Domain mapping with LAR *

Alberto Paoluzzi April 15, 2014

Abstract

In this module a first implementation (no optimisations) is done of several LAR operators, reproducing the behaviour of the plasm STRUCT and MAP primitives, but with better handling of the topology, including the stitching of decomposed (simplicial domains) about their possible sewing. A definition of specialised classes Model, Mat and Verts is also contained in this module, together with the design and the implementation of the *traversal* algorithms for networks of structures.

Contents

1 Introduction

The mapper module, introduced here, aims to provide the tools needed to apply both dimension-independent affine transformations and general simplicial maps to geometric objects and assemblies developed within the LAR scheme.

For this purpose, a simplicial decomposition of the $[0,1]^d$ hypercube $(d \ge 1)$ with any possible shape is firstly given, followed by its scaled version with any according $\mathtt{size} \in \mathbb{E}^d$, being its position vector the mapped image of the point $\mathbf{1} \in \mathbb{E}^d$. A general mapping mechanism is specified, to map any domain decomposition (either simplicial or not) with a given set of ccordinate functions, providing a piecewise-linear approximation of any curved embedding of a d-dimensional domain in any \mathbb{E}^n space, with $n \ge d$. A suitable function is therefore given to identify corresponding vertices when mapping a domain decomposition of the fundamental polygon of a closed manifold.

The geometric tools given in this chapter employ a normalised homogeneous representation of vertices of the represented shapes, where the added coordinate is the *last* of the ordered list of vertex coordinates. The homogeneous representation of vertices is used *implicitly*, by inserting the extra coordinate only when needed by the operation at hand,

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mainly for computing the product of the object's vertices times the matrix of an affine tensor.

A set of primitive surface and solid shapes is also provided, via the mapping mechanism of a simplicial decomposition of a d-dimensional chart. A simplified version of the PLaSM specification of dimension-independent elementary affine transformation is given as well.

The second part of this module is dedicated to the development of a complete framework for the implementation of hierarchical assemblies of shapes and scene graphs, by using the simplest possible set of computing tools. In this case no hierarchical graphs or multigraph are employed, i.e. no specialised data structures are produced. The ordered list model of hierarchical structures, inherited from PHIGS and PLaSM, is employed in this context. A recursive traversal is used to transform all the component parts of a hierarchical assembly into the reference frame of the first object of the assembly, i.e. in world coordinates.

2 Piecewise-linear mapping of topological spaces

A very simple but foundational software subsystem is developed in this section, by giving a general mechanism to produce curved maps of topological spaces, via the simplicial decomposition of a chart, i.e. of a planar embedding of the fundamental polygon of a d-dimensional manifold, and the definition of coordinate functions to be applied to its vertices (0-cells of the decomposition) to generate an embedding of the manifold.

2.1 Domain decomposition

A simplicial map is a map between simplicial complexes with the property that the images of the vertices of a simplex always span a simplex. Simplicial maps are thus determined by their effects on vertices, and provide a piecewise-linear approximation of their underlying polyhedra.

Standard and scaled decomposition of unit domain The larDomain of given shape is decomposed by larSimplexGrid as an hypercube of dimension $d \equiv len(shape)$, where the shape tuple provides the number or row, columns, pages, etc. of the decomposition.

@D Generate a simplicial decomposition of the $[0,1]^d$ domain @""" simplicial decomposition of the unit d-cube """ def larDomain(shape): V,CV = larSimplexGrid(shape) V = scalePoints(V, [1./d for d in shape]) return V,CV @

A scaled simplicial decomposition is provided by the second-order larIntervals function, with len(shape) and len(size) parameters, where the d-dimensionale vector len(size) is assumed as the scaling vector to be applied to the point $\mathbf{1} \in \mathbb{E}^d$.

@D Scaled simplicial decomposition of the $[0,1]^d$ domain @def larIntervals(shape): def larIntervals0(size): V,CV = larDomain(shape) V = scalePoints(V, [scaleFactor for scaleFactor in size]) return V,CV return larIntervals0 @

2.2 Mapping domain vertices

The second-order textttlarMap function is the LAR implementation of the PLaSM primitive MAP. It is applied to the array coordFuncs of coordinate functions and to the simplicially decomposed domain, returning an embedded and/or curved domain instance.

@D Primitive mapping function @def larMap(coordFuncs): def larMap0(domain): V,CV = domain V = TRANS(CONS(coordFuncs)(V)) plasm CONStruction return checkModel((V,CV)) return larMap0 @

2.3 Identify close or coincident points

The function checkModel, applied to a model parameter, i.e. to a (vertices, cells) pair, returns the model after identification of vertices with coincident or very close position vectors. The checkModel function works as follows: first a dictionary vertDict is created, with key a suitably approximated position converted into a string by the vcode converter (given in the Appendix), and with value the list of vertex indices with the same (approximated) position. Then, an invertedindex array is created, associating each original vertex index with the new index produced by enumerating the (distinct) keys of the dictionary. Finally, a new list CV of cells is created, by substituting the new vertex indices for the old ones.

@D Create a dictionary with key the point location @def checkModel(model): V,CV = model; n = len(V) vertDict = defaultdict(list) for k,v in enumerate(V): vertDict[vcode(v)].append(k) points,verts = TRANS(vertDict.items()) invertedindex = [None]*n <math>V = [] for k,v alue in enumerate(verts): V.append(eval(points[k])) for i in value: invertedindex[i]=k CV = [[invertedindex[<math>v] for v in cell] for cell in v0 filter out degenerate cells v0 for cell in v1 filter out degenerate cells v2 for v3 for cell in v4 for cell in v5 filter out degenerate cells v6 for cell in v7 filter out degenerate cells v8 for cell in v9 for cell in v9 filter out degenerate cells v9 for cell in v9 for cell in v9 filter out degenerate cells v9 for cell in v9 for cell in v9 filter out degenerate cells v9 for cell in v9 for cell in v9 filter out degenerate cells v9 for cell in v9 filter out degenerate cells v9 for cell in v9 for cell in v9 filter out degenerate cells v9 for cell in v9 filter out degenerate cells v9 for cell in v9 filter out degenerate cells v9 filter out degenerate ce

2.4 Embedding or projecting LAR models

In order to apply 3D transformations to a two-dimensional LAR model, we must embed it in 3D space, by adding one more coordinate to its vertices.

Embedding or projecting a geometric model This task is performed by the function larembed with parameter k, that inserts its d-dimensional geometric argument in the $x_{d+1}, \ldots, x_{d+k} = 0$ subspace of \mathbb{E}^{d+k} . A projection transformation, that removes the last k coordinate of vertices, without changing the object topology, is performed by the function larembed with negative integer parameter.

@D Embedding and projecting a geometric model @def larEmbed(k): def larEmbed0(model): V,CV = model if k $_i$ 0: V = [v+[0.]*k for v in V] elif k $_i$ 0: V = [v[:-k] for v in V] return V,CV return larEmbed0 @

3 Primitive objects

A large number of primitive surfaces or solids is defined in this section, using the larMap mechanism and the coordinate functions of a suitable chart.

3.1 1D primitives

 $\begin{array}{lll} \textbf{Circle} & @D \ Circle \ centered \ in \ the \ origin \ @def \ lar Circle (radius=1.): \ def \ lar Circle 0 (shape=36): \\ domain = \ lar Intervals ([shape])([2*PI]) \ V, CV = \ domain \ x = \ lambda \ V : [radius*COS(p[0]) \\ for \ p \ in \ V] \ y = \ lambda \ V : [radius*SIN(p[0]) \ for \ p \ in \ V] \ return \ lar Map([x,y])(domain) \\ return \ lar Circle 0 \ @ \\ \end{array}$

3.2 2D primitives

Some useful 2D primitive objects either in \mathbb{E}^2 or embedded in \mathbb{E}^3 are defined here, including 2D disks and rings, as well as cylindrical, spherical and toroidal surfaces.

 $\begin{array}{ll} \textbf{Disk} & @D \ \text{Disk centered in the origin @def larDisk(radius=1.): def larDisk0(shape=[36,1]):} \\ domain = larIntervals(shape)([2*PI,radius]) \ V,CV = domain \ x = lambda \ V : [p[1]*COS(p[0]) \\ for \ p \ in \ V] \ y = lambda \ V : [p[1]*SIN(p[0]) \ for \ p \ in \ V] \ return \ larMap([x,y])(domain) \ return \ larDisk0 \ @ \\ \end{array}$

Ring @D Ring centered in the origin @def larRing(params): r1,r2 = params def larRing0(shape=[36,1]): V,CV = larIntervals(shape)([2*PI,r2-r1]) V = translatePoints(V,[0,r1]) domain = V,CV x = lambda V : [p[1] * COS(p[0]) for p in V] y = lambda V : [p[1] * SIN(p[0]) for p in V] return larMap([x,y])(domain) return larRing0 @

 $\begin{array}{ll} \textbf{Cylinder surface} & @D \ \text{Cylinder surface with } z \ \text{axis} \ @from \ \text{scipy.linalg import det }""' \ \text{def} \\ \text{makeOriented(model): } V, CV = \text{model out} = [] \ \text{for cell in CV: mat} = \text{scipy.array}([V[v]+[1] \ \text{for } v \ \text{in cell}] + [[0,0,0,1]]) \ \text{if } \ \text{det}(\text{mat}) \ \text{; } 0.0 \ \text{out.append(cell)} \ \text{else: out.append([cell[1]]+[cell[0]]+cell[2:])} \\ \text{return } V, \text{out }""" \ \text{def} \ \text{larCylinder(params): radius, height= params } \ \text{def} \ \text{larCylinder0(shape=[36,1]):} \\ \text{domain} = \text{larIntervals(shape)}([2*PI,1]) \ V, CV = \text{domain } x = \text{lambda } V : [\text{radius*COS(p[0])} \ \text{for p in } V] \ y = \text{lambda } V : [\text{radius*SIN(p[0])} \ \text{for p in } V] \ z = \text{lambda } V : [\text{height*p[1]} \ \text{for p in } V] \ \text{mapping} = [x,y,z] \ \text{model} = \text{larMap(mapping)(domain)} \ \text{model} = \text{makeOriented(model)} \\ \text{return model return } \text{larCylinder0} \ @ \\ \end{array}$

Spherical surface of given radius @D Spherical surface of given radius @def lar-Sphere(radius=1): def larSphere0(shape=[18,36]): V,CV = larIntervals(shape)([PI,2*PI]) V = translatePoints(V,[-PI/2,-PI]) domain = V,CV x = lambda V : [radius*COS(p[0])*COS(p[1]) for p in V] y = lambda V : [radius*COS(p[0])*SIN(p[1]) for p in V] z = lambda V : [radius*SIN(p[0]) for p in V] return larMap([x,y,z])(domain) return larSphere0 @

Toroidal surface @D Toroidal surface of given radiuses @def larToroidal(params): r,R = params def larToroidal0(shape=[24,36]): domain = larIntervals(shape)([2*PI,2*PI]) V,CV = domain x = lambda V : [(R + r*COS(p[0])) * COS(p[1]) for p in V] y = lambda V : [(R + r*COS(p[0])) * SIN(p[1]) for p in V] z = lambda V : [-r * SIN(p[0]) for p in V] return larMap([x,y,z])(domain) return larToroidal0 @

Crown surface @D Half-toroidal surface of given radiuses @def larCrown(params): r,R = params def larCrown0(shape=[24,36]): V,CV = larIntervals(shape)([PI,2*PI]) V = translatePoints(V,[-PI/2,0]) domain = V,CV x = lambda V : [(R + r*COS(p[0])) * COS(p[1]) for p in V] y = lambda V : [(R + r*COS(p[0])) * SIN(p[1]) for p in V] z = lambda V : [-r * SIN(p[0]) for p in V] return larMap([x,y,z])(domain) return larCrown0 @

3.3 3D primitives

Ball @D Solid Sphere of given radius @def larBall(radius=1): def larBall0(shape=[18,36]): V,CV = checkModel(larSphere(radius)(shape)) return V,[range(len(V))] return larBall0 @

Solid cylinder @D Solid cylinder of given radius and height @def larRod(params): radius,height= params def larRod0(shape=[36,1]): V,CV = checkModel(larCylinder(params)(shape)) return V,[range(len(V))] return larRod0 @

Solid torus @D Solid torus of given radiuses @def larTorus(params): r,R = params def larTorus(shape=[24,36,1]): domain = larIntervals(shape)([2*PI,2*PI,r]) V,CV = domain x = lambda V : [(R + p[2]*COS(p[0])) * COS(p[1]) for p in V] y = lambda V : [(R + p[2]*COS(p[0])) * SIN(p[1]) for p in V] z = lambda V : [-p[2] * SIN(p[0]) for p in V] return larMap([x,y,z])(domain) return larTorus0 @

Solid pizza @D Solid pizza of given radiuses @def larPizza(params): r,R= params def larPizza0(shape=[24,36]): V,CV= checkModel(larCrown(params)(shape)) return V,[range(len(V))] return larPizza0 @

4 Affine transformations

4.1 Design decision

First we state the general rules that will be satisfied by the matrices used in this module, mainly devoted to apply affine transformations to vertices of models in structure environments:

1. assume the scipy ndarray as the type of vertices, stored in row-major order;

- 2. use the last coordinate as the homogeneous coordinate of vertices, but do not store it explicitly;
- 3. store explicitly the homogeneous coordinate of transformation matrices.
- 4. use labels 'verts' and 'mat' to distinguish between vertices and transformation matrices.
- 5. transformation matrices are dimension-independent, and their dimension is computed as the length of the parameter vector passed to the generating function.

4.2 map

@D Apply an affine transformation to a LAR model @def larApply(affineMatrix): def larApply0(model): if isinstance(model,Model): V = scipy.dot([v.tolist()+[1.0] for v in model.verts], affineMatrix.T).tolist() $V = [v[:-1] \text{ for } v \text{ in } V] \text{ CV} = copy(model.cells) \text{ d} = copy(model.d) \text{ return } Model((V,CV),d) \text{ elif isinstance}(model,tuple): } V,CV = model V = scipy.dot([v+[1.0] \text{ for } v \text{ in } V], \text{ affineMatrix.T}).tolist() \text{ return } [v[:-1] \text{ for } v \text{ in } V],CV \text{ return } \text{ larApply0} @$

4.3 Elementary matrices

Elementary matrices for affine transformation of vectors in any dimensional vector space are defined here. They include translation, scaling, rotation and shearing.

Translation @D Translation matrices @def t(*args): d = len(args) mat = scipy.identity(d+1) for k in range(d): mat[k,d] = args[k] return mat.view(Mat) @

Scaling @D Scaling matrices @def s(*args): d = len(args) mat = scipy.identity(d+1) for k in range(d): mat[k,k] = args[k] return mat.view(Mat) @

Rotation @D Rotation matrices @def r(*args): args = list(args) n = len(args) @; plane rotation (in 2D) @; space rotation (in 3D) @; return mat.view(Mat) @ @D plane rotation (in 2D) @if n == 1: rotation in 2D angle = args[0]; cos = COS(angle); sin = SIN(angle) mat = scipy.identity(3) mat[0,0] = cos; mat[0,1] = -sin; mat[1,0] = sin; mat[1,1] = cos; @ @D space rotation (in 3D) @if n == 3: rotation in 2D mat = scipy.identity(4) angle = VECTNORM(args); axis = UNITVECT(args) cos = COS(angle); sin = SIN(angle) @; elementary rotations (in 3D) @; @; general rotations (in 3D) @; @ @D elementary rotations (in 3D) @if axis[1]==axis[2]==0.0: rotation about x mat[1,1] = cos; mat[1,2] = -sin; mat[2,1] = sin; mat[2,2] = cos; elif axis[0]==axis[2]==0.0: rotation about z mat[0,2] = sin; mat[2,0] = -sin; mat[1,0] = sin; mat[1,1] = cos; @ @D rotation about z mat[0,0] = cos; mat[0,1] = -sin; mat[1,0] = sin; mat[1,1] = cos; @ @D

general rotations (in 3D) @else: general 3D rotation (Rodrigues' rotation formula) I = scipy.identity(3); u = axis Ux = scipy.array([[0, -u[2], u[1]], [u[2], 0, -u[0]], [-u[1], u[0], 0]) UU = scipy.array([[u[0]*u[0], u[0]*u[1], u[0]*u[2]], [u[1]*u[0], u[1]*u[1], u[1]*u[2]], [u[2]*u[0], u[2]*u[1], u[2]*u[2]])) mat[:3,:3] = cos*I + sin*Ux + (1.0-cos)*UU @

5 Hierarchical complexes

Hierarchical models of complex assemblies are generated by an aggregation of subassemblies, each one defined in a local coordinate system, and relocated by affine transformations of coordinates. This operation may be repeated hierarchically, with some subassemblies defined by aggregation of simpler parts, and so on, until one obtains a set of elementary components, which cannot be further decomposed.

Two main advantages can be found in a hierarchical modeling approach. Each elementary part and each assembly, at every hierarchical level, are defined independently from each other, using a local coordinate frame, suitably chosen to make its definition easier. Furthermore, only one copy of each component is stored in the memory, and may be instanced in different locations and orientations how many times it is needed.

5.1 Traversal of hierarchical structures

Of course, the main algorithm with hierarchical structures is the *traversal* of the structure network, whose aim is to transform every encountered object from local to global coordinates, where the global coordinates are those of the network root (the only node with indegree zero).

A structure network can be modelled using a directed acyclic multigraph, i.e. a triple (N, A, f) made by a set N of nodes, a set A of arcs, and a function $f: A \to N^2$ from arcs to ordered pairs of nodes. Conversely that in standard oriented graphs, in this kind of structure more than one oriented arc is allowed between a pair on nodes.

A simple modification of a DFS (Depth First Search) visit of a graph can be used to traverse the structure network This algorithm is given in Figure ?? from Reference [?].

5.1.1 Traversal of nested lists

The representation chosen for structure networks with LAR is the serialised one, consisting in ordered sequences (lists) of either LAR models, or affine transformations, or references to other structures, either directly nested within some given structure, or called by reference (name) from within the list.

The aim of a structure network traversal is, of course, to transform every component structure, usually defined in a local coordinate system, into the reference frame of the structure as a whole, normally corresponding with the reference system of the structure's root, called the *world coordinate* system.

```
Script 8.3.1 (Traversal of a multigraph)
algorithm Traversal ((N, A, f) : multigraph) {
   CTM := identity matrix;
   TraverseNode (root)
proc TraverseNode (n:node) {
   foreach a \in A outgoing from n do TraverseArc (a);
   ProcessNode (n)
}
proc TraverseArc (a = (n, m) : arc) {
   Stack.push (CTM);
   CTM := CTM * a.mat;
   TraverseNode (m);
   CTM := Stack.pop()
}
proc ProcessNode (n : node) {
   foreach object \in n do Process( CTM * object)
```

Figure 1: Traversal algorithm of an acyclic multigraph.

The pattern of calls and returned values In order to better understand the task of the traversal algorithm, where every transformation is applied to all the following models, — only if included in the same list (i.e. structure) — it is very useful to start with some algorithm emulation. In particular, the recursive script below discriminate between three different cases (number, string, or sequence) as the actual traversal must do with Models, Matrices, and Structures, respectively.

@D Emulation of scene multigraph traversal @from pyplasm import * def $_{traverse(CTM,stack,o):foriinrange(len(o))}$ def algorithm(data): CTM,stack = ["I"],[] $_{traverse(CTM,stack,data)}$ @

Some use example of the above algorithm are provided below. The printout produced at run time is shown from the emulation of traversal algorithm macro. @D Examples of multigraph traversal @data = [1, "A", 2, 3, "B", [4, "C", 5], [6, "D", "E", 7, 8], 9] print algorithm(data) data = [1, "A", [2, 3, "B", 4, "C", 5, 6, "D"], "E", 7, 8, 9] print algorithm(data) data = [2, 3, "B", 4, "C", 5, 6, "D"] print algorithm(data) dat = [2, 3, "B", 4, "C", 5, 6, "D"] print algorithm(data) dat = [2, 3, "B", 4, "C", 5, 6, "D"] data = [1, "A", dat, "E", 7, 8, 9] print algorithm(data) @ @D Emulation of traversal algorithm @1 ['1'] 2 ['A', '1'] 3 ['A', '1'] 4 ['B', 'A', '1'] 5 ['C', 'B', 'A', '1'] 9 ['B', 'A', '1'] None 1 ['1'] 2 ['A', '1'] 3 ['A', '1'] 4 ['B', 'A', '1'] 8 ['E', 'A', '1'] 6 ['C', 'B', 'A', '1'] 7 ['E', 'A', '1'] None 1 ['1'] 2 ['A', '1'] 3 ['A', '1'] 4 ['B', 'A', '1'] 5 ['C', 'B', 'A', '1'] 5 ['C', 'B', 'A', '1'] 7 ['E', 'A', '1'] None 1 ['1'] 2 ['A', '1'] 3 ['A', '1'] 4 ['B', 'A', '1'] 5 ['C', 'B', 'A', '1'] 6 ['C', 'B', 'A', '1'] 7 ['E', 'A', '1'] 8 ['E', 'A', '1'] 9 ['E', 'A', '1'] 4 ['B', 'A', '1'] 5 ['C', 'B', 'A', '1'] 6 ['C', 'B', 'A', '1'] 7 ['E', 'A', '1'] 8 ['E', 'A', '1'] 9 ['E'

Traversal of a scene multigraph The previous traversal algorithm is here customised for scene multigraph, where the objects are LAR models, i.e. pairs of vertices of type 'Verts and cells, and where the transformations are matrix transformations of type 'Mat'.

Check models for common dimension The input list of a call to larStruct primitive is preliminary checked for uniform dimensionality of the enclosed LAR models and transformations. The common dimension dim of models and matrices is returned by the function checkStruct, within the class definition Struct in the module lar2psm. Otherwise, an exception is generated (TODO).

Initialization and call of the algorithm @D Traversal of a scene multigraph @""" Traversal of a scene multigraph """ @; Structure traversal algorithm @;

def evalStruct(struct): dim = struct.n CTM, stack = scipy.identity(dim+1), [] scene = traverse(CTM, stack, struct) return scene @

Structure traversal algorithm @D Structure traversal algorithm @def traverse(CTM, stack, o, scene=[]): for i in range(len(o)): if isinstance(o[i],Model): scene += [larAp-ply(CTM)(o[i])] elif isinstance(o[i],Mat): CTM = scipy.dot(CTM, o[i]) elif isinstance(o[i],Struct): stack.append(CTM) traverse(CTM, stack, o[i], scene) CTM = stack.pop() return scene @

5.2 Example

Some simple examples of structures as combinations of LAR models and affine transformations are given in this section.

Global coordinates We start with a simple 2D example of a non-nested list of translated 2D object instances and rotation about the origin.

@O test/py/mapper/test04.py @""" Example of non-nested structure with translation and rotations """ @; Initial import of modules @; from mapper import * square = lar-Cuboids([1,1]) table = larApply(t(-.5,-.5))(square) chair = larApply(s(.35,.35))(table) chair1 = larApply(t(.75, 0))(chair) chair2 = larApply(r(PI/2))(chair1) chair3 = larApply(r(PI/2))(chair2) chair4 = larApply(r(PI/2))(chair3) VIEW(SKEL_1(STRUCT(MKPOLS(table) + MKPOLS(chair1) + MKPOLS(chair2) + MKPOLS(chair3) + MKPOLS(chair4))))@

Local coordinates A different composition of transformations, from local to global coordinate frames, is used in yie following example.

@O test/py/mapper/test05.py @""" Example of non-nested structure with translation and rotations """ @; Initial import of modules @; from mapper import * square = larCuboids([1,1]) square = Model(square,2) table = larApply(t(-.5,-.5))(square) chair = larApply(s(.35,.35))(table) chair = larApply(t(.75, 0))(chair) struct = Struct([table] + 4*[chair, r(PI/2)]) scene = evalStruct(struct) VIEW(SKEL₁(STRUCT(CAT(AA(MKPOLS)(scene)))))@ @O test/py/mapper/test06.py @""" Example of nested structures with translation and rotations """ @; Initial import of modules @; from mapper import * square = larCuboids([1,1]) square = Model(square,2) table = larApply(t(-.5,-.5))(square) chair = Struct([t(.75, 0), s(.35,.35), table]) struct = Struct([t(2,1)] + [table] + 4*[r(PI/2), chair]) scene = evalStruct(struct) VIEW(SKEL₁(STRUCT(CAT(AA(MKPOLS)(scene)))))@

6 Computational framework

6.1 Exporting the library

@O lib/py/mapper.py @""" Mapping functions and primitive objects """ @; Initial import of modules @; @; Generate a simplicial decomposition of the $[0,1]^d$ domain @; @; Scaled simplicial decomposition of the $[0,1]^d$ domain @; @; Create a dictionary with key the point location @; @; Primitive mapping function @; @; Basic tests of mapper module @; @; Circle centered in the origin @; @; Disk centered in the origin @; @; Ring centered in the origin @; @; Spherical surface of given radius @; @; Cylinder surface with z axis @; @; Toroidal surface of given radiuses @; @; Half-toroidal surface of given radiuses @; @; Solid Sphere of given radius @; @; Solid cylinder of given radius and height @; @; Solid torus of given radiuses @; @; Solid pizza of given radiuses @; @; Translation matrices @; @; Scaling matrices @; @; Rotation matrices @; @; Embedding and projecting a geometric

model @; @; Apply an affine transformation to a LAR model @; @; Traversal of a scene multigraph @; @

6.2 Examples

3D rotation about a general axis @O test/py/mapper/test02.py @""" General 3D rotation of a toroidal surface """ @; Initial import of modules @; from mapper import * model = checkModel(larToroidal([0.2,1])()) a,b,c = SCALARVECTPROD([PI/2,UNITVECT([1,1,0])]) model = larApply(r(a,b,c))(model) VIEW(STRUCT(MKPOLS(model))) @

3D rotation of a 2D circle @O test/py/mapper/test03.py @""" Elementary 3D rotation of a 2D circle """ @; Initial import of modules @; from mapper import * model = checkModel(larCircle(1)()) model = larEmbed(1)(model) model = larApply(r(PI/2,0,0))(model) VIEW(STRUCT(MKPOLS(model))) @

6.3 Tests

```
@D Basic tests of mapper module @if _{name_{=="main,:V,EV=larDomain([5])VIEW(EXPLODE(1.5,1.5,1.5)(MKPOLS((V,EV))))V,EV=larDomain([5,3])} VIEW(EXPLODE(1.5,1.5,1.5)(MKPOLS((V,FV)))) V,FV = larIntervals([36,3])([2*PI,1.]) VIEW(EXPLODE(1.5,1.5,1.5)(MKPOLS((V,FV)))) V,CV = larDomain([5,3,1]) VIEW(EXPLODE(1.5,1.5,1.5)(MKPOLS((V,CV)))) V,CV = larIntervals([36,2,3])([2*PI,1.,1.]) VIEW(EXPLODE(1.5,1.5,1.5)(MKPOLS((V,CV)))) @
```

Circumference of unit radius @O test/py/mapper/test01.py @""" Circumference of unit radius """ @; Initial import of modules @; from mapper import * model = check-Model(larCircle(1)()) VIEW(EXPLODE(1.2,1.2,1.2)(MKPOLS(model))) model = check-Model(larDisk(1)([36,4])) VIEW(EXPLODE(1.2,1.2,1.2)(MKPOLS(model))) model = check-Model(larRing([.9, 1.])([36,2])) VIEW(EXPLODE(1.2,1.2,1.2)(MKPOLS(model))) model = checkModel(larCylinder([.5,2.])([32,1])) VIEW(STRUCT(MKPOLS(model))) model = larBall(1)() VIEW(STRUCT(MKPOLS(model))) model = larBall(1)() VIEW(STRUCT(MKPOLS(model))) model = checkModel(larToroidal([0.5,1]) VIEW(STRUCT(MKPOLS(model))) model = checkModel(larToroidal([0.5,1]) VIEW(STRUCT(MKPOLS(model))) model = checkModel(larToroidal([0.5,1]) VIEW(STRUCT(MKPOLS(model))) model = check-Model(larTorus([0.5,1])([0.5,1])([0.5,1])() VIEW(STRUCT(MKPOLS(model))) @

A Utility functions

@D Initial import of modules @from pyplasm import * from scipy import * import os,sys """ import modules from larcc/lib """ sys.path.insert(0, 'lib/py/') @; Import the module @(lar2psm@) @; @; Import the module @(simplexn@) @; @; Import the module

@(larcc@) @; @; Import the module @(largrid@) @; @; Import the module @(boolean2@) @; @

@D Import the module @import @1 from @1 import * @

A.1 Numeric utilities

A small set of utility functions is used to transform a point representation as array of coordinates into a string of fixed format to be used as point key into python dictionaries.

@D Symbolic utility to represent points as strings @""" TODO: use package Decimal (http://docs.python.org/2/library/decimal.html) """ ROUND $_ZERO=1E-07defround_or_zero(x,prec=7):$ """ Decision procedure to approximate a small number to zero. Returne ither the input number or zero.""" definition to the first procedure to approximate a small number to zero. Returne ither the input number or zero."" definition to the first procedure to approximate a small number to zero. The first part of the first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first procedure to approximate a small number to zero. The first proced

def fixedPrec(value): if abs(value - int(value)); ROUND $_ZERO$: value = int(value)out = ('ifout == '-0.': out = '0.'returnout

def vcode (vect): """ To generate a string representation of a number array. Used to generate the vertex keys in PointSet dictionary, and other similar operations. """ return prepKey(AA(fixedPrec)(vect)) @