BIODEG

Biodegradation and Corrosion Simulation using Finite Element

User Manual

Version 0.8

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KU LEUVEN

BioDeg website

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1 Introduction

BIODEG is an open source software written in FreeFEM (a domain-specific language for finite element programming), C++, and Python for modeling the degradation of metallic biomaterials and simulating the biodegradation behavior of medical devices, implants and scaffolds in corrosion experiments. It can handle any geometry of desire and supports parallel computing to simulate models with a high level of complexity.

1.1 Authors

BIODEG is developed by the Biomechanics Research group at KU Leuven and University of Liege. The code is currently maintained by its principal developer, who manage the development of the mathematical models and the core functionalities.

Principal developer

• Mojtaba Barzegari (University of Leuven, Belgium).

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1.2 Acknowledgments

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1.3 Referencing BioDeg

Please refer to BIODEG repository, section "Publications and referencing" to properly cite the use of BIODEG in your scientific work.

2 Useful background information

You may refer to the following articles for a background of the methods and algorithms implemented in BioDeg.

In addition, below are some useful references on finite element method and some online resources that provide a background of finite elements and their application to the solution of partial differential equations.

3 Installation

All the underlying installation instructions assume a Linux operating system. We assume standard tools and libraries like CMake, compilers- (C, C++ and Fortran), and MPI libraries are pre-installed. Most high-performance computers would have the latest version of these libraries in the default environment.

3.1 Compiling and installing external libraries

- 3.1.1 Instructions for PETSc and Qt
- 3.1.2 Instructions for FreeFEM

3.2 Obtaining and Compiling BioDeg

Assuming that you have already installed the above external dependencies, next follow the steps below to obtain and compile BioDeg.

4 Running BioDeg

- 4.1 Configuring the simulation
- 4.1.1 Example 1
- 4.1.2 Example 2

5 Future extensions to BioDeg

The future versions of BioDeg will focus on the implementation of the following methodologies/features.

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6 Finding answers to more questions

If you have questions that go beyond this manual, there are a number of resources:

- For questions/suggestions about BIODEG installation, bugs, or similar stuff please use the BIODEG issue tracker.
- BIODEG is primarily based on the FreeFEM. If you have particular questions about FreeFEM, contact the community at https://community.freefem.org/.
- If you have specific questions about BioDeg that are not suitable for public and archived mailing lists, you can contact the primary developer and mentor:
 - Mojtaba Barzegari: mojtaba.barzegari@kuleuven.be.
 - Liesbet Geris: liesbet.geris@kuleuven.be (Mentor).

A Run-time input parameters

The underlying description of the input parameters also includes a "Standard/Advanced" label, which signifies whether an input parameter is a standard one or an advanced level parameter. The default values of the "Advanced" parameters are good enough for almost all cases. However, in some cases user may need to use "Advanced" labeled parameters. For user convenience, all input parameters are also indexed at the end of this manual in Section A.4.

A.1 Geometry and mesh parameters

• Parameter name: mesh_file

Default: Should be provided

Description: [Standard]

 $Possible\ values:\ {\bf Any\ string\ value}$

• Parameter name: label_scaffold

Default: 1

Description: [Standard]

Possible values: Any positive integer value

• Parameter name: label_medium

Default: 2

Description: [Standard]

Possible values: Any positive integer value

• Parameter name: label_wall

Default: 3

Description: [Advanced]

Possible values: Any positive integer value

• Parameter name: label_inlet

Default: 4

Description: [Advanced]

Possible values: Any positive integer value

• Parameter name: label_outlet

Default: 5

Description: [Advanced]

Possible values: Any positive integer value

• Parameter name: box_length

Default: 20.0

Description: [Standard]

Possible values: Any positive floating point number

• Parameter name: cube_size_x

Default: 13.0

Description: [Standard]

Possible values: Any positive floating point number

• Parameter name: cube_size_y

Default: 13.0

Description: [Standard]

• Parameter name: cube_size_z

Default: 4.0

Description: [Standard]

Possible values: Any positive floating point number

• Parameter name: mesh_size

Default: 32

Description: [Standard]

Possible values: Any positive integer number

• Parameter name: refine_initial_mesh

Default: false (0)

Description: [Advanced]

Possible values: A boolean value (1 or 0)

• Parameter name: mshmet_error

Default: 0.01

Description: [Advanced]

Possible values: Any floating point number

• Parameter name: mesh_size_min

Default: 0.04

Description: [Advanced]

Possible values: Any floating point number

• Parameter name: mesh_size_max

Default: 0.8

Description: [Advanced]

Possible values: Any floating point number

A.2 Materials and boundary conditions parameters

• Parameter name: material_density

Default: 1.735e-3

Description: [Standard]

Possible values: Any floating point number

• Parameter name: film_density

 $Default:\ 2.3446\text{e--}3$

Description: [Standard]

Possible values: Any floating point number

• Parameter name: material_satur

Default: 0.134e-3

Description: [Advanced]

• Parameter name: material_eps

Default: 0.55

Description: [Advanced]

Possible values: Any floating point number between 0 and 1

• Parameter name: material_tau

Default: 1.0

Description: [Advanced]

Possible values: Any floating point number

• Parameter name: d_mg

 $Default:\ 0.05$

Description: [Standard]

Possible values: Any floating point number

• Parameter name: d_cl

Default: 0.05

Description: [Standard]

Possible values: Any floating point number

• Parameter name: d_oh

Default: 25.2

Description: [Standard]

Possible values: Any floating point number

• Parameter name: k1

Default: 7.0

Description: [Standard]

Possible values: Any floating point number

• Parameter name: k2

Default: 1e15

Description: [Standard]

Possible values: Any floating point number

• Parameter name: fluid_nu

Default: 0.85

Description: [Advanced]

Possible values: Any floating point number

• Parameter name: fluid_in_x

Default: 0.1

Description: [Advanced]

• Parameter name: fluid_in_y

Default: 0

Description: [Advanced]

Possible values: Any floating point number

• Parameter name: fluid_in_z

Default: 0

Description: [Advanced]

Possible values: Any floating point number

• Parameter name: initial_cl

Default: 5.175e-6

Description: [Standard]

Possible values: Any floating point number

• Parameter name: initial_oh

Default: 1e-7

Description: [Standard]

Possible values: Any floating point number

A.3 Solver parameters

• Parameter name: solve_mg

Default: true (1)

Description: [Standard]

Possible values: Any boolean value (1 or 0)

• Parameter name: solve_film

Default: true (1)

Description: [Standard]

Possible values: Any boolean value (1 or 0)

• Parameter name: solve_cl

Default: true (1)

Description: [Standard]

Possible values: Any boolean value (1 or 0)

• Parameter name: solve_oh

Default: true (1)

Description: [Standard]

Possible values: Any boolean value (1 or 0)

• Parameter name: solve_ls

Default: true (1)

Description: [Standard]

Possible values: Any boolean value (1 or 0)

• Parameter name: solve_fluid

Default: false (0)

Description: [Standard]

Possible values: Any boolean value (1 or 0)

• Parameter name: solve_full_ns

Default: true (1)

Description: [Advanced]

 $Possible\ values:$ Any boolean value (1 or 0)

• Parameter name: write_fluid_output

Default: true (1)

Description: [Advanced]

Possible values: Any boolean value (1 or 0)

• Parameter name: solve_fluid_each

Default: 10

Description: [Advanced]

Possible values: Any positive integer number

• Parameter name: time_step

Default: 0.025

Description: [Advanced]

Possible values: Any floating point number

• Parameter name: final_time

Default: 21.0

Description: [Standard]

Possible values: Any floating point number

• Parameter name: do_redistance

Default: true (1)

Description: [Advanced]

Possible values: Any boolean value (1 or 0)

• Parameter name: redistance_time

Default: 1.0

Description: [Advanced]

A.4 Output parameters

• Parameter name: text_output_file

Default: "output/result.txt"
Description: [Standard]

Possible values: Any string value referring to a valid path

• Parameter name: write_vtk

Default: true (1)

Description: [Standard]

Possible values: Any boolean value (1 or 0)

• Parameter name: vtk_output_name

Default: "output/output" Description: [Standard]

Possible values: Any string value referring to a valid path

• Parameter name: save_each

Default: 0.25

Description: [Standard]

Possible values: Any floating point number

• Parameter name: save_last_state

Default: true (1)

Description: [Standard]

Possible values: Any boolean value (1 or 0)

• Parameter name: output_per_area

Default: false (0)

Description: [Advanced]

Possible values: Any boolean value (1 or 0)

• Parameter name: save_multiplier

Default: 1.0

Description: [Advanced]

Possible values: Any floating point number

• Parameter name: export_scaffold

Default: false (0)

Description: [Advanced]

Possible values: Any boolean value (1 or 0)

• Parameter name: export_scaffold_each

Default: 1.0

Description: [Advanced]

• Parameter name: export_scaffold_volume

Default: true (1)

Description: [Advanced]

Possible values: Any boolean value (1 or 0)

• Parameter name: export_scaffold_surface

Default: true (1)

Description: [Advanced]

Possible values: Any boolean value (1 or 0)

• Parameter name: save_initial_mesh

Default: false (0)

Description: [Advanced]

Possible values: Any boolean value (1 or 0)

ullet Parameter name: save_initial_partitioned_mesh

Default: false (0)

Description: [Advanced]

Possible values: Any boolean value (1 or 0)

Index of run-time parameters with section names

The following is a listing of all run-time parameters, sorted by the section in which they appear.

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