

# Solfec User Manual

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# 1 Introduction

## 2 Input language

Solfec input file is essentially a Python source code. Python interpreter is embedded in Solfec. At the same time Solfec extends Python by adding a number of objects and routines. There are few general principles to remember:

- Zero based indexing is observed in routine arguments.
- Parameters after the bar | are optional. For example *FUNCTION* (*a*, *b* | *c*, *d*) has two optional parameters *c*, *d*.
- Passing Solfec objects to some routines *empties* them. This means that a variable, that was passed as an argument, no longer stores data. For example: let  $x = \text{CREATE1}()$  create an object  $x$ , and let  $y = \text{CREATE2}(x)$  create an object  $y$ , using  $x$ . If  $\text{CREATE2}(x)$  empties  $x$ , then after the call  $x$  becomes an empty placeholder. One can use it to assign value,  $x = \text{CREATE1}()$ , but using it as an argument,  $z = \text{CREATE2}(x)$ , will cause an abnormal termination. One can create a copy of an object by calling  $z = \text{COPY}(x)$ , hence using  $y = \text{CREATE2}(\text{COPY}(x))$  leaves  $x$  intact.

Sections below document Solfec objects and routines used for their manipulation.

### 2.1 CONVEX

An object of type CONVEX is either an arbitrary convex polyhedron, or it is a collection of such polyhedrons.

**obj = CONVEX (vertices, faces, valid | convex)**

This routine creates a CONVEX object from a detailed input data.

- **obj** - created CONVEX object
- **vertices** - list of vertices: [ $x0, y0, z0, x1, y1, z1, \dots$ ]
- **faces** - list of faces: [ $n1, v1, v2, \dots, vn1, s1, n2, v1, v2, \dots, vn2, s2, \dots$ ], where  $n1$  is the number of vertices of the first face,  $v1, v2, \dots, vn1$  enumerate the vertices in the CCW order when looking from the outside, and  $s1$  is the surface identifier of the face. Similarly for the second face and so on.
- **valid** - volume identifier

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- **convex** (emptied) - collection of CONVEX objects appending **obj**

Some parameters can also be accessed as members and methods of a CONVEX object. These are

Read-only members and methods
<i><b>obj.nver</b> - number of convex vertices</i>
<i><b>obj.vertex</b> (<b>n</b>) - returns a (<i>x</i>, <i>y</i>, <i>z</i>) tuple storing coordinates of <i>nth</i> vertex</i>

### **obj = HULL (points, valid, surfid | convex)**

This routine creates a CONVEX object as a convex hull of a point set.

- **obj** - created CONVEX object
- **points** - list of points: [*x0*, *y0*, *z0*, *x1*, *y1*, *z1*, ...]
- **valid** - volume identifier
- **surfid** - surface identifier common to all faces
- **convex** (emptied) - collection of CONVEX objects appending **obj**

## 2.2 MESH

An object of type MESH describes an arbitrary volumetric mesh, comprising tetrahedrons, pyramids, wedges, and hexahedrons (Figure 2.1).

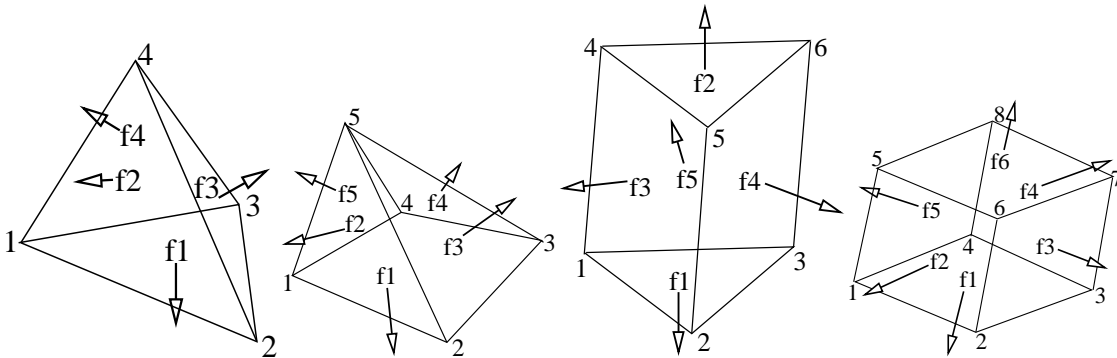


Figure 2.1: Element types in Solfec

### **obj = MESH (nodes, elements, surfids)**

This routine creates a MESH object from a detailed input data.

- **obj** - created MESH object
- **nodes** - list of nodes:  $[x0, y0, z0, x1, y1, z1, \dots]$
- **elements** - list of elements:  $[e1, n1, n2, \dots, ne1, v1, e2, n1, n2, \dots, ne2, v2, \dots]$ , where  $e1$  is the number of nodes of the first element,  $n1, n2, \dots, ne1$  enumerate the element nodes, and  $v1$  is the volume identifier of the element. Similarly for the second element and so on.
- **surfids** - list of surface identifiers:  $[gid, f1, n1, n2, \dots, nf1, s1, f2, n1, n2, \dots, nf2, s2, \dots]$ , where  $gid$  is the global surface identifier for all not specified faces,  $f1$  is the number of nodes in the first specified face,  $n1, n2, \dots, nf1$  enumerate the face nodes, and  $s1$  is the surface identifier of the face. Similarly for other specified faces. If only the  $gid$  is given, this can be done either as  $[gid]$  or as  $gid$  alone.

Some parameters can also be accessed as members and methods of a MESH object. These are

<b>Read-only</b> members and methods
<i>obj.nnod</i> - number of mesh nodes
<i>obj.node (n)</i> - returns a $(x, y, z)$ tuple storing coordinates of $n$ th node

### **obj = HEX (nodes, i, j, k, volid, surfids | dx, dy, dz)**

This routine creates a MESH object corresponding to a hexahedral shape (hexahedral elements are used).

- **obj** - created MESH object
- **nodes** - list of 8 nodes:  $[x0, y0, z0, x1, y1, z1, \dots, x7, y7, z7]$ . The hexahedral shape will be stretched between those nodes using a linear interpolation.
- **i, j, k** - numbers of subdivisions along the local  $x, y, z$  directions.
- **volid** - volume identifier
- **surfids** - list of six surface identifiers:  $[s1, s2, \dots, s6]$ , corresponding to the faces of the hexahedral shape
- **dx, dy, dz** - lists of subdivision schemes along local  $x, y, z$  directions. By default a subdivision is uniform. When  $dx = [1, 1, 5, 5, 1, 1]$  is present, then this scheme will be normalized (actual numbers do not matter, but their ratios) and applied to the local  $x$  direction of the generated shape.

### **obj = ROUGH\_HEX (shape, i, j, k | dx, dy, dz)**

This routine creates a hexahedral MESH object corresponding to a given shape. The resultant mesh properly contains the input shape and with its orientation (which is based on the inertia properties of the shape).

- **obj** - created MESH object
- **shape** - an input shape defined by a collection of CONVEX objects; a list of CONVEX objects (or their collections) *[cvx1, cvx2, cvx3, ....]* is as well accepted.
- **i, j, k** - numbers of subdivisions along the local *x, y, z* directions of the principal inertia axes
- **dx, dy, dz** - lists of subdivision schemes along local *x, y, z* directions. By default a subdivision is uniform. When *dx = [1, 1, 5, 5, 1, 1]* is present, then this scheme will be normalized (actual numbers do not matter, but their ratios) and applied to the local x direction of the generated shape.

## 2.3 SPHERE

An object of type SPHERE is either a sphere, or it is a collections of spheres.

### **obj = SPHERE (center, radius, volid, surfid | sphere)**

This routine creates a SPHERE object.

- **obj** - created SPHERE object
- **center** - tuple *(x, y, z)* defining the center
- **radius** - sphere radius
- **volid** - volume identifier
- **surfid** - surface identifier
- **sphere** (emptied) - collection of SPHERE objects appending **obj**

Some parameters can also be accessed as members of a MESH object. These are

Read-only members and methods
<i>obj.center, obj.radius</i>

## 2.4 SOLFEC

An object of type SOLFEC represents the Solfec algorithm. One can use several SOLFEC objects to run several analyzes from a single input file.

## **obj = SOLFEC (analysis, step, output)**

This routine creates a SOLFEC object.

- **obj** - created SOLFEC object
- **analysis** - 'DYNAMIC' or 'QUASI\_STATIC' analysis kind
- **step** - initially assumed time step, regarded as an upper bound
- **output** - defines the output **directory** path (**Important note:** if this directory exists and contains valid output data SOLFEC is created in 'READ' mode, otherwise SOLFEC is created in 'WRITE' mode)

Some parameters can also be accessed as members of a SOLFEC object. These are

<b>Read-only</b> members
<i>obj.analysis</i>
<i>obj.time</i> - current time
<i>obj.mode</i> - either 'READ' or 'WRITE' as described above
<i>obj.constraints</i> - list of constraints (cf. Section 2.11)
<i>obj.ncon</i> - number of constraints
<i>obj.bodies</i> - list of bodies (cf. Section 2.7)
<i>obj.nbod</i> - number of bodies
<b>Read/write</b> members
<i>obj.step</i>

## 2.5 SURFACE\_MATERIAL

An object of type SURFACE\_MATERIAL represents material properties on the interface between two surfaces. Surfaces identifiers were included in definitions of all geometric objects.

### **obj = SURFACE\_MATERIAL (solfec| surf1, surf2, model, label, friction, cohesion, restitution, spring, dashpot)**

This routine creates a SURFACE\_MATERIAL object.

- **obj** - created SURFACE\_MATERIAL object
- **solfec** - **obj** is created for this SOLFEC object
- **surf1** - first surface identifier (default: 0)
- **surf2** - second surface identifier (default: 0). If **surf1** or **surf2** (or both) are not specified, a *default* surface material is being defined (one used when a specific surface pairing cannot be found for a new contact point).

## 2 Input language

Model name	Employs variables
'SIGNORINI_COULOMB'	friction, cohesion, restitution
'SPRING_DASHPOT'	spring, dashpot, friction, cohesion

Table 2.1: Surface material models.

- **model** - material model name (default: 'SIGNORINI\_COULOMB'), see Table 2.1 and Chapter 4
- **label** - label string (default: 'SURFACE\_MATERIAL\_*i*', where *i* is incremented for each call)
- **friction** - friction coefficient (default: 0.0)
- **cohesion** - cohesion per unit area (default: 0.0)
- **restitution** - velocity restitution (default: 0.0)
- **spring** - spring stiffness (default: 0.0)
- **dashpot** - dashpot stiffness (default: 0.0)

Some parameters can also be accessed as members of a SURFACE\_MATERIAL object. These are

<b>Read-only</b> members
<i>obj.surf1, obj.surf2, obj.label</i>
<b>Read/write</b> members
<i>obj.model, obj.friction, obj.cohesion, obj.restitution, obj.spring, obj.dashpot</i>

## 2.6 BULK\_MATERIAL

An object of type BULK\_MATERIAL represents material properties of a volume.

**obj = BULK\_MATERIAL (solfec| model, label, young, poisson, density)**

This routine creates a BULK\_MATERIAL object.

- **obj** - created BULK\_MATERIAL object
- **solfec** - **obj** is created for this SOLFEC object
- **model** - material model name (default: 'KIRCHHOFF'), see Table 2.2 and Chapter 4



## 2 Input language

Model name	Employs variables
'KIRCHHOFF'	young, poisson, density

Table 2.2: Bulk material models.

- **label** - label string (default: 'BULK\_MATERIAL\_*i*', where *i* is incremented for each call)
- **young** - Young's modulus (default: 1E6)
- **poisson** - Poisson's coefficient (default: 0.25)
- **density** - material density (default: 1E3)

Some parameters can also be accessed as members of a BULK\_MATERIAL object. These are

<b>Read-only</b> members
<i>obj.model</i> , <i>obj.label</i>
<b>Read/write</b> members
<i>obj.young</i> , <i>obj.poisson</i> , <i>obj.density</i>

## 2.7 BODY

An object of type BODY represents a solid body.

**obj = BODY (solfec, kind, shape, material | label, formulation, mesh)**

This routine creates a body.

- **obj** - created BODY object
- **solfec** - **obj** is created for this SOLFEC object
- **kind** - a string: 'RIGID', 'PSEUDO\_RIGID', 'FINITE\_ELEMENT' or 'OBSTACLE' describing the kinematic model
- **shape** (emptied) - this is can be a CONVEX/MESH/SPHERE object, or a list [*obj1*, *obj2*, ...], where each object is of type CONVEX/MESH/SPHERE. If the **kind** is 'FINITE\_ELEMENT', then two cases are possible:
  - **shape** is a single MESH object: the mesh describes both the shape and the discretisation of the motion of a body
  - **shape** is solely composed of CONVEX objects: here a separate **mesh** must be given to discretise motion of a body (see the **mesh** argument below)

## 2 Input language

Formulation	Remarks
'FEM_O1'	Use first order elements
'FEM_O2'	Use second order elements

Table 2.3: Bulk material models.

- **material** - a BULK\_MATERIAL object or a label of a bulk material (specifies an initial body-wise material, see also the **MATERIAL (...)** routine in Section 2.14)
- **label** - a label string (no label is assigned by default)
- **formulation** - valid when **kind** equals 'FINITE\_ELEMENT', ignored otherwise (default: 'FEM\_O1'). This argument specifies a formulation of the finite element method. See Table 2.3.
- **mesh** - optional when **kind** equals 'FINITE\_ELEMENT', ignored otherwise. This variable must be a MESH object describing a finite element mesh properly containing the **shape** composed solely of CONVEX objects. This way the 'FINITE\_ELEMENT' model allows to handle complicated shapes with less finite elements, e.g. an arbitrary shape could be contained in just one hexahedron.

Some parameters can also be accessed as members of a BODY object. These are

Read-only members
<i>obj.kind</i> , <i>obj.label</i>
<i>obj.conf</i> - tuple ( $q1, q2, \dots, qN$ ) storing configuration of the body. See Table 2.4.
<i>obj.velo</i> - tuple ( $u1, u2, \dots, uN$ ) storing velocity of the body. See Table 2.5.

Read/write members
<i>obj.selfcontact</i> - self-contact detection flag (default: 'OFF') taking values 'ON' or 'OFF'.
<i>obj.scheme</i> - time integration scheme (default: 'DEFAULT') used to integrate motion. See Table 2.6.
<i>obj.damping</i> - mass proportional damping coefficient (default: 0.0) for the dynamic case.

## 2.8 TIME\_SERIES

An object of type TIME\_SERIES is a linear spline based on a series of 2-points.

Body kind	Configuration description
'RIGID'	Column-wise rotation matrix followed by the current mass center.
'PSEUDO_RIGID'	Column-wise deformation gradient followed by the current mass center.
'FINITE_ELEMENT'	Current coordinates x, y, z of mesh nodes.
'OBSTACLE'	Python <i>None</i> object.

Table 2.4: Types of configurations.

Body kind	Velocity description
'RIGID'	Referential angular velocity followed by the spatial velocity of mass center.
'PSEUDO_RIGID'	Deformation gradient velocity followed by the spatial velocity of mass center.
'FINITE_ELEMENT'	Components x, y, z of spatial velocities of mesh nodes.
'OBSTACLE'	Python <i>None</i> object.

Table 2.5: Types of velocities.

### **obj = TIME\_SERIES (points)**

This routine creates a TIME\_SERIES object.

- **obj** - created TIME\_SERIES object
- **points** - either a list  $[t0, v0, t1, v1, \dots]$  of points (where  $t_i < t_j$ , when  $i < j$ ), or a path to a file storing times and values pairs

## **2.9 GAUSS\_SEIDEL\_SOLVER**

An object of type GAUSS\_SEIDEL\_SOLVER represents a nonlinear block Gauss-Seidel solver, employed for the calculation of constraint reactions.

### **obj = GAUSS\_SEIDEL\_SOLVER (epsilon, maxiter | failure, diagepsilon, diagmaxiter, diagsolver, data, callback)**

This routine creates a GAUSS\_SEIDEL\_SOLVER object.

- **obj** - created GAUSS\_SEIDEL\_SOLVER object
- **epsilon** - relative accuracy of constraint reactions sufficient for termination
- **maxiter** - maximal number of iterations before termination

## 2 Input language

Scheme	Kinematics	Remarks
'DEFAULT'	all	Use a default time integrator regardless of underlying kinematics.
'RIG_POS'	rigid	NEW1 in [3]: explicit, positive energy drift, no momentum conservation
'RIG_NEG'	rigid	NEW2 in [3]: explicit, negative energy drift, exact momentum conservation; <b>default</b> for rigid kinematics
'RIG_IMP'	rigid	NEW3 in [3]: semi-explicit, no energy drift and exact momentum conservation
'DEF_EXP'	pseudo-rigid, finite element	Explicit trapezoidal-like scheme described in Chapter 5 of [2]; <b>default</b> for deformable kinematics
'DEF_IMP'	pseudo-rigid, finite element	Linearly implicit trapezoidal scheme similar to [1]; <b>TODO</b> : make it default after more testing

Table 2.6: Time integration schema.

- **failure** - failure (lack of convergence) action (default: 'CONTINUE'). Available failure actions are: 'CONTINUE' (simulation is continued), 'EXIT' (simulation is stopped and Solfec exits), 'CALLBACK' (a callback function is called if it was set or otherwise the 'EXIT' scenario is executed). In all cases *obj.error* variable is set up, cf. Table 2.7.
- **diagepsilon** - diagonal block solver relative accuracy of constraint reactions (default:  $\epsilon / 100$ )
- **diagmaxiter** - diagonal block solver maximal number of iterations (default:  $\max(100, \maxiter / 100)$ )
- **diagsolver** - diagonal block solver kind (default: 'SEMISMOOTH\_NEWTON'). Available diagonal solvers are 'SEMISMOOTH\_NEWTON', 'PROJECTED\_GRADIENT', 'DE\_SAXE\_AND\_FENG', cf. Chapter 5.
- **data** - data passed to the failure callback function (if this is a tuple it will accordingly expand the parameter list of the callback routine)
- **callback** - failure callback function of form:  $value = callback(obj, data)$ , where for the returned value equal zero Solfec run is stopped

Some parameters can also be accessed as members of a GAUSS\_SEIDEL\_SOLVER object. These are

<b>Read-only members</b>
<i>obj.failure</i>
<i>obj.error</i> - current error code, cf. Table 2.7
<i>obj.itors</i> - number of iterations during a last run of solver
<i>obj.rerhist</i> - if history recording is on, this is a list of relative error values for each iteration of the last run. Otherwise a <i>None</i> object is returned.

<b>Read/write members</b>
<i>obj.epsilon</i> , <i>obj.maxiter</i> , <i>obj.diagepsilon</i> , <i>obj.diagmaxiter</i> , <i>obj.diamsolver</i>
<i>obj.history</i> - 'ON' or 'OFF' flag switching history recording (default is 'OFF')
<i>obj.reverse</i> - 'ON' or 'OFF' flag switching iteration reversion modes (whether to alternate backward and forward or not, default is 'OFF')
<i>obj.variant</i> - variant of parallel Gauss-Seidel update (default: 'FULL'), cf. Table 2.8. Ignored in sequential mode.

'OK'	No error has occurred
'DIVERGED'	Global iteration loop divergence
'DIAGONAL_DIVERGED'	Diagonal solver iteration loop divergence
'DIAGONAL_FAILED'	Failure of a diagonal solver (e.g. singularity)

Table 2.7: Error codes of GAUSS\_SEIDEL\_SOLVER object

## 2.10 EXPLICIT\_SOLVER

An object of type EXPLICIT\_SOLVER represents a penalty based constraint solver. When in use, all 'SIGNORONI\_COULOMB' type contact interfaces are regarded as 'SPRING\_DASHPOT' ones. One should then remember about specifying the *spring* value for those.

**obj = EXPLICIT\_SOLVER ()**

- **obj** - created EXPLICIT\_SOLVER object

## 2.11 CONSTRAINT

An object of type CONSTRAINT represents a constraint and some of its associated data (e.g. constraint reaction). Both user prescribed constraints and contact constraints are represented by an object of the same type.

**obj = FIX\_POINT (solfec, body, point)**

This routine creates a fixed point constraint.

## 2 Input language

'FULL'	Full Gauss-Seidel update as in sequential case. Although the slowest, it works in all cases. It should be noted, that all of the below variants will usually fail for all-rigid-body models.
'MIDDLE_JACOBI'	Jacobi update for off-processor data of <b>W</b> matrix blocks communicating with processors of higher and lower colors. Of use for deformable kinematics, where off-diagonal interactions are weaker. The Gauss-Seidel run-time should be halved for large numbers of processors.
'BOUNDARY_JACOBI'	Use Jacobi update for all off-processor data. This approach will fail in most cases. It servers as illustration.
'SIMPLIFIED'	A single sweep over contacts is done with previous values of off-processor data. This is followed by local Gauss-Seidel iterations for all non-contact constraints. This is the fastest and least consistent approach, of use for deformable kinematics dominated models.

Table 2.8: Variants of parallel Gauss-Seidel update.

- **obj** - created CONSTRAINT object
- **solfec** - **obj** is created for this SOLFEC object
- **body** - BODY object whose motion is constrained
- **point** -  $(x, y, z)$  tuple with referential point coordinates

**obj = FIX\_DIRECTION (solfec, body, point, direction)**

This routine fixes the motion of a referential point along a specified spatial direction.

- **obj** - created CONSTRAINT object
- **solfec** - **obj** is created for this SOLFEC object
- **body** - BODY object whose motion is constrained
- **point** -  $(x, y, z)$  tuple with referential point coordinates
- **direction** -  $(vx, vy, vz)$  tuple with spatial direction components

**obj = FIX\_SURFACE (solfec, body, surfid, direction)TODO**

This routine fixes the motion of a referential surface along a specified spatial direction.

- **obj** - list of created point CONSTRAINT objects
- **solfec** - **obj** is created for this SOLFEC object
- **body** - BODY object whose motion is constrained
- **surfid** - surface identifier
- **direction** -  $(vx, vy, vz)$  tuple defining the direction of load, or 'NORMAL' if load is normal to the surface, or 'TANGENT1' if load acts along the first tangent direction, or 'TANGENT2' if it acts along the second tangent direction. The normal direction ('NORMAL') is outward. The first tangent direction ('TANGENT1') is the one of the steepest descent, or a global  $x$  direction if the surface is horizontal. The second tangent direction ('TANGENT2') is such that the local ('TANGENT1', 'TANGENT2', 'NORMAL') coordinate system is right-handed.

**obj = SET\_DISPLACEMENT (solfec, body, point, direction, tms)**

This routine prescribes a displacement history of a referential point along a specified spacial direction.

- **obj** - created CONSTRAINT object
- **solfec** - **obj** is created for this SOLFEC object
- **body** - BODY object whose motion is constrained
- **point** -  $(x, y, z)$  tuple with referential point coordinates
- **direction** -  $(vx, vy, vz)$  tuple with spatial direction components
- **tms** - TIME\_SERIES object with the displacement history

**obj = SET\_VELOCITY (solfec, body, point, direction, value)**

This routine prescribes a velocity history of a referential point along a specified spacial direction.

- **obj** - created CONSTRAINT object
- **solfec** - **obj** is created for this SOLFEC object
- **body** - BODY object whose motion is constrained
- **point** -  $(x, y, z)$  tuple with referential point coordinates
- **direction** -  $(vx, vy, vz)$  tuple with spatial direction components
- **value** - a constant value or a TIME\_SERIES object with the velocity history

**obj = SET\_ACCELERATION (solfec, body, point, direction, tms)**

This routine prescribes an acceleration history of a referential point along a specified spacial direction.

- **obj** - created CONSTRAINT object
- **solfec** - **obj** is created for this SOLFEC object
- **body** - BODY object whose motion is constrained
- **point** -  $(x, y, z)$  tuple with referential point coordinates
- **direction** -  $(vx, vy, vz)$  tuple with spatial direction components
- **tms** - TIME\_SERIES object with the acceleration history

**obj = PUT\_RIGID\_LINK (solfec, body1, body2, point1, point2)**

This routine creates a rigid link constraints between two referential points of two distinct bodies.

- **obj** - created CONSTRAINT object
- **solfec** - **obj** is created for this SOLFEC object
- **body1** - BODY object one whose motion is constrained (could be *None* when **body2** is not *None* - then one of the points is fixed “in the air”)
- **body2** - BODY object two whose motion is constrained (could be *None* when **body1** is not *None*)
- **point1** -  $(x1, y1, z1)$  tuple with the first referential point coordinates
- **point2** -  $(x2, y2, z2)$  tuple with the second referential point coordinates

Some parameters can also be accessed as members of a CONSTRAINT object. These are



Read-only members
<b>obj.kind</b> - kind of constraint: 'CONTACT', 'FIXPNT' (fixed point), 'FIXDIR' (fixed direction), 'VELODIR' (prescribed velocity; note that prescribed displacement and acceleration are converted into this case), 'RIGLNK' (rigid link)
<b>obj.R</b> - current average constraint reaction in a form of a tuple: $(RT1, RT2, RN)$ given with respect to a local base stored at <b>obj.base</b>
<b>obj.base</b> - current spatial coordinate system in a form of a tuple: $(eT1x, eT2x, eNx, eT1y, eT2y, eNy, eT1z, eT2z, eNz)$ where $x, y, z$ components are global
<b>obj.point</b> - current spatial point where the constraint force acts. This is a $(x, y, z)$ tuple for all constraint types, but 'RIGLNK' for which this is a $(x1, y1, z1, x2, y2, z2)$ tuple.
<b>obj.adjbod</b> - adjacent bodies. This is a tuple (body1, body2) of BODY objects for 'CONTACT' and 'RIGLNK' or a single BODY object otherwise.
<b>obj.matlab</b> - surface material label for constraints of kind 'CONTACT', or a <i>None</i> object otherwise.

## 2.12 Applying loads

Routines listed in this section apply loads.

### GRAVITY (solfec, direction, value)

This routine sets up the gravitational acceleration.

- **solfec** - SOLFEC object for which the acceleration is set up
- **direction** -  $(vx, vy, vz)$  tuple defining the direction
- **value** - a number or a TIME\_SERIES object defining the value of the acceleration

### FORCE (body, kind, point, direction, value| data)

This routine applies a point force to a body.

- **body** - BODY object to which the force is applied
- **kind** - either 'SPATIAL' or 'CONVECTED'; the *spatial* direction remains fixed, while the *convected* one follows deformation
- **point** -  $(x, y, z)$  tuple with the referential point where the force is applied
- **direction** -  $(vx, vy, vz)$  tuple defining the direction of force

## 2 Input language

- **value** - a number, a TIME\_SERIES object or a callback routine defining the value of the applied force. In case of a callback routine, the following format is assumed:

$$force = value\_callback (data, q, u, time, step)$$

where: **data** is the optional user data passed to **FORCE** routine (if **data** is a tuple it will expand the list of parameters to the callback), **q** is the configuration of the body passed to the callback, **u** is the velocity of the body passed to the callback, **time** is the current time passed to the callback and **step** is the current time step passed to the callback. The callback returns a **force** tuple. For rigid body the force reads (*spatial force, spatial torque, referential torque*), while for other kinds of bodies this is a generalised force of the same dimension as the velocity **u** (power conjugate to it).

- **data** - callback routine user data

### TORQUE (body, kind, direction, value)

This routine applies a torque to a *rigid* body.

- **body** - BODY object of kind 'RIGID' to which the torque is applied
- **kind** - either 'SPATIAL' or 'CONVECTED'; the *spatial* direction remains fixed, while the *convected* one follows deformation
- **direction** - (*vx, vy, vz*) tuple defining the direction of torque
- **value** - a number or a TIME\_SERIES object defining the value of the applied torque

### LOAD (body, kind, surfid, direction, value) **TODO**

This routine applies a surface load.

- **body** - BODY object to which the load is applied
- **kind** - either 'SPATIAL' or 'CONVECTED'; the *spatial* direction remains fixed, while the *convected* one follows deformation
- **surfid** - the integer surface identifier
- **direction** - (*vx, vy, vz*) tuple defining the direction of load, or 'NORMAL' if load is normal to the surface, or 'TANGENT1' if load acts along the first tangent direction, or 'TANGENT2' if it acts along the second tangent direction. The normal direction ('NORMAL') is outward. The first tangent direction ('TANGENT1') is the one of the steepest descent, or a global *x* direction if the surface is horizontal. The second tangent direction ('TANGENT2') is such that the local ('TANGENT1', 'TANGENT2', 'NORMAL') coordinate system is right-handed.

- **value** - a number or a TIME\_SERIES object defining the value of the applied load

### 2.13 Running simulations

Routines listed in this section control the solution process.

#### **RUN (solfec, solver, duration)**

This routine runs a simulation.

- **solfec** - SOLFEC object
- **solver** - constraint solver object (e.g. GAUSS\_SEIDEL\_SOLVER, EXPLICIT\_SOLVER)
- **duration** - duration of analysis

#### **OUTPUT (solfec, interval | compression)**

This routine specifies the frequency of writing to the output file.

- **solfec** - SOLFEC object
- **interval** - length of the time interval elapsing before consecutive output file writes
- **compression** - output compression mode: 'OFF' (default) or 'FASTLZ'. Compressed output files are smaller, although they might not be portable between hardware platforms.

#### **EXTENTS (solfec, extents)**

This routine bounds the simulation space. Bodies falling outside of the extents are deleted from the simulation.

- **solfec** - SOLFEC object
- **extents** -  $(x_{min}, y_{min}, z_{min}, x_{max}, y_{max}, z_{max})$  tuple

#### **CALLBACK (solfec, interval, data, callback)**

This routine defines a callback function, invoked during a run of Solfec every interval of time. A callback routine can interrupt the course of **RUN** command by returning 0.

- **solfec** - SOLFEC object
- **interval** - length of the time interval elapsing before consecutive callback calls
- **data** - data passed to the callback function
- **callback** - callback function of form:  $value = callback(data)$ , where for the returned value equal zero Solfec run is stopped

### UNPHYSICAL\_PENETRATION (solfec, depth)

This routine sets a depth of an unphysical interpenetration. Once it is exceeded, the simulation is stopped and a suitable error message printed out.

- **solfec** - SOLFEC object
- **depth** - interpenetration depth bound (default:  $\infty$ )

## 2.14 Utilities

Various utility routines are listed below.

### IMBALANCE\_TOLERANCE (solfec, tolerance)

This routine sets the imbalance tolerance for parallel balancing of Solfec data. A ratio of maximal to minimal per processor count of objects used. Hence, 1.0 indicates perfect balance, while any ratio  $> 1.0$  indicates an imbalance. Initially imbalance tolerances are all set to 1.3. This routine is ignored during sequential runs.

- **solfec** - SOLFEC object
- **tolerance** - data imbalance tolerance (default: 1.3)

### num = NCPU ()

This routine returns the number CPUs used in the analysis.

- **num** - the number of CPUs

### ret = HERE (solfec, object)

This routine tests whether an object is located on the current processor. During parallel runs objects migrate between processors. When calling a function (or a member) for an object not present on the current processor, the call will usually return `None` or be ignored. Hence, it is convenient to check whether an object resides on the current processor.

- **ret** - *True* or *False*
- **solfec** - SOLFEC object
- **object** - BODY or CONSTRAINT object

### obj = VIEWER ()

This routine tests whether the viewer is enabled.

- **obj** - *True* or *False* depending on whether the viewer (*-v* command line option) was enabled

### **BODY\_CHARS (body, mass, volume, center, tensor)**

This routine overwrites referential characteristics of a body.

- **body** - BODY object
- **mass** - body mass
- **volume** - body volume
- **center** -  $(x, y, z)$  mass center
- **tensor** -  $(t_{11}, t_{21}, \dots, t_{33})$  column-wise inertia tensor for a rigid body or Euler tensor otherwise

### **INITIAL\_VELOCITY (body, linear, angular)**

This routine applies initial (at time zero) linear and angular (in the sense of rigid motion) velocity to a body.

- **body** - BODY object
- **linear** - linear velocity  $(v_x, v_y, v_z)$
- **angular** - angular velocity  $(\omega_x, \omega_y, \omega_z)$

### **MATERIAL (solfec, body, valid, material)**

This routine applies material to a subset of geometric objects with the given volume identifier.

- **solfec** - SOLFEC object
- **body** - BODY object
- **valid** - volume identifier
- **material** - MATERIAL object or material label

### **DELETE (solfec, object)**

This routine deletes a BODY object or a CONSTRAINT object from a SOLFEC object.

- **solfec** - SOLFEC object
- **object** (emptied) - BODY or CONSTRAINT object

### **obj = SCALE (shape, coefs)**

This routine scales a geometrical object or a collection of such objects.

- **obj** - when **shape** is not  $(x, y, z)$  tuple: same as **shape**, returned for convenience. Otherwise the  $(x \cdot \text{coefs}[0], y \cdot \text{coefs}[1], z \cdot \text{coefs}[2])$  tuple.
- **shape** - object, collection of objects, or a list  $[a, b, c, \dots]$  of objects of type CONVEX, MESH, SPHERE. Alternately this can be a single  $(x, y, z)$  tuple.
- **coefs** -  $(cx, cy, cz)$  tuple of scaling factors along each axis

### **obj = TRANSLATE (shape, vector)**

This routine translates a geometrical object or a collection of such objects.

- **obj** - when **shape** is not  $(x, y, z)$  tuple: same as **shape**, returned for convenience. Otherwise the  $(x + \text{vector}[0], y + \text{vector}[1], z + \text{vector}[2])$  tuple.
- **shape** - object, collection of objects, or a list  $[a, b, c, \dots]$  of objects of type CONVEX, MESH, SPHERE. Alternately this can be a single  $(x, y, z)$  tuple.
- **vector** -  $(vx, vy, vz)$  tuple defining the translation

### **obj = ROTATE (shape, point, vector, angle)**

This routine rotates a geometrical object or a collection of such objects.

- **obj** - when **shape** is not  $(x, y, z)$  tuple: same as **shape**, returned for convenience. Otherwise the rotated  $(x1, y1, z1)$  image of  $(x, y, z)$ .
- **shape** - object, collection of objects, or a list  $[a, b, c, \dots]$  of objects of type CONVEX, MESH, SPHERE. Alternately this can be a single  $(x, y, z)$  tuple.
- **point** -  $(px, py, pz)$  tuple defining a point passed by the rotation axis
- **vector** -  $(vx, vy, vz)$  tuple defining a direction of the rotation axis
- **angle** - rotation angle in degrees

### **(one, two) = SPLIT (shape, point, normal)**

This routine splits a geometrical object (or a collection of objects) by a plane passing by a point.

- **one** - objects placed below the splitting plane (*None* if no objects were placed below)
- **two** - objects placed above the splitting plane (*None* if no objects were placed above)

## 2 Input language

- **shape** (emptied) - object, collection of objects, or a list  $[a, b, c, \dots]$  of objects of type CONVEX or SPHERE
- **point** -  $(px, py, pz)$  tuple defining a point passed by the splitting plane
- **normal** -  $(nx, ny, nz)$  tuple defining the splitting plane normal

### **obj = COPY (shape)**

This routine makes a copy of input objects.

- **obj** - created collection of copied objects
- **shape** - object, collection of objects, or a list  $[a, b, c, \dots]$  of objects of type CONVEX, MESH, SPHERE

### **obj = BYLABEL (solfec, kind, label)**

This routine finds a labeled object inside of a SOLFEC object.

- **obj** - returned object (*None* if a labeled object was not found)
- **solfec** - SOLFEC object
- **kind** - labeled object: 'SURFACE\_MATERIAL', 'BULK\_MATERIAL', 'BODY'
- **label** - the label string

### **obj = MASS\_CENTER (shape)**

This routine calculates the mass center of a geometrical object or a collection of such objects.

- **obj** -  $(x, y, z)$  tuple storing the mass center
- **shape** - object, collection of objects, or a list  $[a, b, c, \dots]$  of objects of type CONVEX, MESH, SPHERE. Alternately this can be a single BODY tuple.

### **CONTACT\_EXCLUDE\_BODIES (solfec, body1, body2)**

This routine disables contact detection for a specific pair of bodies. By default contact detection is enabled for all possible body pairs. **NOTE:** *must be invoked on all processors during a parallel run (do not use from within a callback).*

- **solfec** - SOLFEC object
- **body1** - first BODY object
- **body2** - second BODY object

### **CONTACT\_EXCLUDE\_OBJECTS (solfec, body1, point1, body2, point2)**

This routine disables contact detection for a specific pair of geometric objects (e.g. elements, convices, sheres). By default, between different bodies, contact detection is enabled for all possible object pairs. **NOTE:** *must be invoked on all processors during a parallel run (do not use from within a callback).*

- **solfec** - SOLFEC object
- **body1** - first BODY object
- **point1** - referential point properly contained in the 1st geometric object
- **bod2** - second BODY object
- **point2** - referential point properly contained in the 2nd geometric object

### **CONTACT\_SPARSIFY (solfec, threshold)**

This routine modifies contact filtering (sparsification) behaviour. Generally speaking, some contact points are filtered out in order to avoid unnecessary dense contact point clusters. If a pair of bodies is connected by two or more contact points, one of the points generated by topologically adjacent entities (elements, convices) will be removed (sparsified) if the ratio of contact areas of is smaller than the prescribed threshold.

- **solfec** - SOLFEC object
- **threshold** - sparsification threshold (default: 0.01) from within the interval [0, 1]. Zero corresponds to the lack of sparsification.

## **2.15 Results access**

Results can be accessed either in the 'READ' mode of a SOLFEC object, or in the 'WRITE' mode once some analysis has been run.

### **value = DURATION (solfec)**

This routine returns the duration of a simulation in SOLFEC's 'READ' mode, or *solfec.time* in the 'WRITE' mode.

- **value** - (*t0*, *t1*) duration limits of the simulation in 'READ' mode or current *time* in 'WRITE' mode
- **solfec** - SOLFEC object



### **FORWARD (solfec, steps)**

This routine steps forward within the simulation output file. Ignored in SOLFEC's 'WRITE' mode.

- **solfec** - SOLFEC object
- **steps** - numbers of steps forward

### **BACKWARD (solfec, steps)**

This routine steps backward within the simulation output file. Ignored in SOLFEC's 'WRITE' mode.

- **solfec** - SOLFEC object
- **steps** - number of steps backward

### **SEEK (solfec, time)**

This routine to a specific time within the simulation output file. Ignored in SOLFEC's 'WRITE' mode.

- **solfec** - SOLFEC object
- **time** - time to start reading at

### **disp = DISPLACEMENT (body, point)**

This routine outputs the displacement of a referential point.

- **disp** -  $(dx, dy, dz)$  tuple storing the displacement
- **body** - BODY object
- **point** -  $(x, y, z)$  tuple storing the referential point

### **velo = VELOCITY (body, point)**

This routine outputs the velocity of a referential point.

- **velo** -  $(vx, vy, vz)$  tuple storing the velocity
- **body** - BODY object
- **point** -  $(x, y, z)$  tuple storing the referential point

### **stre = STRESS (body, point)**

This routine outputs the Cauchy stress of a referential point.

- **stre** -  $(sx, sy, sz, sxy, sxz, syz, mises)$  tuple storing the Cauchy stress and the von Mises norm of it
- **body** - BODY object
- **point** -  $(x, y, z)$  tuple storing the referential point

### **ene = ENERGY (solfec | object)**

The routine outputs the value of energy of a specific object.

- **ene** -  $(kinetic, internal, external, conact, friction)$  tuple of energy values; *internal* energy corresponds to the work of internal forces, *external* energy corresponds to the work of external forces (including constraint reactions), *contact* energy corresponds to the work of normal contact reactions, *friction* energy corresponds to the work of tangential contact reactions
- **solfec** - SOLFEC object
- **object** - SOLFEC object, BODY object or a list of BODY objects

### **tim = TIMING (solfec, kind)**

The routine outputs the value of a specific action timing per time step.

- **tim** - value of timing
- **solfec** - SOLFEC object
- **kind** - this is one of: 'TIMINT' (time integration), 'CONDET' (contact detection), 'LOCDYN' (local dynamics setup), 'CONSOL' (constraints solution), 'PARBAL' (parallel load balancing). The load balancing timing is non-zero only for parallel runs.

### **hist = HISTORY (solfec, list, t0, t1 | skip)**

This routine outputs time histories of entities.

- **hist** - a tuple of list objects storing the histories:  $(times, values1, values2, ..., valuesN)$
- **solfec** - SOLFEC object
- **list** - list of objects  $[object1, object2, ..., objectN]$  indicating requested values. The valid objects are:

## 2 Input language

- a tuple  $(body, point, entity)$  where *body* is a BODY object, *point* is a  $(x, y, z)$  tuple storing the referential point, and *entity* is one of: 'DX', 'DY', 'DZ' (displacement), 'VX', 'VY', 'VZ' (velocity), 'SX', 'SY', 'SZ', 'SXY', 'SXZ', 'SYZ' (stress), 'MISES' (von Mises norm of stress)
  - a tuple  $(object, kind)$  where *object* is a SOLFE object, a BODY object or a list of BODY objects, and *kind* is a string 'KINETIC', 'INTERNAL', 'EXTERNAL', 'CONTACT', 'FRICTION' and it corresponds to the energy kind
  - a string 'TIMINT', 'CONDET', 'LOCDDYN', 'CONSOL', 'PARBAL' for timing histories
- **t0** - time interval start
  - **t1** - time interval end
  - **skip** - number of steps to skip between two time instants

## 3 Tutorials

## 4 Materials

## 5 Solvers

# Bibliography

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- [3] Tomasz Koziara and Nenad Bićanić. Simple and efficient integration of rigid rotations suitable for constraint solvers. *To appear in the International Journal for Numerical Methods in Engineering*, 2009.