Boundary operators on LAR *

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Abstract

The various versions of boundary operators on Linear Algebraic Representation of cellular complexes are developed in this module, in order to maintain under focus their proper development, including the possible special cases.

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1 Introduction

In the current LarLib implementation, we have to distinguish between between dimension-independent, dimension-dependent, signed and non-signed operators. Therefore a code refactoring of LarLib—related to boundary/coboundary operators—started here, with the aim of both providing a precise mathematical definition within the LAR framework, and to simplify and generalise the implemented algorithms.

2 Implementation

We start this section by making a distinction between the (matrices of) boundary operators for the linear spaces C_k of chains over the field $\mathbb{Z}_2 = \{0,1\}$ and over the field \mathbb{Z} of integer numbers. We call either non-signed or signed the corresponding boundary operators, respectively, since the matrix elements take values within the sets $\{-1,0,+1\}$ or $\{0,1\}$, correspondingly. Of course, the associated matrices of coboundary operators are their transpose matrices.

2.1 Non-signed operators

For several computations, the knowledge of the matrices of non-signed boundary operators is sufficient. Therefore we will use such tool wherever possible, since its computation is much faster in term of computing time.

In the following we provide be binary operator matrices provided by two implementations, respectively named boundary and boundary1. The first one works correctly only with convex cells; the second one works also with non-convex but path-connected cells.

2.1.1 Dimension-independence

As we show in the following, in order to compute the non-signed boundary operator ∂_d , it it sufficient to have knowledge of the M_d and M_{d-1} characteristic matrices of d-cells and their (d-1)-facets, at least in the case of cellular complexes with convex cells. Conversely, for more general non-convex but simply-connected cells, also the M_{d-2} matrix is needed.

Convex-cells The algorithm used is pretty easy to present. The compressed characteristic matrices of d-cells and (d-1)-cells, denoted as **cells** and **facets**, respectively, are first put in **csr** format as **csrCV** and **csrFV**. Then the incidence matrix **csrFC** in compressed sparse row format is computed by matrix product of the compressed characteristic matrices.

The element (i, j) of this matrix provides the number of vertices in the intersection of facet i and cell j, whereas the number of non-zero elements in each csrFV row gives the number of vertices of the facet represented by the row, and is stored in facetLengths.

The boundary function—to be used only with dimension-independent LAR convex cells—is written efficiently in the following script, by using only the standard functions and attributes of the scipy.sparse module.

The variable facetCoboundary stores in a list, for every facet (for h in range(m)) the list of cells in its *coboundary*, to be stored in the output csr_matrix boundary matrix as column indices of elements with non-zero (i.e. 1) value.

Notice that both the computation of facetCoboundary contents, and the output of the compressed boundary matrix, are performed in the most efficient way—according to the internal design of the scipy's csr sparse data structure.

```
\langle \text{ convex-cells boundary operator } 3 \rangle \equiv
     """ convex-cells boundary operator --- best implementation """
     def boundary(cells,facets):
         lenV = max(CAT(cells))+1
         csrCV = csrCreate(cells,lenV)
         csrFV = csrCreate(facets,lenV)
         csrFC = csrFV * csrCV.T
         facetLengths = [csrFacet.getnnz() for csrFacet in csrFV]
         m,n = csrFC.shape
         facetCoboundary = [[csrFC.indices[csrFC.indptr[h]+k]
              for k,v in enumerate(csrFC.data[csrFC.indptr[h]:csrFC.indptr[h+1]])
                  if v==facetLengths[h]] for h in range(m)]
         indptr = [0]+list(cumsum(AA(len)(facetCoboundary)))
         indices = CAT(facetCoboundary)
         data = [1]*len(indices)
         return csr_matrix((data,indices,indptr),shape=(m,n),dtype='b')
```

Macro referenced in 5a.

Non-convex LAR cells A more general boundary1 operator is given in the following, aiming at compute the boundary matrix for general non-convex cellular decompositions, including multiply connected LAR models. Notice that in this case an input triple made by CV, FV, and EV is needed, where—more in general embedded in \mathbf{E}^d —they stand for the (binary compressed) characteristic matrices M_d , M_{d-1} , and M_{d-2} .

First the boundary operator for the convex case is computed within the out variable of $\mathtt{csr_matrix}$ type. Then every out row (i.e. every (d-1)-facet of the d-complex) is tested for $\mathit{reliability}$, since every (d-1)-face can be shared by $\mathit{at\ most\ two\ d}$ -cells in a d-complex. When this condition is not satisfied, deeper tests are needed to understand what row elements must be forced to value 1, since the (d-1)-face itself is a subset, but not actually a facet, of the corresponding d-cell.

In presence of some "unreliable" facets, the matrix csrBBMat of the operator $\partial_{d-1} \circ \partial_d$ and the relation FE between faces of dimensions d-1 and d-2 are computed. Now, let us notice that the columns of csrBBMat report the number of incidences of the d-2 faces (as

belonging to (d-1)-facets embedded on the boundary) and d-cells (that are associated to such matrix columns). Hence, in a regular (convex) d-complex, such numbers are always even, and in \mathbb{Z}_2 arithmetic are reduced to zero, in order to satisfy the fundaments equation $\partial \partial = 0$.

Conversely, with non-convex LAR cells, some incidence numbers may get odd values, due to the non-strict coincidence between cell facets and vertex subsets. Therefore, for "unreliable" h rows (facets) the csrBBMat columns tracked by ones in $[\partial_d]$ are checked, looking for elements of (h, k) indices with value greater that 2.

```
\langle path-connected-cells boundary operator 4a \rangle \equiv
     """ path-connected-cells boundary operator """
     import larlib
     import larcc
     from larcc import *
     def csrBoundaryFilter1(unreliable,out,csrBBMat,cells,FE):
         for row in unreliable:
              for j in range(len(cells)):
                  if out[row,j] == 1:
                      cooCE = csrBBMat.T[j].tocoo()
                      flawedCells = [cooCE.col[k] for k,datum in enumerate(cooCE.data)
                           if datum>2]
                      if all([facet in flawedCells for facet in FE[row]]):
                           out[row,j]=0
         return out
     def boundary1(CV,FV,EV):
         out = boundary(CV,FV)
         def csrRowSum(h):
              return sum(out.data[out.indptr[h]:out.indptr[h+1]])
         unreliable = [h for h in range(len(FV)) if csrRowSum(h) > 2]
         if unreliable != ∏:
              csrBBMat = boundary(FV,EV) * boundary(CV,FV)
              print "\ncsrBBMat =",csrBBMat.todense(),"\n"
              lenV = max(CAT(CV))+1
              FE = larcc.crossRelation(lenV,FV,EV)
              out = csrBoundaryFilter1(unreliable,out,csrBBMat,CV,FE)
         return out
Macro referenced in 5a.
\langle From cells and facets to boundary cells 4b\rangle \equiv
     def totalChain(cells):
         return csr_matrix(len(cells)*[[1]])
```

```
def boundaryCells(cells,facets):
   csrBoundaryMat = boundary(cells,facets)
   csrChain = csr_matrix(totalChain(cells))
   csrBoundaryChain = csrBoundaryMat * csrChain
   for k,value in enumerate(csrBoundaryChain.data):
        if value % 2 == 0: csrBoundaryChain.data[k] = 0
   out = [k for k,val in enumerate(csrBoundaryChain.data.tolist()) if val == 1]
   return out
def boundaryCells1(cells,facets,faces):
    csrBoundaryMat = boundary1(cells,facets,faces)
    csrChain = csr_matrix(totalChain(cells))
   csrBoundaryChain = csrBoundaryMat * csrChain
   for k,value in enumerate(csrBoundaryChain.data):
        if value % 2 == 0: csrBoundaryChain.data[k] = 0
   out = [k for k,val in enumerate(csrBoundaryChain.data.tolist()) if val == 1]
   return out
```

Macro referenced in 5a.

2.2 Signed operators

3 Exporting

```
"larlib/larlib/boundary.py" 5a =

""" boundary operators """

from larlib import *

⟨convex-cells boundary operator 3⟩

⟨path-connected-cells boundary operator 4a⟩

⟨From cells and facets to boundary cells 4b⟩

⋄
```

4 Testing

5 Non-signed operators

Correct boundary extraction example The boundary.boundary() operator is applied here to a cellular 2-complex of convex cells, producing correct result.

```
"test/py/boundary/test01.py" 5b \equiv """ testing boundary operators (correct result) """ from larlib import *
```

```
filename = "test/svg/inters/boundarytest0.svg"
lines = svg2lines(filename)
VIEW(STRUCT(AA(POLYLINE)(lines)))

V,FV,EV,polygons = larFromLines(lines)
VV = AA(LIST)(range(len(V)))
submodel = STRUCT(MKPOLS((V,EV)))
VIEW(larModelNumbering(1,1,1)(V,[VV,EV,FV],submodel,0.2))
VIEW(EXPLODE(1.2,1.2,1.2)(MKPOLS((V,[EV[e] for e in boundaryCells(FV,EV)],))))
VIEW(EXPLODE(1.2,1.2,1.2)(MKTRIANGLES((V,FV,EV))))

boundaryOp = boundary1(FV,EV,VV)

for k in range(1,len(FV)+1):
    faceChain = k*[1]
    BF = chain2BoundaryChain(boundaryOp)(faceChain)
    VIEW(STRUCT(MKPOLS((V,[EV[e] for e in BF]))))
```

Wrong boundary extraction example The boundary() operator, applied to a cellular 2-complex wih some non-convex cells, produces incorrect results. In such cases a correct result may be produced only by chance (sometimes this happens). So, be careful to use it only when the precondition (of cell convexity) is everywhere verified. In order to get always a correct result, use the boundary1 operator.

```
VIEW(EXPLODE(1.2,1.2,1.2)(MKFACES((V,FV,EV))))
VIEW(SKEL_1(EXPLODE(1.2,1.2,1.2)(MKTRIANGLES((V,FV,EV)))))
for k in range(1,len(FV)+1):
    faceChain = k*[1]
    boundaryChain = chain2BoundaryChain(boundaryOp)(faceChain)
    VIEW(STRUCT(MKPOLS((V,[EV[e] for e in boundaryChain]))))
```

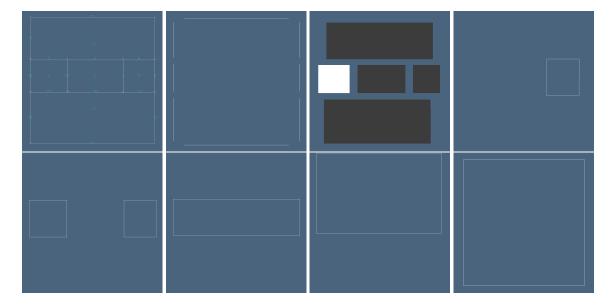


Figure 1: Convex-cell 2-complex. (a) Indexing of 0-,1-,and 2-cells; (b) exploded 2-boundary cells; (c) exploded 2-cells; (d) boundary of a singleton 2-chain; (e-h) boundaries of some 2-chains.

Example Comparison of two implementations of the ∂ operator. Notice the difference between the penultimate rows. In particular, the penultimate row of the matrix generated by boundary(FV,EV) is plain wrong. It means that the edge e_{10} is shared by all the (three) 2-cells of the complex. Conversely, it is well known that, for a solid complex, i.e. a d-complex embedded in \mathbb{E}^d , every (d-1)-facet may be shared by no more than 2 d-cells. The resulting boundary of the total chain $[f_0, f_1, f_2]$ codified in coordinates as [1, 1, 1], and shown in Figure 2d, is sonsequently incorrect.

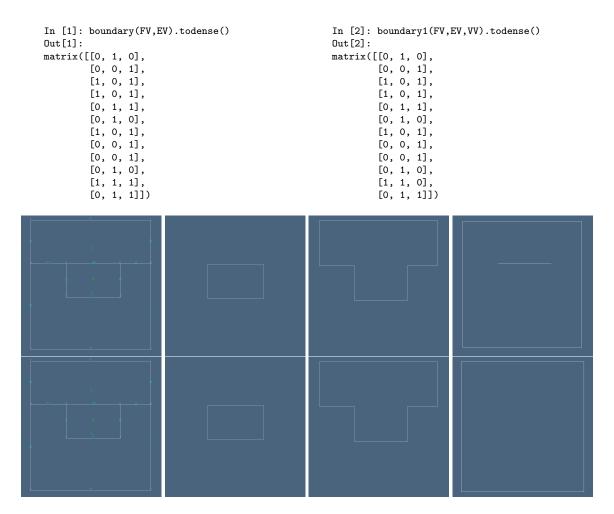


Figure 2: Non-working (i.e. *wrong*) example with boundary. (a) Indexing of 0-,1-,and 2-cells; (b) boundary of a singleton 2-chain; (c) exploded 2-cells; (d) boundary of a singleton 2-chain. Working (i.e. *exact*) example using boundary1: (e-h) as above.

3D non-convex LAR cells

```
"test/py/boundary/test03.py" 8 =
    """ 3D non-convex LAR cells """
    from larlib import *
    # LAR model with non-contractible faces
    # -------

V = [[0.25, 0.25, 0.0], [0.25, 0.75, 0.0], [0.75, 0.75, 0.0], [0.75, 0.25, 0.0], [1.0, 0.0, 0.0], [0.0, 0.0, 0.0], [1.0, 1.0, 0.0], [0.0, 1.0, 0.0], [0.25, 0.25, 1.0], [0.25,
```

```
0.25, 2.0, [0.25, 0.75, 2.0], [0.25, 0.75, 1.0], [0.25, 0.75, -1.0], [0.25, 0.25, 0.25, 0.25]
-1.0], [0.75, 0.75, -1.0], [0.75, 0.25, -1.0], [0.75, 0.25, 1.0], [0.75, 0.75, 1.0],
[1.0, 0.0, 1.0], [0.0, 0.0, 1.0], [1.0, 1.0, 1.0], [0.0, 1.0, 1.0], [0.75, 0.75, 2.0],
[0.75, 0.25, 2.0]
CV = [(0,1,2,3,4,5,6,7,8,11,16,17,18,19,20,21),
(0,1,2,3,8,11,16,17),
(0,1,2,3,12,13,14,15),
(8,9,10,11,16,17,22,23)
FV = [(2,3,16,17),(6,7,20,21),(12,13,14,15),(0,1,8,11),(1,2,11,17),(0,1,12,13),
(4,6,18,20),(5,7,19,21),(0,3,13,15),(0,3,8,16),(0,1,2,3),
(10,11,17,22),(2,3,14,15),(8,9,16,23),(8,11,16,17),
(1,2,12,14),(16,17,22,23),(4,5,18,19),(8,9,10,11),(
9,10,22,23),(0,1,2,3,4,5,6,7),(8,11,16,17,18,19,20,21)
EV = [(3,15),(7,21),(10,11),(4,18),(12,13),(5,19),(8,9),(18,19),(22,23),(0,3),(1,11),
(16,17),(0,8),(6,7),(20,21),(3,16),(10,22),(18,20),(19,21),(1,2),(12,14),(4,5),(12,14),(12,14),(13,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,14),(14,1
8,11),(13,15),(16,23),(14,15),(11,17),(17,22),(2,14),(2,17),(0,1),(9,10),(8,16),
(4,6),(1,12),(5,7),(0,13),(9,23),(6,20),(2,3)
VV = AA(LIST)(range(len(V)))
hpc = STRUCT(MKPOLS((V,EV)))
VIEW(larModelNumbering(1,1,1)(V,[VV,EV,FV,CV],hpc,0.6))
BF = boundaryCells1(CV,FV,EV)
VIEW(EXPLODE(1.2,1.2,1.2)(MKTRIANGLES((V,[FV[f] for f in BF],EV))))
```

References

[CL13] CVD-Lab, *Linear algebraic representation*, Tech. Report 13-00, Roma Tre University, October 2013.

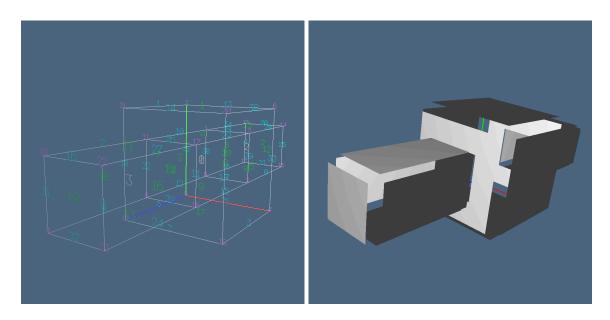


Figure 3: Non-convex 3-complex. (a) Indexing of 0-,1-,2- and 3-cells; (b) exploded 2-boundary cells. Notice that two faces are multiply-connected.