

Qualitative Spatial Reasoning about Line Segments

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Abstract. Representing and reasoning about orientation information is an important aspect of qualitative spatial reasoning. We present a novel approach for dealing with intrinsic orientation information by specifying qualitative relations between oriented line segments, the simplest possible spatial entities being extended and having an intrinsic direction. We identify a set of 24 atomic relations which form a relation algebra and for which we compute relational compositions based on their algebraic semantics. Reasoning over the full algebra turns out to be NP-hard. Potential applications of the calculus are motivated with a small example which shows the reasoning capabilities of the dipole calculus using constraint-based reasoning methods.

1 Introduction

Qualitative representation of space abstracts from the physical world and enables computers to make predictions about spatial relations, even when precise quantitative information is not available [2]. Different aspects of space can be represented in a qualitative way. The most important of these are topological information and orientation information about physical objects which are usually spatially extended. While it is common for representing topological information to use extended spatial regions as the basic entities, most approaches to qualitatively representing and reasoning about orientation information deal with points as the basic entities. Those orientation approaches that use extended spatial regions as the basic entities mostly approximate regions by using, for instance, minimal bounded rectangles whose sides are parallel to the axes of the global reference frame. This, however, does not account for representing intrinsic orientation information.

In this paper we develop one of the simplest possible calculi for representing intrinsic orientation information, namely, by using oriented line segments represented by their start and end points as the basic entities.

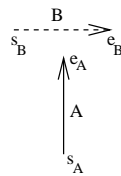


Figure 1. Orientation relation between two dipoles

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We propose calculi on different levels of granularity which all form relation algebras and as such allow for using standard constraint based reasoning mechanisms originally developed for temporal reasoning. Even on the coarsest level of granularity our calculi enable to represent polygonal lines which are particularly interesting for applications such as cognitive robotics [10] or spatial information systems [6].

2 The Basic Representation of the Dipole Relations

The basic entities we are using are dipoles, i.e., oriented line segments formed by a pair of two points, a start point and an end point. Dipoles are denoted by A, B, C, \dots , the start point by s_A , the end point by e_A , respectively (see Figure 1). These dipoles are used for representing spatial objects with an intrinsic orientation. Given a set of dipoles it is possible to specify many different relations of different arity, e.g., depending on the length of dipoles, the angle between different dipoles, or the dimension and nature of the underlying space. The goal of identifying different relations is to obtain a set of jointly exhaustive and pairwise disjoint *atomic* relations, i.e., between any two dipoles exactly one relation holds. If these relations form a *relation algebra* it is possible to apply standard constraint-based reasoning mechanisms which were originally developed for temporal reasoning and which have also proved valuable for spatial reasoning. In order to enable efficient reasoning, it should be tried to keep the number of different base relations relatively small.

For this reason, we will restrict for now to using two-dimensional continuous space, in particular \mathbb{R}^2 , and distinguish the location and orientation of the different dipoles only according to whether a point lies to the left, to the right, or on the straight line through the referring dipole. Then s_B can either lie to the left of A (see figure 1), on the straight line through A or to the right of A , expressed as $A \mid s_B$, $A \circ s_B$ or $A r s_B$, respectively. Using these three relations between a dipole and a point it is possible to specify the relations between two dipoles with the following four relationships:

$$A R s_B \wedge A R e_B \wedge B R s_A \wedge B R e_A,$$

where R is one of $\{r, l, o\}$. Since this still leads to a very large number of different atomic relations, we require in the first version of our algebra all points to be in *general position*, i.e., no more than two points are on a line (the extended version of the algebra is described in section 5). This gives us the following 14 relations that hold if the four points s_B, e_B, s_A, e_A are distinct:

$$A r r r r B := A r s_B \wedge A r e_B \wedge B r s_A \wedge B r e_A$$

$$A r r r l B := A r s_B \wedge A r e_B \wedge B r s_A \wedge B l e_A$$

$$A r r l r B := A r s_B \wedge A r e_B \wedge B l s_A \wedge B r e_A$$

$$A r r l l B := A r s_B \wedge A r e_B \wedge B l s_A \wedge B l e_A$$

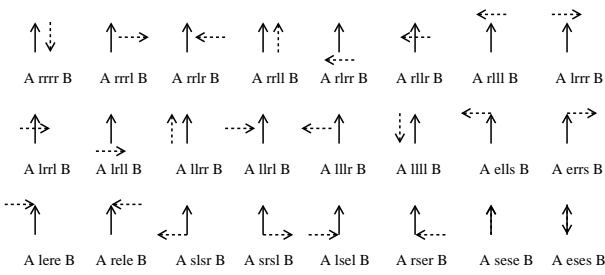


Figure 2. The 24 atomic relations of the dipole calculus

$$\begin{aligned}
A \text{ rrrr } B &:= A \text{ r } \mathbf{s}_B \wedge A \text{ l } \mathbf{e}_B \wedge B \text{ r } \mathbf{s}_A \wedge B \text{ r } \mathbf{e}_A \\
A \text{ rllr } B &:= A \text{ r } \mathbf{s}_B \wedge A \text{ l } \mathbf{e}_B \wedge B \text{ l } \mathbf{s}_A \wedge B \text{ r } \mathbf{e}_A \\
A \text{ rlll } B &:= A \text{ r } \mathbf{s}_B \wedge A \text{ l } \mathbf{e}_B \wedge B \text{ l } \mathbf{s}_A \wedge B \text{ l } \mathbf{e}_A \\
A \text{ lrrr } B &:= A \text{ l } \mathbf{s}_B \wedge A \text{ r } \mathbf{e}_B \wedge B \text{ r } \mathbf{s}_A \wedge B \text{ r } \mathbf{e}_A \\
A \text{ lrrl } B &:= A \text{ l } \mathbf{s}_B \wedge A \text{ r } \mathbf{e}_B \wedge B \text{ r } \mathbf{s}_A \wedge B \text{ l } \mathbf{e}_A \\
A \text{ lrll } B &:= A \text{ l } \mathbf{s}_B \wedge A \text{ r } \mathbf{e}_B \wedge B \text{ l } \mathbf{s}_A \wedge B \text{ l } \mathbf{e}_A \\
A \text{ llrr } B &:= A \text{ l } \mathbf{s}_B \wedge A \text{ l } \mathbf{e}_B \wedge B \text{ r } \mathbf{s}_A \wedge B \text{ r } \mathbf{e}_A \\
A \text{ llrl } B &:= A \text{ l } \mathbf{s}_B \wedge A \text{ l } \mathbf{e}_B \wedge B \text{ r } \mathbf{s}_A \wedge B \text{ l } \mathbf{e}_A \\
A \text{ llrr } B &:= A \text{ l } \mathbf{s}_B \wedge A \text{ l } \mathbf{e}_B \wedge B \text{ l } \mathbf{s}_A \wedge B \text{ r } \mathbf{e}_A \\
A \text{ llrr } B &:= A \text{ l } \mathbf{s}_B \wedge A \text{ l } \mathbf{e}_B \wedge B \text{ l } \mathbf{s}_A \wedge B \text{ l } \mathbf{e}_A
\end{aligned}$$

The cases $A \text{ r } \mathbf{s}_B \wedge A \text{ l } \mathbf{e}_B \wedge B \text{ r } \mathbf{s}_A \wedge B \text{ l } \mathbf{e}_A$ and $A \text{ l } \mathbf{s}_B \wedge A \text{ r } \mathbf{e}_B \wedge B \text{ l } \mathbf{s}_A \wedge B \text{ r } \mathbf{e}_A$ cannot be realized on the plane. These 14 relations are similar to the relations between line segments derived by Schlieder [13]. However, in order to obtain a relation algebra, we also have to consider those relations where two dipoles share common points. Then \mathbf{s}_B can be equivalent to the start point of A or to the end point of A . This is denoted as $A \text{ s } \mathbf{s}_B$ or $A \text{ e } \mathbf{s}_B$, respectively. Using these additional dipole-point relations, we obtain the following ten additional dipole-dipole relations: {ells, errs, lere, rele, slsr, srsl, lsrl, rser, sese, eses}. Altogether we obtain 24 different atomic relations. These relations are jointly exhaustive and pairwise disjoint provided that all points are in general position. The relation sese is the identity relation. We use \mathcal{D}_{24} to refer to the set of 24 atomic relations, and $\mathcal{DR}\mathcal{A}_{24}$ to refer to the powerset of \mathcal{D}_{24} which contains all 2^{24} possible unions of the atomic relations.

The relations which are introduced above in an informal way can be defined in an algebraic way. Every dipole D on the plane \mathbb{R}^2 is an ordered pair of two points \mathbf{s}_D and \mathbf{e}_D , each of them is represented by its Cartesian coordinates x and y , with $x, y \in \mathbb{R}$ and $\mathbf{s}_D \neq \mathbf{e}_D$.

$$D = (\mathbf{s}_D, \mathbf{e}_D), \quad \mathbf{s}_D = ((\mathbf{s}_D)_x, (\mathbf{s}_D)_y)$$

The basic relations are then described as polynomial equations with the coordinates as variables. The set of solutions for a system of equations describes all the possible coordinates for these points. As an example, we will have a more detailed look at the relation $A \text{ rrrr } B$. We need to find an equation, which is solvable iff a point lies to the right of a given line. Then, we can use this equation to express the premises of the relation: $A \text{ r } \mathbf{s}_B, A \text{ r } \mathbf{e}_B, B \text{ r } \mathbf{s}_A, B \text{ r } \mathbf{e}_A$. The equation for “right of” is constructed as follows ($A \text{ r } \mathbf{s}_B$ serves as example):

With $\vec{A} = \begin{pmatrix} (\mathbf{e}_A)_x - (\mathbf{s}_A)_x \\ (\mathbf{e}_A)_y - (\mathbf{s}_A)_y \end{pmatrix}$, hence $\vec{A}^t = \begin{pmatrix} (\mathbf{e}_A)_y - (\mathbf{s}_A)_y \\ (\mathbf{s}_A)_x - (\mathbf{e}_A)_x \end{pmatrix}$ and $\vec{P} = \begin{pmatrix} (\mathbf{s}_B)_x - (\mathbf{s}_A)_x \\ (\mathbf{s}_B)_y - (\mathbf{e}_A)_y \end{pmatrix}$. Whenever \mathbf{s}_B lies on the right of the line $\overline{\mathbf{s}_A \mathbf{e}_A}$, the inequation

$${}^t \vec{A} \cdot \vec{P} > 0 \quad (1)$$

holds. To change this into an equation, we introduce a new variable v .

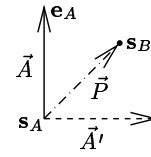


Figure 3. Constructing equations with the coordinates as variables

As v^2 can only take nonnegative values, the resulting equation

$${}^t \vec{A}' \cdot \vec{P} - v^2 = 0 \quad \text{with } v \in \mathbb{R} \setminus \{0\} \quad (2)$$

will have a solution iff the point \mathbf{s}_B lies to the right of the line $\overline{\mathbf{s}_A \mathbf{e}_A}$. The equation 1 is modified in a similar way for the premise “left of” (l):

$$A \text{ l } \mathbf{s}_B : {}^t \vec{A}' \cdot \vec{P} + v^2 = 0 \quad \text{with } v \in \mathbb{R} \setminus \{0\} \quad (3)$$

Note that the equations will only have a solution, when $\mathbf{s}_A \neq \mathbf{e}_A$. Constructing the dipole-point relations s and e is done by using the same variables for the identical points.

For the following substitutions $x_1 = (\mathbf{s}_A)_x, x_2 = (\mathbf{s}_A)_y, x_3 = (\mathbf{e}_A)_x, x_4 = (\mathbf{e}_A)_y, x_5 = (\mathbf{s}_B)_x, x_6 = (\mathbf{s}_B)_y, x_7 = (\mathbf{e}_B)_x, x_8 = (\mathbf{e}_B)_y$, and new introduced variables v_1, \dots, v_4 , the complete set of equations describing relation $A \text{ rrrr } B$ reads as:

$$\begin{aligned}
-x_1 x_4 + x_1 x_6 + x_2 x_3 - x_2 x_5 - x_3 x_6 + x_4 x_5 - v_1^2 &= 0 \\
-x_1 x_4 + x_1 x_8 + x_2 x_3 - x_2 x_7 - x_3 x_8 + x_4 x_7 - v_2^2 &= 0 \\
-x_1 x_6 + x_1 x_8 + x_2 x_5 - x_2 x_7 - x_5 x_8 + x_6 x_7 - v_3^2 &= 0 \\
-x_3 x_6 + x_3 x_8 + x_4 x_5 - x_4 x_7 - x_5 x_8 + x_6 x_7 - v_4^2 &= 0
\end{aligned}$$

with $x_1, \dots, x_8 \in \mathbb{R}, v_1, \dots, v_4 \in \mathbb{R} \setminus \{0\}$. The other relations are constructed in an analogous way.

3 Constraint Reasoning with the Dipole Calculus

For reasoning about the dipole relations we apply constraint-based reasoning techniques which were originally introduced for temporal reasoning [1] and which also proved valuable for spatial reasoning [12]. In order to apply these techniques to a set of relations, these relations must form a relation algebra [8], i.e., they must be closed under composition (\circ), intersection (\cap), complement ($\bar{}$), and converse (\smile) and there must be an empty relation, a universal relation, and an identity relation. While the converse (see Table 1), the complement, and the intersection of relations can be computed from the set-theoretic definitions of the relations, the composition of relations must be computed based on the semantics of the relations. The compositions are usually computed only for the atomic relations which are then stored in a composition table. The composition of compound relations can be obtained as the union of the compositions of the corresponding atomic relations.

We computed the compositions of the atomic relations using the algebraic semantics of the relations. For this we apply the method of “Gröbner Bases” using a geometric theorem prover [3]. A possible composition table entry $R_x \circ R_y \mapsto R_z$ is represented (for every combination of R_x, R_y , and R_z) by a set of equations. This set results from the union of three sets, one for each relation as shown in the previous section. $R_x(A, B) \wedge R_y(B, C) \wedge R_z(A, C)$ is a contradiction if and only if the set of equations has no solution. This can happen because of an equation with no solution (e.g. $x_i^2 = -1$) or a violation of the condition $v_1, \dots, v_n \in \mathbb{R} \setminus \{0\}$ (e.g. $v_i^2 + v_j^2 = 0$). By computing the Gröbner Base, equations are generated which do not change the systems solution. These generated equations allow the prover to detect, if there cannot be a solution. For all combinations

of R_x, R_y , and R_z where no contradiction was detected, we have to construct a possible configuration of points in the plane. Instead of generating this configuration from the equations (which can be quite complicated), we simply search for a valid configuration of points on a grid.

R	rrrr	rrrl	rrlr	rrll	rlrr	rllr	rlll	lrrr
R^{\sim}	rrrr	rlrr	lrrr	llrr	rrrl	lrrl	llrl	rrlr
R'	llll	lllr	llrl	llrr	lrlr	lrrl	lrrr	rlll
R	lrrl	lrll	llrr	llrl	lllr	llll	ells	errs
R^{\sim}	rlrr	lllr	rrll	rlll	lrll	llll	lsel	rser
R'	rlrr	rlrr	rrll	rrlr	rrrl	rrrr	errs	ells
R	lere	rele	slsr	srsr	lsel	rser	sese	eses
R^{\sim}	rele	lere	srsr	srsr	ells	errs	sese	eses
R'	rele	lere	srsr	srsr	rser	lsel	sese	eses

Table 1. Converse and reflection table of the dipole calculus

The composition table for the atomic relations is given in Table 2⁴. We use * to mark places which can be filled with r or l. In order to reduce the size of the table, trivial cases (sese,eses) for the columns are omitted. Symmetric cases can be derived using the converse operation and a reflection operation (reflection on an axis, denoted R' , see also Table 1). The missing entries can be calculated using the following equation:

$$R_1 \circ R_2 = (R_2^{\sim} \circ R_1^{\sim})^{\sim} = (R_1' \circ R_2')' \quad (4)$$

Dipole constraints are written as xRy where x, y are variables for dipoles and R is a \mathcal{DRA}_{24} relation. Given a set Θ of dipole constraints, an important reasoning problem is deciding whether Θ is *consistent*, i.e., whether there is an assignment of all variables of Θ with dipoles such that all constraints are satisfied (a *solution*). We call this problem DSAT. DSAT is a Constraint Satisfaction Problem (CSP) [9] and can be solved using the standard methods developed for CSP's with infinite domains (see, e.g. [8]).

A partial method for determining inconsistency of a set of constraints Θ is the *path-consistency method* which enforces path-consistency on Θ [9]. A set of constraints is path-consistent if and only if for any two variables, there exists an instantiation of any third variable such that the three values taken together are consistent. It is necessary but not sufficient for the consistency of a set of constraints that path-consistency can be enforced. A naive way to enforce path-consistency is to strengthen relations by successively applying the following operation until a fixed point is reached:

$$\forall i, j, k : R_{ij} \leftarrow R_{ij} \cap (R_{ik} \circ R_{kj})$$

where i, j, k are nodes and R_{ij} is the relation between i and j . The resulting set of constraints is equivalent to the original set, i.e., it has the same set of solutions. If the empty relation occurs while performing this operation Θ is inconsistent, otherwise the resulting set is path-consistent. In Section 6 we use the path-consistency method to solve a small navigation problem with the dipole calculus.

4 Computational Properties of the Dipole Calculus

Although we restricted the possible binary relations between dipoles to 24 atomic relations, \mathcal{DRA}_{24} is very expressive. For instance, it is

⁴ An electronic version of the table can be obtained at <http://www.informatik.uni-hamburg.de/WSV/DRA>

possible to express directed and undirected graphs and their properties such as planarity or (convex) cycles. Hence, it is not surprising that DSAT(\mathcal{DRA}_{24}) is NP-hard which can be shown by reduction of the BETWEENNESS problem (Instance: Finite set A , collection C of ordered triples (a, b, c) of distinct elements from A , Question: Is there a one-to-one function f from A to $1, 2, \dots, |A|$ such that for each (a, b, c) in C , $f(a) < f(b) < f(c)$ or $f(c) < f(b) < f(a)$). [5])

Theorem 1 DSAT(\mathcal{DRA}_{24}) is NP-hard

Proof. Reduction from BETWEENNESS. Given a finite set A and a collection C of ordered triples (a, b, c) of distinct elements from A . For every element a of A introduce two dipoles a_1 and a_2 such that $a_1\{ells, errs\}a_2$ holds. For every pair a, b of distinct elements of A we require that $a_1\{slsr, srsr\}b_1$, $a_1\{lere, rele\}b_1$, and $a_i\{rllr, lrll\}b_i$ (for $i = 1, 2$) holds. The latter constraint guarantees that the graph formed by the dipoles $a_1, a_2, b_1, b_2, \dots$ is planar (see Figure 4).

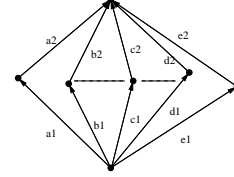


Figure 4. Reduction of a set A to a graph of dipoles

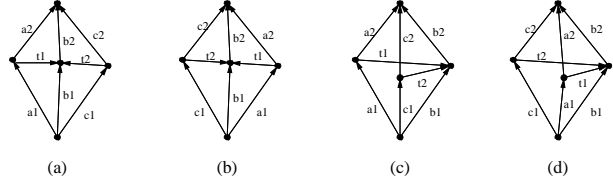


Figure 5. Reduction of a triple (a, b, c) to dipole constraints. If b is between a and c , the constraints are satisfied (see (a),(b)), if b is between a and c , then either t_1 overlaps c_1 or c_2 or t_2 overlaps a_1 or a_2 which contradicts the constraints (see (c),(d)).

For every ordered triple $t = (a, b, c)$ we introduce the two dipoles t_1, t_2 and the constraints $a_1\{ells, errs\}t_1$, $b_1\{lere, rele\}t_1$, $b_1\{lere, rele\}t_2$, $c_1\{ells, errs\}t_2$, $a_i\{rllr, lrll\}t_2$, and $b_i\{rllr, lrll\}t_1$. As it can be seen in Figure 5, these constraints guarantee that the set of dipole constraints Θ is consistent iff there is a one-to-one function f from A to $1, 2, \dots, |A|$ such that for each (a, b, c) in C , $f(a) < f(b) < f(c)$ or $f(c) < f(b) < f(a)$. \square

We have so far neither been able to prove that DSAT(\mathcal{DRA}_{24}) is a member of NP nor whether reasoning over the atomic relations is tractable. However, it follows from the algebraic semantics of the relations that DSAT(\mathcal{DRA}_{24}) is a member of PSPACE. This is because all relations can be expressed as equalities over polynomials with integer coefficients. Systems of such equalities can be solved using polynomial space [11].

5 An Extended Version of the Dipole Calculus

In certain domains we might want to represent spatial arrangements in which more than two start or end points of dipoles are on a straight line. Then we need three more dipole-point relations. The additional relations describe the cases when the point is straight behind the dipole (b), in the interior of the dipole (i) or straight in front of the dipole (f). The corresponding regions are shown on Figure 6. Such a set of relations was proposed by Freksa [4].

