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# Metamodel Matching for Automatic Model Transformation Generation<sup>\*</sup>

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**Abstract.** Applying Model-Driven Engineering (MDE) leads to the creation of a large number of metamodels, since MDE recommends an intensive use of models defined by metamodels. Metamodels with similar objectives are then inescapably created. A recurrent issue is thus to turn compatible models conforming to similar metamodels, for example to use them in the same tool. The issue is classically solved developing ad hoc model transformations. In this paper, we propose an approach that automatically detects mappings between two metamodels and uses them to generate an alignment between those metamodels. This alignment needs to be manually checked and can then be used to generate a model transformation. Our approach is built on the Similarity Flooding algorithm used in the fields of schema matching and ontology alignment. Experimental results comparing the effectiveness of the application of various implementations of this approach on real-world metamodels are given.

## 1 Introduction

The large success of Model-Driven Engineering leads to a relative abundance of metamodels, for various domains. Tools are developed to handle models conforming to the metamodels : code generators, model transformations and graphical editors. Inescapably, several metamodels for the same kind of applications are independently created. It is then quite natural to wish to use a tool-set adapted to a given metamodel for models conforming to another metamodel with the same purpose. For example, several metamodels exist to represent class models, and several tools exist to automatically refactor class models. The problem is to use a tool dedicated to a metamodel with a class model conforming to another metamodel allowing to express the same kind of class models. Even for the same application domain, various versions of the same metamodels usually exist, and the compatibility between the models conforming to the different versions has to be ensured. Usually, this problem is solved using manually written, ad hoc model transformations. Such model transformations are not difficult to write, but are numerous, and thus represent a large work load. In this paper, we propose to generate automatically an alignment between two similar metamodels. This

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alignment represents the correspondences between the elements of the source metamodel and those of the target metamodel. This alignment can then be used to generate the code of a model transformation.

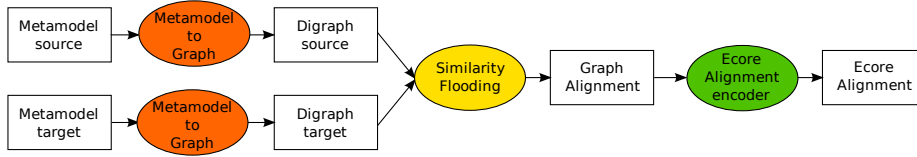
The match operation is well-known in many application domains, such as semantic web, schema and ontology integration, data warehouses, e-commerce [1,2]. It takes as input two schemas and produces as output a set of relations (*e.g.*, equivalence and subsumption) between the entities of the two input schemas. Such an output is often called an alignment of two schemas. The notion of schema can appear as unclear since it highly depends on the concerned domain. Concretely, a schema can be an XML Schema, an ontology, a database schema or an object-oriented class model. Despite the variation between those formats, the mechanisms involved to perform the match operation are highly similar.

In this paper, we report experiments which evaluate and adapt the Similarity Flooding algorithm [3] to compute metamodel alignments. The context of our approach is the automated generation of model transformations. The input of the Similarity Flooding algorithm being two labeled directed graphs, part of our contribution is the proposal and assessment of several ways of encoding a given metamodel into a labeled directed graph (Section 3). In order to ensure the reproducibility of our experiments by readers of this paper, we present in detail the implementation choices we have taken for the Similarity Flooding algorithm (Section 4). Finally, the output is filtered and transformed into a metamodel alignment (Section 5).

## 2 Overview of the approach

In this paper, we aim at generating a model transformation from an input model conforming to a metamodel  $MM_{source}$  to an output model conforming to a metamodel  $MM_{target}$ , aligning  $MM_{source}$  on  $MM_{target}$ . Such a transformation can be obtained using a schema matching approach.

*The Similarity Flooding algorithm* Several surveys [1,2] describe the existing approaches concerning the matching problem. Those approaches could roughly be classified in two categories: schema level approaches that produce alignments using the schemas (*e.g.* an XML Schema), and instance level approaches that produce alignments using instances of the schemas (*e.g.* an XML document conforming to a given Schema). Our objective being to generate model transformations based on the metamodels, we focused on the first category. Among the many such approaches found in the literature, we chose to use the Similarity Flooding algorithm [3], for two main reasons. First, it takes as input directed labeled graphs representing the schemas to match, and converting metamodels to directed labeled graphs is an easy operation. Second, its generic design allows easy tuning in order to produce better results using information specific to the input format. The similarity algorithm works on directed labeled graphs. Those graphs are used in a fix-point computation process whose result tells us what nodes in one graph are similar to nodes in the second graph. This process is



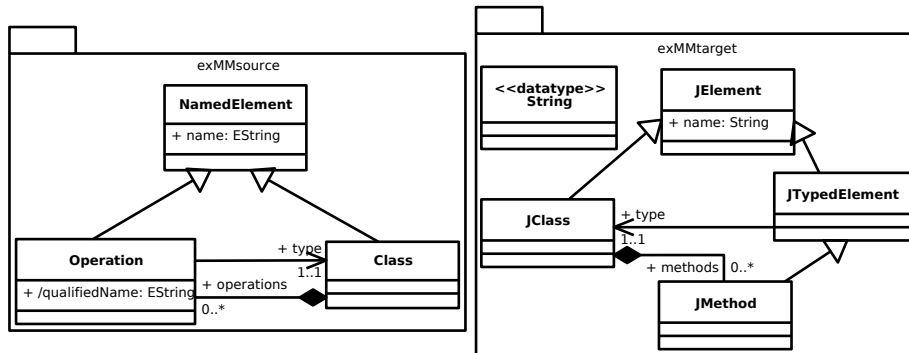
**Fig. 1.** Overview of the approach

based on the following intuition: if a node is similar to another node, the neighbours of this node may be similar to the neighbours of the other node. We have the same intuition concerning metamodel alignment.

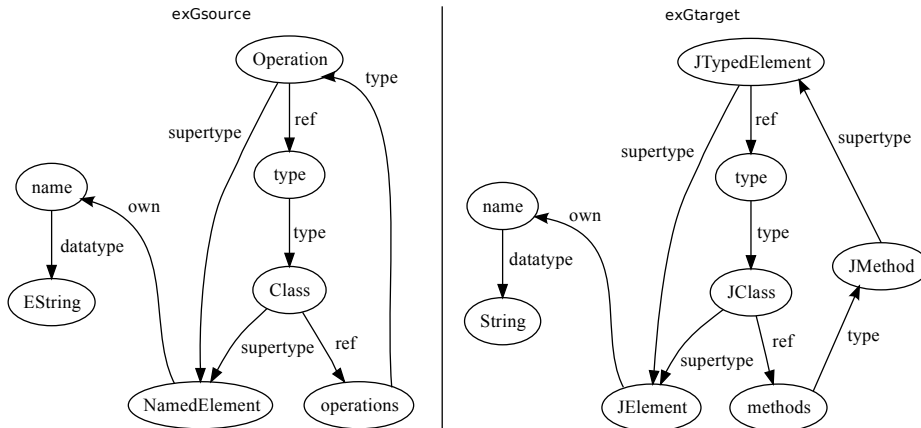
*Overall approach* The approach we propose to compute a metamodel alignment between two metamodels  $MM_{source}$  and  $MM_{target}$  is three-phased (see Fig. 1):

1. Transformation of  $MM_{source}$  and  $MM_{target}$  to the directed labeled graphs  $G_{source}$  and  $G_{target}$ . There are several strategies to encode a metamodel into such a graph, and thus several graphs can be generated in a deterministic way from the same metamodel, depending on the chosen strategy. This step will be detailed in Section 3.
2. Application of the Similarity Flooding algorithm. To keep this paper self-contained, we explain this algorithm in Section 4.
3. Generation of an alignment between  $MM_{source}$  and  $MM_{target}$ . This step is described in Section 5.

We illustrate these three steps using the two metamodels  $exMM_{source}$  and  $exMM_{target}$  shown in Fig. 2. These two metamodels are both used to express class models, but the names of the elements and their structure are slightly different (one is UML-like and the other is Java-like).



**Fig. 2.** Example  $exMM_{source}$  (left) and  $exMM_{target}$  (right) metamodels



**Fig. 3.** Graphs generated using the example metamodels and the Minimal configuration

### 3 From metamodels to directed labeled graphs

In this section, we explain how to transform a metamodel into a directed labeled graph that can be exploited by the Similarity Flooding algorithm. This transformation has a crucial importance: it is in charge of extracting from the metamodels the information that will be used to compute the mappings between the metamodel elements. It also structures this chosen information: the structure of the graph is important since the Similarity Flooding algorithm computes the alignment based on the structure of the input graphs. In the following, we refer to the choice of the elements in the metamodel and the way to structure them in a directed labeled graph as a configuration.

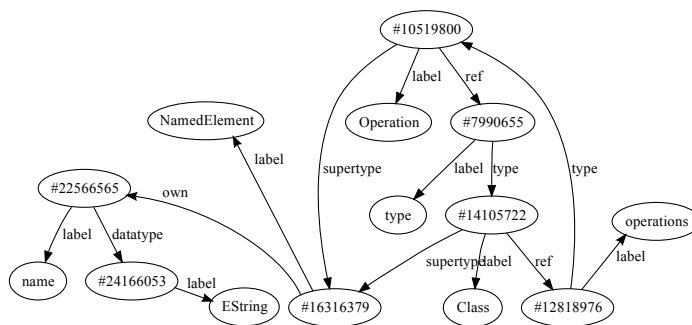
To determine the most adequate configuration, we designed six possible configurations that explore the effects of derived elements (that may introduce redundancy and increases the complexity of the Similarity Flooding algorithm), of abstract classes, and of transitive closure on relations based on the inheritance relation. We describe below those six configurations, and how they impact the resulting alignment in the case study presented in Section 6. Our approach uses **Ecore** [4] as the meta-metamodel, thus the configurations refer to **Ecore** elements (*e.g.*, **EClass**, **EReference**).

The first configuration is called **Minimal**: it is the simplest and most intuitive one. It uses a small subset of elements included in the metamodel: it does not use any meta-attribute of the meta-classes. One labeled node is created for each **EClass**, each non-derived **EAttribute**, each non-derived **EReference**, each **EDataType**, each **EEnum** and each **EEnumLiteral** contained in the metamodel. The labels of those nodes are the names of the corresponding elements. A directed arc labeled *supertype* is created between two nodes that represent an **EClass** whenever those two **EClasses** are in an inheritance relationship. An *own* arc is created when an **EClass** owns an **EAttribute**. A *ref* arc is created when an **EClass** owns an **EReference**. A *type* arc is created when an **EReference** is

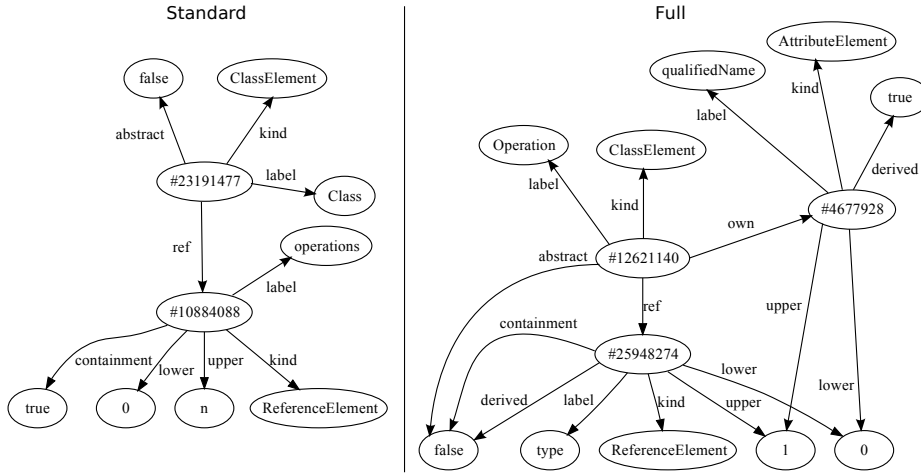
typed by an `EClass`. A *datatype* arc is created when an `EAttribute` is typed by an `EDataType` or an `EEnum`. A *literal* arc is created when an `EEnum` owns an `EEnumLiteral`. Fig. 3 shows the graphs  $exG_{source}$  and  $exG_{target}$ , corresponding respectively to  $exMM_{source}$  and  $exMM_{target}$  metamodels, generated with this configuration.

The second configuration, called **Basic**, differs from the Minimal configuration by the way it represents the names of the elements. With the Minimal configuration, an element  $E$  (for example an `EClass`) with name  $n$  is represented by a node  $N$  with label  $n$ , while with the Basic configuration,  $E$  is represented by a node  $N$  labeled by a unique identifier  $\#ID$ , and linked to a node labeled  $n$  with an arc *label*. Fig. 4 shows the graph  $exG_{source}$  generated with this configuration. Such a configuration allows us to directly have the information on the frequency a name is used (we just have to count the number of arcs *label* entering a given node). This frequency can then be exploited by the Similarity Flooding algorithm.

The third configuration, **Standard**, extends the Basic configuration to obtain similarities not only inspecting the names of the elements, but also their types, and main attributes (`abstract` for the `EClass`, `lowerBound` and `upperBound` from `EAttribute` and `EReference`, and `containment` from `EReference`). In that purpose, a node representing an element  $E$  is linked by an arc labeled *kind* to a node  $N$  representing the type of the element  $E$ .  $N$  is labeled according to the type of the element  $E$ , removing the prefix “ $E$ ” for all the elements of `Ecore`, and adding the suffix “*Element*” (e.g. `EClass` is turned into `ClassElement`). To deal with the main attributes, when an element  $E$  has for type a meta-class with an attribute  $A$ , then the node corresponding to  $E$  is linked with an arc labeled  $A$  to a node whose label is the value of the attribute  $A$  for  $E$ . For example, to represent an abstract `EClass` `myEClass`, the node representing `myEClass` is linked with an arc labeled `abstract` to a node labeled `true`. The left side of Fig. 5 shows an excerpt of the graph  $exG_{source}$  generated using the Standard configuration. The **Full** configuration extends the Standard configuration taking into account the derived `EAttributes` and `EReferences`. All the nodes representing an `EAttribute` or



**Fig. 4.** Graph  $exG_{source}$  corresponding to  $exMM_{source}$  using the Basic configuration (the name of the elements are in a dedicated node)



**Fig. 5.** Extracts of graph  $exG_{source}$  generated with the Standard and Full configurations

an **EReference** are added an arc labeled *derived* and leading to a node labeled *true* or *false* depending whether the element is derived. The right side of Fig. 5 shows an extract of the graph  $exG_{source}$  generated using this configuration.

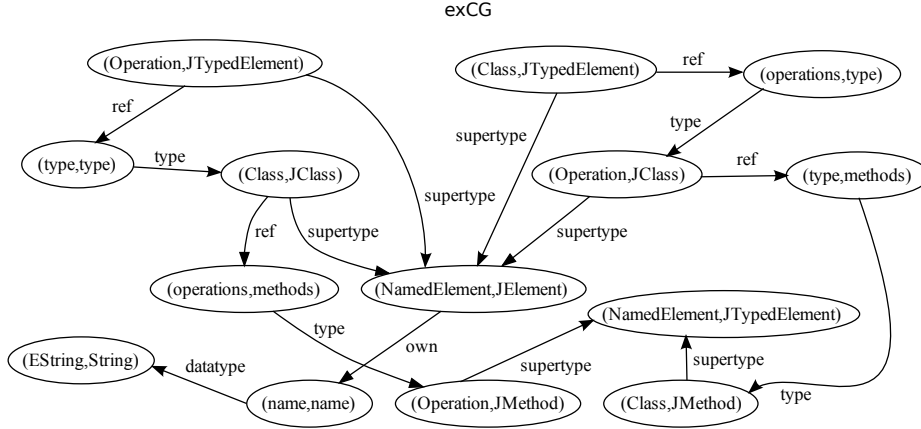
The **Flattened** configuration is based on the Standard one, but with flattened inheritance. The nodes representing abstract **EClasses** and the arcs labeled *supertype* are deleted. Instead, arcs labeled *own* (resp. *ref*) connect nodes representing an **EClass**  $Ecl$  to nodes representing the **EAttribute** (resp. **EReference**) defined by  $Ecl$  and all its superclasses. Moreover, when an **EReference**  $Eref$  is typed by an abstract **EClass**  $Ecl$ , a *type* arc is created from the node corresponding to  $Eref$  to each non-abstract sub-class of  $Ecl$ .

The last configuration is called **Saturated** and is still based on the Standard one. Here, the transitive relations are saturated, like in [5]. **EClass** nodes are now connected by a *supertype* arc to the nodes representing all the super-classes of this **EClass**. **EClass** nodes are also connected by *own* (resp. *ref*) arcs to the nodes representing the **EAttribute** (resp. **EReference**) introduced and inherited by the **EClass**. Finally, for a node representing an **EReference**, *type* arcs are created to the node representing the **EClass** that types the **EReference** as well as all the nodes representing the super-classes of the **EClass**.

## 4 The Similarity Flooding algorithm

This section describes the steps performed by the Similarity Flooding algorithm [3] to compute an alignment between two graphs  $G_{source}$  and  $G_{target}$ .

*Step 1* The first step of this algorithm is the computation of a compatibility graph (called pairwise connectivity graph in [3]) between  $G_{source}$  and  $G_{target}$ .



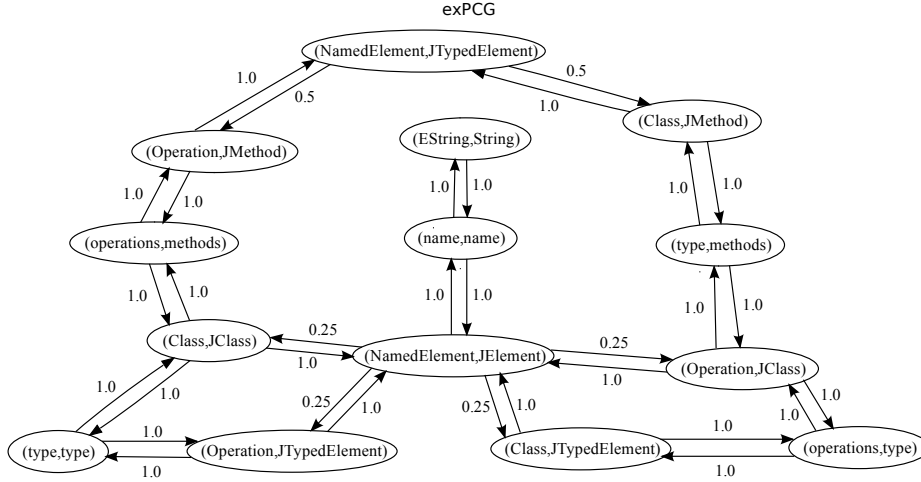
**Fig. 6.** The compatibility graph  $exCG$  between  $exG_{source}$  and  $exG_{target}$

A graph  $G_i$  is here a set of triples  $(x, r, y)$  with  $x$  and  $y$  two labeled nodes and  $r$  a labeled arc. The compatibility graph  $CG$  between two graphs  $G_{source}$  and  $G_{target}$  is built this way:  $((x_1, x_2), r, (y_1, y_2)) \in CG$  iff  $(x_1, r, y_1) \in G_{source}$  and  $(x_2, r, y_2) \in G_{target}$ . Therefore, a compatibility graph is a set of triples  $(s, r, t)$  such as  $s = (x_1, y_1)$  and  $t = (x_2, y_2)$  with  $x_1$  and  $x_2$  labeled nodes of  $G_{source}$ ,  $y_1$  and  $y_2$  labeled nodes of  $G_{target}$ , and  $r$  a labeled arc. Nodes  $s$  and  $t$  are called compatibility nodes. Fig. 6 shows the compatibility graph built using the two sample graphs of Fig. 3.

*Step 2* In this step, we compute a propagation graph using this compatibility graph. The labeled arcs of the compatibility graph created during the previous step are replaced by several weighted arcs. Different ways of computing the weights are compared in [3]. Here, we have chosen to use one that has proved to give good results. First, for each arc  $(s, r, t)$  of the compatibility graph, a reverse arc  $(t, r, s)$  is also created. Then, for each compatibility node  $n$  we compute  $O_l^n$  the number of arcs that start from  $n$  and are labeled by  $l$ . For instance, in Fig. 6,  $O_{supertype}^{(NamedElement, JElement)} = 4$  and  $O_{own}^{(NamedElement, JElement)} = 1$ . For each triple  $(s, r, t)$  of the compatibility graph, a  $(s, v, t)$  triple is created in the propagation graph with  $v = 1/O_r^s$ . In Fig. 6 we can see that  $(n_1, supertype, n_2)$  is replaced by  $(n_1, 0.25, n_2)$  because  $O_{supertype}^{n_1} = 4$ , with  $n_1 = (NamedElement, JElement)$  and  $n_2 = (Class, JClass)$ . Similarly,  $(n_1, own, n_3)$  is replaced by  $(n_1, 1, n_3)$  because  $O_{own}^{n_1} = 1$ , with  $n_3 = (name, name)$ . Fig. 7 shows the propagation graph computed using the compatibility graph of Fig. 6.

*Step 3* We dispose now of a propagation graph that indicates how the similarity values of the nodes will propagate through the graph. Before starting the fix-point computation process, it is necessary to assign an initial similarity value to each node of the propagation graph. We chose to assign the initial value  $s_n^0$  of a compatibility node  $n = (x, y)$  with the following formula, similar to the one used





**Fig. 7.** The propagation graph *exPCG*

in [3]: if  $x$  or  $y$  is an identifier  $s_n^0 = 0$ , else  $s_n^0 = 1 - lev(x, y) / \max(len(x), len(y))$ , with  $len(x)$  the length of label  $x$  and  $lev(x, y)$  the Levenshtein distance [6] between labels  $x$  and  $y$ . The Levenshtein distance between two labels is given by the minimum number of operations needed to transform one label into the other, where an operation is an insertion, deletion, or substitution of a single character. Table 1 shows the initial similarity value computed for each node of the propagation graph shown in Fig. 7. An iterative fix-point computation process will propagate these initial similarity values through the propagation graph. There are again many ways to propagate those similarity values. We chose the fix-point formula that has proved to give the best result and has the best convergence properties in [3]. This formula has the following definition. Let  $n$  be a compatibility node and  $I_n$  be the set of compatibility nodes that are connected to  $n$  by a weighted arc that ends at  $n$ , and let  $w(s, t)$  be the weight of the arc between  $s$  and  $t$ . The similarity value  $s_n^{i+1}$  of node  $n$  at step  $i + 1$  is computed according to the following formula:  $s_n^{i+1} = s_n^i + s_n^0 + \sum_{m \in I_n} w(m, n) \times (s_m^0 + s_m^i)$ .

At the end of a step of this iterative process, the similarity value of each compatibility node is normalized; i.e. divided by the greatest similarity value computed in the step. For instance, in the sample propagation graph shown in Fig. 7,

$$\begin{aligned}
 s_{n_4}^1 &= s_{n_4}^0 + s_{n_4}^0 + w(n_5, n_4) \times (s_{n_5}^0 + s_{n_5}^0) + w(n_6, n_4) \times (s_{n_6}^0 + s_{n_6}^0) \\
 &\simeq 0.10 + 0.10 + 1 \times (0.11 + 0.11) + 1 \times (0.08 + 0.08) \simeq 0.58 \\
 \text{with } n_4 &= (\text{operations}, \text{type}), n_5 = (\text{Operation}, \text{JClass}), \\
 n_6 &= (\text{Class}, \text{JTypedElement}).
 \end{aligned}$$

For this sample propagation graph, the best computed similarity value during the first step is  $s_{(\text{NamedElement}, \text{JElement})}^1 \simeq 5.67$  hence at the end of the first

Compatibility node	Initial similarity value	Final similarity value
(NamedElement, JElement)	0.5833334	1.0
(name, name)	1.0	0.7771561
(EString, String)	0.85714287	0.5186626
(NamedElement, JTypedElement)	0.6923077	0.2985012
(Operation, JTypedElement)	0.23076922	0.44552472
(Class, JTypedElement)	0.07692307	0.13148126
(type, type)	1.0	0.63979846
(operations, type)	0.100000024	0.11389989
(Operation, JClass)	0.111111104	0.1868067
(Class, JClass)	0.8333333	0.7449572
(type, methods)	0.14285713	0.11238139
(operations, methods)	0.39999998	0.52174574
(Operation, JMethod)	0.3333333	0.3444618
(Class, JMethod)	0.0	0.14695258

**Table 1.** Initial and final similarity value of each compatibility node

step,  $s_{(operations, type)}^1 \simeq 0, 10$  and  $s_{(NamedElement, JElement)}^1 = 1$ . Let  $(S^i, S^{i+1})$  be a vector of the similarity values computed at two successive steps  $i$  and  $i + 1$ . This iterative fix-point computation process ends whenever the euclidean norm of this vector becomes less than  $\varepsilon$  for some  $i > 0$ . If the computation does not converge (it means that the euclidean norm of  $(S^i, S^{i+1})$  never becomes less than  $\varepsilon$ ), the process terminates after a maximal number of steps. More information on the complexity and the convergence properties of this algorithm are given in [3]. Table 1 shows the final similarity values of the compatibility nodes of the sample propagation graph shown in Fig. 7, with  $\varepsilon = 0.05$ . The set of tuples  $(n, v)$  with  $n$  a compatibility node and  $v$  a similarity value, corresponding to the columns *Compatibility node* and *Final similarity value* of Table 1, is called the multimapping between  $G_{source}$  and  $G_{target}$ .

*Step 4* This last step filters the multimapping to produce a final alignment. We call  $R$  the multimapping and  $A$  the final alignment, with  $A \subseteq R$ . The method we use to filter the produced multimapping is called *SelectThreshold* in [3]. In the multimapping, a node of the source graph can be associated to several nodes of the target graph. Fig. 8 shows an extract of the multimapping produced with the sample graphs. In this extract, we can clearly see that the *type* node of the source graph can be associated to the *type* node or the *methods* node of the target graph. The *SelectThreshold* filter works as follows. For each couple  $(n, v)$  of the multimapping, with  $n = (x, y)$ , the best match for  $x$  (resp  $y$ ) is looked for in the multimapping. The best match for  $x$  (resp  $y$ ) is a couple  $(m, v) \in R$  such as  $m = (x, z)$  (resp  $m = (w, y)$ ) with  $v$  the best similarity value for a node like  $m$ . We call  $s_x^{max}$  (resp  $s_y^{max}$ ) the similarity value for the best match for  $x$  (resp  $y$ ). For the previous couple  $(n, v)$ , we compute  $s_x^n = v/s_x^{max}$  and  $s_y^n = v/s_y^{max}$ . The couple  $(n, v) \in A$  iff  $s_x^n \geq threshold$  and  $s_y^n \geq threshold$ , with  $threshold \in$

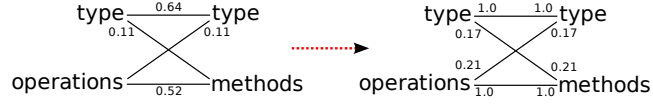


Fig. 8. Extract of the multimapping (left) and of the mutual multimapping (right)

$[0, 1]$ . In the example of Fig. 8, we have  $s_{operations}^{max} \simeq 0.52$  and  $s_{type}^{max} \simeq 0.64$ . Therefore,  $s_{operations}^{(operations, type)} \simeq \frac{0.11}{0.52} \simeq 0.21$ . Similarly,  $s_{type}^{(operations, type)} \simeq \frac{0.11}{0.64} \simeq 0.17$ . With  $threshold = 0.2$ , the couple  $(operations, type)$  is discarded while with  $threshold = 0.15$ , it is included in the alignment. The choice of  $threshold$  determines the alignment: a high value of  $threshold$  will lead to  $[0 - 1] \times [0 - 1]$  cardinality mappings (one element from the source metamodel corresponds at most to one element of the target metamodel). On the other hand, a low value of  $threshold$  will create  $[0 - n] \times [0 - n]$  cardinality mappings (a set of elements of the source metamodel corresponds to a set of elements of the target metamodel). Fig. 9 shows the alignment computed on the example graphs with  $threshold = 0.95$ . Each mapping of this alignment is correct, except the mapping between *Operation* and *JTypedElement*.

## 5 Computation of the metamodel alignment

In our approach, we consider a model transformation from a source metamodel  $MM_{source}$  to a target metamodel  $MM_{target}$  as a procedure that produces objects conforming to concrete classes of  $MM_{target}$  by the analysis of objects conforming to concrete classes of  $MM_{source}$ . Data used when analysing an object of the source model correspond to the non-derived attributes and references, because Ecore does not allow to express how an operation is computed or a derived feature. In the MDE technological space, many transformation languages exist. We did not want our approach to be tied to a particular language. To avoid that, we chose to build a language-independent alignment metamodel, inspired by the

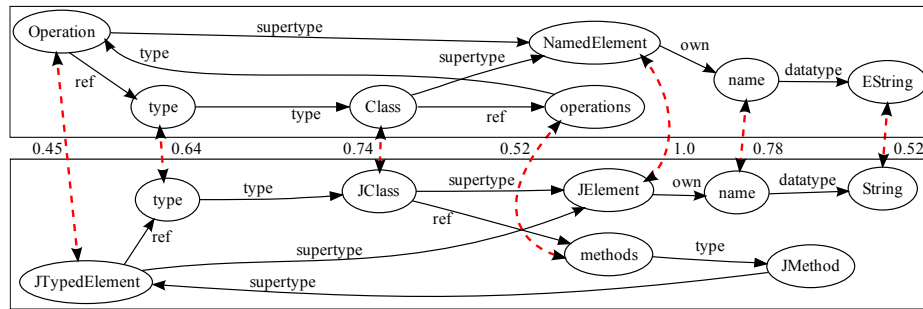


Fig. 9. Computed alignment between  $exG_{source}$  and  $exG_{target}$

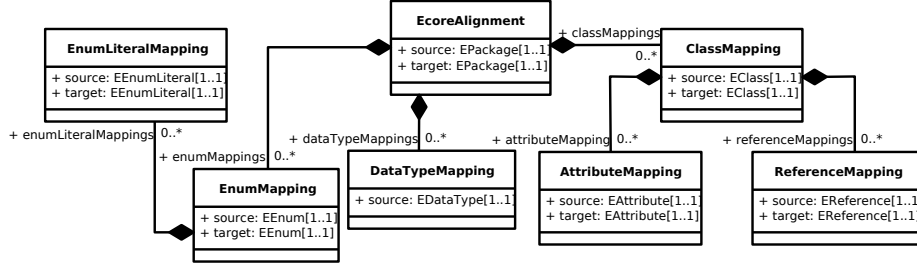


Fig. 10. Ecore alignment metamodel

schema-matching metamodel given in [7]. Models conforming to this metamodel show how a metamodel is aligned to another metamodel. Such a model is built by analysing and reorganizing the alignment produced by the Similarity Flooding algorithm. A transformation generator then reads an alignment model and generates the code of the transformation in a target language. Such an architecture allows to generate transformations in various languages. However, such code generators are out of the scope of this paper. The alignment metamodel is shown in Fig. 10. An Ecore alignment is composed of several class mappings, data type mappings and enumeration mappings. A class mapping is composed of several attribute and reference mappings. An enumeration mapping is composed of several enumeration literal mappings. Because we did not want to tackle complex element transformations (such as Class to Attribute or Attribute to Class mappings) in this paper, our metamodel only allows to express mappings between elements of the same kind. However, such cross-kind mappings can be found in the alignment produced by Similarity Flooding. According to our definition of a model transformation, we add the following constraints on our alignment metamodel: two mapped `EClass` are not abstract and two mapped `EReference` or `EAttribute` are not derived. The steps required to build such an Ecore alignment from a Similarity Flooding (SF) alignment are:

1. Search in the SF alignment all the mappings between two concrete `EClass` to build the `ClassMapping` elements. Then for each such mapping:
  - (a) When an `EAttribute` from the source `EClass` and its super-classes is mapped to an `EAttribute` contained by the target `EClass` or its super-classes, add an `AttributeMapping` element in the `ClassMapping`;
  - (b) When an `EReference` from the source `EClass` and its super-classes is mapped to an `EReference` contained by the target `EClass` or its super-classes, add a `ReferenceMapping` element in the `ClassMapping`;
2. Search in the SF alignment all the mappings from a source `EEnum` to a target `EEnum` to build the `EnumMapping` elements. For each literal of a mapping source `EEnum` which matches on a literal of the target `EEnum`, add an `EnumLiteralMapping` element.
3. Search in the SF alignment all the mappings from a source `EDataType` to a target `EDataType` to build the `DataTypeMapping` elements.

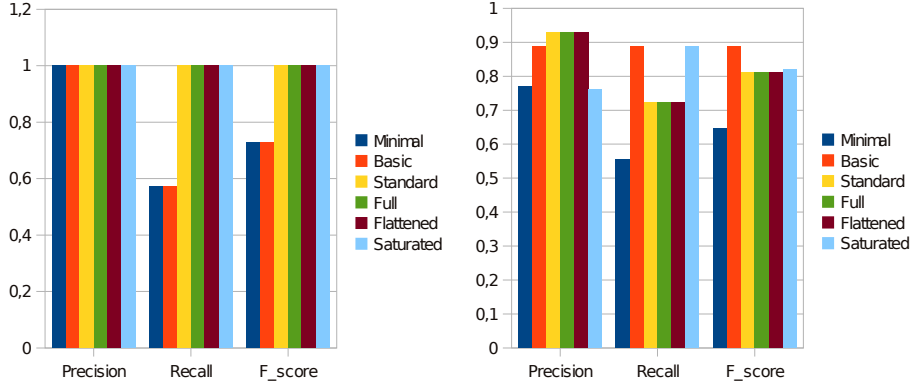


Fig. 11. Results of  $exMM_{source} \leftrightarrow exMM_{target}$  (left) and  $Ecore \leftrightarrow Minjava$  (right)

## 6 Case study

This case study compares the results obtained on various metamodels with the different configurations proposed in Section 3. The parameters of the similarity flooding algorithm used in this case study are:  $\varepsilon = 0.05$  and  $threshold = 0.95$ . Since many of the real world metamodels are designed to express class models, we chose to use those metamodels as a benchmark. The chosen alignment scenarios are:  $exMM_{source} \leftrightarrow exMM_{target}$ ,  $Ecore \leftrightarrow Minjava$ ,  $Ecore \leftrightarrow Kermeta$  and  $Ecore \leftrightarrow UML$ . Ecore [4] is the meta-metamodel and hence a metamodel. Minjava [8] is a simple metamodel for the Java language, that allows to represent the structure of a Java program. Since Minjava imports the models from Java bytecode, it does not provide a way to describe how the methods of the different classes of a program are structured and thus is very similar to Ecore. Kermeta [9] is an extension of Ecore that allows to give a behavioral description of the operations and derived properties of a class model in addition to the structural description. UML [10] is a multi-purpose software modeling language that deals with class models (including behavioral description), use cases or statecharts. Since the elements from the metamodel used in the matching process are not the same for the different configurations of Section 3, we needed a common denominator between those configurations to compare them. This common denominator is the Ecore alignment model produced by the Ecore alignment encoder described in Section 5. This model only shows mappings between non abstract classes, non derived attributes and references, datatypes and enumerations. Those elements are included in all the different configurations. We compare the Ecore alignment model produced by the alignment encoder to an ideal alignment model designed by an expert (standard transformations such as Ecore to Kermeta have been used to build those ideal models). We evaluate our results with common metrics from the information retrieval field [11]. For each alignment, we compute  $precision = \frac{\#correct\_found\_mappings}{\#total\_found\_mappings}$ ,  $recall = \frac{\#correct\_found\_mappings}{\#total\_correct\_mappings}$

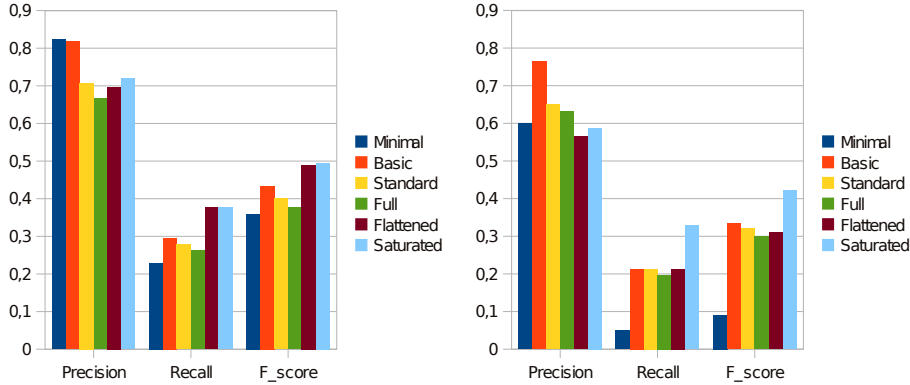


Fig. 12. Results of  $Ecore \leftrightarrow Kermeta$  (left) and  $Ecore \leftrightarrow UML$  (right)

and  $f\_score = \frac{2 \times recall \times precision}{recall + precision}$ . We have  $precision \in [0, 1]$ . The higher is the  $precision$  value, the smaller is the set of wrong mappings.  $recall \in [0, 1]$ . The higher is the  $recall$  value, the smaller is the set of the mappings that have not been found.  $f\_score$  is the harmonic mean of  $precision$  and  $recall$ .  $f\_score \in [0, 1]$ . It can be considered as a global measure of the alignment quality. A high value of  $f\_score$  is obtained when the produced alignment has a good quality.

The results of our experiments are shown in Fig. 11 and 12. The results show that the best configuration is generally Saturated. On the other hand, the Minimal configuration gives very bad results. Full and Standard configurations, despite using more information from the metamodels, seem to produce slightly poorer results than the Basic configuration. The metamodels we used to match on Ecore vary in size. Minjava has more or less the same size as Ecore, Kermeta is a little bigger, and UML is very large. We can clearly see from the results that the alignment quality decreases when the size difference between the two matched metamodels is increasing. This is partly caused by the *SelectThreshold* filter that filters the alignment produced by Similarity Flooding. Results on  $Ecore \leftrightarrow Minjava$  and  $Ecore \leftrightarrow Kermeta$  show that good quality alignments can be produced by our approach.

## 7 Related work and conclusion

*Related Work* Automating the discovery of mappings between schemas, ontologies, documents or (meta)models has been thoroughly investigated [1,2]. In the context of Model-Driven Engineering, several approaches for semi-automatic generation of transformations based on mapping have recently been proposed. In [12], model transformations are generated based on ontological information. The two metamodels are supposed to have a semantics given using a mapping onto

a known ontology. Reasoning on the ontology then allows to generate a model transformation, adapting a bootstrap transformation that is whether automatically generated or existing. The transformation is written in the paper in QVT Relational. [13] studies the semi-automatic derivation of graph transformation rules based on manual set-up of prototype mappings between models. In another Model Transformation By-Example approach [14], users define a mapping directly on models in concrete syntax (M1 level), then, using a predefined known mapping between concrete and abstract syntaxes, ATL rules can be generated. Compared to these two approaches that are based on a given mapping (onto a pivot ontology or between examples), we focus on automatically determining the mapping on which is based the transformation generation.

In [7], the matching operation focuses on classes, datatypes and enumerations. The proposed class matching algorithm iterates until a fix-point is reached on evaluating if two leaf classes match. Characteristics including attributes, references and relationships (based on closure and composition of inheritance, association, agregation, type-of, etc.) are then used: if the number of matching characteristics is greater than a threshold average, the two classes are considered as similar. The exact computation of characteristic similarity is not much detailed in [7], but it is said that they used name, type and graph matching similarity. For datatypes, the possibility of representing a value of a type (*e.g.* int) by another type (*e.g.* float) is used. For enumeration, the numbers of identical literals are compared. Experiments on partial UML and Java metamodels are shown in [15]. In [16], an architecture is proposed for semi-automating the generation of model transformations. A weaving metamodel describes the kinds of links generated by the matching engine and current transformation patterns. Matching is based on element to element and on structural similarities. Structural similarities computation is inspired by the Similarity Flooding approach. A configuration metamodel is introduced to tune the matching phase. Several case studies are conducted in [16]. Here we go deeper into the encoding of metamodels for applying Similarity Flooding. We detail and compare several configurations on known, available, full metamodels. These configurations are helpful for future case tool constructions. The whole process is detailed enough in order to be reproducible and compared to future metamodel matching approaches.

*Conclusions and perspectives* In this paper, we have presented an approach that produces an alignment between two metamodels in an automated way. This approach is based on the application on metamodels of a well-known schema matching algorithm called Similarity Flooding. Our main contribution is the study on the various ways of encoding a given metamodel into a directed labeled graph that can then be exploited by the Similarity Flooding algorithm. We compared six strategies in a case study in terms of precision and recall of the computed mappings between metamodels. This case study points out that the more intuitive encoding strategy (Minimal) leads to poor results, and that the Saturated strategy is the best-suited in the majority of the cases. The generation of an Ecore alignment model (conforming to an Ecore alignment metamodel) from which model transformation code can easily be derived is our second main

contribution. Future work will consist in designing code generators both for imperative and declarative languages. We also currently follow another perspective: we study how the Similarity Flooding algorithm can be tuned to our particular domain (metamodel matching). We are working on the way the initial similarity value of a compatibility node is computed, and on the way we compute the weights of the arcs in the propagation graph.

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