,1]]	inertia tensor [SI:kgm <sup>2</sup> ] of rigid body w.r.t. to the reference point of the body, auto-computed from the mass matrix <b>M<sub>ff</sub></b>
physicsCenterOfMass	Vector3D	3	[0.,0.,0.]	local position of center of mass (COM); auto-computed from mass matrix <b>M</b>
PHItTM	NumpyMatrix		Matrix[]	projector matrix; may be removed in future
tempUserFunctionForce	NumpyVector		[]	temporary vector for UF force
tempRefPosSkew	NumpyMatrix		Matrix[]	matrix with skew symmetric local (deformed) node positions
tempVelSkew	NumpyMatrix		Matrix[]	matrix with skew symmetric local node velocities
visualization	VObjectFFRFReducedOrder			parameters for visualization of item

The item VObjectFFRFReducedOrder has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown; use visualizationSettings.bodies.deformationScaleFactor to draw scaled (local) deformations; the reference frame node is shown with additional letters RF
color	Float4	4	[-1.,-1.,-1.,-1.]	RGBA color for object; 4th value is alpha-transparency; R=-1.f means, that default color is used
triangleMesh	NumpyMatrixI		MatrixI[]	a matrix, containing node number triples in every row, referring to the node numbers of the GenericODE2 object; the mesh uses the nodes to visualize the underlying object; contour plot colors are still computed in the local frame!

showNodes	bool		False	set true, nodes are drawn uniquely via the mesh, eventually using the floating reference frame, even in the visualization of the node is show=False; node numbers are shown with indicator 'NF'
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#### Detailed information on ObjectFFRFReducedOrder:

input parameter	symbol	description see tables above
nodeNumbers	$\mathbf{n} = [n_0, n_1]^T$	
massMatrixReduced	$\mathbf{M}_{red} \in \mathbb{R}^{n_{cf} \times n_{cf}}$	
stiffnessMatrixReduced	$\mathbf{K}_{red} \in \mathbb{R}^{n_{cf} \times n_{cf}}$	
dampingMatrixReduced	$\mathbf{D}_{red} \in \mathbb{R}^{n_{cf} \times n_{cf}}$	
forceUserFunction	$\mathbf{f}_{user} \in \mathbb{R}^{n_c}$	
massMatrixUserFunction	$\mathbf{M}_{user} \in \mathbb{R}^{n_c \times n_c}$	
modeBasis	$\boldsymbol{\psi} \in \mathbb{R}^{n_{cf} \times n_m}$	
outputVariableModeBasis	$\boldsymbol{\psi}_{OV} \in \mathbb{R}^{n_m \times (n_m \cdot s_{OV})}$	
referencePositions	${}^b \mathbf{r}_f \in \mathbb{R}^{n_f}$	
physicsMass	$m$	
physicsInertia	$\mathbf{J}_r \in \mathbb{R}^{3 \times 3}$	
physicsCenterOfMass	${}^b \mathbf{p}_{COM}$	
PHItTM	$\Phi_i^T \in \mathbb{R}^{n_{cf} \times 3}$	
tempUserFunctionForce	$\mathbf{v}_{temp} \in \mathbb{R}^{n_c}$	
tempRefPosSkew	$\tilde{\mathbf{p}}_f \in \mathbb{R}^{n_{cf} \times 3}$	
tempVelSkew	$\dot{\tilde{\mathbf{c}}}_f \in \mathbb{R}^{n_{cf} \times 3}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Coordinates		all ODE2 coordinates
Coordinates_t		all ODE2 velocity coordinates
Force		generalized forces for all coordinates (residual of all forces except mass*acceleration; corresponds to ComputeODE2LHS)

#### Description of Item:

, 'Stress': allows to compute linearized, corotational nodal stresses (in mesh nodes, in body frame) based on modal stress values provided in outputVariableModeBasis; the flag outputVariableTypeModeBasis must be set in this case to exu.Outputvariable.Stress', 'Strain': allows to compute linearized, corotational nodal strains (in mesh nodes, in body frame) based on modal strain values provided in outputVariableModeBasis; the flag outputVariableTypeModeBasis must be set in this case to exu.Outputvariable.Strain' The object additionally provides the following output variables for mesh nodes (use textttmbs.GetObjectOutputSuperElement(...) or SensorSuperElement):

mesh node output variables	symbol	description
Position	${}^0 \mathbf{r}_{n_i}$	position of node with mesh node number $n_i$ in global coordinates

Position	${}^0\mathbf{r}_{n_i}$	position of node with mesh node number $n_i$ in global coordinates
DisplacementLocal (mesh node $i$ )	${}^b\mathbf{u}_{f,i}$	local nodal mesh displacement in reference (body) frame
VelocityLocal (mesh node $i$ )	${}^b\dot{\mathbf{u}}_{f,i}$	local nodal mesh velocity in reference (body) frame
Displacement (mesh node $i$ )	${}^0\mathbf{u}_{i,config} = {}^0\mathbf{q}_{t,config} + {}^{0b}\mathbf{A}_{config} {}^b\mathbf{p}_{f,i,config} - ({}^0\mathbf{q}_{t,ref} + {}^{0b}\mathbf{A}_{ref} {}^b\mathbf{r}_{f,i})$	nodal mesh displacement in global coordinates
Position (mesh node $i$ )	${}^0\mathbf{p}_i = {}^0\mathbf{p}_t + {}^{0b}\mathbf{A} {}^b\mathbf{p}_{f,i}$	nodal mesh displacement in global coordinates
Velocity (mesh node $i$ )	${}^0\dot{\mathbf{u}}_i = {}^0\dot{\mathbf{q}}_t + {}^{0b}\mathbf{A} ({}^b\dot{\mathbf{u}}_{f,i} + {}^b\tilde{\omega} {}^b\mathbf{u}_{f,i})$	nodal mesh velocity in global coordinates
Stress (mesh node $i$ )	${}^b\boldsymbol{\sigma}_i = (\boldsymbol{\psi}_{OV}\boldsymbol{\zeta})_{3 \dots 3+i+5}$	linearized stress components of mesh node $i$ in reference frame; $\boldsymbol{\sigma} = [\sigma_{xx}, \sigma_{yy}, \sigma_{zz}, \sigma_{yz}, \sigma_{xz}, \sigma_{xy}]^T$ ; ONLY available, if $\boldsymbol{\psi}_{OV}$ is provided and outputVariableTypeModeBasis==exu.OutputVariableT
Strain (mesh node $i$ )	${}^b\boldsymbol{\varepsilon}_i = (\boldsymbol{\psi}_{OV}\boldsymbol{\zeta})_{3 \dots 3+i+5}$	linearized stress components of mesh node $i$ in reference frame; $\boldsymbol{\sigma} = [\sigma_{xx}, \sigma_{yy}, \sigma_{zz}, \sigma_{yz}, \sigma_{xz}, \sigma_{xy}]^T$ ; ONLY available, if $\boldsymbol{\psi}_{OV}$ is provided and outputVariableTypeModeBasis==exu.OutputVariableT

intermediate variables	symbol	description
flexible coordinates transformation matrix	${}^{0b}\mathbf{A}_{bd} = \text{diag}([{}^{0b}\mathbf{A}, \dots, {}^{0b}\mathbf{A}])$	block diagonal transformation matrix, which transforms all flexible coordinates from local to global coordinates
coordinate vector	$\mathbf{q} = [{}^0\mathbf{q}_t, \boldsymbol{\psi}, \boldsymbol{\zeta}]$	vector of object coordinates; $\mathbf{q}_t$ and $\boldsymbol{\psi}$ are the translation and rotation part of displacements of the reference frame, provided by the rigid body node (node number 0)
reference frame position	${}^0\mathbf{p}_{t,config} = {}^0\mathbf{q}_{t,config} + {}^0\mathbf{q}_{t,ref}$	reference frame position in any configuration except reference
reference frame rotation	$\boldsymbol{\theta}_{config} = \boldsymbol{\theta}_{config} + \boldsymbol{\theta}_{ref}$	reference frame rotation parameters in any configuration except reference
vector of modal coordinates	$\boldsymbol{\zeta} = [\zeta_0, \dots, \zeta_{n_m-1}]^T$	vector of modal coordinates
vector of mesh coordinates	${}^b\mathbf{q}_f = \boldsymbol{\psi}\boldsymbol{\zeta}$	vector of alternating x,y, an z coordinates of local (in body frame) mesh displacements reconstructed from modal coordinates $\boldsymbol{\zeta}$
local mesh displacements	${}^b\mathbf{u}_{f,i} = \begin{bmatrix} {}^b\mathbf{q}_{f,i:3} \\ {}^b\mathbf{q}_{f,i:3+1} \\ {}^b\mathbf{q}_{f,i:3+2} \end{bmatrix}$	nodal mesh displacement in local coordinates (body frame)
local mesh position	${}^b\mathbf{p}_{f,i} = \begin{bmatrix} {}^b\mathbf{q}_{f,i:3} \\ {}^b\mathbf{q}_{f,i:3+1} \\ {}^b\mathbf{q}_{f,i:3+2} \end{bmatrix} + \begin{bmatrix} {}^b\mathbf{r}_{f,i:3} \\ {}^b\mathbf{r}_{f,i:3+1} \\ {}^b\mathbf{r}_{f,i:3+2} \end{bmatrix}$	(deformed) nodal mesh position in local coordinates (body frame)

Some definitions:

- body frame (b) = reference frame
- $n_n$  ... number of mesh nodes

- $n_c = 3 \cdot n_n$  ... number of mesh coordinates
- $n_{rigid}$  ... number of rigid body node coordinates: 6 in case of Euler angles and 7 in case of Euler parameters
- $n_{ODE2} = n_c + n_{rigid}$  ... total number of object coordinates
- $n_m$  ... number of modal coordinates; computed from number of columns in modeBasis
- $\psi$  ... mode basis, containing eigenmodes and static modes
- ${}^b\mathbf{r}_f$  ... node reference coordinates for mesh nodes

**Equations of motion**, in case that `computeFFRFterms = True` (NEEDS TO BE UPDATED FOR FFRF!!!!):

$$\left( \mathbf{M}_{user}(t, \mathbf{c}, \dot{\mathbf{c}}) + \begin{bmatrix} \mathbf{M}_{tt} & \mathbf{M}_{tr} & \mathbf{M}_{tf} \\ & \mathbf{M}_{rr} & \mathbf{M}_{rf} \\ \text{sym.} & & \mathbf{M}_{ff} \end{bmatrix} \right) \ddot{\mathbf{c}} + \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & \mathbf{D}_{ff} \end{bmatrix} \dot{\mathbf{c}} + \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & \mathbf{K}_{ff} \end{bmatrix} \mathbf{c} = \mathbf{f}_Q(\mathbf{c}, \dot{\mathbf{c}}) + \mathbf{f}_{user}(t, \mathbf{c}, \dot{\mathbf{c}}) \quad (6.49)$$

In case that `computeFFRFterms = False`, the mass terms  $\mathbf{M}_{tt} \dots \mathbf{M}_{ff}$  are zero (not computed) and the quadratic velocity vector  $\mathbf{f}_Q = \mathbf{0}$ . Note that the user functions  $\mathbf{f}_{user}(t, \mathbf{c}, \dot{\mathbf{c}})$  and  $\mathbf{M}_{user}(t, \mathbf{c}, \dot{\mathbf{c}})$  may be empty (=0). The detailed equations of motion for this element can be found in [9].

CoordinateLoads are integrated for each ODE2 coordinate on the RHS of the latter equation.

### 6.3.10 ObjectANCFcable2D

A 2D cable finite element using 2 nodes of type NodePoint2DSlope1.

**Additional information for ObjectANCFcable2D:**

- Requested node type = Position2D + Orientation2D + Point2DSlope1 + Position + Orientation
- **Short name** for Python = **Cable2D**
- **Short name** for Python (visualization object) = **VCable2D**

The item **ObjectANCFcable2D** with type = 'ANCFcable2D' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
physicsLength	UReal		0.	reference length $L$ [SI:m] of beam; such that the total volume (e.g. for volume load) gives $\rho AL$
physicsMassPerLength	UReal		0.	mass $\rho A$ [SI:kg/m <sup>2</sup> ] of beam
physicsBendingStiffness	UReal		0.	bending stiffness $EI$ [SI:Nm <sup>2</sup> ] of beam; the bending moment is $m = EI(\kappa - \kappa_0)$ , in which $\kappa$ is the material measure of curvature
physicsAxialStiffness	UReal		0.	axial stiffness $EA$ [SI:N] of beam; the axial force is $f_{ax} = EA(\varepsilon - \varepsilon_0)$ , in which $\varepsilon =  \mathbf{r}'  - 1$ is the axial strain
physicsBendingDamping	UReal		0.	bending damping $d_{EI}$ [SI:Nm <sup>2</sup> /s] of beam; the additional virtual work due to damping is $\delta W_{\dot{\kappa}} = \int_0^L \dot{\kappa} \delta \kappa dx$
physicsAxialDamping	UReal		0.	axial stiffness $d_{EA}$ [SI:N/s] of beam; the additional virtual work due to damping is $\delta W_{\dot{\varepsilon}} = \int_0^L \dot{\varepsilon} \delta \varepsilon dx$
physicsReferenceAxialStrain	UReal		0.	reference axial strain of beam (pre-deformation) $\varepsilon_0$ [SI:1] of beam; without external loading the beam will statically keep the reference axial strain value
physicsReferenceCurvature	UReal		0.	reference curvature of beam (pre-deformation) $\kappa_0$ [SI:1/m] of beam; without external loading the beam will statically keep the reference curvature value
nodeNumbers	NodeIndex2		[MAXINT, MAXINT]	two node numbers ANCF cable element
useReducedOrderIntegration	Bool		False	false: use Gauss order 9 integration for virtual work of axial forces, order 5 for virtual work of bending moments; true: use Gauss order 7 integration for virtual work of axial forces, order 3 for virtual work of bending moments
visualization	VObjectANCFcable2D			parameters for visualization of item

The item VObjectANCFcable2D has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawHeight	float		0.	if beam is drawn with rectangular shape, this is the drawing height
color	Float4		[-1.,-1.,-1.,-1.]	RGBA color of the object; if R=-1, use default color

**Detailed information on ObjectANCF Cable2D:** The following output parameters are available as **Output-VariableType** in sensors and other functions:

output parameter	symbol	description
Position		global position vector of local axis (1) and cross section (2) position
Displacement		global displacement vector of local axis (1) and cross section (2) position
Velocity		global velocity vector of local axis (1) and cross section (2) position
Director1		(axial) slope vector of local axis position
Strain		axial strain (scalar)
Curvature		axial strain (scalar)
Force		(local) section normal force (scalar)
Torque		(local) bending moment (scalar)

**Description of Item:** A 2D cable finite element using 2 nodes of type NodePoint2DSlope1. The element has 8 coordinates and uses cubic polynomials for position interpolation. The Bernoulli-Euler beam is capable of large deformation as it employs the material measure of curvature for the bending.

For a detailed description of this beam element, see Gerstmayr and Irschik [5].

### 6.3.11 ObjectALEANCFCable2D

A 2D cable finite element using 2 nodes of type NodePoint2DSlope1 and a axially moving coordinate of type NodeGenericODE2.

**Additional information for ObjectALEANCFCable2D:**

- Requested node type: read detailed information of item
- **Short name** for Python = **ALECable2D**
- **Short name** for Python (visualization object) = **VALECable2D**

The item **ObjectALEANCFCable2D** with type = 'ALEANCFCable2D' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
physicsLength	UReal		0.	reference length $L$ [SI:m] of beam; such that the total volume (e.g. for volume load) gives $\rho A L$
physicsMassPerLength	UReal		0.	mass $\rho A$ [SI:kg/m <sup>2</sup> ] of beam
physicsMovingMassFactor	UReal		1.	this factor denotes the amount of $\rho A$ which is moving; physicsMovingMassFactor=1 means, that all mass is moving; physicsMovingMassFactor=0 means, that no mass is moving; factor can be used to simulate e.g. pipe conveying fluid, in which $\rho A$ is the mass of the pipe+fluid, while $physicsMovingMassFactor \cdot \rho A$ is the mass per unit length of the fluid
physicsBendingStiffness	UReal		0.	bending stiffness $EI$ [SI:Nm <sup>2</sup> ] of beam; the bending moment is $m = EI(\kappa - \kappa_0)$ , in which $\kappa$ is the material measure of curvature
physicsAxialStiffness	UReal		0.	axial stiffness $EA$ [SI:N] of beam; the axial force is $f_{ax} = EA(\varepsilon - \varepsilon_0)$ , in which $\varepsilon =  \mathbf{r}'  - 1$ is the axial strain
physicsBendingDamping	UReal		0.	bending damping $d_{EI}$ [SI:Nm <sup>2</sup> /s] of beam; the additional virtual work due to damping is $\delta W_{\dot{\kappa}} = \int_0^L \dot{\kappa} \delta \kappa dx$
physicsAxialDamping	UReal		0.	axial stiffness $d_{EA}$ [SI:N/s] of beam; the additional virtual work due to damping is $\delta W_{\dot{\varepsilon}} = \int_0^L \dot{\varepsilon} \delta \varepsilon dx$
physicsReferenceAxialStrain	UReal		0.	reference axial strain of beam (pre-deformation) $\varepsilon_0$ [SI:1] of beam; without external loading the beam will statically keep the reference axial strain value
physicsReferenceCurvature	UReal		0.	reference curvature of beam (pre-deformation) $\kappa_0$ [SI:1/m] of beam; without external loading the beam will statically keep the reference curvature value

physicsUseCouplingTerms	bool		True	true: correct case, where all coupling terms due to moving mass are respected; false: only include constant mass for ALE node coordinate, but deactivate other coupling terms (behaves like ANCF Cable2D then)
nodeNumbers	NodeIndex3		[MAXINT, MAXINT, MAXINT]	two node numbers ANCF cable element, third node=ALE GenericODE2 node
useReducedOrderIntegration	Bool		False	false: use Gauss order 9 integration for virtual work of axial forces, order 5 for virtual work of bending moments; true: use Gauss order 7 integration for virtual work of axial forces, order 3 for virtual work of bending moments
visualization	VObjectALEANCF	Cable2D		parameters for visualization of item

The item VObjectALEANCF Cable2D has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawHeight	float		0.	if beam is drawn with rectangular shape, this is the drawing height
color	Float4		[-1.,-1.,-1.,-1.]	RGBA color of the object; if R=-1, use default color

**Detailed information on ObjectALEANCF Cable2D:** The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position		global position vector of local axis (1) and cross section (2) position
Displacement		global displacement vector of local axis (1) and cross section (2) position
Velocity		global velocity vector of local axis (1) and cross section (2) position
Director1		(axial) slope vector of local axis position
Strain		axial strain (scalar)
Curvature		axial strain (scalar)
Force		(local) section normal force (scalar)
Torque		(local) bending moment (scalar)

**Description of Item:** A 2D cable finite element using 2 nodes of type NodePoint2DSlope1 and an axially moving coordinate of type NodeGenericODE2. The element has 8+1 coordinates and uses cubic polynomials for position interpolation. In addition to ANCF Cable2D the element adds an Eulerian axial velocity by the GenericODE2 coordinate. The parameter physicsMovingMassFactor allows to control the amount of mass, which moves with the Eulerian velocity (e.g., the fluid), and which is not moving (the pipe). A factor of physicsMovingMassFactor=1 gives an axially moving beam.



The Bernoulli-Euler beam is capable of large deformation as it employs the material measure of curvature for the bending. Note that damping (`physicsBendingDamping`, `physicsAxialDamping`) only acts on the non-moving part of the beam, as it is the case for the pipe.

A detailed paper on this element is yet under submission, but a similar formulation can be found in [7].

### 6.3.12 ObjectGround

A ground object behaving like a rigid body, but having no degrees of freedom; used to attach body-connectors without an action. For examples see spring dampers and joints.

**Additional information for ObjectGround:**

- The Object has the following types = Ground, Body

The item **ObjectGround** with type = 'Ground' has the following parameters:

Name	type	size	default value	description
name	String		"	objects's unique name
referencePosition	Vector3D	3	[0.,0.,0.]	reference position for ground object; local position is added on top of reference position for a ground object
visualization	VObjectGround			parameters for visualization of item

The item VObjectGround has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
graphicsDataUserFunction	PyFunctionGraphicsData		0	A python function which returns a body-GraphicsData object, which is a list of graphics data in a dictionary computed by the user function
color	Float4		[-1.,-1.,-1.,-1.]	RGB node color; if R=-1, use default color
graphicsData	BodyGraphicsData			Structure contains data for body visualization; data is defined in special list / dictionary structure

**Detailed information on ObjectGround:** The following output parameters are available as **OutputVariable-Type** in sensors and other functions:

output parameter	symbol	description
Position		global position vector of rotated and translated local position
Displacement		global displacement vector of local position
Velocity		global velocity vector of local position
AngularVelocity		angular velocity of body
RotationMatrix		rotation matrix in vector form (stored in row-major order)

### 6.3.13 ObjectConnectorSpringDamper

An simple spring-damper element with additional force; connects to position-based markers.

**Additional information for ObjectConnectorSpringDamper:**

- The Object has the following types = Connector
- Requested marker type = Position
- **Short name** for Python = **SpringDamper**
- **Short name** for Python (visualization object) = **VSpringDamper**

The item **ObjectConnectorSpringDamper** with type = 'ConnectorSpringDamper' has the following parameters:

Name	type	size	default value	description
name	String		"	connector's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
referenceLength	UReal		0.	reference length [SI:m] of spring
stiffness	UReal		0.	stiffness [SI:N/m] of spring; acts against (length-initialLength)
damping	UReal		0.	damping [SI:N/(m s)] of damper; acts against d/dt(length)
force	UReal		0.	added constant force [SI:N] of spring; scalar force; f=1 is equivalent to reducing initialLength by 1/stiffness; f > 0: tension; f < 0: compression
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
springForceUserFunction	PyFunctionScalar6		0	A python function which defines the spring force with parameters (time, deltaL, deltaL_t, Real stiffness, Real damping, Real springForce); the parameters are provided to the function using the current values of the SpringDamper object; The python function will only be evaluated, if activeConnector is true, otherwise the SpringDamper is inactive; Example for python function: def f(t, u, v, k, d, F0): return k*u + d*v + F0
visualization	VObjectConnectorSpringDamper			parameters for visualization of item

The item **VObjectConnectorSpringDamper** has the following parameters:

Name	type	size	default value	description
show	Bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = diameter of spring; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R=-1, use default color

Detailed information on ObjectConnectorSpringDamper: The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Distance		distance between both points
Displacement		relative displacement between both points
Velocity		relative velocity between both points
Force		spring-damper force

**Description of Item:**

**Definition of quantities:**

input parameter	symbol	description
referenceLength	$L_0$	
stiffness	$k$	
damping	$d$	
force	$f_a$	additional force (e.g., actuator force)
markerNumbers[0]	$m0$	global marker number m0
markerNumbers[1]	$m1$	global marker number m1

intermediate variables	symbol	description
marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker m0
marker m1 position	${}^0\mathbf{p}_{m1}$	
marker m0 velocity	${}^0\mathbf{v}_{m0}$	current global velocity which is provided by marker m0
marker m1 velocity	${}^0\mathbf{v}_{m1}$	

output variables	symbol	formula
Distance	$L$	$ \Delta^0\mathbf{p} $
Displacement	$\Delta^0\mathbf{p}$	${}^0\mathbf{p}_{m1} - {}^0\mathbf{p}_{m0}$
Velocity	$\Delta^0\mathbf{v}$	${}^0\mathbf{v}_{m1} - {}^0\mathbf{v}_{m0}$
Force	$\mathbf{f}$	see below

**Connector equations:** The unit vector in force direction reads (raises SysError if  $L = 0$ ),

$$\mathbf{v}_f = \frac{1}{L} \Delta^0\mathbf{p} \quad (6.50)$$

If `activeConnector = True`, the scalar spring force is computed as

$$f_{SD} = k \cdot (L - L_0) + d \cdot \Delta^0\mathbf{v}^T \mathbf{v}_f + f_a \quad (6.51)$$

If the springForceUserFunction UF is defined,  $\mathbf{f}$  instead becomes ( $t$  is current time)

$$f_{SD} = \text{UF}(t, L - L_0, \Delta^0\mathbf{v}^T \mathbf{v}_f, k, d, f_a) \quad (6.52)$$

if `activeConnector = False`,  $f_{SD}$  is set to zero.: The vector of the spring force applied at both markers finally reads

$$\mathbf{f} = f_{SD} \mathbf{v}_f \quad (6.53)$$

**Example** for ObjectConnectorSpringDamper:

---

```

node = mbs.AddNode(NodePoint(referenceCoordinates = [1.05,0,0]))
oMassPoint = mbs.AddObject(MassPoint(nodeNumber = node, physicsMass=1))

m0 = mbs.AddMarker(MarkerBodyPosition(bodyNumber=oGround, localPosition=[0,0,0]))
m1 = mbs.AddMarker(MarkerBodyPosition(bodyNumber=oMassPoint, localPosition=[0,0,0]))

mbs.AddObject(ObjectConnectorSpringDamper(markerNumbers=[m0,m1],
                                           referenceLength = 1, #shorter than initial
                                           distance
                                           stiffness = 100,
                                           damping = 1))

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result at default integration time
testError = mbs.GetNodeOutput(node, exu.OutputVariableType.Position)[0] -
0.9736596225944887

```

---

### 6.3.14 ObjectConnectorCartesianSpringDamper

An 3D spring-damper element acting accordingly in three (global) directions (x,y,z) which connects to position-based markers.

**Additional information for ObjectConnectorCartesianSpringDamper:**

- The Object has the following types = Connector
- Requested marker type = Position
- **Short name** for Python = **CartesianSpringDamper**
- **Short name** for Python (visualization object) = **VCartesianSpringDamper**

The item **ObjectConnectorCartesianSpringDamper** with type = 'ConnectorCartesianSpringDamper' has the following parameters:

Name	type	size	default value	description
name	String		"	connector's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
stiffness	Vector3D		[0.,0.,0.]	stiffness [SI:N/m] of springs; act against relative displacements in 0, 1, and 2-direction
damping	Vector3D		[0.,0.,0.]	damping [SI:N/(m s)] of dampers; act against relative velocities in 0, 1, and 2-direction
offset	Vector3D		[0.,0.,0.]	offset between two springs
springForceUserFunction	PyFunctionVector3DScalar5Vector3D		0	A python function which computes the 3D force vector between the two marker points, if activeConnector=True; The function takes the relative displacement (3D) vector (m1.position-m0.position, etc.) and the relative velocity vector (3D), the spring striffness vector 3D, damping and offset parameter vectors (3D): f(time, displacement, velocity, stiffness, damping, offset); Example for python function: def f(t, u, v, k, d, offset): return [u[0]*k[0],u[1]*k[1],u[2]*k[2]]
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temorarily) a connector or constraint
visualization	VObjectConnectorCartesianSpringDamper			parameters for visualization of item

The item **VObjectConnectorCartesianSpringDamper** has the following parameters:

Name	type	size	default value	description
show	Bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = diameter of spring; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R== -1, use default color

### Detailed information on ObjectConnectorCartesianSpringDamper:

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	
stiffness	$\mathbf{k}$	
damping	$\mathbf{d}$	
offset	$\mathbf{v}_{\text{off}}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Displacement	$\Delta^0 \mathbf{p} = {}^0 \mathbf{p}_{m1} - {}^0 \mathbf{p}_{m0}$	relative displacement in global coordinates
Distance	$L =  \Delta^0 \mathbf{p} $	scalar distance between both marker points
Velocity	$\Delta^0 \mathbf{v} = {}^0 \mathbf{v}_{m1} - {}^0 \mathbf{v}_{m0}$	relative translational velocity in global coordinates
Force	$\mathbf{f}_{SD}$	joint force in global coordinates, see equations

### Description of Item:

### Definition of quantities:

intermediate variables	symbol	description
marker m0 position	${}^0 \mathbf{p}_{m0}$	current global position which is provided by marker m0
marker m1 position	${}^0 \mathbf{p}_{m1}$	
marker m0 velocity	${}^0 \mathbf{v}_{m0}$	current global velocity which is provided by marker m0
marker m1 velocity	${}^0 \mathbf{v}_{m1}$	

**Connector equations:** Displacement between marker m0 to marker m1 positions,

$$\Delta^0 \mathbf{p} = {}^0 \mathbf{p}_{m1} - {}^0 \mathbf{p}_{m0} \quad (6.54)$$

and relative velocity,

$$\Delta^0 \mathbf{v} = {}^0 \mathbf{v}_{m1} - {}^0 \mathbf{v}_{m0} \quad (6.55)$$

If activeConnector = True, the spring force vector is computed as

$$\mathbf{f}_{SD} = (\mathbf{k} \cdot (\Delta^0 \mathbf{p} - \mathbf{v}_{\text{off}}) + \mathbf{d} \Delta^0 \mathbf{v}) \quad (6.56)$$

If the springForceUserFunction UF is defined,  $\mathbf{f}_{SD}$  instead becomes ( $t$  is current time)

$$\mathbf{f}_{SD} = \text{UF}(t, \Delta^0 \mathbf{p}, \Delta^0 \mathbf{v}, \mathbf{k}, \mathbf{d}, \mathbf{v}_{\text{off}}) \quad (6.57)$$

if activeConnector = False,  $\mathbf{f}_{SD}$  is set to zero. **Example** for ObjectConnectorCartesianSpringDamper:

---

```
#example with mass at [1,1,0], 5kg under load 5N in -y direction
k=5000
nMass = mbs.AddNode(NodePoint(referenceCoordinates=[1,1,0]))
oMass = mbs.AddObject(MassPoint(physicsMass = 5, nodeNumber = nMass))

mMass = mbs.AddMarker(MarkerNodePosition(nodeNumber=nMass))
```

```

mGround = mbs.AddMarker(MarkerBodyPosition(bodyNumber=oGround, localPosition =
    [1,1,0]))
mbs.AddObject(CartesianSpringDamper(markerNumbers = [mGround, mMass],
    stiffness = [k,k,k],
    damping = [0,k*0.05,0], offset = [0,0,0]))
mbs.AddLoad(Force(markerNumber = mMass, loadVector = [0, -5, 0])) #static solution
    =-5/5000=-0.001m

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result at default integration time
testError = mbs.GetNodeOutput(nMass, exu.OutputVariableType.Displacement)[1] -
    (-0.000999999999997058)

```

---



### 6.3.15 ObjectConnectorRigidBodySpringDamper

An 3D spring-damper element acting on relative displacements and relative rotations of two rigid body (position+orientation) markers; connects to (position+orientation)-based markers and represents a penalty-based rigid joint (or prismatic, revolute, etc.)

**Additional information for ObjectConnectorRigidBodySpringDamper:**

- The Object has the following types = Connector
- Requested marker type = Position + Orientation
- **Short name** for Python = **RigidBodySpringDamper**
- **Short name** for Python (visualization object) = **VRigidBodySpringDamper**

The item **ObjectConnectorRigidBodySpringDamper** with type = 'ConnectorRigidBodySpringDamper' has the following parameters:

Name	type	size	default value	description
name	String		"	connector's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
stiffness	Matrix6D		np.zeros([6,6])	stiffness [SI:N/m or Nm/rad] of translational, torsional and coupled springs; act against relative displacements in x, y, and z-direction as well as the relative angles (calculated as Euler angles); in the simplest case, the first 3 diagonal values correspond to the local stiffness in x,y,z direction and the last 3 diagonal values correspond to the rotational stiffness around x,y and z axis
damping	Matrix6D		np.zeros([6,6])	damping [SI:N/(m/s) or Nm/(rad/s)] of translational, torsional and coupled dampers; very similar to stiffness, however, the rotational velocity is computed from the angular velocity vector
rotationMarker0	Matrix3D		[[1,0,0], [0,1,0], [0,0,1]]	local rotation matrix for marker 0; stiffness, damping, etc. components are measured in local coordinates relative to rotationMarker0
rotationMarker1	Matrix3D		[[1,0,0], [0,1,0], [0,0,1]]	local rotation matrix for marker 1; stiffness, damping, etc. components are measured in local coordinates relative to rotationMarker1
offset	Vector6D		[0.,0.,0.,0.,0.,0.]	translational and rotational offset considered in the spring force calculation
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectConnectorRigidBodySpringDamper			parameters for visualization of item

The item **VObjectConnectorRigidBodySpringDamper** has the following parameters:

Name	type	size	default value	description
show	Bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = diameter of spring; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R== -1, use default color

**Detailed information on ObjectConnectorRigidBodySpringDamper: The following output parameters are available as OutputVariableType in sensors and other functions:**

output parameter	symbol	description
DisplacementLocal	${}^{J0}\Delta\mathbf{p}$	relative displacement in local joint0 coordinates
VelocityLocal	${}^{J0}\Delta\mathbf{v}$	relative translational velocity in local joint0 coordinates
Rotation	${}^{J0}\boldsymbol{\theta} = [\theta_0, \theta_1, \theta_2]^T$	relative rotation parameters (Tait Bryan Rxyz); these are the angles used for calculation of joint torques (e.g. if cX is the diagonal rotational stiffness, the moment for axis X reads $mX=cX*\phi X$ , etc.)
AngularVelocityLocal	${}^{J0}\Delta\boldsymbol{\omega}$	relative angular velocity in local joint0 coordinates
ForceLocal	${}^{J0}\mathbf{f}$	joint force in local joint0 coordinates
TorqueLocal	${}^{J0}\mathbf{m}$	joint torque in in local joint0 coordinates

**Description of Item:**

**Definition of quantities:**

input parameter	symbol	description
stiffness	$\mathbf{k} \in \mathbb{R}^{6 \times 6}$	stiffness in $J0$ coordinates
damping	$\mathbf{d} \in \mathbb{R}^{6 \times 6}$	damping in $J0$ coordinates
offset	${}^{J0}\mathbf{v}_{\text{off}} \in \mathbb{R}^6$	offset in $J0$ coordinates
rotationMarker0	${}^{m0,J0}\mathbf{A}$	rotation matrix which transforms from joint 0 into marker 0 coordinates
rotationMarker1	${}^{m1,J1}\mathbf{A}$	rotation matrix which transforms from joint 1 into marker 1 coordinates
markerNumbers[0]	$m0$	global marker number m0
markerNumbers[1]	$m1$	global marker number m1

intermediate variables	symbol	description
marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker m0
marker m0 orientation	${}^{0,m0}\mathbf{A}$	current rotation matrix provided by marker m0
marker m1 position	${}^0\mathbf{p}_{m1}$	accordingly
marker m1 orientation	${}^{0,m1}\mathbf{A}$	current rotation matrix provided by marker m1

marker m0 velocity	${}^0\mathbf{v}_{m0}$	current global velocity which is provided by marker m0
marker m1 velocity	${}^0\mathbf{v}_{m1}$	accordingly
marker m0 velocity	${}^b\boldsymbol{\omega}_{m0}$	current local angular velocity vector provided by marker m0
marker m1 velocity	${}^b\boldsymbol{\omega}_{m1}$	current local angular velocity vector provided by marker m1
Displacement	${}^0\Delta\mathbf{p}$	${}^0\mathbf{p}_{m1} - {}^0\mathbf{p}_{m0}$
Velocity	${}^0\Delta\mathbf{v}$	${}^0\mathbf{v}_{m1} - {}^0\mathbf{v}_{m0}$
DisplacementLocal	${}^{J0}\Delta\mathbf{p}$	$\left({}^{0,m0}\mathbf{A} \quad {}^{m0,J0}\mathbf{A}\right)^T {}^0\Delta\mathbf{p}$
VelocityLocal	${}^{J0}\Delta\mathbf{v}$	$\left({}^{0,m0}\mathbf{A} \quad {}^{m0,J0}\mathbf{A}\right)^T {}^0\Delta\mathbf{v}$
AngularVelocityLocal	${}^{J0}\Delta\boldsymbol{\omega}$	$\left({}^{0,m0}\mathbf{A} \quad {}^{m0,J0}\mathbf{A}\right)^T \left({}^{0,m1}\mathbf{A} \quad {}^{m1}\boldsymbol{\omega} - {}^{0,m0}\mathbf{A} \quad {}^{m0}\boldsymbol{\omega}\right)$

**Connector equations:** If `activeConnector = True`, the vector spring force is computed as

$$\begin{bmatrix} {}^{J0}\mathbf{f}_{SD} \\ {}^{J0}\mathbf{m}_{SD} \end{bmatrix} = \mathbf{k} \left( \begin{bmatrix} {}^{J0}\Delta\mathbf{p} \\ {}^{J0}\boldsymbol{\theta} \end{bmatrix} - {}^{J0}\mathbf{v}_{\text{off}} \right) + \mathbf{d} \begin{bmatrix} {}^{J0}\Delta\mathbf{v} \\ {}^{J0}\Delta\boldsymbol{\omega} \end{bmatrix} \quad (6.58)$$

For the application of joint forces to markers,  $[{}^{J0}\mathbf{f}_{SD}, {}^{J0}\mathbf{m}_{SD}]^T$  is transformed into global coordinates. if `activeConnector = False`,  ${}^{J0}\mathbf{f}_{SD}$  and  ${}^{J0}\mathbf{m}_{SD}$  are set to zero.

### 6.3.16 ObjectConnectorCoordinateSpringDamper

A 1D (scalar) spring-damper element acting on single ODE2 coordinates; connects to coordinate-based markers; NOTE that the coordinate markers only measure the coordinate (=displacement), but the reference position is not included as compared to position-based markers!; the spring-damper can also act on rotational coordinates.

**Additional information for ObjectConnectorCoordinateSpringDamper:**

- The Object has the following types = Connector
- Requested marker type = Coordinate
- **Short name** for Python = **CoordinateSpringDamper**
- **Short name** for Python (visualization object) = **VCoordinateSpringDamper**

The item **ObjectConnectorCoordinateSpringDamper** with type = 'ConnectorCoordinateSpringDamper' has the following parameters:

Name	type	size	default value	description
name	String		"	connector's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
stiffness	Real		0.	stiffness [SI:N/m] of spring; acts against relative value of coordinates
damping	Real		0.	damping [SI:N/(m s)] of damper; acts against relative velocity of coordinates
offset	Real		0.	offset between two coordinates (reference length of springs), see equation
dryFriction	Real		0.	dry friction force [SI:N] against relative velocity; assuming a normal force $f_N$ , the friction force can be interpreted as $f_\mu = \mu f_N$
dryFrictionProportionalZone	Real		0.	limit velocity [m/s] up to which the friction is proportional to velocity (for regularization / avoid numerical oscillations)
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
springForceUserFunction	PyFunctionScalar8		0	A python function which defines the spring force with 8 parameters, see equations section; Example for python function: def f(t, u, v, k, d, offset, frictionForce, frictionProportionalZone): return k*(u-offset) + d*v
visualization	VObjectConnectorCoordinateSpringDamper			parameters for visualization of item

The item **VObjectConnectorCoordinateSpringDamper** has the following parameters:

Name	type	size	default value	description
show	Bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = diameter of spring; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R==-1, use default color

### Detailed information on ObjectConnectorCoordinateSpringDamper:

input parameter	symbol	description see tables above
stiffness	$k$	
damping	$d$	
offset	$l_{\text{off}}$	
dryFriction	$f_{\mu}$	
dryFrictionProportionalZone	$v_{\mu}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Displacement	$\Delta q$	relative scalar displacement of marker coordinates
Velocity	$\Delta v$	difference of scalar marker velocity coordinates
Force	$f_{SD}$	scalar spring force

### Description of Item:

### Definition of quantities:

intermediate variables	symbol	description
marker m0 coordinate	$q_{m0}$	current displacement coordinate which is provided by marker m0; does NOT include reference coordinate!
marker m1 coordinate	$q_{m1}$	
marker m0 velocity coordinate	$v_{m0}$	current velocity coordinate which is provided by marker m0
marker m1 velocity coordinate	$v_{m1}$	

**Connector equations:** Displacement between marker m0 to marker m1 coordinates (does NOT include reference coordinates),

$$\Delta q = q_{m1} - q_{m0} \quad (6.59)$$

and relative velocity,

$$\Delta v = v_{m1} - v_{m0} \quad (6.60)$$

If  $f_{\mu} > 0$ , the friction force is computed as

$$f_{\text{friction}} = \begin{cases} \text{Sgn}(\Delta v) \cdot f_{\mu} & \text{if } |\Delta v| \geq v_{\mu} \\ \frac{\Delta v}{v_{\mu}} f_{\mu} & \text{if } |\Delta v| < v_{\mu} \end{cases} \quad (6.61)$$

If `activeConnector = True`, the scalar spring force vector is computed as

$$f_{SD} = k(\Delta q - l_{\text{off}}) + d \cdot \Delta v + f_{\text{friction}} \quad (6.62)$$

If the springForceUserFunction UF is defined,  $f_{SD}$  instead becomes ( $t$  is current time)

$$f_{SD} = \text{UF}(t, \Delta q, \Delta v, k, d, l_{\text{off}}, f_{\mu}, v_{\mu}) \quad (6.63)$$

if `activeConnector = False`,  $f_{SD}$  is set to zero. **Example** for ObjectConnectorCoordinateSpringDamper:

---

```

def springForce(t, u, v, k, d, offset, frictionForce, frictionProportionalZone):
    return 0.1*k*u+k*u**3+v*d

nMass=mbs.AddNode(Point(referenceCoordinates = [2,0,0]))
massPoint = mbs.AddObject(MassPoint(physicsMass = 5, nodeNumber = nMass))

groundMarker=mbs.AddMarker(MarkerNodeCoordinate(nodeNumber= nGround, coordinate = 0))
nodeMarker =mbs.AddMarker(MarkerNodeCoordinate(nodeNumber= nMass, coordinate = 0))

#Spring-Damper between two marker coordinates
mbs.AddObject(CoordinateSpringDamper(markerNumbers = [groundMarker, nodeMarker],
                                     stiffness = 5000, damping = 80,
                                     springForceUserFunction = springForce))
loadCoord = mbs.AddLoad(LoadCoordinate(markerNumber = nodeMarker, load = 1)) #static
linear solution:0.002

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result at default integration time
testError = mbs.GetNodeOutput(nMass, exu.OutputVariableType.Displacement)[0] -
0.0019995158325691875

```

---

### 6.3.17 ObjectConnectorDistance

Connector which enforces constant or prescribed distance between two bodies/nodes.

**Additional information for ObjectConnectorDistance:**

- The Object has the following types = Connector, Constraint
- Requested marker type = Position
- **Short name** for Python = **DistanceConstraint**
- **Short name** for Python (visualization object) = **VDistanceConstraint**

The item **ObjectConnectorDistance** with type = 'ConnectorDistance' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
distance	UReal		0.	prescribed distance [SI:m] of the used markers
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectConnectorDistance			parameters for visualization of item

The item **VObjectConnectorDistance** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = link size; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R== -1, use default color

**Detailed information on ObjectConnectorDistance:**

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	
distance	$d_0$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Displacement	${}^0\Delta\mathbf{p}$	relative displacement in global coordinates
Velocity	${}^0\Delta\mathbf{v}$	relative translational velocity in global coordinates
Distance	$ {}^0\Delta\mathbf{p} $	distance between markers (should stay constant; shows constraint deviation)
Force	$\lambda_0$	joint force in global coordinates

### Description of Item: Definition of quantities:

intermediate variables	symbol	description
marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker m0
marker m1 position	${}^0\mathbf{p}_{m1}$	accordingly
marker m0 velocity	${}^0\mathbf{v}_{m0}$	current global velocity which is provided by marker m0
marker m1 velocity	${}^0\mathbf{v}_{m1}$	accordingly
relative displacement	${}^0\Delta\mathbf{p}$	${}^0\mathbf{p}_{m1} - {}^0\mathbf{p}_{m0}$
relative velocity	${}^0\Delta\mathbf{v}$	${}^0\mathbf{v}_{m1} - {}^0\mathbf{v}_{m0}$
algebraicVariable	$\lambda_0$	Lagrange multiplier = force in constraint

**Algebraic constraint equations:** If activeConnector = True, the index 3 algebraic equation reads

$$|{}^0\Delta\mathbf{p}| - d_0 = 0 \quad (6.64)$$

The index 2 (velocity level) algebraic equation reads

$$\left( \frac{{}^0\Delta\mathbf{p}}{|{}^0\Delta\mathbf{p}|} \right)^T \Delta\mathbf{v} = 0 \quad (6.65)$$

if activeConnector = False, the algebraic equation reads

$$\lambda_0 = 0 \quad (6.66)$$

### Example for ObjectConnectorDistance:

---

```

#example with 1m pendulum, 50kg under gravity
nMass = mbs.AddNode(NodePoint2D(referenceCoordinates=[1,0]))
oMass = mbs.AddObject(MassPoint2D(physicsMass = 50, nodeNumber = nMass))

mMass = mbs.AddMarker(MarkerNodePosition(nodeNumber=nMass))
mGround = mbs.AddMarker(MarkerBodyPosition(bodyNumber=oGround, localPosition =
    [0,0,0]))
oDistance = mbs.AddObject(DistanceConstraint(markerNumbers = [mGround, mMass],
    distance = 1))

mbs.AddLoad(Force(markerNumber = mMass, loadVector = [0, -50*9.81, 0]))

#assemble and solve system for default parameters
mbs.Assemble()

sims=exu.SimulationSettings()
sims.timeIntegration.generalizedAlpha.spectralRadius=0.7
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', sims)

#check result at default integration time
testError = mbs.GetNodeOutput(nMass, exu.OutputVariableType.Position)[0] -
    (-0.9845225086606828)

```

---



### 6.3.18 ObjectConnectorCoordinate

A coordinate constraint which constrains two (scalar) coordinates of Marker[Node|Body]Coordinates attached to nodes or bodies. The constraint acts directly on coordinates, but does not include reference values, e.g., of nodal values.

**Additional information for ObjectConnectorCoordinate:**

- The Object has the following types = Connector, Constraint
- Requested marker type = Coordinate
- **Short name** for Python = **CoordinateConstraint**
- **Short name** for Python (visualization object) = **VCoordinateConstraint**

The item **ObjectConnectorCoordinate** with type = 'ConnectorCoordinate' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
offset	UReal		0.	An offset between the two values
factorValue1	UReal		1.	An additional factor multiplied with value1 used in algebraic equation
velocityLevel	bool		False	If true: connector constrains velocities (only works for ODE2 coordinates!); offset is used between velocities; in this case, the offsetUserFunction_t is considered and offsetUserFunction is ignored
offsetUserFunction	PyFunctionScalar2		0	A python function which defines the time-dependent offset; it is highly RECOMMENDED to use sufficiently smooth functions, having consistent initial offsets with initial configuration of bodies, zero or compatible initial offset-velocity, and no accelerations; Example for python function: def UF(t, l_offset): return l_offset*(1-np.cos(t*10*2*np.pi))
offsetUserFunction_t	PyFunctionScalar2		0	time derivative of offsetUserFunction; needed for 'velocityLevel=True', or for index2 time integration and for computation of initial accelerations in SecondOrderImplicit integrators
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectConnectorCoordinate			parameters for visualization of item

The item **VObjectConnectorCoordinate** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

drawSize	float		-1.	drawing size = link size; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R== -1, use default color

#### Detailed information on ObjectConnectorCoordinate:

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	
offset	$l_{\text{off}}$	
factorValue1	$k_{m1}$	
offsetUserFunction	$UF(t, l_{\text{off}})$	
offsetUserFunction_t	$UF_t(t, l_{\text{off}})$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Displacement	$\Delta q$	relative scalar displacement of marker coordinates, not including factorValue1
Velocity	$\Delta v$	difference of scalar marker velocity coordinates, not including factorValue1
ConstraintEquation	<b>c</b>	(residuum of) constraint equation
Force	$\lambda_0$	scalar constraint force (Lagrange multiplier)

#### Description of Item: Definition of quantities:

intermediate variables	symbol	description
marker m0 coordinate	$q_{m0}$	current displacement coordinate which is provided by marker m0; does NOT include reference coordinate!
marker m1 coordinate	$q_{m1}$	
marker m0 velocity coordinate	$v_{m0}$	current velocity coordinate which is provided by marker m0
marker m1 velocity coordinate	$v_{m1}$	
difference of coordinates	$\Delta q = q_{m1} - q_{m0}$	Displacement between marker m0 to marker m1 coordinates (does NOT include reference coordinates)
difference of velocity coordinates	$\Delta v = v_{m1} - v_{m0}$	

**Algebraic constraint equations:** If activeConnector = True, the index 3 algebraic equation reads

$$\mathbf{c}(q_{m0}, q_{m1}) = k_{m1} \cdot q_{m1} - q_{m0} - l_{\text{off}} = 0 \quad (6.67)$$

If the offsetUserFunction UF is defined, **c** instead becomes ( $t$  is current time)

$$\mathbf{c}(q_{m0}, q_{m1}) = k_{m1} \cdot q_{m1} - q_{m0} - UF(t, l_{\text{off}}) = 0 \quad (6.68)$$

The activeConnector = True, index 2 (velocity level) algebraic equation reads

$$\dot{\mathbf{c}}(\dot{q}_{m0}, \dot{q}_{m1}) = k_{m1} \cdot \dot{q}_{m1} - \dot{q}_{m0} - d = 0 \quad (6.69)$$

The factor  $d$  in velocity level equations is zero, except if `parameters.velocityLevel = True`, then  $d = l_{\text{off}}$ . If velocity level constraints are active and the velocity level offsetUserFunction\_t  $UF_t$  is defined,  $\dot{c}$  instead becomes  $\dot{c}(t)$  ( $t$  is current time)

$$\dot{c}(\dot{q}_{m0}, \dot{q}_{m1}) = k_{m1} \cdot \dot{q}_{m1} - \dot{q}_{m0} - UF_t(t, l_{\text{off}}) = 0 \quad (6.70)$$

Note that the index 2 equations are used, if the solver uses index 2 formulation OR if the flag `parameters.velocityLevel = True` (or both). The user functions include dependency on time  $t$ , but this time dependency is not respected in the computation of initial accelerations. Therefore, it is recommended that  $UF$  and  $UF_t$  does not include initial accelerations.

If `activeConnector = False`, the (index 1) algebraic equation reads for ALL cases:

$$c(\lambda_0) = \lambda_0 = 0 \quad (6.71)$$

**Example** for `ObjectConnectorCoordinate`:

---

```
def OffsetUF(t, lOffset): #gives 0.05 at t=1
    return 0.5*(1-np.cos(2*3.141592653589793*0.25*t))*lOffset

nMass=mbs.AddNode(Point(referenceCoordinates = [2,0,0]))
massPoint = mbs.AddObject(MassPoint(physicsMass = 5, nodeNumber = nMass))

groundMarker=mbs.AddMarker(MarkerNodeCoordinate(nodeNumber= nGround, coordinate = 0))
nodeMarker =mbs.AddMarker(MarkerNodeCoordinate(nodeNumber= nMass, coordinate = 0))

#Spring-Damper between two marker coordinates
mbs.AddObject(CoordinateConstraint(markerNumbers = [groundMarker, nodeMarker],
                                  offset = 0.1, offsetUserFunction = OffsetUF))

#assemble and solve system for default parameters
mbs.Assemble()
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', exu.SimulationSettings())

#check result at default integration time
testError = mbs.GetNodeOutput(nMass, exu.OutputVariableType.Displacement)[0] -
0.0499999999999272404
```

---

### 6.3.19 ObjectConnectorCoordinateVector

A constraint which constrains the coordinate vectors of two markers Marker[Node|Object|Body]Coordinates attached to nodes or bodies. The marker uses the objects LTG-lists to build the according coordinate mappings.

**Additional information for ObjectConnectorCoordinateVector:**

- The Object has the following types = Connector, Constraint
- Requested marker type = Coordinate
- **Short name** for Python = **CoordinateVectorConstraint**
- **Short name** for Python (visualization object) = **VCoordinateVectorConstraint**

The item **ObjectConnectorCoordinateVector** with type = 'ConnectorCoordinateVector' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
scalingMarker0	NumpyMatrix		Matrix[]	linear scaling matrix for coordinate vector of marker 0; matrix provided in python numpy format
scalingMarker1	NumpyMatrix		Matrix[]	linear scaling matrix for coordinate vector of marker 1; matrix provided in python numpy format
offset	NumpyVector		[]	offset added to constraint equation; only active, if no userFunction is defined
velocityLevel	bool		False	If true: connector constrains velocities (only works for ODE2 coordinates!); offset is used between velocities; in this case, the offsetUserFunction_t is considered and offsetUserFunction is ignored
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectConnectorCoordinateVector			parameters for visualization of item

The item **VObjectConnectorCoordinateVector** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R=-1, use default color

**Detailed information on ObjectConnectorCoordinateVector:**

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	

scalingMarker0	$\mathbf{X}_{m0} \in \mathbb{R}^{n_{ae} \times n_{q_{m0}}}$	
scalingMarker1	$\mathbf{X}_{m1} \in \mathbb{R}^{n_{ae} \times n_{q_{m1}}}$	
offset	$\mathbf{v}_{\text{off}} \in \mathbb{R}^{n_{ae}}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Displacement	$\Delta \mathbf{q}$	relative scalar displacement of marker coordinates, not including scaling matrices
Velocity	$\Delta \mathbf{v}$	difference of scalar marker velocity coordinates, not including scaling matrices
ConstraintEquation	$\mathbf{c}$	(residuum of) constraint equations
Force	$\lambda$	constraint force vector (vector of Lagrange multipliers), resulting from action of constraint equations

**Description of Item: Definition of quantities:**

intermediate variables	symbol	description
marker m0 coordinate vector	$\mathbf{q}_{m0} \in \mathbb{R}^{n_{q_{m0}}}$	coordinate vector provided by marker $m0$ ; depending on the marker, the coordinates may or may not include reference coordinates
marker m1 coordinate vector	$\mathbf{q}_{m1} \in \mathbb{R}^{n_{q_{m1}}}$	coordinate vector provided by marker $m1$ ; depending on the marker, the coordinates may or may not include reference coordinates
marker m0 velocity coordinate vector	$\dot{\mathbf{q}}_{m0} \in \mathbb{R}^{n_{q_{m0}}}$	velocity coordinate vector provided by marker $m0$
marker m1 velocity coordinate vector	$\dot{\mathbf{q}}_{m1} \in \mathbb{R}^{n_{q_{m1}}}$	velocity coordinate vector provided by marker $m1$
number of algebraic equations	$n_{ae}$	number of algebraic equations must be same as number of rows in $\mathbf{X}_{m0}$ and $\mathbf{X}_{m1}$
difference of coordinates	$\Delta \mathbf{q} = \mathbf{q}_{m1} - \mathbf{q}_{m0}$	Displacement between marker $m0$ to marker $m1$ coordinates
difference of velocity coordinates	$\Delta \mathbf{v} = \dot{\mathbf{q}}_{m1} - \dot{\mathbf{q}}_{m0}$	

**Remarks:** The number of algebraic equations depends on the number of rows in  $\mathbf{X}_{m0}$ , which must be same as the number of rows in  $\mathbf{X}_{m1}$ . The number of columns in  $\mathbf{X}_{m0}$  must agree with the length of the coordinate vector  $\mathbf{q}_{m0}$  and the number of columns in  $\mathbf{X}_{m1}$  must agree with the length of the coordinate vector  $\mathbf{q}_{m1}$ . If one marker  $k$  is a ground marker (node/object), the length of  $\mathbf{q}_{m,k}$  is zero and also the according matrix  $\mathbf{X}_{m,k}$  has zero size and will not be considered in the computation of the constraint equations.

If the number of rows of  $\mathbf{X}_{m0}$  plus the number of rows of  $\mathbf{X}_{m1}$  is larger than the total number of coordinates ( $\mathbf{q}_{m0}$  and  $\mathbf{q}_{m1}$ ), the algebraic equations are underdetermined and probably not solvable.

**Algebraic constraint equations:** If activeConnector = True, the index 3 algebraic equations

$$\mathbf{c}(\mathbf{q}_{m0}, \mathbf{q}_{m1}) = \mathbf{X}_{m1} \cdot \mathbf{q}_{m1} - \mathbf{X}_{m0} \mathbf{q}_{m0} - \mathbf{v}_{\text{off}} = 0 \quad (6.72)$$

If the offsetUserFunction UF is defined,  $\mathbf{c}$  instead becomes ( $t$  is current time)

$$\mathbf{c}(\mathbf{q}_{m0}, \mathbf{q}_{m1}) = \mathbf{X}_{m1} \cdot \mathbf{q}_{m1} - \mathbf{X}_{m0} \mathbf{q}_{m0} - \text{UF}(t, \mathbf{v}_{\text{off}}) = 0 \quad (6.73)$$

The activeConnector = True, index 2 (velocity level) algebraic equation reads

$$\dot{\mathbf{c}}(\dot{\mathbf{q}}_{m0}, \dot{\mathbf{q}}_{m1}) = \mathbf{X}_{m1} \cdot \dot{\mathbf{q}}_{m1} - \mathbf{X}_{m0} \dot{\mathbf{q}}_{m0} - \mathbf{d}_{\text{off}} = 0 \quad (6.74)$$

The vector  $dv$  in velocity level equations is zero, except if `parameters.velocityLevel = True`, then  $\mathbf{d} = \mathbf{v}_{\text{off}}$ .

If velocity level constraints are active and the velocity level `offsetUserFunction_t`  $\text{UF}_t$  is defined,  $\dot{\mathbf{c}}$  instead becomes ( $t$  is current time)

$$\dot{\mathbf{c}}(\dot{\mathbf{q}}_{m0}, \dot{\mathbf{q}}_{m1}) = \mathbf{X}_{m1} \cdot \dot{\mathbf{q}}_{m1} - \mathbf{X}_{m0} \dot{\mathbf{q}}_{m0} - \text{UF}_t(t, \mathbf{v}_{\text{off}}) = 0 \quad (6.75)$$

Note that the index 2 equations are used, if the solver uses index 2 formulation OR if the flag `parameters.velocityLevel = True` (or both). The user functions include dependency on time  $t$ , but this time dependency is not respected in the computation of initial accelerations. Therefore, it is recommended that  $\text{UF}$  and  $\text{UF}_t$  does not include initial accelerations.

If `activeConnector = False`, the (index 1) algebraic equation reads for ALL cases:

$$\mathbf{c}(\lambda) = \lambda = 0 \quad (6.76)$$

### 6.3.20 ObjectConnectorRollingDiscPenalty

A (flexible) connector representing a rolling rigid disc (marker 1) on a flat surface (marker 0, ground body) in global  $x$ - $y$  plane. The connector is based on a penalty formulation and adds friction and slipping. The constraints works for discs as long as the disc axis and the plane normal vector are not parallel. Parameters may need to be adjusted for better convergence (e.g., dryFrictionProportionalZone). The formulation is still under development and needs further testing.

**Additional information for ObjectConnectorRollingDiscPenalty:**

- The Object has the following types = Connector
- Requested marker type = Position + Orientation
- Requested node type = GenericData
- **Short name** for Python = **RollingDiscPenalty**
- **Short name** for Python (visualization object) = **VRollingDiscPenalty**

The item **ObjectConnectorRollingDiscPenalty** with type = 'ConnectorRollingDiscPenalty' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex	2	[ MAXINT, MAX-INT ]	list of markers used in connector
nodeNumber	NodeIndex		MAXINT	node number of a NodeGenericData (size=3) for 3 dataCoordinates
dryFrictionAngle	Real		0.	angle [SI:1 (rad)] which defines a rotation of the local tangential coordinates dry friction; this allows to model Mecanum wheels with specified roll angle
contactStiffness	Real		0.	normal contact stiffness [SI:N/m]
contactDamping	Real		0.	normal contact damping [SI:N/(m s)]
dryFriction	Vector2D		[0,0]	dry friction coefficients [SI:1] in local marker 1 joint /1 coordinates; if $\alpha_t == 0$ , lateral direction $l = x$ and forward direction $f = y$ ; assuming a normal force $f_n$ , the local friction force can be computed as ${}^{J1}\begin{bmatrix} f_{t,x} \\ f_{t,y} \end{bmatrix} = \begin{bmatrix} \mu_x f_n \\ \mu_y f_n \end{bmatrix}$
dryFrictionProportionalZone	Real		0.	limit velocity [m/s] up to which the friction is proportional to velocity (for regularization / avoid numerical oscillations)
rollingFrictionViscous	Real		0.	rolling friction [SI:1], which acts against the velocity of the trail on ground and leads to a force proportional to the contact normal force; currently, only implemented for disc axis parallel to ground!
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
discRadius	Real		0	defines the disc radius
visualization	VObjectConnectorRollingDiscPenalty			parameters for visualization of item

The item VObjectConnectorRollingDiscPenalty has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
discWidth	float		0.1	width of disc for drawing
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R=-1, use default color

Detailed information on ObjectConnectorRollingDiscPenalty:

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	
nodeNumber	$n_d$	
dryFrictionAngle	$\alpha_t$	
contactStiffness	$k_c$	
contactDamping	$d_c$	
dryFriction	$[\mu_x, \mu_y]^T$	
dryFrictionProportionalZone	$v_\mu$	
rollingFrictionViscous	$\mu_r$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_G$	current global position of contact point between rolling disc and ground
VelocityLocal	${}^D\mathbf{v}_G$	current velocity of the trail (contact) point in disc coordinates; this is the velocity with which the contact moves over the ground plane
ForceLocal	${}^1\mathbf{f}$	disc-ground force in special marker 1 joint coordinates, $f_0$ being the lateral force, $f_1$ being the longitudinal force and $f_2$ being the normal force

Description of Item:

Definition of quantities:

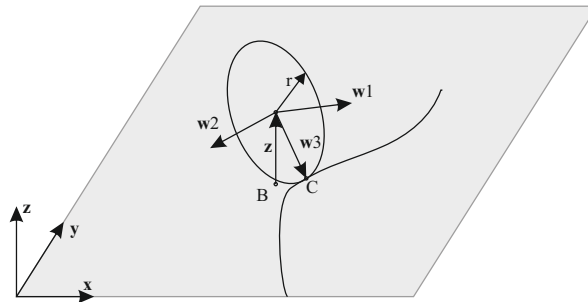
intermediate variables	symbol	description
marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker m0, any ground reference position; currently unused
marker m0 orientation	${}^{0,m0}\mathbf{A}$	current rotation matrix provided by marker m0; currently unused
marker m1 position	${}^0\mathbf{p}_{m1}$	center of disc
marker m1 orientation	${}^{0,m1}\mathbf{A}$	current rotation matrix provided by marker m1



data coordinates	$\mathbf{x} = [x_0, x_1, x_2]^T$	data coordinates for $[x_0, x_1]$ : hold the sliding velocity in lateral and longitudinal direction of last discontinuous iteration; $x_2$ : represents gap of last discontinuous iteration (in contact normal direction)
marker m1 velocity	${}^0\mathbf{v}_{m1}$	accordingly
marker m1 angular velocity	${}^0\boldsymbol{\omega}_{m1}$	current angular velocity vector provided by marker m1
ground normal vector	${}^0\mathbf{v}_{PN}$	normalized normal vector to the ground, currently $[0,0,1]$
ground position B	${}^0\mathbf{p}_B$	disc center point projected on ground (normal projection)
ground position C	${}^0\mathbf{p}_C$	contact point of disc with ground
ground velocity C	${}^0\mathbf{v}_C$	velocity of disc at ground contact point (must be zero at end of iteration)
wheel axis vector	${}^0\mathbf{w}_1 = {}^{0,m1}\mathbf{A} \cdot [1,0,0]^T$	normalized disc axis vector, currently $[1,0,0]^T$ in local coordinates
longitudinal vector	${}^0\mathbf{w}_2$	vector in longitudinal (motion) direction
lateral vector	${}^0\mathbf{w}_l = {}^0\mathbf{v}_{PN} \times {}^0\mathbf{w}_2 = [-\mathbf{w}_{2,y}, \mathbf{w}_{2,x}, 0]$	vector in lateral direction, lies in ground plane
contact point vector	${}^0\mathbf{w}_3$	normalized vector from disc center point in direction of contact point C
connector forces	${}^J\mathbf{f} = [f_{t,x}, f_{t,y}, f_n]^T$	joint force vector at contact point in joint 1 coordinates: $x$ =lateral direction, $y$ =longitudinal direction, $z$ =plane normal (contact normal)

### Geometric relations:

The main geometrical setup is shown in the following figure:



First, the contact point  ${}^0\mathbf{p}_C$  must be computed. With the helper vector,

$${}^0\mathbf{x} = {}^0\mathbf{w}_1 \times {}^0\mathbf{v}_{PN} \quad (6.77)$$

we obtain a disc coordinate system, representing the longitudinal direction,

$${}^0\mathbf{w}_2 = \frac{1}{|{}^0\mathbf{x}|} {}^0\mathbf{x} \quad (6.78)$$

and the vector to the contact point,

$${}^0\mathbf{w}_3 = {}^0\mathbf{w}_1 \times {}^0\mathbf{w}_2 \quad (6.79)$$

The contact point can be computed from

$${}^0\mathbf{p}_C = {}^0\mathbf{p}_{m1} + r \cdot {}^0\mathbf{w}_3 \quad (6.80)$$

The velocity of the contact point at the disc is computed from,

$${}^0\mathbf{v}_C = {}^0\mathbf{v}_{m1} + {}^0\boldsymbol{\omega}_{m1} \times (r \cdot {}^0\mathbf{w}_3) \quad (6.81)$$

The connector forces at the contact point C are computed as follows. The normal contact force reads

$$f_n = k_c \cdot {}^0\mathbf{p}_{C,z} + d_c \cdot {}^0\mathbf{v}_{C,z} \quad (6.82)$$

The tangential forces are computed from the inplane velocity  $\mathbf{v}_t = [{}^0\mathbf{v}_{C,x}, {}^0\mathbf{v}_{C,y}]^T$

$$\mathbf{f}_t = \boldsymbol{\mu} \cdot \phi(|\mathbf{v}_t|, v_\mu) \cdot f_n \cdot \mathbf{e}_t \quad (6.83)$$

with the regularization function:

$$\phi(v, v_\mu) = \begin{cases} (2 - \frac{v}{v_\mu})\frac{v}{v_\mu} & \text{if } v \leq v_\mu \\ 1 & \text{if } v > v_\mu \end{cases} \quad (6.84)$$

and the direction of tangential slip

$$\mathbf{e}_t = \frac{\mathbf{v}_t}{|\mathbf{v}_t|} \quad (6.85)$$

The friction coefficient matrix  $\boldsymbol{\mu}$  is computed from

$$\boldsymbol{\mu} = \begin{bmatrix} \mu_x & 0 \\ 0 & \mu_y \end{bmatrix} \quad (6.86)$$

where for isotropic behaviour of surface and wheel, it will give a diagonal matrix with the friction coefficient in the diagonal. In case that the dry friction angle  $\alpha_t$  is not zero, the  $\boldsymbol{\mu}$  changes to

$$\boldsymbol{\mu} = \begin{bmatrix} \cos(\alpha_t) & \sin(\alpha_t) \\ -\sin(\alpha_t) & \cos(\alpha_t) \end{bmatrix} \begin{bmatrix} \mu_x & 0 \\ 0 & \mu_y \end{bmatrix} \begin{bmatrix} \cos(\alpha_t) & -\sin(\alpha_t) \\ \sin(\alpha_t) & \cos(\alpha_t) \end{bmatrix} \quad (6.87)$$

Finally, the connector forces read in joint coordinates

$${}^I\mathbf{f} = \begin{bmatrix} f_{t,x} \\ f_{t,y} \\ f_n \end{bmatrix} \quad (6.88)$$

and in global coordinates, they are computed from

$${}^0\mathbf{f} = f_{t,x}\mathbf{w}_1 + f_{t,y}\mathbf{w}_2 + f_n\mathbf{w}_{PN} \quad (6.89)$$

The moment caused by the contact forces are given as

$${}^0\mathbf{f} = (r \cdot {}^0\mathbf{w}_3) \times {}^0\mathbf{f} \quad (6.90)$$

if `activeConnector = False`,

$${}^I\mathbf{f} = \mathbf{0} \quad (6.91)$$

### 6.3.21 ObjectContactCoordinate

A penalty-based contact condition for one coordinate; the contact gap  $g$  is defined as  $g = \text{marker.value}[1] - \text{marker.value}[0] - \text{offset}$ ; the contact force  $f_c$  is zero for  $gap > 0$  and otherwise computed from  $f_c = g * \text{contactStiffness} + \dot{g} * \text{contactDamping}$ ; during Newton iterations, the contact force is activated only, if  $\text{dataCoordinate}[0] \leq 0$ ;  $\text{dataCoordinate}$  is set equal to  $gap$  in nonlinear iterations, but not modified in Newton iterations.

**Additional information for ObjectContactCoordinate:**

- The Object has the following types = Connector
- Requested marker type = Coordinate
- Requested node type = GenericData

The item **ObjectContactCoordinate** with type = 'ContactCoordinate' has the following parameters:

Name	type	size	default value	description
name	String		"	connector's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	markers define contact gap
nodeNumber	NodeIndex		MAXINT	node number of a NodeGenericData for 1 dataCoordinate (used for active set strategy ==> holds the gap of the last discontinuous iteration)
contactStiffness	UReal		0.	contact (penalty) stiffness [SI:N/m]; acts only upon penetration
contactDamping	UReal		0.	contact damping [SI:N/(m s)]; acts only upon penetration
offset	UReal		0.	offset [SI:m] of contact
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectContactCoordinate			parameters for visualization of item

The item **VObjectContactCoordinate** has the following parameters:

Name	type	size	default value	description
show	Bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = diameter of spring; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R==-1, use default color

### 6.3.22 ObjectContactCircleCable2D

A very specialized penalty-based contact condition between a 2D circle (=marker0, any Position-marker) on a body and an ANCF Cable2DShape (=marker1, Marker: BodyCable2DShape), in xy-plane; a node NodeGenericData is required with the number of coordinates according to the number of contact segments; the contact gap  $g$  is integrated (piecewise linear) along the cable and circle; the contact force  $f_c$  is zero for  $gap > 0$  and otherwise computed from  $f_c = g * contactStiffness + \dot{g} * contactDamping$ ; during Newton iterations, the contact force is activated only, if  $dataCoordinate[0] \leq 0$ ; dataCoordinate is set equal to gap in nonlinear iterations, but not modified in Newton iterations.

#### Additional information for ObjectContactCircleCable2D:

- The Object has the following types = Connector
- Requested marker type = \_None
- Requested node type = GenericData

The item **ObjectContactCircleCable2D** with type = 'ContactCircleCable2D' has the following parameters:

Name	type	size	default value	description
name	String		"	connector's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAXINT ]	markers define contact gap
nodeNumber	NodeIndex		MAXINT	node number of a NodeGenericData for nSegments dataCoordinates (used for active set strategy ==> hold the gap of the last discontinuous iteration and the friction state)
numberOfContactSegments	Index		3	number of linear contact segments to determine contact; each segment is a line and is associated to a data (history) variable; must be same as in according marker
contactStiffness	UReal		0.	contact (penalty) stiffness [SI:N/(m s)/(contact segment)]; the stiffness is per contact segment; specific contact forces (per length) $f_N$ act in contact normal direction only upon penetration
contactDamping	UReal		0.	contact damping [SI:N/(m s)/(contact segment)]; the damping is per contact segment; acts in contact normal direction only upon penetration
circleRadius	UReal		0.	radius [SI:m] of contact circle
offset	UReal		0.	offset [SI:m] of contact, e.g. to include thickness of cable element
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectContactCircleCable2D			parameters for visualization of item

The item VObjectContactCircleCable2D has the following parameters:

Name	type	size	default value	description
show	Bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = diameter of spring; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R==-1, use default color

### 6.3.23 ObjectContactFrictionCircleCable2D

A very specialized penalty-based contact/friction condition between a 2D circle in the local x/y plane (=marker0, a Rigid-Body Marker) on a body and an ANCF Cable2DShape (=marker1, Marker: BodyCable2DShape), in xy-plane; a node NodeGenericData is required with  $3 \times (\text{number of contact segments})$  – containing per segment: [contact gap, stick/slip (stick=1), last friction position]; the contact gap  $g$  is integrated (piecewise linear) along the cable and circle; the contact force  $f_c$  is zero for  $gap > 0$  and otherwise computed from  $f_c = g * contactStiffness + \dot{g} * contactDamping$ ; during Newton iterations, the contact force is activated only, if  $dataCoordinate[0] \leq 0$ ; dataCoordinate is set equal to gap in nonlinear iterations, but not modified in Newton iterations.

#### Additional information for ObjectContactFrictionCircleCable2D:

- The Object has the following types = Connector
- Requested marker type = \_None
- Requested node type = GenericData

The item **ObjectContactFrictionCircleCable2D** with type = 'ContactFrictionCircleCable2D' has the following parameters:

Name	type	size	default value	description
name	String		"	connector's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	markers define contact gap
nodeNumber	NodeIndex		MAXINT	node number of a NodeGenericData with $3 \times nSegments$ dataCoordinates (used for active set strategy ==> hold the gap of the last discontinuous iteration and the friction state)
numberOfContactSegments	Index		3	number of linear contact segments to determine contact; each segment is a line and is associated to a data (history) variable; must be same as in according marker
contactStiffness	UReal		0.	contact (penalty) stiffness [SI:N/m/(contact segment)]; the stiffness is per contact segment; specific contact forces (per length) $f_N$ act in contact normal direction only upon penetration
contactDamping	UReal		0.	contact damping [SI:N/(m s)/(contact segment)]; the damping is per contact segment; acts in contact normal direction only upon penetration
frictionVelocityPenalty	UReal		0.	velocity dependent penalty coefficient for friction [SI:N/(m s)/(contact segment)]; the coefficient causes tangential (contact) forces against relative tangential velocities in the contact area

frictionStiffness	UReal		0.	CURRENTLY NOT IMPLEMENTED: displacement dependent penalty/stiffness coefficient for friction [SI:N/m/(contact segment)]; the coefficient causes tangential (contact) forces against relative tangential displacements in the contact area
frictionCoefficient	UReal		0.	friction coefficient $\mu$ [SI: 1]; tangential specific friction forces (per length) $f_T$ must fulfill the condition $f_T \leq \mu f_N$
circleRadius	UReal		0.	radius [SI:m] of contact circle
offset	UReal		0.	offset [SI:m] of contact, e.g. to include thickness of cable element
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectContactFrictionCircleCable2D			parameters for visualization of item

The item VObjectContactFrictionCircleCable2D has the following parameters:

Name	type	size	default value	description
show	Bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = diameter of spring; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R==-1, use default color

### 6.3.24 ObjectJointGeneric

A generic joint in 3D; constrains components of the absolute position and rotations of two points given by PointMarkers or RigidMarkers; an additional local rotation can be used to define three rotation axes and/or sliding axes

#### Additional information for ObjectJointGeneric:

- The Object has the following types = Connector, Constraint
- Requested marker type = Position + Orientation
- **Short name** for Python = **GenericJoint**
- **Short name** for Python (visualization object) = **VGenericJoint**

The item **ObjectJointGeneric** with type = 'JointGeneric' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex	2	[ MAXINT, MAX-INT ]	list of markers used in connector
constrainedAxes	ArrayIndex	6	[1,1,1,1,1,1]	flag, which determines which translation (0,1,2) and rotation (3,4,5) axes are constrained; for $j_i$ , two values are possible: 0=free axis, 1=constrained axis
rotationMarker0	Matrix3D		[[1,0,0], [0,1,0], [0,0,1]]	local rotation matrix for marker $m0$ ; translation and rotation axes for marker $m0$ are defined in the local body coordinate system and additionally transformed by rotation-Marker0
rotationMarker1	Matrix3D		[[1,0,0], [0,1,0], [0,0,1]]	local rotation matrix for marker $m1$ ; translation and rotation axes for marker $m1$ are defined in the local body coordinate system and additionally transformed by rotation-Marker1
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
offsetUserFunctionParameters	Vector6D		[0.,0.,0.,0.,0.,0.]	vector of 6 parameters for joint's offsetUser-Function



offsetUserFunction	PyFunctionVector6D	Scalar	Vector6D 0	A python function which defines the time-dependent (fixed) offset of translation (indices 0,1,2) and rotation (indices 3,4,5) joint coordinates with parameters (t, offsetUserFunctionParameters); the offset represents the current value of the object; it is highly RECOMMENDED to use sufficiently smooth functions, having consistent initial offsets with initial configuration of bodies, zero or compatible initial offset-velocity, and no accelerations; Example for python function: def f(t, offsetUserFunctionParameters): return [offsetUserFunctionParameters[0]*(1 - np.cos(t*10*2*np.pi)), 0,0,0,0,0]
offsetUserFunction_t	PyFunctionVector6D	Scalar	Vector6D 0	(NOT IMPLEMENTED YET)time derivative of offsetUserFunction using the same parameters; needed for 'velocityLevel=True', or for index2 time integration and for computation of initial accelerations in SecondOrderImplicit integrators
visualization	VObjectJointGeneric			parameters for visualization of item

The item VObjectJointGeneric has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
axesRadius	float		0.1	radius of joint axes to draw
axesLength	float		0.4	length of joint axes to draw
color	Float4		[-1,-1,-1,-1.]	RGBA connector color; if R== -1, use default color

#### Detailed information on ObjectJointGeneric:

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	
constrainedAxes	$\mathbf{j} = [j_0, \dots, j_5]$	
rotationMarker0	${}^{m0,j0}\mathbf{A}$	
rotationMarker1	${}^{m1,j1}\mathbf{A}$	
offsetUserFunctionParameters	$\mathbf{p}_{par}$	
offsetUserFunction	$UF(t, \mathbf{p}_{par})$	
offsetUserFunction_t	$UF_t(t, \mathbf{p}_{par})$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{m0}$	current global position of position marker $m0$

Velocity	${}^0\mathbf{v}_{m0}$	current global velocity of position marker $m0$
DisplacementLocal	${}^{J0}\Delta\mathbf{p}$	relative displacement in local joint0 coordinates; uses local J0 coordinates even for spherical joint configuration
VelocityLocal	${}^{J0}\Delta\mathbf{v}$	relative translational velocity in local joint0 coordinates
Rotation	${}^{J0}\boldsymbol{\theta} = [\theta_0, \theta_1, \theta_2]^T$	relative rotation parameters (Tait Bryan Rxyz); if all axes are fixed, this output represents the rotational drift; for a revolute joint, it contains the rotation of this axis
AngularVelocityLocal	${}^{J0}\Delta\boldsymbol{\omega}$	relative angular velocity in local joint0 coordinates; if all axes are fixed, this output represents the angular velocity constraint error; for a revolute joint, it contains the angular velocity of this axis
ForceLocal	${}^{J0}\mathbf{f}$	joint force in local J0 coordinates
TorqueLocal	${}^{J0}\mathbf{m}$	joint torque in local J0 coordinates; depending on joint configuration, the result may not be the according torque vector

#### Description of Item:

#### Definition of quantities:

intermediate variables	symbol	description
marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker m0
marker m0 orientation	${}^{0,m0}\mathbf{A}$	current rotation matrix provided by marker m0
joint J0 orientation	${}^{0,J0}\mathbf{A} = {}^{0,m0}\mathbf{A} {}^{m0,J0}\mathbf{A}$	joint J0 rotation matrix
joint J0 orientation vectors	${}^{0,J0}\mathbf{A} = [\mathbf{v}_{x0}, \mathbf{v}_{y0}, \mathbf{v}_{z0}]^T$	orientation vectors used for definition of constraint equations
marker m1 position	${}^0\mathbf{p}_{m1}$	accordingly
marker m1 orientation	${}^{0,m1}\mathbf{A}$	current rotation matrix provided by marker m1
joint J1 orientation	${}^{0,J1}\mathbf{A} = {}^{0,m1}\mathbf{A} {}^{m1,J1}\mathbf{A}$	joint J1 rotation matrix
joint J1 orientation vectors	${}^{0,J1}\mathbf{A} = [\mathbf{v}_{x1}, \mathbf{v}_{y1}, \mathbf{v}_{z1}]^T$	orientation vectors used for definition of constraint equations
marker m0 velocity	${}^0\mathbf{v}_{m0}$	current global velocity which is provided by marker m0
marker m1 velocity	${}^0\mathbf{v}_{m1}$	accordingly
marker m0 velocity	${}^b\boldsymbol{\omega}_{m0}$	current local angular velocity vector provided by marker m0
marker m1 velocity	${}^b\boldsymbol{\omega}_{m1}$	current local angular velocity vector provided by marker m1
Displacement	${}^0\Delta\mathbf{p} = {}^0\mathbf{p}_{m1} - {}^0\mathbf{p}_{m0}$	used, if all translational axes are constrained
Velocity	${}^0\Delta\mathbf{v} = {}^0\mathbf{v}_{m1} - {}^0\mathbf{v}_{m0}$	used, if all translational axes are constrained (velocity level)
DisplacementLocal	${}^{J0}\Delta\mathbf{p}$	$({}^{0,m0}\mathbf{A} {}^{m0,J0}\mathbf{A})^T {}^0\Delta\mathbf{p}$

VelocityLocal	${}^{J0}\Delta \mathbf{v}$	$\left( {}^{0,m0}\mathbf{A} \quad {}^{m0,J0}\mathbf{A} \right)^T {}^0\Delta \mathbf{v}$ ... note that this is the global relative velocity projected into the local $J0$ coordinate system
AngularVelocityLocal	${}^{J0}\Delta \omega$	$\left( {}^{0,m0}\mathbf{A} \quad {}^{m0,J0}\mathbf{A} \right)^T \left( {}^{0,m1}\mathbf{A} \quad {}^{m1}\omega - {}^{0,m0}\mathbf{A} \quad {}^{m0}\omega \right)$
algebraic variables	$\mathbf{z} = [\lambda_0, \dots, \lambda_5]^T$	vector of algebraic variables (Lagrange multipliers) according to the algebraic equations

**Connector equations for translational part (activeConnector = True):**

If  $[j_0, \dots, j_2] = [1, 1, 1]^T$ , meaning that all translational coordinates are fixed, the translational index 3 constraints read  $(UF_{0,1,2}(t, \mathbf{p}_{par}))$  is the translational part of the user function  $UF$ ,

$${}^0\mathbf{p}_{m1} - {}^0\mathbf{p}_{m0} - UF_{0,1,2}(t, \mathbf{p}_{par}) = \mathbf{0} \quad (6.92)$$

and the translational index 2 constraints read

$${}^0\mathbf{v}_{m1} - {}^0\mathbf{v}_{m0} - UF_{t;0,1,2}(t, \mathbf{p}_{par}) = \mathbf{0} \quad (6.93)$$

If  $[j_0, \dots, j_2] \neq [1, 1, 1]^T$ , meaning that at least one translational coordinate is free, the translational index 3 constraints read for every component  $k \in [0, 1, 2]$  of the vector  ${}^{J0}\Delta \mathbf{p}$

$${}^{J0}\Delta p_k - UF_k(t, \mathbf{p}_{par}) = 0 \quad \text{if } j_k = 1 \quad \text{and} \quad (6.94)$$

$$\lambda_k = 0 \quad \text{if } j_k = 0 \quad (6.95)$$

$$(6.96)$$

and the translational index 2 constraints read for every component  $k \in [0, 1, 2]$  of the vector  ${}^{J0}\Delta \mathbf{v}$

$${}^{J0}\Delta v_k - UF_{t;k}(t, \mathbf{p}_{par}) = 0 \quad \text{if } j_k = 1 \quad \text{and} \quad (6.97)$$

$$\lambda_k = 0 \quad \text{if } j_k = 0 \quad (6.98)$$

$$(6.99)$$

**Connector equations for rotational part (activeConnector = True):**

The following equations are exemplarily for certain constrained rotation axes configurations, which shall represent all other possibilities. Equations are only given for the index 3 case; the index 2 case can be derived from these equations easily (see C++ code...). In case of user functions, the additional rotation matrix  ${}^{J0,J0U}\mathbf{A}(UF_{3,4,5}(t, \mathbf{p}_{par}))$ , in which the three components of  $UF_{3,4,5}$  are interpreted as Tait-Bryan angles that are added to the joint frame.

If **3 rotation axes are constrained**,  $[j_3, \dots, j_5] = [1, 1, 1]^T$ , the index 3 constraint equations read

$$\mathbf{v}_{z0}^T \mathbf{v}_{y1} = 0 \quad (6.100)$$

$$\mathbf{v}_{z0}^T \mathbf{v}_{x1} = 0 \quad (6.101)$$

$$\mathbf{v}_{x0}^T \mathbf{v}_{y1} = 0 \quad (6.102)$$

If **2 rotation axes are constrained**, e.g.,  $[j_3, \dots, j_5] = [0, 1, 1]^T$ , the index 3 constraint equations read

$$\lambda_3 = 0 \quad (6.103)$$

$$\mathbf{v}_{x0}^T \mathbf{v}_{y1} = 0 \quad (6.104)$$

$$\mathbf{v}_{x0}^T \mathbf{v}_{z1} = 0 \quad (6.105)$$

If **1 rotation axis is constrained**, e.g.,  $[j_3, \dots, j_5] = [1, 0, 0]^T$ , the index 3 constraint equations read

$$\mathbf{v}_{y0}^T \mathbf{v}_{z1} = 0 \quad (6.106)$$

$$\lambda_4 = 0 \quad (6.107)$$

$$\lambda_5 = 0 \quad (6.108)$$

if `activeConnector = False`,

$$\mathbf{z} = \mathbf{0} \quad (6.109)$$

### 6.3.25 ObjectJointSpherical

A spherical joint, which constrains the relative translation between two position based markers.

**Additional information for ObjectJointSpherical:**

- The Object has the following types = Connector, Constraint
- Requested marker type = Position
- **Short name** for Python = **SphericalJoint**
- **Short name** for Python (visualization object) = **VSphericalJoint**

The item **ObjectJointSpherical** with type = 'JointSpherical' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex	2	[ MAXINT, MAX-INT ]	list of markers used in connector
constrainedAxes	ArrayIndex	3	[1,1,1]	flag, which determines which translation (0,1,2) and rotation (3,4,5) axes are constrained; for $j_i$ , two values are possible: 0=free axis, 1=constrained axis
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectJointSpherical			parameters for visualization of item

The item **VObjectJointSpherical** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
jointRadius	float		0.1	radius of joint to draw
color	Float4		[-1,-1,-1,-1.]	RGBA connector color; if R=-1, use default color

**Detailed information on ObjectJointSpherical:**

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	
constrainedAxes	$\mathbf{j} = [j_0, \dots, j_2]$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{m0}$	current global position of position marker $m0$
Velocity	${}^0\mathbf{v}_{m0}$	current global velocity of position marker $m0$
Displacement	${}^0\Delta\mathbf{p} = {}^0\mathbf{p}_{m1} - {}^0\mathbf{p}_{m0}$	constraint drift or relative motion, if not all axes fixed

Force	${}^0\mathbf{f}$	joint force in global coordinates
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**Description of Item:**

**Definition of quantities:**

intermediate variables	symbol	description
marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker $m0$
marker m1 position	${}^0\mathbf{p}_{m1}$	current global position which is provided by marker $m1$
marker m0 velocity	${}^0\mathbf{v}_{m0}$	current global velocity which is provided by marker $m0$
marker m1 velocity	${}^0\mathbf{v}_{m1}$	current global velocity which is provided by marker $m1$
relative velocity	${}^0\Delta\mathbf{v} = {}^0\mathbf{v}_{m1} - {}^0\mathbf{v}_{m0}$	constraint velocity error, or relative velocity if not all axes fixed
algebraic variables	$\mathbf{z} = [\lambda_0, \dots, \lambda_2]^T$	vector of algebraic variables (Lagrange multipliers) according to the algebraic equations

**Connector equations (activeConnector = True):**

If  $[j_0, \dots, j_2] = [1, 1, 1]^T$ , meaning that all translational coordinates are fixed, the translational index 3 constraints read

$${}^0\mathbf{p}_{m1} - {}^0\mathbf{p}_{m0} = \mathbf{0} \quad (6.110)$$

and the translational index 2 constraints read

$${}^0\mathbf{v}_{m1} - {}^0\mathbf{v}_{m0} = \mathbf{0} \quad (6.111)$$

If  $[j_0, \dots, j_2] \neq [1, 1, 1]^T$ , meaning that at least one translational coordinate is free, the translational index 3 constraints read for every component  $k \in [0, 1, 2]$  of the vector  ${}^0\Delta\mathbf{p}$

$${}^0\Delta p_k = 0 \quad \text{if } j_k = 1 \quad \text{and} \quad (6.112)$$

$$\lambda_k = 0 \quad \text{if } j_k = 0 \quad (6.113)$$

$$(6.114)$$

and the translational index 2 constraints read for every component  $k \in [0, 1, 2]$  of the vector  ${}^0\Delta\mathbf{v}$

$${}^0\Delta v_k = 0 \quad \text{if } j_k = 1 \quad \text{and} \quad (6.115)$$

$$\lambda_k = 0 \quad \text{if } j_k = 0 \quad (6.116)$$

$$(6.117)$$

if activeConnector = False,

$$\mathbf{z} = \mathbf{0} \quad (6.118)$$

### 6.3.26 ObjectJointRollingDisc

A joint representing a rolling rigid disc (marker 1) on a flat surface (marker 0, ground body) in global  $x$ - $y$  plane. The constraint is based on an idealized rolling formulation with no slip. The constraints works for discs as long as the disc axis and the plane normal vector are not parallel. It must be assured that the disc has contact to ground in the initial configuration (adjust  $z$ -position of body accordingly).

**Additional information for ObjectJointRollingDisc:**

- The Object has the following types = Connector, Constraint
- Requested marker type = Position + Orientation
- **Short name** for Python = **RollingDiscJoint**
- **Short name** for Python (visualization object) = **VRollingDiscJoint**

The item **ObjectJointRollingDisc** with type = 'JointRollingDisc' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex	2	[ MAXINT, MAX-INT ]	list of markers used in connector
constrainedAxes	ArrayIndex	3	[1,1,1]	flag, which determines which constraints are active, in which $j_0, j_1$ represent the tangential motion and $j_2$ represents the normal (contact) direction
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
discRadius	Real		0	defines the disc radius
visualization	VObjectJointRollingDisc			parameters for visualization of item

The item **VObjectJointRollingDisc** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
discWidth	float		0.1	width of disc for drawing
color	Float4		[-1,-1,-1,-1.]	RGBA connector color; if R=-1, use default color

**Detailed information on ObjectJointRollingDisc:**

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	
constrainedAxes	$\mathbf{j} = [j_0, \dots, j_2]$	

The following output parameters are available as **OutputVariableType** in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_G$	current global position of contact point between rolling disc and ground

VelocityLocal	${}^D\mathbf{v}_G$	current velocity of the trail (contact) point in disc coordinates; this is the velocity with which the contact moves over the ground plane
ForceLocal	${}^D\mathbf{f} = [-\mathbf{z}^T \mathbf{w}_l, -\mathbf{z}^T \mathbf{w}_2, -z_z]^T$	contact forces acting on disc, in special disc coordinates, $f_x$ being the lateral force, $f_y$ being the longitudinal force and $f_z$ being the normal force

#### Description of Item:

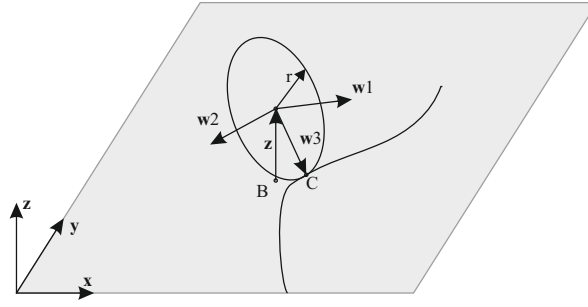
#### Definition of quantities:

intermediate variables	symbol	description
marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker m0, any ground reference position; currently unused
marker m0 orientation	${}^{0,m0}\mathbf{A}$	current rotation matrix provided by marker m0; currently unused
marker m1 position	${}^0\mathbf{p}_{m1}$	center of disc
marker m1 orientation	${}^{0,m1}\mathbf{A}$	current rotation matrix provided by marker m1
marker m1 velocity	${}^0\mathbf{v}_{m1}$	accordingly
marker m1 angular velocity	${}^0\boldsymbol{\omega}_{m1}$	current angular velocity vector provided by marker m1
ground normal vector	${}^0\mathbf{v}_{PN}$	normalized normal vector to the ground plane, currently [0,0,1]
ground position B	${}^0\mathbf{p}_B$	disc center point projected on ground (normal projection)
ground position C	${}^0\mathbf{p}_C$	contact point of disc with ground
ground velocity C	${}^0\mathbf{v}_C$	velocity of disc at ground contact point (must be zero at end of iteration)
wheel axis vector	${}^0\mathbf{w}_1 = {}^{0,m1}\mathbf{A} \cdot [1, 0, 0]^T$	normalized disc axis vector, currently $[1, 0, 0]^T$ in local coordinates
longitudinal vector	${}^0\mathbf{w}_2$	vector in longitudinal (motion) direction
lateral vector	${}^0\mathbf{w}_l = {}^0\mathbf{v}_{PN} \times {}^0\mathbf{w}_2 = [-\mathbf{w}_{2,y}, \mathbf{w}_{2,x}, 0]$	vector in lateral direction, lies in ground plane
contact point vector	${}^0\mathbf{w}_3$	normalized vector from disc center point in direction of contact point C
algebraic variables	$\mathbf{z} = [\lambda_0, \lambda_1, \lambda_2]^T$	vector of algebraic variables (Lagrange multipliers) according to the algebraic equations

#### Geometric relations:

The main geometrical setup is shown in the following figure:





First, the contact point  ${}^0\mathbf{p}_C$  must be computed. With the helper vector,

$${}^0\mathbf{x} = {}^0\mathbf{w}_1 \times {}^0\mathbf{v}_{PN} \quad (6.119)$$

we obtain a disc coordinate system, representing the longitudinal direction,

$${}^0\mathbf{w}_2 = \frac{1}{|{}^0\mathbf{x}|} {}^0\mathbf{x} \quad (6.120)$$

and the vector to the contact point,

$${}^0\mathbf{w}_3 = {}^0\mathbf{w}_1 \times {}^0\mathbf{w}_2 \quad (6.121)$$

The contact point can be computed from

$${}^0\mathbf{p}_C = {}^0\mathbf{p}_{m1} + r \cdot {}^0\mathbf{w}_3 \quad (6.122)$$

The velocity of the contact point at the disc is computed from,

$${}^0\mathbf{v}_C = {}^0\mathbf{v}_{m1} + {}^0\boldsymbol{\omega}_{m1} \times (r \cdot {}^0\mathbf{w}_3) \quad (6.123)$$

#### Connector constraint equations (**activeConnector = True**):

The non-holonomic, index 2 constraints for the tangential motion follow from (an index 3 formulation would be possible, but is not implemented yet because of mixing different jacobians)

$$\begin{bmatrix} {}^0\mathbf{v}_{C,x} \\ {}^0\mathbf{v}_{C,y} \end{bmatrix} = \mathbf{0} \quad (6.124)$$

In the index 2 (velocity level) case, the constraint for the normal direction reads

$${}^0\mathbf{v}_{C,z} = 0 \quad (6.125)$$

if **activeConnector = False**,

$$\mathbf{z} = 0 \quad (6.126)$$

### 6.3.27 ObjectJointRevolute2D

A revolute joint in 2D; constrains the absolute 2D position of two points given by PointMarkers or RigidMarkers

**Additional information for ObjectJointRevolute2D:**

- The Object has the following types = Connector, Constraint
- Requested marker type = Position
- **Short name** for Python = **RevoluteJoint2D**
- **Short name** for Python (visualization object) = **VRevoluteJoint2D**

The item **ObjectJointRevolute2D** with type = 'JointRevolute2D' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectJointRevolute2D			parameters for visualization of item

The item **VObjectJointRevolute2D** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = radius of revolute joint; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R=-1, use default color

### 6.3.28 ObjectJointPrismatic2D

A prismatic joint in 2D; allows the relative motion of two bodies, using two RigidMarkers; the vector  $\mathbf{t}_0$  = axisMarker0 is given in local coordinates of the first marker's (body) frame and defines the prismatic axis; the vector  $\mathbf{n}_1$  = normalMarker1 is given in the second marker's (body) frame and is the normal vector to the prismatic axis; using the global position vector  $\mathbf{p}_0$  and rotation matrix  $\mathbf{A}_0$  of marker0 and the global position vector  $\mathbf{p}_1$  rotation matrix  $\mathbf{A}_1$  of marker1, the equations for the prismatic joint follow as

$$(\mathbf{p}_1 - \mathbf{p}_0)^T \cdot \mathbf{A}_1 \cdot \mathbf{n}_1 = 0 \quad (6.127)$$

$$(\mathbf{A}_0 \cdot \mathbf{t}_0)^T \cdot \mathbf{A}_1 \cdot \mathbf{n}_1 = 0 \quad (6.128)$$

The lagrange multipliers follow for these two equations  $[\lambda_0, \lambda_1]$ , in which  $\lambda_0$  is the transverse force and  $\lambda_1$  is the torque in the joint.

**Additional information for ObjectJointPrismatic2D:**

- The Object has the following types = Connector, Constraint
- Requested marker type = Position + Orientation
- **Short name** for Python = **PrismaticJoint2D**
- **Short name** for Python (visualization object) = **VPrismaticJoint2D**

The item **ObjectJointPrismatic2D** with type = 'JointPrismatic2D' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	list of markers used in connector
axisMarker0	Vector3D		[1.,0.,0.]	direction of prismatic axis, given as a 3D vector in Marker0 frame
normalMarker1	Vector3D		[0.,1.,0.]	direction of normal to prismatic axis, given as a 3D vector in Marker1 frame
constrainRotation	bool		True	flag, which determines, if the connector also constrains the relative rotation of the two objects; if set to false, the constraint will keep an algebraic equation set equal zero
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectJointPrismatic2D			parameters for visualization of item

The item **VObjectJointPrismatic2D** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = radius of revolute joint; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R=-1, use default color



### 6.3.29 ObjectJointSliding2D

A specialized sliding joint (without rotation) in 2D between a Cable2D (marker1) and a position-based marker (marker0); the data coordinate x[0] provides the current index in slidingMarkerNumbers, and x[1] the local position in the cable element at the beginning of the timestep.

#### Additional information for ObjectJointSliding2D:

- The Object has the following types = Connector, Constraint
- Requested marker type = `_None`
- Requested node type = `GenericData`
- **Short name** for Python = **SlidingJoint2D**
- **Short name** for Python (visualization object) = **VSlidingJoint2D**

The item **ObjectJointSliding2D** with type = 'JointSliding2D' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	marker m0: position-marker of mass point or rigid body; marker m1: updated marker to Cable2D element, where the sliding joint currently is attached to; must be initialized with an appropriate (global) marker number according to the starting position of the sliding object; this marker changes with time (PostNewtonStep)
slidingMarkerNumbers	ArrayMarkerIndex		[]	these markers are used to update marker m1, if the sliding position exceeds the current cable's range; the markers must be sorted such that marker $m_{si}$ at $x=cable(i).length$ is equal to marker(i+1) at $x=0$ of cable(i+1)
slidingMarkerOffsets	Vector		[]	this list contains the offsets of every sliding object (given by slidingMarkerNumbers) w.r.t. to the initial position (0): marker m0: offset=0, marker m1: offset=Length(cable0), marker m2: offset=Length(cable0)+Length(cable1), ...
nodeNumber	NodeIndex		MAXINT	node number of a NodeGenericData for 1 dataCoordinate showing the according marker number which is currently active and the start-of-step (global) sliding position
classicalFormulation	bool		True	uses a formulation with 3 equations, including the force in sliding direction to be zero; forces in global coordinates, only index 3; alternatively: use local formulation, which only needs two equations and can be used with index 2 formulation
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint

visualization	VObjectJointSliding2D		parameters for visualization of item
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The item VObjectJointSliding2D has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = radius of revolute joint; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R==-1, use default color

#### Detailed information on ObjectJointSliding2D:

input parameter	symbol	description see tables above
markerNumbers	$[m_0, m_1]^T$	
slidingMarkerNumbers	$[m_{s0}, \dots, m_{sn}]^T$	
slidingMarkerOffsets	$[d_{s0}, \dots, d_{sn}]$	
nodeNumber	$n_{GD}$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position		position vector of joint given by marker0
Velocity		velocity vector of joint given by marker0
SlidingCoordinate		global sliding coordinate along all elements; the maximum sliding coordinate is equivalent to the reference lengths of all sliding elements
Force		joint force vector (3D)

#### Description of Item:

#### Definition of quantities:

intermediate variables	symbol	description
data node	$\mathbf{x} = [x_{data0}, x_{data1}]^T$	coordinates of node with node number $n_{GD}$
data coordinate 0	$x_{data0}$	the current index in slidingMarkerNumbers
data coordinate 1	$x_{data1}$	the global sliding coordinate (ranging from 0 to the total length of all sliding elements) at <b>start-of-step</b> - beginning of the timestep
marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker m0
marker m0 velocity	${}^0\mathbf{v}_{m0}$	current global velocity which is provided by marker m0
cable coordinates	$\mathbf{q}_{ANCF,m1}$	current coordinates of the ANCF cable element with the current marker $m1$ is referring to

sliding position	${}^0\mathbf{r}_{ANCF} = \mathbf{S}(s_{el})\mathbf{q}_{ANCF,m1}$	current global position at the ANCF cable element, evaluated at local sliding position $s_{el}$
sliding position slope	${}^0\mathbf{r}'_{ANCF} = \mathbf{S}'(s_{el})\mathbf{q}_{ANCF,m1} = [r'_0, r'_1]^T$	current global slope vector of the ANCF cable element, evaluated at local sliding position $s_{el}$
sliding velocity	${}^0\mathbf{v}_{ANCF} = \mathbf{S}(s_{el})\dot{\mathbf{q}}_{ANCF,m1}$	current global velocity at the ANCF cable element, evaluated at local sliding position $s_{el}$ ( $s_{el}$ not differentiated!!!)
sliding velocity slope	${}^0\mathbf{v}'_{ANCF} = \mathbf{S}'(s_{el})\dot{\mathbf{q}}_{ANCF,m1}$	current global slope velocity vector of the ANCF cable element, evaluated at local sliding position $s_{el}$
sliding normal vector	${}^0\mathbf{n} = [-r'_1, r'_0]$	2D normal vector computed from slope $\mathbf{r}' = {}^0\mathbf{r}'_{ANCF}$
sliding normal velocity vector	${}^0\dot{\mathbf{n}} = [-\dot{r}'_1, \dot{r}'_0]$	time derivative of 2D normal vector computed from slope velocity $\dot{\mathbf{r}}' = {}^0\dot{\mathbf{r}}'_{ANCF}$
algebraic coordinates	$\mathbf{z} = [\lambda_0, \lambda_1, s]^T$	algebraic coordinates composed of Lagrange multipliers $\lambda_0$ and $\lambda_1$ (in local cable coordinates: $\lambda_0$ is in axis direction) and the current sliding coordinate $s$ , which is local in the current cable element.
local sliding coordinate	$s$	local incremental sliding coordinate $s$ : the (algebraic) sliding coordinate <b>relative to the start-of-step value</b> . Thus, $s$ only contains small local increments.

output variables	symbol	formula
Position	${}^0\mathbf{p}_{m0}$	current global position of position marker $m0$
Velocity	${}^0\mathbf{v}_{m0}$	current global velocity of position marker $m0$
SlidingCoordinate	$s_g = s + x_{data1}$	current value of the global sliding coordinate
Force	$\mathbf{f}$	see below

Assume we have given the sliding coordinate  $s$  (e.g., as a guess of the Newton method or beginning of the time step). The element sliding coordinate (in the local coordinates of the current sliding element) is computed as

$$s_{el} = s + x_{data1} - d_{m1} = s_g - d_{m1}. \quad (6.129)$$

The vector (=difference; error) between the marker  $m0$  and the marker  $m1$  ( $=\mathbf{r}_{ANCF}$ ) positions reads

$${}^0\Delta\mathbf{p} = {}^0\mathbf{r}_{ANCF} - {}^0\mathbf{p}_{m0} \quad (6.130)$$

The vector (=difference; error) between the marker  $m0$  and the marker  $m1$  velocities reads

$${}^0\Delta\mathbf{v} = {}^0\dot{\mathbf{r}}_{ANCF} - {}^0\mathbf{v}_{m0} \quad (6.131)$$

**Algebraic constraint equations (classicalFormulation=True):** The 2D sliding joint is implemented having 3

equations, using the special algebraic coordinates  $\mathbf{z}$ . The algebraic equations read

$${}^0\Delta\mathbf{p} = \mathbf{0}, \quad \dots \text{2 index 3 equations, ensuring the sliding body to stay at the cable} \quad (6.132)$$

$$[\lambda_0, \lambda_1] \cdot {}^0\mathbf{r}'_{ANCF} = 0, \quad \dots \text{1 index 1 equation, ensuring the force in sliding direction} = 0 \quad (6.133)$$

$$(6.134)$$

No index 2 case exists, because no time derivative exists for  $s_{el}$ . The jacobian matrices for algebraic and ODE2 coordinates read

$$J_{AE} = \begin{bmatrix} 0 & 0 & r'_0 \\ 0 & 0 & r'_1 \\ r'_0 & r'_1 & r''_0\lambda_0 + r''_1\lambda_1 \end{bmatrix} \quad (6.135)$$

$$J_{ODE2} = \begin{bmatrix} -J_{pos,m0} & \mathbf{S}(s_{el}) \\ \mathbf{0}^T & [\lambda_0, \lambda_1] \cdot \mathbf{S}'(s_{el}) \end{bmatrix} \quad (6.136)$$

if `activeConnector = False`, the algebraic equations are changed to:

$$\lambda_0 = 0, \quad (6.137)$$

$$\lambda_1 = 0, \quad (6.138)$$

$$s = 0 \quad (6.139)$$

**Algebraic constraint equations (classicalFormulation=False):** The 2D sliding joint is implemented having 3 equations (first equation is dummy and could be eliminated), using the special algebraic coordinates  $\mathbf{z}$ . The algebraic equations read

$$\lambda_0 = 0, \quad \dots \text{this equation is not necessary, but can be used for switching to other mode} \quad (6.140)$$

$${}^0\Delta\mathbf{p}^T {}^0\mathbf{n} = 0, \quad \dots \text{equation ensures that sliding body stays at cable centerline; index3 equation} \quad (6.141)$$

$${}^0\Delta\mathbf{p}^T {}^0\mathbf{r}'_{ANCF} = 0. \quad \dots \text{resolves the sliding coordinate } s; \text{ index1 equation!} \quad (6.142)$$

In the index 2 case, the second equation reads

$${}^0\Delta\mathbf{v}^T {}^0\mathbf{n} + {}^0\Delta\mathbf{p}^T {}^0\dot{\mathbf{n}} = 0 \quad (6.143)$$

if `activeConnector = False`, the algebraic equations are changed to:

$$\lambda_0 = 0, \quad (6.144)$$

$$\lambda_1 = 0, \quad (6.145)$$

$$s = 0 \quad (6.146)$$

**Post Newton Step:** After the Newton solver has converged, a PostNewtonStep is performed for the element, which updates the marker  $m1$  index if necessary.

$$\begin{aligned} s_{el} < 0 & \rightarrow x_{data0} -= 1 \\ s_{el} > L & \rightarrow x_{data0} += 1 \end{aligned} \quad (6.147)$$

Furthermore, it is checked, if  $x_{data0}$  becomes smaller than zero, which raises a warning and keeps  $x_{data0} = 0$ . The same results if  $x_{data0} \geq sn$ , then  $x_{data0} = sn$ . Finally, the data coordinate is updated in order to provide the starting value for the next step,

$$x_{data1} += s. \quad (6.148)$$

**Examples:** see TestModels!



### 6.3.30 ObjectJointALEMoving2D

A specialized axially moving joint (without rotation) in 2D between a ALE Cable2D (marker1) and a position-based marker (marker0); ALE=Arbitrary Lagrangian Eulerian; the data coordinate x[0] provides the current index in slidingMarkerNumbers, and the ODE2 coordinate q[0] provides the (given) moving coordinate in the cable element.

#### Additional information for ObjectJointALEMoving2D:

- The Object has the following types = Connector, Constraint
- Requested marker type = \_None
- Requested node type: read detailed information of item
- **Short name** for Python = **ALEMovingJoint2D**
- **Short name** for Python (visualization object) = **VALEMovingJoint2D**

The item **ObjectJointALEMoving2D** with type = 'JointALEMoving2D' has the following parameters:

Name	type	size	default value	description
name	String		"	constraints's unique name
markerNumbers	ArrayMarkerIndex		[ MAXINT, MAX-INT ]	marker m0: position-marker of mass point or rigid body; marker m1: updated marker to ANCF Cable2D element, where the sliding joint currently is attached to; must be initialized with an appropriate (global) marker number according to the starting position of the sliding object; this marker changes with time (PostNewtonStep)
slidingMarkerNumbers	ArrayMarkerIndex		[]	a list of sn (global) marker numbers which are used to update marker1
slidingMarkerOffsets	Vector		[]	this list contains the offsets of every sliding object (given by slidingMarkerNumbers) w.r.t. to the initial position (0): marker0: offset=0, marker1: offset=Length(cable0), marker2: offset=Length(cable0)+Length(cable1), ...
slidingOffset	Real		0.	sliding offset list [SI:m]: a list of sn scalar offsets, which represent the (reference arc) length of all previous sliding cable elements
nodeNumbers	ArrayNodeIndex		[ MAXINT, MAX-INT ]	node number of NodeGenericData (GD) with one data coordinate and of NodeGenericODE2 (ALE) with one ODE2 coordinate
usePenaltyFormulation	bool		False	flag, which determines, if the connector is formulated with penalty, but still using algebraic equations (IsPenaltyConnector() still false)
penaltyStiffness	Real		0.	penalty stiffness [SI:N/m] used if usePenaltyFormulation=True
activeConnector	bool		True	flag, which determines, if the connector is active; used to deactivate (temporarily) a connector or constraint
visualization	VObjectJointALEMoving2D			parameters for visualization of item

The item VObjectJointALEMoving2D has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
drawSize	float		-1.	drawing size = radius of revolute joint; size == -1.f means that default connector size is used
color	Float4		[-1.,-1.,-1.,-1.]	RGBA connector color; if R== -1, use default color

Detailed information on ObjectJointALEMoving2D:

input parameter	symbol	description see tables above
markerNumbers	$[m0, m1]^T$	
slidingMarkerNumbers	$[m_{s0}, \dots, m_{sn}]^T$	
slidingMarkerOffsets	$[d_{s0}, \dots, d_{sn}]$	
slidingOffset	$s_{off}$	
nodeNumbers	$[n_{GD}, n_{ALE}]$	
penaltyStiffness	$k$	

The following output parameters are available as OutputVariableType in sensors and other functions:

output parameter	symbol	description
Position	${}^0\mathbf{p}_{m0}$	current global position of position marker $m0$
Velocity	${}^0\mathbf{v}_{m0}$	current global velocity of position marker $m0$
SlidingCoordinate	$s_g = q_{ALE} + s_{off}$	current value of the global sliding ALE coordinate, including offset; note that reference coordinate of $q_{ALE}$ is ignored!
Coordinates	$[x_{data0}, q_{ALE}]^T$	provides two values: [0] = current sliding marker index, [1] = ALE sliding coordinate
Coordinates_t	$[\dot{q}_{ALE}]^T$	provides ALE sliding velocity
Force	$\mathbf{f}$	joint force vector (3D)

Description of Item:

intermediate variables	symbol	description
generic data node	$\mathbf{x} = [x_{data0}]^T$	coordinates of node with node number $n_{GD}$
generic ODE2 node	$\mathbf{q} = [q_0]^T$	coordinates of node with node number $n_{ALE}$ , which is shared with all ALE-ANCF and ALE sliding joint objects
data coordinate	$x_{data0}$	the current index in slidingMarkerNumbers
ALE coordinate	$q_{ALE} = q_0$	current ALE coordinate (in fact this is the Eulerian coordinate in the ALE formulation); note that reference coordinate of $q_{ALE}$ is ignored!

marker m0 position	${}^0\mathbf{p}_{m0}$	current global position which is provided by marker m0
marker m0 velocity	${}^0\mathbf{v}_{m0}$	current global velocity which is provided by marker m0
cable coordinates	$\mathbf{q}_{ANCF,m1}$	current coordiantes of the ANCF cable element with the current marker $m1$ is referring to
sliding position	${}^0\mathbf{r}_{ANCF} = \mathbf{S}(s_{el})\mathbf{q}_{ANCF,m1}$	current global position at the ANCF cable element, evaluated at local sliding position $s_{el}$
sliding position slope	${}^0\mathbf{r}'_{ANCF} = \mathbf{S}'(s_{el})\mathbf{q}_{ANCF,m1}$	current global slope vector of the ANCF cable element, evaluated at local sliding position $s_{el}$
sliding velocity	${}^0\mathbf{v}_{ANCF} = \mathbf{S}(s_{el})\dot{\mathbf{q}}_{ANCF,m1} + \dot{q}_{ALE} {}^0\mathbf{r}'_{ANCF}$	current global velocity at the ANCF cable element, evaluated at local sliding position $s_{el}$ , including convective term
sliding normal vector	${}^0\mathbf{n} = [-r'_1, r'_0]$	2D normal vector computed from slope $\mathbf{r}' = {}^0\mathbf{r}'_{ANCF}$
algebraic variables	$\mathbf{z} = [\lambda_0, \lambda_1]^T$	algebraic variables (Lagrange multipliers) according to the algebraic equations

The element sliding coordinate (in the local coordinates of the current sliding element) is computed from the ALE coordinate

$$s_{el} = q_{ALE} + s_{off} - d_{m1} = s_g - d_{m1}. \quad (6.149)$$

The vector (=difference; error) between the marker  $m0$  and the marker  $m1$  ( $=\mathbf{r}_{ANCF}$ ) positions reads

$${}^0\Delta\mathbf{p} = {}^0\mathbf{r}_{ANCF} - {}^0\mathbf{p}_{m0} \quad (6.150)$$

The vector (=difference; error) between the marker  $m0$  and the marker  $m1$  velocities reads

$${}^0\Delta\mathbf{v} = {}^0\mathbf{v}_{ANCF} - {}^0\mathbf{v}_{m0} \quad (6.151)$$

**Algebraic constraint equations:** The 2D sliding joint is implemented having 2 equations, using the Lagrange multipliers  $\mathbf{z}$ . The algebraic (index 3) equations read

$${}^0\Delta\mathbf{p} = 0 \quad (6.152)$$

Note that the Lagrange multipliers  $[\lambda_0, \lambda_1]^T$  are the global forces in the joint. In the index 2 case the algebraic equations read

$${}^0\Delta\mathbf{v} = 0 \quad (6.153)$$

If `usePenalty = True`, the algebraic equations are changed to:

$${}^0\Delta\mathbf{p} - \frac{1}{k}\mathbf{z} = 0. \quad (6.154)$$

If `activeConnector = False`, the algebraic equations are changed to:

$$\lambda_0 = 0, \quad (6.155)$$

$$\lambda_1 = 0. \quad (6.156)$$

**Post Newton Step:** After the Newton solver has converged, a `PostNewtonStep` is performed for the element, which updates the marker  $m1$  index if necessary.

$$\begin{aligned} s_{el} < 0 &\rightarrow x_{data0} -= 1 \\ s_{el} > L &\rightarrow x_{data0} += 1 \end{aligned} \quad (6.157)$$

Furthermore, it is checked, if  $x_{data0}$  becomes smaller than zero, which raises a warning and keeps  $x_{data0} = 0$ . The same results if  $x_{data0} \geq sn$ , then  $x_{data0} = sn$ . Finally, the data coordinate is updated in order to provide the starting value for the next step,

$$x_{data1} += s. \quad (6.158)$$

**Examples:** see TestModels!

## 6.4 Markers

### 6.4.1 MarkerBodyMass

A marker attached to the body mass; use this marker to apply a body-load (e.g. gravitational force).

**Additional information for MarkerBodyMass:**

- The Marker has the following types = Object, Body, BodyMass

The item **MarkerBodyMass** with type = 'BodyMass' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body number to which marker is attached to
visualization	VMarkerBodyMass			parameters for visualization of item

The item **VMarkerBodyMass** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.4.2 MarkerBodyPosition

A position body-marker attached to local position (x,y,z) of the body.

**Additional information for MarkerBodyPosition:**

- The Marker has the following types = Object, Body, Position

The item **MarkerBodyPosition** with type = 'BodyPosition' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body number to which marker is attached to
localPosition	Vector3D	3	[0.,0.,0.]	local body position of marker; e.g. local (body-fixed) position where force is applied to
visualization	VMarkerBodyPosition			parameters for visualization of item

The item VMarkerBodyPosition has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.4.3 MarkerBodyRigid

A rigid-body (position+orientation) body-marker attached to local position (x,y,z) of the body.

**Additional information for MarkerBodyRigid:**

- The Marker has the following types = Object, Body, Position, Orientation

The item **MarkerBodyRigid** with type = 'BodyRigid' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body number to which marker is attached to
localPosition	Vector3D	3	[0.,0.,0.]	local body position of marker; e.g. local (body-fixed) position where force is applied to
visualization	VMarkerBodyRigid			parameters for visualization of item

The item VMarkerBodyRigid has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.4.4 MarkerNodePosition

A node-Marker attached to a position-based node.

**Additional information for MarkerNodePosition:**

- The Marker has the following types = Node, Position

The item **MarkerNodePosition** with type = 'NodePosition' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
nodeNumber	NodeIndex		MAXINT	node number to which marker is attached to
visualization	VMarkerNodePosition			parameters for visualization of item

The item VMarkerNodePosition has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown



### 6.4.5 MarkerNodeRigid

A rigid-body (position+orientation) node-marker attached to a rigid-body node.

**Additional information for MarkerNodeRigid:**

- The Marker has the following types = Node, Position, Orientation

The item **MarkerNodeRigid** with type = 'NodeRigid' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
nodeNumber	NodeIndex		MAXINT	node number to which marker is attached to
visualization	VMarkerNodeRigid			parameters for visualization of item

The item VMarkerNodeRigid has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.4.6 MarkerNodeCoordinate

A node-Marker attached to a ODE2 coordinate of a node; for other coordinates (ODE1,...) other markers need to be defined.

**Additional information for MarkerNodeCoordinate:**

- The Marker has the following types = Node, Coordinate

The item **MarkerNodeCoordinate** with type = 'NodeCoordinate' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
nodeNumber	NodeIndex		MAXINT	node number to which marker is attached to
coordinate	Index		MAXINT	coordinate of node to which marker is attached to
visualization	VMarkerNodeCoordinate			parameters for visualization of item

The item VMarkerNodeCoordinate has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.4.7 MarkerNodeRotationCoordinate

A node-Marker attached to a node containing rotation; the Marker measures a rotation coordinate (Tait-Bryan angles) or angular velocities on the velocity level

**Additional information for MarkerNodeRotationCoordinate:**

- The Marker has the following types = Node, Orientation, Coordinate

The item **MarkerNodeRotationCoordinate** with type = 'NodeRotationCoordinate' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
nodeNumber	NodeIndex		MAXINT	node number to which marker is attached to
rotationCoordinate	Index		MAXINT	rotation coordinate: 0=x, 1=y, 2=z
visualization	VMarkerNodeRotationCoordinate			parameters for visualization of item

The item VMarkerNodeRotationCoordinate has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

## 6.4.8 MarkerSuperElementPosition

A position marker attached to a SuperElement, such as ObjectFFRF, ObjectGenericODE2 and ObjectFFRFreducedOrder (for which it is inefficient for large number of meshNodeNumbers). The marker acts on the mesh (interface) nodes, not on the underlying nodes of the object.

**Additional information for MarkerSuperElementPosition:**

- The Marker has the following types = Object, Body, Position

The item **MarkerSuperElementPosition** with type = 'SuperElementPosition' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body number to which marker is attached to
meshNodeNumbers	ArrayIndex		[]	a list of $n_m$ mesh node numbers of superelement (=interface nodes) which are used to compute the body-fixed marker position; the related nodes must provide 3D position information, such as NodePoint, NodePoint2D, NodeRigidBody[..]; in order to retrieve the global node number, the generic body needs to convert local into global node numbers
weightingFactors	Vector		[]	a list of $n_m$ weighting factors per node to compute the final local position
visualization	VMarkerSuperElementPosition			parameters for visualization of item

The item VMarkerSuperElementPosition has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
showMarkerNodes	bool		True	set true, if all nodes are shown (similar to marker, but with less intensity)

**Detailed information on MarkerSuperElementPosition:**

input parameter	symbol	description see tables above
bodyNumber	$n_b$	
meshNodeNumbers	$[k_0, \dots, k_{n_m-1}]^T$	
weightingFactors	$[w_0, \dots, w_{n_m-1}]^T$	

**Description of Item:**

**Definition of marker quantities:**

intermediate variables	symbol	description
------------------------	--------	-------------

number of mesh nodes	$n_m$	size of meshNodeNumbers and weightingFactors which are marked; this must not be the number of mesh nodes in the marked object
mesh node number	$i = k_i$	abbreviation
mesh node points	${}^0\mathbf{p}_i$	position of mesh node $k_i$ in object $n_b$
mesh node velocities	${}^0\mathbf{v}_i$	velocity of mesh node $i$ in object $n_b$
marker position	${}^0\mathbf{p}_m = \sum_{i=0}^{n-1} w_i \cdot {}^0\mathbf{p}_i$	current global position which is provided by marker
marker velocity	${}^0\mathbf{v}_m = \sum_{i=0}^{n-1} w_i \cdot {}^0\mathbf{v}_i$	current global velocity which is provided by marker

### Marker quantities:

The marker provides a 'position' jacobian, which is the derivative of the marker velocity w.r.t. the object velocity coordinates  $\dot{\mathbf{q}}_{n_b}$ ,

$$\mathbf{J}_{m,pos} = \frac{\partial {}^0\mathbf{v}_m}{\partial \dot{\mathbf{q}}_{n_b}} = \sum_{i=0}^{n-1} w_i \cdot \mathbf{J}_{i,pos} \quad (6.159)$$

in which  $\mathbf{J}_{i,pos}$  denotes the position jacobian of mesh node  $i$ ,

$$\mathbf{J}_{i,pos} = \frac{\partial {}^0\mathbf{v}_i}{\partial \dot{\mathbf{q}}_{n_b}} \quad (6.160)$$

The jacobian  $\mathbf{J}_{i,pos}$  usually contains mostly zeros for ObjectGenericODE2, because the jacobian only affects one single node. In ObjectFFRFReducedOrder, the jacobian may affect all reduced coordinates.

Note that  $\mathbf{J}_{m,pos}$  is actually computed by the ObjectSuperElement within the function GetAccessFunctionSuperElement.

**Example** for MarkerSuperElementPosition:

---

```

#set up a mechanical system with two nodes; it has the structure: |~~M0~~M1
#==>further examples see objectGenericODE2Test.py, objectFFRFTest2.py, etc.
nMass0 = mbs.AddNode(NodePoint(referenceCoordinates=[0,0,0]))
nMass1 = mbs.AddNode(NodePoint(referenceCoordinates=[1,0,0]))
mGround = mbs.AddMarker(MarkerBodyPosition(bodyNumber=oGround, localPosition =
    [1,0,0]))

mass = 0.5 * np.eye(3)      #mass of nodes
stif = 5000 * np.eye(3)    #stiffness of nodes
damp = 50 * np.eye(3)      #damping of nodes
Z = 0. * np.eye(3)         #matrix with zeros
#build mass, stiffness and damping matrices (:
M = np.block([[mass,      0.*np.eye(3)],
               [0.*np.eye(3), mass      ] ])
K = np.block([[2*stif, -stif],
               [-stif,  stif] ])
D = np.block([[2*damp, -damp],
               [-damp,  damp] ])

oGenericODE2 = mbs.AddObject(ObjectGenericODE2(nodeNumbers=[nMass0,nMass1],
                                                massMatrix=M,
                                                stiffnessMatrix=K,
```

```

dampingMatrix=D))

#EXAMPLE for single node marker on super element body, mesh node 1; compare results
to ObjectGenericODE2 example!!!
mSuperElement = mbs.AddMarker(MarkerSuperElementPosition(bodyNumber=oGenericODE2,
    meshNodeNumbers=[1], weightingFactors=[1]))
mbs.AddLoad(Force(markerNumber = mSuperElement, loadVector = [10, 0, 0]))

#assemble and solve system for default parameters
mbs.Assemble()

sims=exu.SimulationSettings()
sims.timeIntegration.generalizedAlpha.spectralRadius=1
SC.TimeIntegrationSolve(mbs, 'GeneralizedAlpha', sims)

#check result at default integration time
testError = mbs.GetNodeOutput(nMass1, exu.OutputVariableType.Position)[0] -
    (1.0039999999354785)

```

---

## 6.4.9 MarkerSuperElementRigid

A position and orientation (rigid-body) marker attached to a SuperElement, such as ObjectFFRF, ObjectGenericODE2 and ObjectFFRFreducedOrder (for which it may be inefficient). The marker acts on the mesh nodes, not on the underlying nodes of the object. Note that in contrast to the MarkerSuperElementPosition, this marker needs a set of interface nodes which are not aligned at one line, such that they can represent rigid body motion. Note that definitions of marker positions are slightly different from MarkerSuperElementPosition.

### Additional information for MarkerSuperElementRigid:

- The Marker has the following types = Object, Body, Position, Orientation

The item **MarkerSuperElementRigid** with type = 'SuperElementRigid' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body number to which marker is attached to
referencePosition	Vector3D	3	[0.,0.,0.]	local marker SuperElement reference position used to compute average displacement and average rotation; currently, this must be the center of weighted nodes of the marker
meshNodeNumbers	ArrayIndex		[]	a list of $n_m$ mesh node numbers of superelement (=interface nodes) which are used to compute the body-fixed marker position and orientation; the related nodes must provide 3D position information, such as NodePoint, NodePoint2D, NodeRigidBody[.]; in order to retrieve the global node number, the generic body needs to convert local into global node numbers
weightingFactors	Vector		[]	a list of $n_m$ weighting factors per node to compute the final local position and orientation; these factors could be based on surface integrals of the constrained mesh faces
visualization	VMarkerSuperElementRigid			parameters for visualization of item

The item VMarkerSuperElementRigid has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown
showMarkerNodes	bool		True	set true, if all nodes are shown (similar to marker, but with less intensity)

### Detailed information on MarkerSuperElementRigid:

input parameter	symbol	description see tables above
bodyNumber	$n_b$	

referencePosition	${}^r\mathbf{p}_{0,ref}$	
meshNodeNumbers	$[k_0, \dots, k_{n_m-1}]^T$	
weightingFactors	$[w_0, \dots, w_{n_m-1}]^T$	

**Description of Item:**

**Definition of marker quantities:**

intermediate variables	symbol	description
number of mesh nodes	$n_m$	size of meshNodeNumbers and weightingFactors which are marked; this must not be the number of mesh nodes in the marked object
mesh node number	$i = k_i$	abbreviation
mesh node local position	${}^r\mathbf{p}_i$	current local (within reference frame $r$ ) position of mesh node $k_i$ in object $n_b$
mesh node local reference position	${}^r\mathbf{p}_{ref,i}$	local (within reference frame $r$ ) reference position of mesh node $k_i$ in object $n_b$
mesh node local displacement	${}^r\mathbf{u}_i$	current local (within reference frame $r$ ) displacement of mesh node $k_i$ in object $n_b$
mesh node local velocity	${}^r\mathbf{v}_i$	current local (within reference frame $r$ ) velocity of mesh node $k_i$ in object $n_b$
super element reference position	${}^0\mathbf{p}_r$	current reference position of super element's floating frame (r), which is zero, if the object does not provide a reference frame (such as GenericODE2)
super element rotation matrix	${}^{0r}\mathbf{A}$	current rigid body transformation matrix of super element's floating frame (r), which is the identity matrix, if the object does not provide a reference frame (such as GenericODE2)
super element angular velocity	${}^r\boldsymbol{\omega}_r$	current local angular velocity of super element's floating frame (r), which is zero, if the object does not provide a reference frame (such as GenericODE2)
marker reference position	${}^0\mathbf{p}_0 = {}^0\mathbf{p}_r + {}^{0r}\mathbf{A} {}^r\mathbf{p}_{0,ref}$	current global marker reference position; note that ${}^0\mathbf{p}_0 = {}^{0r}\mathbf{I} {}^r\mathbf{p}_{0,ref}$ , if the object does not provide a reference frame (such as GenericODE2)
marker position	${}^0\mathbf{p}_m = {}^0\mathbf{p}_0 + {}^{0r}\mathbf{A} \sum_{i=0}^{n-1} w_i \cdot {}^r\mathbf{u}_i$	current global position which is provided by marker
marker velocity	${}^0\mathbf{v}_m = {}^0\dot{\mathbf{p}}_r + {}^{0r}\mathbf{A} {}^r\boldsymbol{\omega}_r {}^r\mathbf{p}_{0,ref} + {}^{0r}\mathbf{A} \left( \sum_{i=0}^{n-1} (w_i \cdot {}^r\mathbf{v}_i) + {}^r\boldsymbol{\omega}_r \sum_{i=0}^{n-1} (w_i {}^r\mathbf{u}_i) \right)$	current global velocity which is provided by marker
marker local rotation	${}^r\boldsymbol{\theta}_m = \frac{\sum_{i=0}^{n-1} w_i {}^r\mathbf{p}_{ref,i} \times {}^r\mathbf{u}_i}{\sum_{i=0}^{n-1} w_i  {}^r\mathbf{p}_{i,ref} ^2}$	current local (within reference frame $r$ ) linearized rotation parameters
marker rotation matrix	${}^{0b}\mathbf{A}_m = {}^{0r}\mathbf{A} \begin{bmatrix} 1 & -\theta_2 & \theta_1 \\ \theta_2 & 1 & -\theta_0 \\ -\theta_1 & \theta_0 & 1 \end{bmatrix}$	current rotation matrix, which transforms the local marker coordinates and adds the rigid body transformation of floating frames ${}^{0r}\mathbf{A}$ ; only valid for small (linearized rotations)!



marker local angular velocity	${}^r\omega_m = {}^r\dot{\theta}_m = \frac{\sum_{i=0}^{n-1} w_i {}^r\tilde{\mathbf{p}}_{ref,i} {}^r\mathbf{v}_i}{\sum_{i=0}^{n-1} w_i  {}^r\tilde{\mathbf{p}}_{i,ref} ^2}$	local (within reference frame $r$ ) angular velocity due to mesh node velocity only
marker inertial angular velocity	${}^0\omega_m = {}^0\omega_r + {}^{0r}\mathbf{A} {}^r\omega_m$	current inertial angular velocity

#### Marker quantities:

The marker provides a 'position' jacobian, which is the derivative of the marker velocity w.r.t. the object velocity coordinates  $\dot{\mathbf{q}}_{n_b}$ ,

$$\mathbf{J}_{m,pos} = \frac{\partial {}^0\mathbf{v}_m}{\partial \dot{\mathbf{q}}_{n_b}} = \dots \quad (6.161)$$

In `ObjectFFRFRducedOrder`, the jacobian may affect all reduced coordinates.

The marker also provides a 'rotation' jacobian, which is the derivative of the marker angular velocity  ${}^0\omega_m$  w.r.t. the object velocity coordinates  $\dot{\mathbf{q}}_{n_b}$ ,

$$\mathbf{J}_{m,rot} = \frac{\partial {}^0\omega_m}{\partial \dot{\mathbf{q}}_{n_b}} = \frac{\partial {}^{0r}\mathbf{A} ({}^r\omega_r + {}^r\omega_m)}{\partial \dot{\mathbf{q}}_{n_b}} = {}^{0r}\mathbf{A} (\mathbf{G}_{local} + \frac{\sum_{i=0}^{n-1} w_i {}^r\tilde{\mathbf{p}}_{ref,i} \mathbf{J}_{i,pos}}{\sum_{i=0}^{n-1} w_i |{}^r\tilde{\mathbf{p}}_{i,ref}|^2}) \quad (6.162)$$

with the matrix  $\mathbf{G}_{local} = \frac{\partial {}^r\omega_r}{\partial \dot{\mathbf{q}}_{n_b}}$ .

#### Alternative computation of rotation (under development):

In the alternative mode, where the weighting matrix  $\mathbf{W}$  has the interpretation of an inertia tensor of built from nodes using weights = node mass, and the local angular momentum reads

$$\mathbf{W} {}^r\omega_m = \sum_{i=0}^{n-1} w_i {}^r\tilde{\mathbf{p}}_{ref,i} {}^r\mathbf{v}_i = - \sum_{i=0}^{n-1} (w_i {}^r\tilde{\mathbf{p}}_{i,ref} {}^r\tilde{\mathbf{p}}_{i,ref}) {}^r\omega_m \quad (6.163)$$

and the marker local rotations and marker local angular velocity are defined as

$${}^r\theta_m = \mathbf{W}^{-1} \sum_{i=0}^{n-1} w_i {}^r\tilde{\mathbf{p}}_{ref,i} {}^r\mathbf{u}_i, \quad {}^r\omega_m = \mathbf{W}^{-1} \sum_{i=0}^{n-1} w_i {}^r\tilde{\mathbf{p}}_{ref,i} {}^r\mathbf{v}_i, \quad (6.164)$$

with the weighting matrix (which must be invertable)

$$\mathbf{W} = - \sum_{i=0}^{n-1} w_i {}^r\tilde{\mathbf{p}}_{i,ref} {}^r\tilde{\mathbf{p}}_{i,ref} \quad (6.165)$$

and in the alternative mode, the angular velocity is defined as

$$\mathbf{J}_{m,rot} = \frac{\partial {}^0\omega_m}{\partial \dot{\mathbf{q}}_{n_b}} = \frac{\partial {}^{0r}\mathbf{A} ({}^r\omega_r + {}^r\omega_m)}{\partial \dot{\mathbf{q}}_{n_b}} = {}^{0r}\mathbf{A} (\mathbf{G}_{local} + \mathbf{W}^{-1} \sum_{i=0}^{n-1} w_i {}^r\tilde{\mathbf{p}}_{ref,i} \mathbf{J}_{i,pos}) \quad (6.166)$$

Note that  $\mathbf{J}_{m,rot}$  is computed by the `ObjectSuperElement` within the function `GetAccessFunctionSuperElement`.

#### EXAMPLE for marker on body 4, mesh nodes 10,11,12,13:

`MarkerSuperElementRigid(bodyNumber = 4, meshNodeNumber = [10, 11, 12, 13], weightingFactors = [0.25, 0.25, 0.25, 0.25], referencePosition=[0,0,0])`

For detailed examples, see `TestModels`.

### 6.4.10 MarkerObjectODE2Coordinates

A Marker attached to all coordinates of an object (currently only body is possible), e.g. to apply special constraints or loads on all coordinates. The measured coordinates INCLUDE reference + current coordinates.

**Additional information for MarkerObjectODE2Coordinates:**

- The Marker has the following types = Object, Body, Coordinate

The item **MarkerObjectODE2Coordinates** with type = 'ObjectODE2Coordinates' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
objectNumber	ObjectIndex		MAXINT	body number to which marker is attached to
visualization	VMarkerObjectODE2Coordinates			parameters for visualization of item

The item **VMarkerObjectODE2Coordinates** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.4.11 MarkerBodyCable2DShape

A special Marker attached to a 2D ANCF beam finite element with cubic interpolation and 8 coordinates.

**Additional information for MarkerBodyCable2DShape:**

- The Marker has the following types = Object, Body, Coordinate

The item **MarkerBodyCable2DShape** with type = 'BodyCable2DShape' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body number to which marker is attached to
numberOfSegments	Index		3	number of number of segments; each segment is a line and is associated to a data (history) variable; must be same as in according contact element
visualization	VMarkerBodyCable2DShape			parameters for visualization of item

The item VMarkerBodyCable2DShape has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.4.12 MarkerBodyCable2DCoordinates

A special Marker attached to the coordinates of a 2D ANCF beam finite element with cubic interpolation.

**Additional information for MarkerBodyCable2DCoordinates:**

- The Marker has the following types = Object, Body, Coordinate

The item **MarkerBodyCable2DCoordinates** with type = 'BodyCable2DCoordinates' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body number to which marker is attached to
visualization	VMarkerBodyCable2DCoordinates			parameters for visualization of item

The item VMarkerBodyCable2DCoordinates has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

## 6.5 Loads

### 6.5.1 LoadForceVector

Load with (3D) force vector; attached to position-based marker.

**Additional information for LoadForceVector:**

- Requested marker type = **Position**
- **Short name** for Python = **Force**
- **Short name** for Python (visualization object) = **VForce**

The item **LoadForceVector** with type = 'ForceVector' has the following parameters:

Name	type	size	default value	description
name	String		"	load's unique name
markerNumber	MarkerIndex		MAXINT	marker's number to which load is applied
loadVector	Vector3D		[0.,0.,0.]	vector-valued load [SI:N]
bodyFixed	Bool		False	if bodyFixed is true, the load is defined in body-fixed (local) coordinates, leading to a follower force; if false: global coordinates are used
loadVectorUserFunction	PyFunctionVector3DScalar	Vector3D	0	A python function which defines the time-dependent load with parameters (Real t, Vector3D load); the load represents the current value of the load; WARNING: this factor does not work in combination with static computation (loadFactor); Example for python function: def f(t, loadVector): return [loadVector[0]*np.sin(t*10*2*3.1415),0,0]
visualization	VLoadForceVector			parameters for visualization of item

The item **VLoadForceVector** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

## 6.5.2 LoadTorqueVector

Load with (3D) torque vector; attached to rigidbody-based marker.

**Additional information for LoadTorqueVector:**

- Requested marker type = Orientation
- **Short name** for Python = **Torque**
- **Short name** for Python (visualization object) = **VTorque**

The item **LoadTorqueVector** with type = 'TorqueVector' has the following parameters:

Name	type	size	default value	description
name	String		"	load's unique name
markerNumber	MarkerIndex		MAXINT	marker's number to which load is applied
loadVector	Vector3D		[0.,0.,0.]	vector-valued load [SI:N]
bodyFixed	Bool		False	if bodyFixed is true, the load is defined in body-fixed (local) coordinates, leading to a follower torque; if false: global coordinates are used
loadVectorUserFunction	PyFunctionVector3DScalar		Vector3D 0	A python function which defines the time-dependent load with parameters (Real t, Vector3D load); the load represents the current value of the load; WARNING: this factor does not work in combination with static computation (load-Factor); Example for python function: def f(t, loadVector): return [loadVector[0]*np.sin(t*10*2*3.1415),0,0]
visualization	VLoadTorqueVector			parameters for visualization of item

The item **VLoadTorqueVector** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.5.3 LoadMassProportional

Load attached to BodyMass-based marker, applying a 3D vector load (e.g. the vector [0,-g,0] is used to apply gravitational loading of size g in negative y-direction).

**Additional information for LoadMassProportional:**

- Requested marker type = Body + BodyMass
- **Short name** for Python = **Gravity**
- **Short name** for Python (visualization object) = **VGravity**

The item **LoadMassProportional** with type = 'MassProportional' has the following parameters:

Name	type	size	default value	description
name	String		"	load's unique name
markerNumber	MarkerIndex		MAXINT	marker's number to which load is applied
loadVector	Vector3D		[0.,0.,0.]	vector-valued load [SI:N/kg = m/s <sup>2</sup> ]
loadVectorUserFunction	PyFunctionVector3DScalar		Vector3D 0	A python function which defines the time-dependent load with parameters (Real t, Vector3D load); the load represents the current value of the load; WARNING: this factor does not work in combination with static computation (load-Factor); Example for python function: def f(t, loadVector): return [loadVector[0]*np.sin(t*10*2*3.1415),0,0]
visualization	VLoadMassProportional			parameters for visualization of item

The item **VLoadMassProportional** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

## 6.5.4 LoadCoordinate

Load with scalar value, which is attached to a coordinate-based marker; the load can be used e.g. to apply a force to a single axis of a body, a nodal coordinate of a finite element or a torque to the rotatory DOF of a rigid body.

### Additional information for LoadCoordinate:

- Requested marker type = Coordinate

The item **LoadCoordinate** with type = 'Coordinate' has the following parameters:

Name	type	size	default value	description
name	String		"	load's unique name
markerNumber	MarkerIndex		MAXINT	marker's number to which load is applied
load	Real		0.	scalar load [SI:N]
loadUserFunction	PyFunctionScalar2		0	A python function which defines the time-dependent load with parameters (Real t, Real load); the load represents the current value of the load; WARNING: this factor does not work in combination with static computation (loadFactor); Example for python function: <code>def f(t, load): return load*np.sin(t*10*2*3.1415)</code>
visualization	VLoadCoordinate			parameters for visualization of item

The item **VLoadCoordinate** has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown



## 6.6 Sensors

### 6.6.1 SensorNode

A sensor attached to a node. The sensor measures OutputVariables and outputs values into a file, showing per line [time, sensorValue[0], sensorValue[1], ...]. A user function can be attached to modify sensor values accordingly.

The item **SensorNode** with type = 'Node' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
nodeNumber	NodeIndex		MAXINT	node number to which sensor is attached to
writeToFile	bool		True	true: write sensor output to file
fileName	String		"	directory and file name for sensor file output; default: empty string generates sensor + sensorNumber + outputVariableType; directory will be created if it does not exist
outputVariableType	OutputVariableType		OutputVariableType:None	OutputVariableType for sensor
visualization	VSensorNode			parameters for visualization of item

The item VSensorNode has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

## 6.6.2 SensorObject

A sensor attached to any object except bodies (connectors, constraint, spring-damper, etc). As a difference to other SensorBody, the connector sensor measures quantities without a local position. The sensor measures OutputVariable and outputs values into a file, showing per line [time, sensorValue[0], sensorValue[1], ...]. A user function can be attached to postprocess sensor values accordingly.

The item **SensorObject** with type = 'Object' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
objectNumber	ObjectIndex		MAXINT	object (e.g. connector) number to which sensor is attached to
writeToFile	bool		True	true: write sensor output to file
fileName	String		"	directory and file name for sensor file output; default: empty string generates sensor + sensorNumber + outputVariableType; directory will be created if it does not exist
outputVariableType	OutputVariableType		OutputVariableType	OutputVariableType for sensor
visualization	VSensorObject			parameters for visualization of item

The item VSensorObject has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown; sensors can be shown at the position associated with the object - note that in some cases, there might be no such position (e.g. data object)!

### 6.6.3 SensorBody

A sensor attached to a body-object with local position. As a difference to other ObjectSensors, the body sensor has a local position at which the sensor is attached to. The sensor measures OutputVariableBody and outputs values into a file, showing per line [time, sensorValue[0], sensorValue[1], ...]. A user function can be attached to postprocess sensor values accordingly.

The item **SensorBody** with type = 'Body' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body (=object) number to which sensor is attached to
localPosition	Vector3D	3	[0.,0.,0.]	local (body-fixed) body position of sensor
writeToFile	bool		True	true: write sensor output to file
fileName	String		"	directory and file name for sensor file output; default: empty string generates sensor + sensorNumber + outputVariableType; directory will be created if it does not exist
outputVariableType	OutputVariableType		OutputVariableType:None	OutputVariableType for sensor
visualization	VSensorBody			parameters for visualization of item

The item VSensorBody has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

## 6.6.4 SensorSuperElement

A sensor attached to a SuperElement-object with mesh node number. As a difference to other ObjectSensors, the SuperElement sensor has a mesh node number at which the sensor is attached to. The sensor measures OutputVariableSuperElement and outputs values into a file, showing per line [time, sensorValue[0], sensorValue[1], ...]. A user function can be attached to postprocess sensor values accordingly.

The item **SensorSuperElement** with type = 'SuperElement' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
bodyNumber	ObjectIndex		MAXINT	body (=object) number to which sensor is attached to
meshNodeNumber	Index		-1	mesh node number, which is a local node number with in the object (starting with 0); the node number may represent a real Node in mbs, or may be virtual and reconstructed from the object coordinates such as in ObjectFFRFRducedOrder
writeToFile	bool		True	true: write sensor output to file
fileName	String		"	directory and file name for sensor file output; default: empty string generates sensor + sensorNumber + outputVariableType; directory will be created if it does not exist
outputVariableType	OutputVariableType		OutputVariableType	OutputVariableType for sensor
visualization	VSensorSuperElement			parameters for visualization of item

The item VSensorSuperElement has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown

### 6.6.5 SensorLoad

A sensor attached to a load. The sensor measures the load values and outputs values into a file, showing per line [time, sensorValue[0], sensorValue[1], ...].

The item **SensorLoad** with type = 'Load' has the following parameters:

Name	type	size	default value	description
name	String		"	marker's unique name
loadNumber	LoadIndex		MAXINT	load number to which sensor is attached to
writeToFile	bool		True	true: write sensor output to file
fileName	String		"	directory and file name for sensor file output; default: empty string generates sensor + sensorNumber + outputVariableType; directory will be created if it does not exist
visualization	VSensorLoad			parameters for visualization of item

The item VSensorLoad has the following parameters:

Name	type	size	default value	description
show	bool		True	set true, if item is shown in visualization and false if it is not shown; CURRENTLY NOT AVAILABLE

## 6.7 GraphicsData

Some items may include a 'graphicsData' dictionary structure. GraphicsData dictionaries can be created with functions provided in the utilities module `exudynGraphicsDataUtilities.py`. GraphicsData contains a list of graphicsData items, i.e. `graphicsData = [graphicsItem1, graphicsItem2, ...]`. Every single graphicsItem may be defined as one of the following structures using a specific 'type':

Name	type	default value	description
<b>type = 'Line':</b>			<i>draws a polygonal line between all specified points</i>
color	list	[0,0,0,1]	list of 4 floats to define RGB-color and transparency
data	list	mandatory	list of float triples of x,y,z coordinates of the line floats to define RGB-color and transparency; Example: <code>data=[0,0,0, 1,0,0, 1,1,0, 0,1,0, 0,0,0]</code> ... draws a rectangle with side length 1
<b>type = 'Circle':</b>			<i>draws a circle with center point, normal (defines plane of circle) and radius</i>
color	list	[0,0,0,1]	list of 4 floats to define RGB-color and transparency
radius	float	mandatory	radius
position	list	mandatory	list of float triples of x,y,z coordinates of center point of the circle
normal	list	[0,0,1]	list of float triples of x,y,z coordinates of normal to the plane of the circle; the default value gives a circle in the (x, y)-plane
<b>type = 'Text':</b>			<i>places the given text at position</i>
color	list	[0,0,0,1]	list of 4 floats to define RGB-color and transparency
text	string	mandatory	text to be displayed
position	list	mandatory	list of float triples of [x,y,z] coordinates of the left upper position of the text; e.g. <code>position=[20,10,0]</code>
<b>type = 'TriangleList':</b>			<i>draws a flat triangle mesh for given points and connectivity</i>
points	list	mandatory	list <code>[x0,y0,z0, x1,y1,z1, ...]</code> containing $n \times 3$ floats (grouped <code>x0,y0,z0, x1,y1,z1, ...</code> ) to define x,y,z coordinates of points, $n$ being the number of points (=vertices)
colors	list	empty	list <code>[R0,G0,B0,A0, R1,G2,B1,A1, ...]</code> containing $n \times 4$ floats to define RGB-color and transparency A, where $n$ must be according to number of points; if field 'colors' does not exist, default colors will be used
normals	list	empty	list <code>[n0x,n0y,n0z, ...]</code> containing $n \times 3$ floats to define normal direction of triangles per point, where $n$ must be according to number of points; if field 'normals' does not exist, default normals <code>[0,0,0]</code> will be used
triangles	list	mandatory	list <code>[T0point0, T0point1, T0point2, ...]</code> containing $n_{trig} \times 3$ floats to define point indices of each vertex of the triangles (=connectivity); point indices start with index 0; the maximum index must be $\leq$ <code>points.size()</code>

Examples of GraphicsData can be found in the Python examples and in `exudynUtilities.py`.



## Chapter 7

# EXUDYN Settings

This section includes the reference manual for settings which are available in the python interface, e.g. simulation settings, visualization settings, and others.

### 7.1 Simulation settings

This section includes hierarchical structures for simulation settings, e.g., time integration, static solver, Newton iteration and solution file export.

#### 7.1.1 SolutionSettings

General settings for exporting the solution (results) of a simulation.

SolutionSettings has the following items:

Name	type/function return type	size	default value / function args	description
writeSolutionToFile	bool		True	flag (true/false), which determines if (global) solution vector is written to file
appendToFile	bool		False	flag (true/false); if true, solution and solver-Information is appended to existing file (otherwise created)
writeFileHeader	bool		True	flag (true/false); if true, file header is written (turn off, e.g. for multiple runs of time integration)
writeFileFooter	bool		True	flag (true/false); if true, information at end of simulation is written: convergence, total solution time, statistics
solutionWritePeriod	UReal		0.01	time span (period), determines how often the solution is written during a simulation
sensorsAppendToFile	bool		False	flag (true/false); if true, sensor output is appended to existing file (otherwise created)
sensorsWriteFileHeader	bool		True	flag (true/false); if true, file header is written for sensor output (turn off, e.g. for multiple runs of time integration)



sensorsWritePeriod	UReal		0.01	time span (period), determines how often the sensor output is written during a simulation
exportVelocities	bool		True	solution is written as displacements, velocities[, accelerations] [,algebraicCoordinates] [,DataCoordinates]
exportAccelerations	bool		True	solution is written as displacements, [velocities,] accelerations [,algebraicCoordinates] [,DataCoordinates]
exportAlgebraicCoordinates	bool		True	solution is written as displacements, [velocities,] [accelerations,], algebraicCoordinates (=Lagrange multipliers) [,DataCoordinates]
exportDataCoordinates	bool		True	solution is written as displacements, [velocities,] [accelerations,] [,algebraicCoordinates (=Lagrange multipliers)] ,DataCoordinates
coordinatesSolutionFileName	FileName		'coordinatesSolution.txt'	filename and (relative) path of solution file containing all coordinates versus time; directory will be created if it does not exist
solverInformationFileName	FileName		'solverInformation.txt'	filename and (relative) path of text file showing detailed information during solving; detail level according to yourSolver.verboseModeFile; if solutionSettings.appendToFile is true, the information is appended in every solution step; directory will be created if it does not exist
solutionInformation	String		"	special information added to header of solution file (e.g. parameters and settings, modes, ...)
outputPrecision	Index		10	precision for floating point numbers written to solution and sensor files
recordImagesInterval	Real		-1.	record frames (images) during solving; amount of time to wait until next image (frame) is recorded; set recordImages = -1. if no images shall be recorded; set, e.g., recordImages = 0.01 to record an image every 10 milliseconds (requires that the time steps / load steps are sufficiently small!); for file names, etc., see VisualizationSettings.exportImages

### 7.1.2 NumericalDifferentiationSettings

Settings for numerical differentiation of a function (needed for computation of numerical jacobian e.g. in implicit integration); HOTINT1:  $\text{relativeEpsilon} * \text{Maximum}(\text{minimumCoordinateSize}, \text{fabs}(x(i)))$ .

NumericalDifferentiationSettings has the following items:

Name	type/function return type	size	default value / function args	description
relativeEpsilon	UReal		1e-7	relative differentiation parameter epsilon; the numerical differentiation parameter $\varepsilon$ follows from the formula ( $\varepsilon = \varepsilon_{\text{relative}} * \max(q_{\min},  q_i + [q_i^{\text{Ref}}]  )$ ), with $\varepsilon_{\text{relative}} = \text{relativeEpsilon}$ , $q_{\min} = \text{minimumCoordinateSize}$ , $q_i$ is the current coordinate which is differentiated, and $q_{\text{Ref}_i}$ is the reference coordinate of the current coordinate
minimumCoordinateSize	UReal		1e-2	minimum size of coordinates in relative differentiation parameter
doSystemWideDifferentiation	bool		False	true: system wide differentiation (e.g. all ODE2 equations w.r.t. all ODE2 coordinates); false: only local (object) differentiation
addReferenceCoordinatesToEpsilon	bool		False	true: for the size estimation of the differentiation parameter, the reference coordinate $q_i^{\text{Ref}}$ is added to ODE2 coordinates -> see; false: only the current coordinate is used for size estimation of the differentiation parameter

### 7.1.3 NewtonSettings

Settings for Newton method used in static or dynamic simulation.

NewtonSettings has the following items:

Name	type/function return type	size	default value / function args	description
numericalDifferentiation	NumericalDifferentiationSettings			numerical differentiation parameters for numerical jacobian (e.g. Newton in static solver or implicit time integration)
useNumericalDifferentiation	bool		False	flag (true/false); false = perform direct computation of jacobian, true = use numerical differentiation for jacobian
useNewtonSolver	bool		True	flag (true/false); false = linear computation, true = use Newton solver for nonlinear solution
relativeTolerance	UReal		1e-8	relative tolerance of residual for Newton (general goal of Newton is to decrease the residual by this factor)

absoluteTolerance	UReal		1e-10	absolute tolerance of residual for Newton (needed e.g. if residual is fulfilled right at beginning); condition: $\sqrt{q \cdot q} / \text{numberOfCoordinates} \leq \text{absoluteTolerance}$
weightTolerancePerCoordinate	bool		False	flag (true/false); false = compute error as L2-Norm of residual; true = compute error as $(\text{L2-Norm of residual}) / (\sqrt{\text{number of coordinates}})$ , which can help to use common tolerance independent of system size
newtonResidualMode	Index		0	0 ... use residual for computation of error (standard); 1 ... use change of solution increment for error (set relTol and absTol to same values!) ==> may be advantageous if residual is zero, e.g., in kinematic analysis; TAKE CARE with this flag
adaptInitialResidual	bool		True	flag (true/false); false = standard; true: if initialResidual is very small (or zero), it may increase dramatically in first step; to achieve relativeTolerance, the initialResidual will be updated by a higher residual within the first Newton iteration
modifiedNewtonContractivity	UReal		0.5	maximum contractivity (=reduction of error in every Newton iteration) accepted by modified Newton; if contractivity is greater, a Jacobian update is computed
useModifiedNewton	bool		False	true: compute Jacobian only at first step; no Jacobian updates per step; false: Jacobian computed in every step
modifiedNewtonJacUpdatePerStep	bool		False	true: compute Jacobian at every time step, but not in every iteration (except for bad convergence ==> switch to full Newton)
maxIterations	Index		25	maximum number of iterations (including modified + restart Newton steps); after that iterations, the static/dynamic solver stops with error
maxModifiedNewtonIterations	Index		8	maximum number of iterations for modified Newton (without Jacobian update); after that number of iterations, the modified Newton method gets a Jacobian update and is further iterated
maxModifiedNewtonRestartIterations	Index		7	maximum number of iterations for modified Newton after a Jacobian update; after that number of iterations, the full Newton method is started for this step

maximumSolutionNorm	UReal		1e38	this is the maximum allowed value for solutionU.L2NormSquared() which is the square of the square norm (value= $u_1^2+u_2^2+\dots$ ), and solutionV/A...; if the norm of solution vectors are larger, Newton method is stopped; the default value is chosen such that it would still work for single precision numbers (float)
maxDiscontinuousIterations	Index		5	maximum number of discontinuous (post Newton) iterations
ignoreMaxDiscontinuousIterations	bool		True	continue solver if maximum number of discontinuous (post Newton) iterations is reached (ignore tolerance)
discontinuousIterationTolerance	UReal		1	absolute tolerance for discontinuous (post Newton) iterations; the errors represent absolute residuals and can be quite high
stepInformation	Index		2	0 ... only current step time, 1 ... show time to go, 2 ... show newton iterations (Nit) per step, 3 ... show discontinuous iterations (Dit) and newton jacobians (jac) per step

### 7.1.4 GeneralizedAlphaSettings

Settings for generalized-alpha, implicit trapezoidal or Newmark time integration methods.

GeneralizedAlphaSettings has the following items:

Name	type/function return type	size	default value / function args	description
newmarkBeta	UReal		0.25	value beta for Newmark method; default value beta = $\frac{1}{4}$ corresponds to (undamped) trapezoidal rule
newmarkGamma	UReal		0.5	value gamma for Newmark method; default value gamma = $\frac{1}{2}$ corresponds to (undamped) trapezoidal rule
useIndex2Constraints	bool		False	set useIndex2Constraints = true in order to use index2 (velocity level constraints) formulation
useNewmark	bool		False	if true, use Newmark method with beta and gamma instead of generalized-Alpha
spectralRadius	UReal		0.9	spectral radius for Generalized-alpha solver; set this value to 1 for no damping or to $0 < \text{spectralRadius} < 1$ for damping of high-frequency dynamics; for position-level constraints (index 3), spectralRadius must be $< 1$

computeInitialAccelerations	bool		True	true: compute initial accelerations from system EOM in acceleration form; NOTE that initial accelerations that are following from user functions in constraints are not considered for now! false: use zero accelerations
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### 7.1.5 TimeIntegrationSettings

General parameters used in time integration; specific parameters are provided in the according solver settings, e.g. for generalizedAlpha.

TimeIntegrationSettings has the following items:

Name	type/function return type	size	default value / function args	description
newton	NewtonSettings			parameters for Newton method; used for implicit time integration methods only
startTime	UReal		0	start time of time integration (usually set to zero)
endTime	UReal		1	end time of time integration
numberOfSteps	UInt		100	number of steps in time integration; stepsize is computed from (endTime-startTime)/numberOfSteps
adaptiveStep	bool		True	true: use step reduction if step fails; false: constant step size
minimumStepSize	UReal		1e-8	lower limit of time step size, before integrator stops
verboseMode	Index		0	0 ... no output, 1 ... show short step information every 2 seconds (error), 2 ... show every step information, 3 ... show also solution vector, 4 ... show also mass matrix and jacobian (implicit methods), 5 ... show also Jacobian inverse (implicit methods)
verboseModeFile	Index		0	same behaviour as verboseMode, but outputs all solver information to file
generalizedAlpha	GeneralizedAlphaSettings			parameters for generalized-alpha, implicit trapezoidal rule or Newmark (options only apply for these methods)
preStepPyExecute	String		"	DEPRECATED, use preStepFunction in simulation settings; Python code to be executed prior to every step and after last step, e.g. for postprocessing

## 7.1.6 StaticSolverSettings

Settings for static solver linear or nonlinear (Newton).

StaticSolverSettings has the following items:

Name	type/function return type	size	default value / function args	description
newton	NewtonSettings			parameters for Newton method (e.g. in static solver or time integration)
numberOfLoadSteps	Index		1	number of load steps; if numberOfLoadSteps=1, no load steps are used and full forces are applied at once
loadStepDuration	UReal		1	quasi-time for all load steps (added to current time in load steps)
loadStepStart	UReal		0	a quasi time, which can be used for the output (first column) as well as for time-dependent forces; quasi-time is increased in every step $i$ by $\text{loadStepDuration}/\text{numberOfLoadSteps}$ ; $\text{loadStepTime} = \text{loadStepStart} + i * \text{loadStepDuration}/\text{numberOfLoadSteps}$ , but $\text{loadStepStart}$ untouched ==> increment by user
loadStepGeometric	bool		False	if $\text{loadStepGeometric}=\text{false}$ , the load steps are incremental (arithmetic series, e.g. 0.1,0.2,0.3,...); if true, the load steps are increased in a geometric series, e.g. for $n = 8$ numberOfLoadSteps and $d = 1000$ loadStepGeometricRange, it follows: $1000^{1/8}/1000 = 0.00237$ , $1000^{2/8}/1000 = 0.00562$ , $1000^{3/8}/1000 = 0.0133$ , ..., $1000^{7/8}/1000 = 0.422$ , $1000^{8/8}/1000 = 1$
loadStepGeometricRange	UReal		1000	if $\text{loadStepGeometric}=\text{true}$ , the load steps are increased in a geometric series, see loadStepGeometric
useLoadFactor	bool		True	true: compute a load factor $\in [0,1]$ from static step time; all loads are scaled by the load factor; false: loads are always scaled with 1 – use this option if time dependent loads use a userFunction
stabilizerODE2term	UReal		0	add mass-proportional stabilizer term in ODE2 part of jacobian for stabilization (scaled), e.g. of badly conditioned problems; the diagonal terms are scaled with $\text{stabilizer} = (1 - \text{loadStepFactor}^2)$ , and go to zero at the end of all load steps: $\text{loadStepFactor} = 1 \rightarrow \text{stabilizer} = 0$
adaptiveStep	bool		True	true: use step reduction if step fails; false: fixed step size
minimumStepSize	UReal		1e-8	lower limit of step size, before nonlinear solver stops

verboseMode	Index		1	0 ... no output, 1 ... show errors and load steps, 2 ... show short Newton step information (error), 3 ... show also solution vector, 4 ... show also jacobian, 5 ... show also Jacobian inverse
verboseModeFile	Index		0	same behaviour as verboseMode, but outputs all solver information to file
preStepPyExecute	String		"	Python code to be executed prior to every load step and after last step, e.g. for post-processing

### 7.1.7 SimulationSettings

General Settings for simulation; according settings for solution and solvers are given in subitems of this structure.

SimulationSettings has the following items:

Name	type/function return type	size	default value / function args	description
timeIntegration	TimeIntegrationSettings			time integration parameters
solutionSettings	SolutionSettings			settings for solution files
staticSolver	StaticSolverSettings			static solver parameters
linearSolverType	LinearSolverType		LinearSolverType::EXUdense	selection of numerical linear solver: exu.LinearSolverType.EXUdense (dense matrix inverse), exu.LinearSolverType.EigenSparse (sparse matrix LU-factorization), ... (enumeration type)
cleanUpMemory	bool		False	true: solvers will free memory at exit (recommended for large systems); false: keep allocated memory for repeated computations to increase performance
displayStatistics	bool		False	display general computation information at end of time step (steps, iterations, function calls, step rejections, ...)
displayComputationTime	bool		False	display computation time statistics at end of solving
pauseAfterEachStep	bool		False	pause after every time step or static load step(user press SPACE)
outputPrecision	Index		6	precision for floating point numbers written to console; e.g. values written by solver
numberOfThreads	Index		1	number of threads used for parallel computation (1 == scalar processing); not yet implemented (status: Nov 2019)

## 7.2 Visualization settings

This section includes hierarchical structures for visualization settings, e.g., drawing of nodes, bodies, connectors, loads and markers and furthermore OpenGL, window and save image options.

### 7.2.1 VSettingsGeneral

General settings for visualization.

VSettingsGeneral has the following items:

Name	type/function return type	size	default value / function args	description
graphicsUpdateInterval	float		0.1	interval of graphics update during simulation in seconds; 0.1 = 10 frames per second; low numbers might slow down computation speed
autoFitScene	bool		True	automatically fit scene within first second after StartRenderer()
textSize	float		12.	general text size if not overwritten
minSceneSize	float		0.1	minimum scene size for initial scene size and for autoFitScene, to avoid division by zero; SET GREATER THAN ZERO
backgroundColor	Float4	4	[1.,1.,1.,1.]	red, green, blue and alpha values for background of render window (white=[1,1,1,1]; black = [0,0,0,1])
coordinateSystemSize	float		0.4	size of coordinate system relative to screen
drawCoordinateSystem	bool		True	false = no coordinate system shown
showComputationInfo	bool		True	false = no info about computation (current time, solver, etc.) shown
pointSize	float		0.01	global point size (absolute)
circleTiling	Index		16	global number of segments for circles; if smaller than 2, 2 segments are used (flat)
cylinderTiling	Index		16	global number of segments for cylinders; if smaller than 2, 2 segments are used (flat)
sphereTiling	Index		6	global number of segments for spheres; if smaller than 2, 2 segments are used (flat)
axesTiling	Index		12	global number of segments for drawing axes cylinders and cones (reduce this number, e.g. to 4, if many axes are drawn)

### 7.2.2 VSettingsWindow

Window and interaction settings for visualization; handle changes with care, as they might lead to unexpected results or crashes.

VSettingsWindow has the following items:



Name	type/function return type	size	default value / function args	description
renderWindowSize	Index2	2	[1024,768]	initial size of OpenGL render window in pixel
startupTimeout	Index		5000	OpenGL render window startup timeout in ms (change might be necessary if CPU is very slow)
alwaysOnTop	bool		False	true: OpenGL render window will be always on top of all other windows
maximize	bool		False	true: OpenGL render window will be maximized at startup
showWindow	bool		True	true: OpenGL render window is shown on startup; false: window will be iconified at startup (e.g. if you are starting multiple computations automatically)
keypressRotationStep	float		5.	rotation increment per keypress in degree (full rotation = 360 degree)
mouseMoveRotationFactor	float		1.	rotation increment per 1 pixel mouse movement in degree
keypressTranslationStep	float		0.1	translation increment per keypress relative to window size
zoomStepFactor	float		1.15	change of zoom per keypress (keypad +/-) or mouse wheel increment

### 7.2.3 VSettingsOpenGL

OpenGL settings for 2D and 2D rendering.

VSettingsOpenGL has the following items:

Name	type/function return type	size	default value / function args	description
initialCenterPoint	Float3	3	[0.,0.,0.]	centerpoint of scene (3D) at renderer startup; overwritten if autoFitScene = True
initialZoom	float		1.	initial zoom of scene; overwritten/ignored if autoFitScene = True
initialMaxSceneSize	float		1.	initial maximum scene size (auto: diagonal of cube with maximum scene coordinates); used for 'zoom all' functionality and for visibility of objects; overwritten if autoFitScene = True
initialModelRotation	StdArray33F	3x3	[Matrix3DF[3,3,1.,0.,0.,0.,1.,0.,0.,0.,1.]]	initial model rotation matrix for OpenGL; in python use e.g.: initialModelRotation=[[1,0,0],[0,1,0],[0,0,1]]

multiSampling	Index	1	1	multi sampling turned off ( $\leq 1$ ) or turned on to given values (2, 4, 8 or 16); increases the graphics buffers and might crash due to graphics card memory limitations; only works if supported by hardware; if it does not work, try to change 3D graphics hardware settings!
lineWidth	float	1	1.	width of lines used for representation of lines, circles, points, etc.
lineSmooth	bool	1	True	draw lines smooth
textLineWidth	float	1	1.	width of lines used for representation of text
textLineSmooth	bool	1	False	draw lines for representation of text smooth
showFaces	bool	1	True	show faces of triangles, etc.; using the options showFaces=false and showFaceEdges=true gives wire frame representation
showFaceEdges	bool	1	False	show edges of faces; using the options showFaces=false and showFaceEdges=true gives wire frame representation
shadeModelSmooth	bool	1	True	true: turn on smoothing for shaders, which uses vertex normals to smooth surfaces
materialSpecular	Float4	4	[1.,1.,1.,1.]	4f specular color of material
materialShininess	float	1	60.	shininess of material
enableLight0	bool	1	True	turn on/off light0
light0position	Float4	4	[1.,1.,-10.,0.]	4f position vector of GL light0; 4th value should be 0, otherwise the vector obtains a special interpretation, see opengl manuals
light0ambient	float	1	0.25	ambient value of GL light0
light0diffuse	float	1	0.4	diffuse value of GL light0
light0specular	float	1	0.4	specular value of GL light0
enableLight1	bool	1	True	turn on/off light1
light1position	Float4	4	[0.,3.,2.,0.]	4f position vector of GL light1; 4th value should be 0, otherwise the vector obtains a special interpretation, see opengl manuals
light1ambient	float	1	0.25	ambient value of GL light1
light1diffuse	float	1	0.4	diffuse value of GL light1
light1specular	float	1	0.	specular value of GL light1
drawFaceNormals	bool	1	False	draws triangle normals, e.g. at center of triangles; used for debugging of faces
drawVertexNormals	bool	1	False	draws vertex normals; used for debugging
drawNormalsLength	float	1	0.1	length of normals; used for debugging

## 7.2.4 VSettingsContour

Settings for contour plots; use these options to visualize field data, such as displacements, stresses, strains, etc. for bodies, nodes and finite elements.

VSettingsContour has the following items:

Name	type/function return type	size	default value / function args	description
outputVariableComponent	Index	1	0	select the component of the chosen output variable; e.g., for displacements, 3 components are available: 0 == x, 1 == y, 2 == z component; if this component is not available by certain objects or nodes, no value is drawn
outputVariable	OutputVariableType		OutputVariableType::_None	selected contour plot output variable type; select OutputVariableType::_None to deactivate contour plotting.
minValue	float	1	0	minimum value for contour plot; set manually, if automaticRange == False
maxValue	float	1	1	maximum value for contour plot; set manually, if automaticRange == False
automaticRange	bool		True	if true, the contour plot value range is chosen automatically to the maximum range
reduceRange	bool		True	if true, the contour plot value range is also reduced; better for static computation; in dynamic computation set this option to false, it can reduce visualization artifacts; you should also set minVal to max(float) and maxVal to min(float)
showColorBar	bool		True	show the colour bar with minimum and maximum values for the contour plot
colorBarTiling	Index	1	12	number of tiles (segments) shown in the colorbar for the contour plot

## 7.2.5 VSettingsExportImages

Functionality to export images to files (.tga format) which can be used to create animations; to activate image recording during the solution process, set SolutionSettings.recordImagesInterval accordingly.

VSettingsExportImages has the following items:

Name	type/function return type	size	default value / function args	description
saveImageTimeOut	Index		5000	timeout for saving a frame as image to disk; this is the amount of time waited for re-drawing; increase for very complex scenes

saveImageFileName	FileName		'images/frame'	filename (without extension!) and (relative) path for image file(s) with consecutive numbering (e.g., frame0000.tga, frame0001.tga,...); ; directory will be created if it does not exist
saveImageFileCounter	Index		0	current value of the counter which is used to consecutively save frames (images) with consecutive numbers
saveImageSingleFile	bool		False	true: only save single files with given filename, not adding numbering; false: add numbering to files, see saveImageFileName

### 7.2.6 VSettingsNodes

Visualization settings for nodes.

VSettingsNodes has the following items:

Name	type/function return type	size	default value / function args	description
show	bool		True	flag to decide, whether the nodes are shown
showNumbers	bool		False	flag to decide, whether the node number is shown
drawNodesAsPoint	bool		True	simplified/faster drawing of nodes; uses general->pointSize as drawing size; if drawNodesAsPoint==True, the basis of the node will be drawn with lines
showBasis	bool		False	show basis (three axes) of coordinate system in 3D nodes
basisSize	float		0.2	size of basis for nodes
tiling	Index		4	tiling for node if drawn as sphere; used to lower the amount of triangles to draw each node; if drawn as circle, this value is multiplied with 4
defaultSize	float		-1.	global node size; if -1.f, node size is relative to openGL.initialMaxSceneSize
defaultColor	Float4	4	[0.2,0.2,1.,1.]	default cRGB color for nodes; 4th value is alpha-transparency
showNodalSlopes	Index		False	draw nodal slope vectors, e.g. in ANCF beam finite elements

### 7.2.7 VSettingsBeams

Visualization settings for beam finite elements.

VSettingsBeams has the following items:

Name	type/function return type	size	default value / function args	description
axialTiling	Index		8	number of segments to discretise the beams axis

## 7.2.8 VSettingsBodies

Visualization settings for bodies.

VSettingsBodies has the following items:

Name	type/function return type	size	default value / function args	description
show	bool		True	flag to decide, whether the bodies are shown
showNumbers	bool		False	flag to decide, whether the body(=object) number is shown
defaultSize	Float3	3	[1.,1.,1.]	global body size of xyz-cube
defaultColor	Float4	4	[0.3,0.3,1.,1.]	default cRGB color for bodies; 4th value is
deformationScaleFactor	float		1	global deformation scale factor; also applies to nodes, if drawn; used for scaled drawing of (linear) finite elements, beams, etc.
beams	VSettingsBeams			visualization settings for beams (e.g. AN-CFCable or other beam elements)

## 7.2.9 VSettingsConnectors

Visualization settings for connectors.

VSettingsConnectors has the following items:

Name	type/function return type	size	default value / function args	description
show	bool		True	flag to decide, whether the connectors are shown
showNumbers	bool		False	flag to decide, whether the connector(=object) number is shown
defaultSize	float		0.1	global connector size; if -1.f, connector size is relative to maxSceneSize
showJointAxes	bool		False	flag to decide, whether contact joint axes of 3D joints are shown
jointAxesLength	float		0.2	global joint axes length
jointAxesRadius	float		0.02	global joint axes radius
showContact	bool		False	flag to decide, whether contact points, lines, etc. are shown

springNumberOfWindings	Index		8	number of windings for springs drawn as helical spring
contactPointsDefaultSize	float		0.02	global contact points size; if -1.f, connector size is relative to maxSceneSize
defaultColor	Float4	4	[0.2,0.2,1.,1.]	default cRGB color for connectors; 4th value is alpha-transparency

### 7.2.10 VSettingsMarkers

Visualization settings for markers.

VSettingsMarkers has the following items:

Name	type/function return type	size	default value / function args	description
show	bool		True	flag to decide, whether the markers are shown
showNumbers	bool		False	flag to decide, whether the marker numbers are shown
drawSimplified	bool		True	draw markers with simplified symbols
defaultSize	float		-1.	global marker size; if -1.f, marker size is relative to maxSceneSize
defaultColor	Float4	4	[0.1,0.5,0.1,1.]	default cRGB color for markers; 4th value is alpha-transparency

### 7.2.11 VSettingsLoads

Visualization settings for loads.

VSettingsLoads has the following items:

Name	type/function return type	size	default value / function args	description
show	bool		True	flag to decide, whether the loads are shown
showNumbers	bool		False	flag to decide, whether the load numbers are shown
defaultSize	float		0.2	global load size; if -1.f, load size is relative to maxSceneSize
defaultRadius	float		0.005	global radius of load axis if drawn in 3D
fixedLoadSize	bool		True	if true, the load is drawn with a fixed vector length in direction of the load vector, independently of the load size
drawSimplified	bool		True	draw markers with simplified symbols
loadSizeFactor	float		0.1	if fixedLoadSize=false, then this scaling factor is used to draw the load vector

defaultColor	Float4	4	[0.7,0.1,0.1,1.]	default cRGB color for loads; 4th value is alpha-transparency
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### 7.2.12 VSettingsSensors

Visualization settings for sensors.

VSettingsSensors has the following items:

Name	type/function return type	size	default value / function args	description
show	bool		True	flag to decide, whether the sensors are shown
showNumbers	bool		False	flag to decide, whether the sensor numbers are shown
drawSimplified	bool		True	draw sensors with simplified symbols
defaultSize	float		-1.	global sensor size; if -1.f, sensor size is relative to maxSceneSize
defaultColor	Float4	4	[0.6,0.6,0.1,1.]	default cRGB color for sensors; 4th value is alpha-transparency

### 7.2.13 VisualizationSettings

Settings for visualization.

VisualizationSettings has the following items:

Name	type/function return type	size	default value / function args	description
general	VSettingsGeneral			general visualization settings
window	VSettingsWindow			visualization window and interaction settings
openGL	VSettingsOpenGL			OpenGL rendering settings
contour	VSettingsContour			contour plot visualization settings
exportImages	VSettingsExportImages			settings for exporting (saving) images to files in order to create animations
nodes	VSettingsNodes			node visualization settings
bodies	VSettingsBodies			body visualization settings
connectors	VSettingsConnectors			connector visualization settings
markers	VSettingsMarkers			marker visualization settings
loads	VSettingsLoads			load visualization settings
sensors	VSettingsSensors			sensor visualization settings

## 7.3 Solver substructures

This section includes structures contained in the solver, which can be accessed via the python interface during solution or for building a customized solver in python.

### 7.3.1 CSolverTimer

Structure for timing in solver. Each Real variable is used to measure the CPU time which certain parts of the solver need. This structure is only active if the code is not compiled with the `__FAST_EXUDYN_LINALG` option and if `displayComputationTime` is set True. Timings will only be filled, if `useTimer` is True.

CSolverTimer has the following items:

Name	type/function return type	size	default value / function args	description
useTimer	bool		True	flag to decide, whether the timer is used (true) or not
total	Real		0.	total time measured between start and end of computation (static/dynamics)
factorization	Real		0.	solve or inverse
newtonIncrement	Real		0.	$Jac^{-1} * RHS$ ; backsubstitution
integrationFormula	Real		0.	time spent for evaluation of integration formulas
ODE2RHS	Real		0.	time for residual evaluation of ODE2 right-hand-side
AERHS	Real		0.	time for residual evaluation of algebraic equations right-hand-side
totalJacobian	Real		0.	time for all jacobian computations
jacobianODE2	Real		0.	jacobian w.r.t. coordinates of ODE2 equations (not counted in sum)
jacobianODE2_t	Real		0.	jacobian w.r.t. coordinates_t of ODE2 equations (not counted in sum)
jacobianAE	Real		0.	jacobian of algebraic equations (not counted in sum)
massMatrix	Real		0.	mass matrix computation
reactionForces	Real		0.	$CqT * \lambda$
postNewton	Real		0.	post newton step
writeSolution	Real		0.	time for writing solution
overhead	Real		0.	overhead, such as initialization, copying and some matrix-vector multiplication
python	Real		0.	time spent for python functions
visualization	Real		0.	time spent for visualization in computation thread
Reset(...)	void		useSolverTimer	reset solver timings to initial state by assigning default values; useSolverTimer sets the useTimer flag
Sum()	Real			compute sum of all timers (except for those counted multiple, e.g., jacobians)
StartTimer(...)	void		value	start timer function for a given variable; subtracts current CPU time from value



StopTimer(...)	void		value	stop timer function for a given variable; adds current CPU time to value
ToString()	String			converts the current timings to a string

### 7.3.2 SolverLocalData

Solver local data structure for solution vectors, system matrices and temporary vectors and data structures.

SolverLocalData has the following items:

Name	type/function return type	size	default value / function args	description
nODE2	Index		0	number of second order ordinary diff. eq. coordinates
nODE1	Index		0	number of first order ordinary diff. eq. coordinates
nAE	Index		0	number of algebraic coordinates
nData	Index		0	number of data coordinates
nSys	Index		0	number of system (unknown) coordinates = nODE2+nODE1+nAE
startAE	Index		0	start of algebraic coordinates, but set to zero if nAE==0
systemResidual	ResizableVector			system residual vector (vectors will be linked to this vector!)
newtonSolution	ResizableVector			Newton decrement (computed from residual and jacobian)
tempODE2	ResizableVector			temporary vector for ODE2 quantities; use in initial accelerations and during Newton
temp2ODE2	ResizableVector			second temporary vector for ODE2 quantities; use in static computation
tempODE2F0	ResizableVector			temporary vector for ODE2 Jacobian
tempODE2F1	ResizableVector			temporary vector for ODE2 Jacobian
startOfStepStateAAlgorithmic	ResizableVector			additional term needed for generalized alpha (startOfStep state)
aAlgorithmic	ResizableVector			additional term needed for generalized alpha (current state)
CleanUpMemory()	void			if desired, temporary data is cleaned up to safe memory
SetLinearSolverType(...)	void		linearSolverType	set linear solver type and matrix version: links system matrices to according dense/sparse versions
GetLinearSolverType()	LinearSolverType			return current linear solver type (dense/sparse)

### 7.3.3 SolverIterationData

Solver internal structure for counters, steps, step size, time, etc.; solution vectors, residuals, etc. are SolverLocalData. The given default values are overwritten by the simulationSettings when initializing the solver. SolverIterationData has the following items:

Name	type/function return type	size	default value / function args	description
maxStepSize	Real		0.	constant or maximum stepSize
minStepSize	Real		0.	minimum stepSize for static/dynamic solver; only used, if adaptive step is activated
currentStepSize	Real		0.	stepSize of current step
numberOfSteps	Index		0	number of time steps (if fixed size); $n$
currentStepIndex	Index		0	current step index; $i$
adaptiveStep	bool		True	if true, the step size may be adaptively controlled
currentTime	Real		0.	holds the current simulation time, copy of state.current.time; interval is [start-Time,tEnd]; in static solver, duration is loadStepDuration
startTime	Real		0.	time at beginning of time integration
endTime	Real		0.	end time of static/dynamic solver
discontinuousIteration	Index		0	number of current discontinuous iteration
newtonSteps	Index		0	number of current newton steps
newtonStepsCount	Index		0	count total Newton steps
newtonJacobiCount	Index		0	count total Newton jacobian computations
rejectedModifiedNewtonSteps	Index		0	count the number of rejected modified Newton steps (switch to full Newton)
discontinuousIterationsCount	Index		0	count total number of discontinuous iterations (min. 1 per step)
ToString()	String			convert iteration statistics to string; used for displayStatistics option

### 7.3.4 SolverConvergenceData

Solver internal structure for convergence information: residua, iteration loop errors and error flags. For detailed behavior of these flags, visit the source code!.

SolverConvergenceData has the following items:

Name	type/function return type	size	default value / function args	description
stepReductionFailed	bool		False	true, if iterations over time/static steps failed (finally, cannot be recovered)
discontinuousIterationsFailed	bool		False	true, if discontinuous iterations failed (may be recovered if adaptive step is active)
linearSolverFailed	bool		False	true, if linear solver failed to factorize
newtonConverged	bool		False	true, if Newton has (finally) converged

newtonSolutionDiverged	bool		False	true, if Newton diverged (may be recovered)
jacobianUpdateRequested	bool		True	true, if a jacobian update is requested in modified Newton (determined in previous step)
massMatrixNotInvertible	bool		True	true, if mass matrix is not invertible during initialization or solution (explicit solver)
discontinuousIterationError	Real		0.	error of discontinuous iterations (contact, friction, ...) outside of Newton iteration
residual	Real		0.	current Newton residual
lastResidual	Real		0.	last Newton residual to determine contractivity
contractivity	Real		0.	Newton contractivity = geometric decay of error in every step
errorCoordinateFactor	Real		1.	factor may include the number of system coordinates to reduce the residual
InitializeData()	void			initialize SolverConvergenceData by assigning default values

### 7.3.5 SolverOutputData

Solver internal structure for output modes, output timers and counters.

SolverOutputData has the following items:

Name	type/function return type	size	default value / function args	description
finishedSuccessfully	bool		False	flag is false until solver finished successfully (can be used as external trigger)
verboseMode	Index		0	this is a copy of the solvers verboseMode used for console output
verboseModeFile	Index		0	this is a copy of the solvers verboseMode-File used for file
writeToSolutionFile	bool		False	if false, no solution file is generated and no file is written
writeToSolverFile	bool		False	if false, no solver output file is generated and no file is written
sensorValuesTemp	ResizableVector			temporary vector for per sensor values (overwritten for every sensor; usually contains last sensor)
lastSolutionWritten	Real		0.	simulation time when last solution has been written
lastSensorsWritten	Real		0.	simulation time when last sensors have been written
lastImageRecorded	Real		0.	simulation time when last image has been recorded
cpuStartTime	Real		0.	CPU start time of computation (starts counting at computation of initial conditions)

cpuLastTimePrinted	Real		0.	CPU time when output has been printed last time
InitializeData()	void			initialize SolverOutputData by assigning default values

### 7.3.6 MainSolverStatic

PyBind interface (trampoline) class for static solver. With this interface, the static solver and its substructures can be accessed via python. NOTE that except from SolveSystem(...), these functions are only intended for experienced users and they need to be handled with care, as unexpected crashes may happen if used inappropriate. Furthermore, the functions have a lot of overhead (performance much lower than internal solver) due to python interfaces, and should thus be used for small systems. To access the solver in python, write:

```
solver = MainSolverStatic()
```

and hereafter you can access all data and functions via 'solver'.

MainSolverStatic has the following items:

Name	type/function return type	size	default value / function args	description
timer	CSolverTimer			timer which measures the CPU time of solver sub functions
it	SolverIterationData			all information about iterations (steps, discontinuous iteration, newton,...)
conv	SolverConvergenceData			all information about tolerances, errors and residua
output	SolverOutputData			output modes and timers for exporting solver information and solution
newton	NewtonSettings			copy of newton settings from timeint or staticSolver
loadStepGeometricFactor	Real			multiplicative load step factor; this factor is computed from loadStepGeometric parameters in SolveSystem(...)
CheckInitialized(...)	bool		mainSystem	check if MainSolver and MainSystem are correctly initialized ==> otherwise raise SysError
ComputeLoadFactor(...)	Real		simulationSettings	for static solver, this is a factor in interval [0,1]; MUST be overwritten
GetSolverName()	std::string			get solver name - needed for output file header and visualization window
IsStaticSolver()	bool			return true, if static solver; needs to be overwritten in derived class
GetSimulationEndTime(...)	Real		simulationSettings	compute simulation end time (depends on static or time integration solver)
ReduceStepSize(...)	bool		mainSystem, simulationSettings, severity	reduce step size (1..normal, 2..severe problems); return true, if reduction was successful

IncreaseStepSize(...)	void		mainSystem, simulationSettings	increase step size if convergence is good
InitializeSolver(...)	bool		mainSystem, simulationSettings	initialize solverSpecific,data,it,conv; set/compute initial conditions (solver-specific!); initialize output files
PreInitializeSolverSpecific(...)	void		mainSystem, simulationSettings	pre-initialize for solver specific tasks; called at beginning of InitializeSolver, right after Solver data reset
InitializeSolverOutput(...)	void		mainSystem, simulationSettings	initialize output files; called from InitializeSolver()
InitializeSolverPreChecks(...)	bool		mainSystem, simulationSettings	check if system is solvable; initialize dense/sparse computation modes
InitializeSolverData(...)	void		mainSystem, simulationSettings	initialize all data,it,conv; called from InitializeSolver()
InitializeSolverInitialConditions(...)	void		mainSystem, simulationSettings	set/compute initial conditions (solver-specific!); called from InitializeSolver()
PostInitializeSolverSpecific(...)	void		mainSystem, simulationSettings	post-initialize for solver specific tasks; called at the end of InitializeSolver
SolveSystem(...)	bool		mainSystem, simulationSettings	solve System: InitializeSolver, SolveSteps, FinalizeSolver
FinalizeSolver(...)	void		mainSystem, simulationSettings	write concluding information (timer statistics, messages) and close files
SolveSteps(...)	bool		mainSystem, simulationSettings	main solver part: calls multiple InitializeStep(...)/ DiscontinuousIteration(...)/ FinishStep(...); do step reduction if necessary; return true if success, false else
UpdateCurrentTime(...)	void		mainSystem, simulationSettings	update currentTime (and load factor); MUST be overwritten in special solver class
InitializeStep(...)	void		mainSystem, simulationSettings	initialize static step / time step; python-functions; do some outputs, checks, etc.
FinishStep(...)	void		mainSystem, simulationSettings	finish static step / time step; write output of results to file
DiscontinuousIteration(...)	bool		mainSystem, simulationSettings	perform discontinuousIteration for static step / time step; CALLS ComputeNewtonResidual
Newton(...)	bool		mainSystem, simulationSettings	perform Newton method for given solver method
ComputeNewtonResidual(...)	void		mainSystem, simulationSettings	compute residual for Newton method (e.g. static or time step); store result in system-Residual
ComputeNewtonUpdate(...)	void		mainSystem, simulationSettings	compute update for currentState from newtonSolution (decrement from residual and jacobian)
ComputeNewtonJacobian(...)	void		mainSystem, simulationSettings	compute jacobian for newton method of given solver method; store result in system-Jacobian
WriteSolutionFileHeader(...)	void		mainSystem, simulationSettings	write unique file header, depending on static/ dynamic simulation
WriteCoordinatesToFile(...)	void		mainSystem, simulationSettings	write unique coordinates solution file

IsVerboseCheck(...)	bool		level	return true, if file or console output is at or above the given level
VerboseWrite(...)	void		level, str	write to console and/or file in case of level
GetODE2size()	Index			number of ODE2 equations in solver
GetODE1size()	Index			number of ODE1 equations in solver (not yet implemented)
GetASize()	Index			number of algebraic equations in solver
GetDataSize()	Index			number of data (history) variables in solver
GetSystemJacobian()	NumpyMatrix			get locally stored / last computed system jacobian of solver
GetSystemMassMatrix()	NumpyMatrix			get locally stored / last computed mass matrix of solver
GetSystemResidual()	NumpyVector			get locally stored / last computed system residual
GetNewtonSolution()	NumpyVector			get locally stored / last computed solution (=increment) of Newton
SetSystemJacobian(...)	void		systemJacobian	set locally stored system jacobian of solver; must have size nODE2+nODE1+nAE
SetSystemMassMatrix(...)	void		systemMassMatrix	set locally stored mass matrix of solver; must have size nODE2+nODE1+nAE
SetSystemResidual(...)	void		systemResidual	set locally stored system residual; must have size nODE2+nODE1+nAE
ComputeMassMatrix(...)	void		mainSystem, scalarFactor=1.	compute systemMassMatrix (multiplied with factor) in cSolver and return mass matrix
ComputeJacobianODE2RHS(...)	void		mainSystem, scalarFactor=1.	set systemJacobian to zero and add jacobian (multiplied with factor) of ODE2RHS to systemJacobian in cSolver
ComputeJacobianODE2RHS_t(...)	void		mainSystem, scalarFactor=1.	add jacobian of ODE2RHS_t (multiplied with factor) to systemJacobian in cSolver
ComputeJacobianAE(...)	void		mainSystem, scalarFactor_ODE2=1., scalarFactor_ODE2_t=1., velocityLevel=false	add jacobian of algebraic equations (multiplied with factor) to systemJacobian in cSolver; the scalarFactors are scaling the derivatives w.r.t. ODE2 coordinates and w.r.t. ODE2_t (velocity) coordinates; if velocityLevel==true, the constraints are evaluated at velocity level
ComputeODE2RHS(...)	void		mainSystem	compute the RHS of ODE2 equations in systemResidual in range(0,nODE2)
ComputeAlgebraicEquations(...)	void		mainSystem, velocityLevel=false	compute the algebraic equations in systemResidual in range(nODE2+nODE1, nODE2+nODE1+nAE)

### 7.3.7 MainSolverImplicitSecondOrder

PyBind interface (trampoline) class for dynamic implicit solver. Note that this solver includes the classical Newmark method (set useNewmark True; with option of index 2 reduction) as well as the generalized-alpha method. With the interface, the dynamic implicit solver and its substructures can be accessed via python.

NOTE that except from SolveSystem(...), these functions are only intended for experienced users and they need to be handled with care, as unexpected crashes may happen if used inappropriate. Furthermore, the functions have a lot of overhead (performance much lower than internal solver) due to python interfaces, and should thus be used for small systems. To access the solver in python, write

```
solver = MainSolverImplicitSecondOrder()
```

and hereafter you can access all data and functions via 'solver'. In this solver, user functions are possible to extend the solver at certain parts, while keeping the overall C++ performance.

MainSolverImplicitSecondOrder has the following items:

Name	type/function return type	size	default value / function args	description
timer	CSolverTimer			timer which measures the CPU time of solver sub functions
it	SolverIterationData			all information about iterations (steps, discontinuous iteration, newton,...)
conv	SolverConvergenceData			all information about tolerances, errors and residua
output	SolverOutputData			output modes and timers for exporting solver information and solution
newton	NewtonSettings			copy of newton settings from timeint or staticSolver
newmarkBeta	Real			copy of parameter in timeIntegration.generalizedAlpha
newmarkGamma	Real			copy of parameter in timeIntegration.generalizedAlpha
alphaM	Real			copy of parameter in timeIntegration.generalizedAlpha
alphaF	Real			copy of parameter in timeIntegration.generalizedAlpha
spectralRadius	Real			copy of parameter in timeIntegration.generalizedAlpha
factJacAlgorithmic	Real			locally computed parameter from generalizedAlpha parameters
CheckInitialized(...)	bool		mainSystem	check if MainSolver and MainSystem are correctly initialized ==> otherwise raise SysError
ComputeLoadFactor(...)	Real		simulationSettings	for static solver, this is a factor in interval [0,1]; MUST be overwritten
GetAAAlgorithmic()	NumpyVector			get locally stored / last computed algorithmic accelerations
GetStartOfStepStateAAAlgorithmic()	NumpyVector			get locally stored / last computed algorithmic accelerations at start of step
SetUserFunctionUpdateCurrentTime(...)	void		mainSystem, user-Function	set user function
SetUserFunctionInitializeStep(...)	void		mainSystem, user-Function	set user function
SetUserFunctionFinishStep(...)	void		mainSystem, user-Function	set user function

SetUserFunctionDiscontinuousIteration(...)	void		mainSystem, user-Function	set user function
SetUserFunctionNewton(...)	void		mainSystem, user-Function	set user function
SetUserFunctionComputeNewtonUpdate(...)	void		mainSystem, user-Function	set user function
SetUserFunctionComputeNewtonResidual(...)	void		mainSystem, user-Function	set user function
SetUserFunctionComputeNewtonJacobian(...)	void		mainSystem, user-Function	set user function
GetSolverName()	std::string			get solver name - needed for output file header and visualization window
IsStaticSolver()	bool			return true, if static solver; needs to be overwritten in derived class
GetSimulationEndTime(...)	Real		simulationSettings	compute simulation end time (depends on static or time integration solver)
ReduceStepSize(...)	bool		mainSystem, simulationSettings, severity	reduce step size (1..normal, 2..severe problems); return true, if reduction was successful
IncreaseStepSize(...)	void		mainSystem, simulationSettings	increase step size if convergence is good
InitializeSolver(...)	bool		mainSystem, simulationSettings	initialize solverSpecific,data,it,conv; set/compute initial conditions (solver-specific!); initialize output files
PreInitializeSolverSpecific(...)	void		mainSystem, simulationSettings	pre-initialize for solver specific tasks; called at beginning of InitializeSolver, right after Solver data reset
InitializeSolverOutput(...)	void		mainSystem, simulationSettings	initialize output files; called from InitializeSolver()
InitializeSolverPreChecks(...)	bool		mainSystem, simulationSettings	check if system is solvable; initialize dense/sparse computation modes
InitializeSolverData(...)	void		mainSystem, simulationSettings	initialize all data,it,conv; called from InitializeSolver()
InitializeSolverInitialConditions(...)	void		mainSystem, simulationSettings	set/compute initial conditions (solver-specific!); called from InitializeSolver()
PostInitializeSolverSpecific(...)	void		mainSystem, simulationSettings	post-initialize for solver specific tasks; called at the end of InitializeSolver
SolveSystem(...)	bool		mainSystem, simulationSettings	solve System: InitializeSolver, SolveSteps, FinalizeSolver
FinalizeSolver(...)	void		mainSystem, simulationSettings	write concluding information (timer statistics, messages) and close files
SolveSteps(...)	bool		mainSystem, simulationSettings	main solver part: calls multiple InitializeStep(...)/ DiscontinuousIteration(...)/ FinishStep(...); do step reduction if necessary; return true if success, false else
UpdateCurrentTime(...)	void		mainSystem, simulationSettings	update currentTime (and load factor); MUST be overwritten in special solver class
InitializeStep(...)	void		mainSystem, simulationSettings	initialize static step / time step; python-functions; do some outputs, checks, etc.



FinishStep(...)	void		mainSystem, simulationSettings	finish static step / time step; write output of results to file
DiscontinuousIteration(...)	bool		mainSystem, simulationSettings	perform discontinuousIteration for static step / time step; CALLS ComputeNewtonResidual
Newton(...)	bool		mainSystem, simulationSettings	perform Newton method for given solver method
ComputeNewtonResidual(...)	void		mainSystem, simulationSettings	compute residual for Newton method (e.g. static or time step); store result in system-Residual
ComputeNewtonUpdate(...)	void		mainSystem, simulationSettings	compute update for currentState from newtonSolution (decrement from residual and jacobian)
ComputeNewtonJacobian(...)	void		mainSystem, simulationSettings	compute jacobian for newton method of given solver method; store result in system-Jacobian
WriteSolutionFileHeader(...)	void		mainSystem, simulationSettings	write unique file header, depending on static/ dynamic simulation
WriteCoordinatesToFile(...)	void		mainSystem, simulationSettings	write unique coordinates solution file
IsVerboseCheck(...)	bool		level	return true, if file or console output is at or above the given level
VerboseWrite(...)	void		level, str	write to console and/or file in case of level
GetODE2size()	Index			number of ODE2 equations in solver
GetODE1size()	Index			number of ODE1 equations in solver (not yet implemented)
GetAEsize()	Index			number of algebraic equations in solver
GetDataSize()	Index			number of data (history) variables in solver
GetSystemJacobian()	NumpyMatrix			get locally stored / last computed system jacobian of solver
GetSystemMassMatrix()	NumpyMatrix			get locally stored / last computed mass matrix of solver
GetSystemResidual()	NumpyVector			get locally stored / last computed system residual
GetNewtonSolution()	NumpyVector			get locally stored / last computed solution (=increment) of Newton
SetSystemJacobian(...)	void		systemJacobian	set locally stored system jacobian of solver; must have size nODE2+nODE1+nAE
SetSystemMassMatrix(...)	void		systemMassMatrix	set locally stored mass matrix of solver; must have size nODE2+nODE1+nAE
SetSystemResidual(...)	void		systemResidual	set locally stored system residual; must have size nODE2+nODE1+nAE
ComputeMassMatrix(...)	void		mainSystem, scalarFactor=1.	compute systemMassMatrix (multiplied with factor) in cSolver and return mass matrix
ComputeJacobianODE2RHS(...)	void		mainSystem, scalarFactor=1.	set systemJacobian to zero and add jacobian (multiplied with factor) of ODE2RHS to systemJacobian in cSolver

ComputeJacobianODE2RHS_t(...)	void		mainSystem, scalarFactor=1.	add jacobian of ODE2RHS_t (multiplied with factor) to systemJacobian in cSolver
ComputeJacobianAE(...)	void		mainSystem, scalarFactor_ODE2=1., scalarFactor_ODE2_t=1., velocityLevel=false	add jacobian of algebraic equations (multiplied with factor) to systemJacobian in cSolver; the scalarFactors are scaling the derivatives w.r.t. ODE2 coordinates and w.r.t. ODE2_t (velocity) coordinates; if velocityLevel == true, the constraints are evaluated at velocity level
ComputeODE2RHS(...)	void		mainSystem	compute the RHS of ODE2 equations in systemResidual in range(0,nODE2)
ComputeAlgebraicEquations(...)	void		mainSystem, velocityLevel=false	compute the algebraic equations in systemResidual in range(nODE2+nODE1, nODE2+nODE1+nAE)



## Chapter 8

# 3D Graphics Visualization

The 3D graphics visualization window is kept simple, but useful to see the animated results of the multibody system.

### 8.1 Mouse input

The following table includes the mouse functions.

Button	action	remarks
left mouse button	move model	keep left mouse button pressed to move the model in the current x/y plane
right mouse button	rotate model	keep right mouse button pressed to rotate model around current current $X_1/X_2$ axes
mouse wheel	zoom	use mouse wheel to zoom (on touch screens 'pinch-to-zoom' might work as well)

### 8.2 Keyboard input

The following table includes the keyboard shortcuts available in the window.

Key(s)	action	remarks
1,2,3,4 or 5	visualization update speed	the entered digit controls the visualization update, which can be changed from 1=1 update per 20ms to 5=1 update per 100s
'.' or KEYPAD +	zoom in	zoom one step into scene (additionally press CTRL to perform small zoom step)
'/' or KEYPAD -	zoom out	zoom one step out of scene (additionally press CTRL to perform small zoom step)
CTRL+1 or SHIFT+CTRL+1	change view	set view in 1/2-plane (+SHIFT: view from opposite side)
CTRL+2 or SHIFT+CTRL+2	change view	set view in 1/3-plane (+SHIFT: view from opposite side)

<b>CTRL+3</b> <b>SHIFT+CTRL+3</b>	<b>or</b>	change view	set view in 2/3-plane (+SHIFT: view from opposite side)
<b>CTRL+4</b> <b>SHIFT+CTRL+4</b>	<b>or</b>	change view	set view in 2/1-plane (+SHIFT: view from opposite side)
<b>CTRL+5</b> <b>SHIFT+CTRL+5</b>	<b>or</b>	change view	set view in 3/1-plane (+SHIFT: view from opposite side)
<b>CTRL+6</b> <b>SHIFT+CTRL+6</b>	<b>or</b>	change view	set view in 3/2-plane (+SHIFT: view from opposite side)
<b>A</b>		zoom all	set zoom such that the whole scene is visible
<b>CURSOR UP, DOWN, ...</b>		move scene	use cursor keys to move the scene up, down, left, and right (use CTRL for small movements)
<b>KEYPAD 2/8,4/6,1/9</b>		rotate scene	about 1, 2 and 3-axis (use CTRL for small rotations)
<b>C</b>		show/hide connectors	pressing this key switches the visibility of connectors
<b>CTRL+C</b>		show/hide connector numbers	pressing this key switches the visibility of connector numbers
<b>B</b>		show/hide bodies	pressing this key switches the visibility of bodies
<b>CTRL+B</b>		show/hide body numbers	pressing this key switches the visibility of body numbers
<b>L</b>		show/hide loads	pressing this key switches the visibility of loads
<b>CTRL+L</b>		show/hide load numbers	pressing this key switches the visibility of load numbers
<b>M</b>		show/hide markers	pressing this key switches the visibility of markers
<b>CTRL+M</b>		show/hide marker numbers	pressing this key switches the visibility of marker numbers
<b>N</b>		show/hide nodes	pressing this key switches the visibility of nodes
<b>CTRL+N</b>		show/hide node numbers	pressing this key switches the visibility of node numbers
<b>S</b>		show/hide sensors	pressing this key switches the visibility of sensors
<b>CTRL+S</b>		show/hide sensor numbers	pressing this key switches the visibility of sensor numbers
<b>Q</b>		stop simulation	simulation is stopped and cannot be recovered
<b>X</b>		execute command	open dialog to enter a python command (in global python scope); dialog may appear behind the visualization window! User errors may lead to crash – be careful! Examples: <code>'print(mbs)'</code> , <code>'x=5'</code> , <code>'mbs.GetObject(0)'</code> , etc.
<b>V</b>		visualization settings	open dialog to modify visualization settings; dialog may appear behind the visualization window!
<b>ESCAPE</b>		close render window	stops the simulation and closes the render window
<b>SPACE</b>		continue simulation	if simulation is paused, it can be continued by pressing space; use SHIFT+SPACE to continuously activate 'continue simulation'

## Chapter 9

# System of equations

The general idea of the code is to have objects, which provide equations (ODE2, ODE1, AE). The solver then assembles these equations and solves the static or dynamic problem. The system structure and solver follow partially the previous implementation in HOTINT [6, 1, 3].

### 9.1 LHS-RHS naming conventions in EXUDYN

Functions and variables contain the abbreviations LHS (*left-hand-side*) and RHS (*right-hand-side*), sometimes lower-case, in order to distinguish if terms are computed at the LHS or RHS.

The objects have the following LHS–RHS conventions:

- the acceleration term, e.g.,  $m \cdot \ddot{q}$  is always positive on the LHS
- objects, connectors, etc., use LHS conventions for most terms: mass, stiffness matrix, elastic forces, damping, etc., are computed at LHS of the object equation
- object forces are written at the RHS of the object equation
- in case of constraint or connector equations, there is no LHS or RHS, as there is no acceleration term. Therefore, the computation function evaluates the term as given in the description of the object, adding it to the LHS.

Object equations may read, e.g., for one coordinate  $q$ , mass  $m$ , damping coefficient  $d$ , stiffness  $k$  and applied force  $f$ ,

$$\underbrace{m \cdot \ddot{q} + d \cdot \dot{q} + k \cdot q}_{LHS} = \underbrace{f}_{RHS} \quad (9.1)$$

In this case, the C++ function `ComputeODE2LHS(const Vector& ode2Lhs)` will compute the term  $d \cdot \dot{q} + k \cdot q$  with positive sign. Note that the acceleration term  $m \cdot \ddot{q}$  is computed separately, as it is computed from mass matrix and acceleration.

However, system quantities (e.g. within the solver) are always written on RHS, except for the acceleration and mass matrix:

$$\underbrace{M_{sys} \cdot \ddot{q}_{sys}}_{LHS} = \underbrace{f_{sys}}_{RHS} \quad (9.2)$$

In the case of the object equation

$$m \cdot \ddot{q} + d \cdot \dot{q} + k \cdot q = f, \quad (9.3)$$

the RHS term becomes  $f_{sys} = -(d \cdot \dot{q} + k \cdot q) + f$  and it is computed by the C++ function `ComputeSystemODE2RHS`.

This means, that during computation, terms which appear at the LHS of the object are transferred to the RHS of the system equation. This enables a simpler setup of equations for the solver.



# Chapter 10

## Solver

### 10.1 Jacobian computation

The computation of the global jacobian matrix is time consuming for the static solver or implicit time integration. The equations are split into 2<sup>nd</sup> order differential equations, 1<sup>st</sup> order differential equations and algebraic equations parts. From this structure, in the general non-symmetric case,  $3 \times 3$  submatrices result for the jacobian. Every submatrix of the jacobian has a certain meaning and needs to be computed individually. Specifically, in implicit time integration the 2<sup>nd</sup> order differential equations  $\times$  2<sup>nd</sup> order differential equations term includes the (tangent) stiffness matrix and the mass matrix.

For efficient computation purpose, the elements provide a list of flags, which determine the dependencies as well as available (analytical) functions to compute the local (object) jacobian:

- ODE2\_ODE2 ... derivative of ODE2 equations with respect to ODE2 variables
- ODE2\_ODE2\_t ... derivative of ODE2 equations with respect to ODE2\_t (velocity) variables
- ODE1\_ODE1 ... derivative of ODE1 equations with respect to ODE1 variables (NOT YET AVAILABLE)
- AE\_ODE2 ... derivative of AE (algebraic) equations with respect to ODE2 variables
- AE\_ODE2\_t ... derivative of AE (algebraic) equations with respect to ODE2\_t (velocity) variables (NOT YET AVAILABLE)
- AE\_ODE1 ... derivative of AE (algebraic) equations with respect to ODE1 variables (NOT YET AVAILABLE)
- AE\_AE ... derivative of AE (algebraic) equations with respect to AE variables

If one of these flags is set (binary; e.g. ODE2\_ODE2 + ODE2\_ODE2\_t), then the according local jacobian is computed and assembled into the global jacobian in the static or implicit dynamic solver.

Jacobians can also be supplied in analytical (function) form, which is indicated by an additional flag with the same name but an additional term '\_function', e.g. 'ODE2\_ODE2\_function' indicates that the derivative of ODE2 equations with respect to its ODE2 coordinates is provided in an analytical form (this is the tangent stiffness matrix).

Two **object** functions are used to compute the local jacobians:

- **ComputeJacobianODE2\_ODE2(Matrix& jacobian, Matrix& jacobian\_ODE2\_t):** computes the ODE2\_ODE2 and ODE2\_ODE2\_t jacobians
- **ComputeJacobianAE(Matrix& jacobian, Matrix& jacobian\_AE):** computes the AE\_ODE2 and AE\_AE jacobians of the object ITSELF

Two **connector** functions are used to compute the local jacobians, using **MarkerData**:



- **ComputeJacobianODE2\_ODE2(Matrix& jacobian, Matrix& jacobian\_ODE2\_t, const MarkerDataStructure& markerData):** computes the ODE2\_ODE2 and ODE2\_ODE2\_t jacobians of the connector; e.g. for spring-damper
- **ComputeJacobianAE(Matrix& jacobian, Matrix& jacobian\_AE, const MarkerDataStructure& markerData):** computes the AE\_ODE2 and AE\_AE jacobians of the connector; e.g. for coordinate constraint

The system jacobian has the structure (2= ODE2, 1= ODE1,  $\lambda$  = AE;  $\bar{\mathbf{f}}_2$  = according system residual including dynamic (mass matrix) terms in time integration;  $\mathbf{g}_\lambda$  = algebraic equations):

$$\begin{bmatrix} \frac{\partial \bar{\mathbf{f}}_2}{\partial \mathbf{q}_2} & 0 & \left( \frac{\partial \mathbf{g}_\lambda}{\partial \mathbf{q}_2} \right)^T \\ 0 & \frac{\partial \bar{\mathbf{f}}_1}{\partial \mathbf{q}_1} & \left( \frac{\partial \mathbf{g}_\lambda}{\partial \mathbf{q}_1} \right)^T \\ \frac{\partial \mathbf{g}_\lambda}{\partial \mathbf{q}_2} & \frac{\partial \mathbf{g}_\lambda}{\partial \mathbf{q}_1} & \frac{\partial \mathbf{g}_\lambda}{\partial \mathbf{q}_\lambda} \end{bmatrix} \quad (10.1)$$

Two system jacobian functions are currently available:

- **JacobianODE2RHS(temp, newton, factorODE2, factorODE2\_t, jacobian\_ODE2, jacobian\_ODE2\_t):** compute analytical/numerical differentiation of ODE2RHS w.r.t. ODE2 and ODE2\_t coordinates; if analytical/functional version of jacobian is available and Newton flag 'useNumericalDifferentiation'=false, then the according jacobian is computed by its according function; results are 2 jacobians; the factors 'factor\_ODE2' and 'factor\_ODE2\_t' are used to scale the two jacobians; if a factor is zero, the according jacobian is not computed.
- **JacobianAE(temp, newton, jacobian, factorODE2, velocityLevel, fillIntoSystemMatrix):** compute constraint jacobian of AE with respect to ODE2 ('fillIntoSystemMatrix'=true: also w.r.t. [ODE1] and AE) coordinates → direct computation given by access functions; 'factorODE2' is used to scale the ODE2-part of the jacobian (to avoid postmultiplication); velocityLevel = true: velocityLevel constraints are used, if available; 'fillIntoSystemMatrix'=true: fill in both  $\frac{\partial \bar{\mathbf{f}}_1}{\partial \mathbf{q}_2}$ ,  $\frac{\partial \bar{\mathbf{f}}_1}{\partial \mathbf{q}_2}^T$  AND  $\frac{\partial \bar{\mathbf{f}}_1}{\partial \mathbf{q}_\lambda}$  at according locations into system matrix; 'fillIntoSystemMatrix'=false: (this is a temporary/WORKAROUND function):

The system jacobian functions compute the local jacobians either by means of a provided function or numerically, using the 'NumericalDifferentiation' settings of 'Newton'.

## 10.2 Implicit trapezoidal rule solver

This solver represents a class of solvers, which are based on the implicit trapezoidal rule. This integration includes the start value and the end value of a time step for the interpolation, thus being a trapezoidal integration rule. In some specializations, e.g. the Newmark method, the interpolation might only depend on the start value or the end value.

Most important representations of this rule:

- Trapezoidal rule (= Newmark with  $\beta = \frac{1}{4}$  and  $\gamma = \frac{1}{2}$ )
- Newmark method
- Generalized- $\alpha$  method (= generalized Newmark method with additional parameters)

## 10.3 Representation of coordinates and equations of motion

Nomenclature:

- '2' ... second order equations (usually of a mechanical system)
- '1' ... first order equations (e.g. of a controller, fluid, etc.)
- 'λ' ... algebraic equations (usually of joints)
- **M** ... mass matrix
- **q<sub>2</sub>** ... 'displacement' coordinates of ODE2 equations
- **q̇<sub>2</sub>** ... 'velocity' coordinates of ODE2 equations
- **q̈<sub>2</sub>** ... 'acceleration' coordinates of ODE2 equations
- **q<sub>1</sub>** ... coordinates of ODE1 equations
- **q̇<sub>1</sub>** ... 'velocity' coordinates of ODE1 equations
- **q<sub>λ</sub>** ... Lagrange multipliers
- **f<sub>2</sub>** ... right-hand-side of ODE2 equations (except for action of joint reaction forces)
- **f<sub>1</sub>** ... right-hand-side of ODE1 equations
- **g** ... algebraic equations
- **K** ... (tangent) stiffness matrix
- **D** ... damping/gyroscopic matrix
- **h** ... step size of time integration method

The equations of motion in EXUDYN are represented as

$$\mathbf{M}\ddot{\mathbf{q}}_2 + \frac{\partial \mathbf{g}}{\partial \mathbf{q}_2^T} \mathbf{q}_\lambda = \mathbf{f}_2(\mathbf{q}_2, \dot{\mathbf{q}}_2, t) \quad (10.2)$$

$$\dot{\mathbf{q}}_1 + \frac{\partial \mathbf{g}}{\partial \mathbf{q}_1^T} \mathbf{q}_\lambda = \mathbf{f}_1(\mathbf{q}_1, t) \quad (10.3)$$

$$\mathbf{g}(\mathbf{q}_2, \dot{\mathbf{q}}_2, \mathbf{q}_1, \mathbf{q}_\lambda, t) = 0 \quad (10.4)$$

Note that the term  $\frac{\partial \mathbf{g}}{\partial \mathbf{q}_1^T} \mathbf{q}_\lambda$  is not yet implemented, such that algebraic equations may not yet depend on 1<sup>st</sup> order differential equations coordinates.

It is important to note, that for linear mechanical systems, **f<sub>2</sub>** becomes

$$\mathbf{f}_2^{lin} = \mathbf{f}^a - \mathbf{K}\mathbf{q}_2 - \mathbf{D}\dot{\mathbf{q}}_1 \quad (10.5)$$

in which **f<sup>a</sup>** represents applied forces and **K** and **D** become part of the system Jacobian for time integration.

## 10.4 Newmark method

The Newmark method obtains two parameters  $\beta$  and  $\gamma$ . The main ideas are

- Interpolate the displacements and the velocities linearly using the accelerations of the beginning of the time step (subindex '0') and the end of the time step (subindex 'T').
- Solve the system equations at the end of the time step for the unknown accelerations as well as for 1<sup>st</sup> order differential equations and algebraic equations coordinates.

We abbreviate the unknown accelerations by  $\ddot{\mathbf{q}} = \mathbf{a}$  and the unknown velocities  $\dot{\mathbf{q}} = \mathbf{v}$ . Thus, the equations at the end of the time step read (bring all terms to LHS):

$$\mathbf{f}_2^{\text{Newmark}} = \mathbf{M}\mathbf{a}_2^T + \frac{\partial \mathbf{g}}{\partial \mathbf{q}_2^T} \mathbf{q}_\lambda^T - \mathbf{f}_2(\mathbf{q}_2^T, \dot{\mathbf{q}}_2^T, t) = 0 \quad (10.6)$$

$$\mathbf{f}_1^{\text{Newmark}} = \mathbf{v}_1^T + \frac{\partial \mathbf{g}}{\partial \mathbf{q}_1^T} \mathbf{q}_\lambda^T - \mathbf{f}_1(\mathbf{q}_1^T, t) = 0 \quad (10.7)$$

$$\mathbf{f}_\lambda^{\text{Newmark}} = \mathbf{g}(\mathbf{q}_2^T, \dot{\mathbf{q}}_2^T, \mathbf{q}_1^T, \mathbf{q}_\lambda^T, t) = 0 \quad (10.8)$$

Within Eq. (10.6), the 2<sup>nd</sup> order differential equations displacements and velocities and for 1<sup>st</sup> order differential equations coordinates are given by

$$\begin{aligned}\mathbf{q}_2^T &= \mathbf{q}_2^0 + h\dot{\mathbf{q}}_2^0 + h^2\left(\frac{1}{2} - \beta\right)\mathbf{a}_2^0 + h^2\beta\mathbf{a}_2^T \\ \dot{\mathbf{q}}_2^T &= \dot{\mathbf{q}}_2^0 + h(1 - \gamma)\mathbf{a}_2^0 + h\gamma\mathbf{a}_2^T \\ \mathbf{q}_1^T &= \mathbf{q}_1^0 + h(1 - \gamma)\mathbf{v}_1^0 + h\gamma\mathbf{v}_1^T\end{aligned}\quad (10.9)$$

The unknowns for the Newton method are

$$\mathbf{q}^{\text{Newton}} = \begin{bmatrix} \mathbf{a}_2^T \\ \mathbf{v}_1^T \\ \mathbf{q}_\lambda^T \end{bmatrix} \quad (10.10)$$

For the Newton method, we need to compute an update for the unknowns Eq. (10.10), using the known residual  $\mathbf{r}_{i-1}$  and the inverse of the Jacobian  $\mathbf{J}_{i-1}$  of step  $i - 1$ ,

$$\mathbf{q}_i^{\text{Newton}} = \mathbf{q}_{i-1}^{\text{Newton}} - \mathbf{J}^{-1}(\mathbf{q}_{i-1}^{\text{Newton}}) \mathbf{r}(\mathbf{q}_{i-1}^{\text{Newton}}) \quad (10.11)$$

The Jacobian has the following  $3 \times 3$  structure,

$$\mathbf{J} = \begin{bmatrix} \mathbf{J}_{22} & \mathbf{J}_{21} & \mathbf{J}_{2\lambda} \\ \mathbf{J}_{12} & \mathbf{J}_{11} & \mathbf{J}_{1\lambda} \\ \mathbf{J}_{\lambda 2} & \mathbf{J}_{\lambda 1} & \mathbf{J}_{\lambda\lambda} \end{bmatrix} \quad (10.12)$$

Note that currently, all terms related to '1' are not implemented. The other terms are only evaluated in the specific jacobian computation, if according flags are set in GetAvailableJacobian(). Otherwise, the constraint needs to be implemented as object which can employ all kinds of coordinates, which do not depend on coordinates of markers.

The available Jacobians need to be rewritten in terms of the Newton unknowns (10.10), and thus read

$$\begin{aligned}\mathbf{J}_{22} &= \frac{\partial \mathbf{f}_2^{\text{Newmark}}}{\partial \mathbf{a}_2^T} = \frac{\partial \mathbf{f}_2^{\text{Newmark}}}{\partial \mathbf{q}_2^T} \frac{\mathbf{q}_2}{\mathbf{a}_2^T} + \frac{\partial \mathbf{f}_2^{\text{Newmark}}}{\partial \dot{\mathbf{q}}_2^T} \frac{\dot{\mathbf{q}}_2}{\mathbf{a}_2^T} = h^2\beta\mathbf{K} + h\gamma\mathbf{D} \\ \mathbf{J}_{2\lambda} &= \frac{\partial \mathbf{f}_2^{\text{Newmark}}}{\partial \mathbf{q}_\lambda^T} = \frac{\partial \mathbf{g}}{\partial \mathbf{q}_\lambda^T} \\ \mathbf{J}_{\lambda 2} &= \frac{\partial \mathbf{f}_\lambda^{\text{Newmark}}}{\partial \mathbf{a}_2^T} = \frac{\partial \mathbf{g}}{\partial \mathbf{a}_2^T} = \frac{\partial \mathbf{g}}{\partial \mathbf{q}_2^T} \frac{\mathbf{q}_2}{\mathbf{a}_2^T} + \frac{\partial \mathbf{g}}{\partial \dot{\mathbf{q}}_2^T} \frac{\dot{\mathbf{q}}_2}{\mathbf{a}_2^T} = h^2\beta \frac{\partial \mathbf{g}}{\partial \mathbf{q}_2^T} + h\gamma \frac{\partial \mathbf{g}}{\partial \dot{\mathbf{q}}_2^T} \\ \mathbf{J}_{\lambda\lambda} &= \frac{\partial \mathbf{f}_\lambda^{\text{Newmark}}}{\partial \mathbf{q}_\lambda^T} = \frac{\partial \mathbf{g}}{\partial \mathbf{q}_\lambda^T}\end{aligned}\quad (10.13)$$

Note that the derivative  $\frac{\mathbf{q}_2}{\mathbf{a}_2^T}$  follows from the Newmark interpolation (10.9) using the relation between  $\mathbf{q}_2^T$  and  $\mathbf{a}_2^T$ . The tangent stiffness matrix  $\mathbf{K}$  must also include derivatives of applied forces  $\mathbf{f}^a$ , which is currently not implemented. Furthermore, the Jacobian is not symmetric, which could be obtained by according scaling.

Once an update  $\mathbf{q}_i^{\text{Newton}}$  has been computed, the interpolation formulas (10.9) need to be evaluated before the next residual and Jacobian can be computed.

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# Chapter 11

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