Extraction of a submap from a Cell Designer map (PART 3)

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Reviewer(s):

INTRODUCTION

Given a generic map of metabolism, creat a derivative map by removing a subset of the reactions.

EQUIPMENT SETUP

To visualise the metabolic maps it is necessary to obtain the version 4.4 of CellDesigner. This software can be freely downloaded from:

http://www.celldesigner.org/download.html

PROCEDURE

1. Import a CellDesigner XML file to MATLAB environment

The transformXML2Map function parses an XML file from Cell Designer (CD) into a Matlab structure. This structure is organised similarly to the structure found in the COnstraint-Base and Reconstruction Analysis (COBRA) models.

Read in a map

```
[GlyXml, GlyMap] = transformXML2Map('glycolysisAndTCA.xml');
Elapsed time is 1.834226 seconds.
```

2. Remove some reactions from the map

```
if 1
    rxnRemoveList={'ENO';'PFK';'PGMT';'ABC'};
else
    rxnRemoveList={'PGMT'};
end
printLevel=1;
[GlyXmlStructSubset,GlyMapSubset,rxnNotInMap] =
removeMapReactions(GlyXml,GlyMap,rxnRemoveList,printLevel);
```

ABC not present in the map

There is no reaction 'ABC' in the map, so the function alerts that it has not been removed.

```
rxnRemoveList(rxnNotInMap)
```

```
ans = 1 \times 1 cell array
```

3. Export the modified map

```
transformMap2XML(GlyXmlStructSubset,GlyMapSubset,'glycolysisAndTCA_subset.xml
');
```

Elapsed time is 1.317182 seconds.

4. Import the modified map and compare it with the matlab structures

```
[GlyXmlStructSubset2,GlyMapSubset2] =
transformXML2Map('glycolysisAndTCA_subset.xml');

Elapsed time is 1.667602 seconds.
ans = struct with fields:
    id: 're2'
metaid: 're57'
    name: 'DPGM'
ans = struct with fields:
    id: 're2'
metaid: 're57'
    metaid: 're57'
    name: 'DPGM'
reversible: 'true'
```

Compare xml structure

```
[resultXml, whyXml] = structeq(GlyXmlStructSubset, GlyXmlStructSubset2)
resultXml = logical
whyXml = struct with fields:
   Reason: 'Properties are different <- Unequal Subcell <- Properties are different <- Unequal Subcell <-
    Where: '(1).sbml(1).model(1).listOfReactions(1).reaction{2}(1).Attributes'
GlyXmlStructSubset.sbml(1).model(1).listOfReactions(1).reaction{2}
(1).Attributes
ans = struct with fields:
      id: 're2'
   metaid: 're57'
     name: 'DPGM'
GlyXmlStructSubset2.sbml(1).model(1).listOfReactions(1).reaction{2}
(1).Attributes
ans = struct with fields:
          id: 're2'
       metaid: 're57'
         name: 'DPGM'
   reversible: 'true'
```

Compare map structure

```
[resultMap, whyMap] = structeq(GlyMapSubset, GlyMapSubset2)
```

```
whyMap = struct with fields:
    Reason: ''
    Where: ''

return
```

4. Remove reactions from the map directly

resultMap = logical

```
fileName = 'glycolysisAndTCA.xml';
printLevel=1;
removeCDReactions(fileName,rxnRemoveList,printLevel)
```