

Interval Reachability Analysis using Second-Order Sensitivity^{*}

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Abstract: We propose a new approach to compute an interval over-approximation of the finite time reachable set for a large class of nonlinear systems. This approach relies on the notions of sensitivity matrices, which are the partial derivatives representing the variations of the system trajectories in response to variations of the initial states. Using interval arithmetics, we first over-approximate the possible values of the second-order sensitivity at the final time of the reachability problem. Then we exploit these bounds and the evaluation of the first-order sensitivity matrices at a few sampled initial states to obtain an over-approximation of the first-order sensitivity, which is in turn used to over-approximate the reachable set of the initial system. Unlike existing methods relying only on the first-order sensitivity matrix, this new approach provides guaranteed over-approximations of the first-order sensitivity and can also provide such over-approximations with an arbitrary precision by increasing the number of samples.

Keywords: Reachability analysis, mixed-monotonicity, sensitivity, interval.

1. INTRODUCTION

Reachability analysis is the problem of evaluating the set of all the successor states that can be reached in finite time by a system starting from a given set of initial states (Blanchini and Miani, 2008). Since the reachable set can rarely be computed exactly, we often rely on methods to over-approximate this set. In the literature, we primarily find two classes of reachability approaches. The first class considers complex and flexible set representations, such as zonotopes (Althoff, 2015), zonotope bundles (Althoff and Krogh, 2011) ellipsoids (Kurzahnskiy and Varaiya, 2007), support functions (Girard and Le Guernic, 2008), paving of intervals (Jaulin, 2001). Their main focus is to over-approximate the reachable set as tightly as possible, which is particularly interesting to solve simple verification problems such as those with safety or reachability specifications where the obtained over-approximation is immediately checked against a set of unsafe or target states.

The second class considers a simpler set representation in the form of (multi-dimensional) intervals, using methods based on differential inequalities (Scott and Barton, 2013), Taylor models (Chen et al., 2012), growth bounds (Reissig et al., 2016) or monotonicity (Meyer et al., 2019). Due to the simpler set representation, these methods tend to offer better efficiency and scalability at the cost of the accuracy of the over-approximations, and are thus particularly used in the field of abstraction-based control synthesis (see e.g. Moor and Raisch, 2002; Coogan and Arcak, 2015; Reissig et al., 2016; Meyer and Dimarogonas, 2019) where the number of reachable set over-approximations required for

the creation of an abstraction grows exponentially in the dimension of the state space.

In the subset of monotonicity-based interval reachability approach, the simplest method, used in Moor and Raisch (2002), relies directly on a monotonicity property (Angeli and Sontag, 2003) and guarantees that an interval over-approximation of the reachable set can be computed by evaluating the successors of only two vertices of the interval of initial states. A generalization of this property called mixed-monotonicity was then introduced and used for reachability analysis in Coogan and Arcak (2015), where an auxiliary monotone system can be created by decomposing the initial system into its increasing and decreasing components. A further generalization of mixed-monotonicity to any system with a bounded Jacobian matrix was recently proposed in Yang et al. (2019) and used for reachability analysis in Meyer and Dimarogonas (2019). Finally, another interval reachability method inspired by the notion of mixed-monotonicity and applicable to continuous-time nonlinear systems was proposed in Meyer et al. (2018), where bounds on the sensitivity matrix (partial derivative describing the influence of initial conditions on successor states) are used to compute an over-approximation interval of the reachable set.

While Meyer et al. (2018) considers two approaches to evaluate these sensitivity bounds, both have shortcomings: one provides very conservative bounds by applying the interval arithmetics results from Althoff et al. (2007), the other only computes empirical bounds through a time-consuming sampling procedure which is not guaranteed to result in an over-approximation of the sensitivity values. In this paper, we propose a novel and more flexible algorithm to obtain sensitivity bounds by combining the advantages of these two approaches while overcoming their main drawbacks. In addition to the first-order sensitivity matrix

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used above, the proposed approach also relies on the second-order sensitivity in the following 3-step procedure:

- first over-approximate the reachable tube (over the whole time range) for the first-order sensitivity matrix using interval arithmetics,
- next use these bounds to over-approximate the reachable set (at the final time only) for the second-order sensitivity using interval arithmetics,
- finally combine the second-order sensitivity bounds with the numerical evaluation of the first-order sensitivity on some sampled initial states to obtain an over-approximation of the reachable set of the first-order sensitivity.

This result has two major advantages. Compared to the purely empirical sampling approach from Meyer et al. (2018), the proposed algorithm is sound since for any number of samples we are guaranteed to over-approximate the set of first-order sensitivity values. Compared to the one-step interval arithmetics method from Meyer et al. (2018), which is conservative, we can now obtain arbitrarily tight bounds of the first-order sensitivity by increasing the number of samples. Indeed, the sampling in our third step can be used to tune the desired tradeoff between the computational complexity and the conservativeness of the over-approximation. Compared to methods relying on Taylor models (Chen et al., 2012) which usually require a decomposition of the time range to reduce the accumulation of errors, the proposed approach relying on mixed-monotonicity does not have this problem and all over-approximations can be computed in a single time step.

Due to space limitation, proofs and additional details are provided in the extended version of this paper.¹

2. PRELIMINARIES

Let \mathbb{R} and \mathbb{N} be the sets of reals and positive integers, respectively. Let $I_n \in \mathbb{R}^{n \times n}$ and $\mathbf{0}_{n \times p}, \mathbf{1}_{n \times p} \in \mathbb{R}^{n \times p}$ denote the identity matrix of dimension n and the $n \times p$ matrices filled with zeros and ones, respectively. Given two matrices $A \in \mathbb{R}^{n \times p}$ and $B \in \mathbb{R}^{q \times r}$, we denote their matrix product (if $p = q$) as $A * B = AB \in \mathbb{R}^{n \times r}$ and their Kronecker product as $A \otimes B \in \mathbb{R}^{nq \times pr}$.

Let $\mathcal{I} \subseteq 2^{\mathbb{R}}$ be the set of closed real intervals, i.e., for all $X \in \mathcal{I}$, there exist $\underline{x}, \bar{x} \in \mathbb{R}$ such that $X = [\underline{x}, \bar{x}] = \{x \in \mathbb{R} \mid \underline{x} \leq x \leq \bar{x}\} \subseteq \mathbb{R}$. \mathcal{I}^n and $\mathcal{I}^{n \times p}$ then represent the sets of interval vectors in \mathbb{R}^n and interval matrices in $\mathbb{R}^{n \times p}$, respectively. Given two interval matrices $[\underline{A}, \bar{A}], [\underline{B}, \bar{B}] \in \mathcal{I}^{n \times p}$, their sum is: $[\underline{A}, \bar{A}] + [\underline{B}, \bar{B}] = [\underline{A} + \underline{B}, \bar{A} + \bar{B}]$. From Jaulin (2001), the product of two scalar intervals is defined as

$$[\underline{a}, \bar{a}] * [\underline{b}, \bar{b}] = [\min(\underline{a}\underline{b}, \underline{a}\bar{b}, \bar{a}\underline{b}, \bar{a}\bar{b}), \max(\underline{a}\underline{b}, \underline{a}\bar{b}, \bar{a}\underline{b}, \bar{a}\bar{b})] \in \mathcal{I}.$$

For $[\underline{A}, \bar{A}] \in \mathcal{I}^{n \times p}$ and $[\underline{B}, \bar{B}] \in \mathcal{I}^{p \times q}$, the product $[\underline{C}, \bar{C}] = [\underline{A}, \bar{A}] * [\underline{B}, \bar{B}] \in \mathcal{I}^{n \times q}$ is defined elementwise such that

$$[\underline{C}_{ij}, \bar{C}_{ij}] = \sum_{k=1}^p [\underline{A}_{ik}, \bar{A}_{ik}] * [\underline{B}_{kj}, \bar{B}_{kj}] \in \mathcal{I},$$

and the product of a scalar interval with a matrix interval is defined as $[\underline{C}, \bar{C}] = [\underline{a}, \bar{a}] * [\underline{B}, \bar{B}] \in \mathcal{I}^{p \times q}$ with

$$[\underline{C}_{ij}, \bar{C}_{ij}] = [\underline{a}, \bar{a}] * [\underline{B}_{ij}, \bar{B}_{ij}] \in \mathcal{I}.$$

For $[\underline{A}, \bar{A}] \in \mathcal{I}^{n \times p}$ and $[\underline{B}, \bar{B}] \in \mathcal{I}^{q \times r}$, the interval Kronecker product $[\underline{C}, \bar{C}] = [\underline{A}, \bar{A}] \otimes [\underline{B}, \bar{B}] \in \mathcal{I}^{nq \times pr}$ is defined as a $n \times p$ block interval matrix with (i, j) block

$$[\underline{C}_{ij}, \bar{C}_{ij}] = [\underline{A}_{ij}, \bar{A}_{ij}] * [\underline{B}, \bar{B}] \in \mathcal{I}^{q \times r}.$$

2.1 Functional matrices

In this section, we provide definitions and results on the manipulation of functional matrices used throughout the paper. We first introduce the differential operator D for a scalar differentiable function $f : \mathbb{R}^n \rightarrow \mathbb{R}$ to be:

$$Df(x) = \left(\frac{\partial f(x)}{\partial x_1} \quad \dots \quad \frac{\partial f(x)}{\partial x_n} \right).$$

Then for a functional matrix $A : \mathbb{R}^n \rightarrow \mathbb{R}^{p \times q}$, its differential $DA(x) \in \mathbb{R}^{p \times nq}$ is the $p \times q$ block matrix where each element $A_{ij}(x) \in \mathbb{R}$ of $A(x) \in \mathbb{R}^{p \times q}$ is replaced by the row vector of its differential $DA_{ij}(x) \in \mathbb{R}^{1 \times n}$:

$$DA(x) = \begin{pmatrix} DA_{11}(x) & \dots & DA_{1q}(x) \\ \vdots & \ddots & \vdots \\ DA_{p1}(x) & \dots & DA_{pq}(x) \end{pmatrix} \quad (1)$$

$$= \begin{pmatrix} \frac{\partial A_{11}(x)}{\partial x_1} & \dots & \frac{\partial A_{11}(x)}{\partial x_n} & \dots & \frac{\partial A_{1q}(x)}{\partial x_1} & \dots & \frac{\partial A_{1q}(x)}{\partial x_n} \\ \vdots & & \vdots & & \vdots & & \vdots \\ \frac{\partial A_{p1}(x)}{\partial x_1} & \dots & \frac{\partial A_{p1}(x)}{\partial x_n} & \dots & \frac{\partial A_{pq}(x)}{\partial x_1} & \dots & \frac{\partial A_{pq}(x)}{\partial x_n} \end{pmatrix}.$$

This notation ensures that we only work with 2-dimensional matrices, instead of matrices with more than two dimensions for which cumbersome matrix product definitions would need to be introduced.

For a time-varying functional matrix $A : \mathbb{R} \times \mathbb{R}^n \rightarrow \mathbb{R}^{p \times q}$, its time derivative is denoted with a dot

$$\dot{A}(t, x) = \frac{\partial A(t, x)}{\partial t},$$

and we keep the notation $DA(t, x)$ as in (1) to denote its derivative with respect to the second variable $x \in \mathbb{R}^n$.

For the product of two functional matrices, its differential is obtained as in the following result from (Cheng et al., 2012, Corollary 18.1).

Lemma 1. (Product rule). Given $A : \mathbb{R}^n \rightarrow \mathbb{R}^{p \times q}$, $B : \mathbb{R}^n \rightarrow \mathbb{R}^{q \times r}$, we have $D(A(x)B(x)) \in \mathbb{R}^{p \times nr}$ given by

$$D(A(x)B(x)) = DA(x) * (B(x) \otimes I_n) + A(x) * DB(x).$$

Next, we introduce the chain rule for the composition of a functional vector and functional matrix.

Lemma 2. (Chain rule). Given $A : \mathbb{R}^m \rightarrow \mathbb{R}^{p \times q}$ and $b : \mathbb{R}^n \rightarrow \mathbb{R}^m$, we have $D(A(b(x))) \in \mathbb{R}^{p \times nq}$ given by

$$D(A(b(x))) = DA(y)|_{y=b(x)} * (I_q \otimes Db(x)).$$

2.2 Reachability analysis of interval affine systems

The method presented in this paper partly relies on results from Althoff et al. (2007) which use interval arithmetics to over-approximate the reachable set and reachable tube of affine interval systems. These results are summarized in this section for self-containment of the paper.

Consider an affine interval system of the form

$$\dot{z} \in \mathcal{A}z + \mathcal{B}, \quad (2)$$

¹ Available at: <https://arxiv.org/abs/1911.09775>

with state $z \in \mathbb{R}^{p \times q}$ and interval matrices $\mathcal{A} = [\underline{A}, \overline{A}] \in \mathcal{I}^{p \times p}$ and $\mathcal{B} = [\underline{B}, \overline{B}] \in \mathcal{I}^{p \times q}$. Given an interval matrix of initial states $Z_0 = [z_0, \overline{z}_0] \in \mathcal{I}^{p \times q}$ and a time step $\tau > 0$, we denote the reachable set of (2) as $z(\tau, Z_0) \subseteq \mathbb{R}^{p \times q}$ and its reachable tube as $z([0, \tau], Z_0) = \bigcup_{t \in [0, \tau]} z(t, Z_0) \subseteq \mathbb{R}^{p \times q}$.

The results from Althoff et al. (2007) rely on Taylor series truncated at an order $r \in \mathbb{N}$ which needs to satisfy $r > \|\mathcal{A}\|_\infty \tau - 2$, where the infinity norm of the interval matrix is defined by $\|\mathcal{A}\|_\infty = \|\max(|\underline{A}|, |\overline{A}|)\|_\infty$ using componentwise absolute value and max operators. Then we introduce

$$C(\tau) = [-\mathbf{1}_{p \times p}, \mathbf{1}_{p \times p}] * \frac{(\|\mathcal{A}\|_\infty \tau)^{r+1}}{(r+1)!} \frac{r+2}{r+2 - \|\mathcal{A}\|_\infty \tau},$$

$$D(\tau) = \sum_{i=0}^r \frac{(\mathcal{A}\tau)^i}{i!} + C(\tau),$$

$$E(\tau) = \sum_{i=0}^r \frac{\mathcal{A}^i \tau^{i+1}}{(i+1)!} + C(\tau)\tau,$$

$$F(\tau) = \left[\sum_{i=2}^r \left(i^{\frac{-i}{i-1}} - i^{\frac{-1}{i-1}} \right) \frac{(\mathcal{A}\tau)^i}{i!}, \mathbf{0}_{p \times p} \right] + C(\tau),$$

where all sums and products of interval matrices follow the definitions in Section 2. We also define the interval hull of two interval matrices $[\underline{a}, \overline{a}], [\underline{b}, \overline{b}] \in \mathcal{I}^{p \times q}$ as $H([\underline{a}, \overline{a}], [\underline{b}, \overline{b}]) = [\min(\underline{a}, \underline{b}), \max(\overline{a}, \overline{b})]$ using the componentwise min and max operators.

Lemma 3. (Althoff et al. (2007)). The reachable set of (2) at time $\tau \geq 0$ is over-approximated by an interval in $\mathcal{I}^{p \times q}$ as follows:

$$z(\tau, Z_0) \subseteq D(\tau)Z_0 + E(\tau)\mathcal{B}. \quad (3)$$

If in addition we have $\mathcal{B} = \{\mathbf{0}_{p \times q}\}$, then the reachable tube of (2) over time range $[0, \tau]$ is over-approximated by an interval in $\mathcal{I}^{p \times q}$ as follows:

$$z([0, \tau], Z_0) \subseteq H(Z_0, D(\tau)Z_0) + F(\tau)Z_0. \quad (4)$$

3. PROBLEM FORMULATION

We consider a continuous-time, time-varying system

$$\dot{x} = f(t, x), \quad (5)$$

with state $x \in \mathbb{R}^n$ and vector field $f : \mathbb{R} \times \mathbb{R}^n \rightarrow \mathbb{R}^n$ assumed to be twice differentiable in the state. We denote as $\Phi(t; t_0, x_0) \in \mathbb{R}^n$ the state reached by (5) at time $t \geq t_0$ from initial state x_0 .

Problem 4. Given a time range $[t_0, t_f] \in \mathcal{I}$ and an interval of initial states $X_0 = [\underline{x}, \overline{x}] \in \mathcal{I}^n$, find an interval in \mathcal{I}^n over-approximating the reachable set of system (5) defined as $R(t_f; t_0, X_0) = \{\Phi(t_f; t_0, x_0) \mid x_0 \in X_0\}$.

To solve Problem 4 with the method presented in Section 5, we assume that bounds on both the first-order and second-order Jacobian matrices of (5) are provided by the user. These two Jacobian matrices are defined below using the differential operator D of the vector field $f(t, x)$ with respect to state x as introduced in Section 2.1:

$$J^x(t, x) = Df(t, x) \in \mathbb{R}^{n \times n},$$

$$J^{xx}(t, x) = DJ^x(t, x) \in \mathbb{R}^{n \times n^2}.$$

Assumption 5. Given an invariant state space $X \subseteq \mathbb{R}^n$ for system (5), there exist $[J^x, \overline{J^x}] \in \mathcal{I}^{n \times n}$ and $[J^{xx}, \overline{J^{xx}}] \in$

$\mathcal{I}^{n \times n^2}$ such that for all $t \in [t_0, t_f]$ and $x \in X$ we have $J^x(t, x) \in [J^x, \overline{J^x}]$ and $J^{xx}(t, x) \in [J^{xx}, \overline{J^{xx}}]$.

4. SENSITIVITY EQUATIONS

The method presented in Section 5 to solve Problem 4 relies on the definition of the sensitivity matrices of system (5) representing the differential influence of the initial conditions on the successor $\Phi(t; t_0, x_0)$ at time t . Similarly to the definition of the Jacobian matrices above, we use D to denote the differential operator of the trajectory $\Phi(t; t_0, x_0)$ with respect to initial state x_0 . Then the first-order and second-order sensitivity matrices are defined as:

$$S^x(t; t_0, x_0) = D\Phi(t; t_0, x_0) \in \mathbb{R}^{n \times n}, \quad (6)$$

$$S^{xx}(t; t_0, x_0) = DS^x(t; t_0, x_0) \in \mathbb{R}^{n \times n^2}. \quad (7)$$

Both sensitivity matrices defined in (6) and (7) can also be described by the time-varying affine systems below.

Proposition 6. Using the short-hand notations $S^x := S^x(t; t_0, x_0)$, $S^{xx} := S^{xx}(t; t_0, x_0)$, $J^x := J^x(t, \Phi(t; t_0, x_0))$ and $J^{xx} := J^{xx}(t, \Phi(t; t_0, x_0))$, the sensitivity matrices defined in (6) and (7) follow:

$$\dot{S}^x = J^x * S^x, \quad (8)$$

$$\dot{S}^{xx} = J^x * S^{xx} + J^{xx} * (S^x \otimes S^x), \quad (9)$$

with $S^x(t_0; t_0, x_0) = I_n$ and $S^{xx}(t_0; t_0, x_0) = \mathbf{0}_{n \times n^2}$.

Alternative derivations of second-order sensitivity equations have been obtained in Choi et al. (2016) for differential algebraic equations and Geng and Hiskens (2019) for hybrid systems.

5. REACHABILITY ALGORITHM

The proposed approach to solve Problem 4 is summarized in Algorithm 1 and Figure 1. Below, we briefly explain this algorithm by going backwards from step 4 to step 1.

The end goal in step 4 is to over-approximate the reachable set of the nonlinear system (5) using the recent reachability method in Meyer et al. (2018) that relies on interval bounds on the reachable set of the first-order sensitivity $S^x(t_f; t_0, X_0)$. The method in Meyer et al. (2018) uses either conservative bounds from a direct application of Lemma 3 or empirical bounds from a sampling procedure. In contrast, here we derive guaranteed bounds on S^x in step 3 by combining bounds on the reachable set of the second-order sensitivity $S^{xx}(t_f; t_0, X_0)$ with the numerical evaluation of S^x at time t_f on a finite set of sampled initial states. The resulting bounds on $S^x(t_f; t_0, X_0)$ can be made arbitrarily tight by increasing the number of samples.

The bounds on S^{xx} are computed in step 2 by applying (3) in Lemma 3 to (9), which requires the knowledge of bounds of both Jacobian matrices (from Assumption 5) and on the reachable tube of the first-order sensitivity $S^x([t_0, t_f]; t_0, X_0)$. This reachable tube of S^x is over-approximated in step 1 by applying (4) in Lemma 3 to (8), which requires bounds on J^x taken from Assumption 5.

5.1 Interval arithmetics on the sensitivity systems

For the first step of Algorithm 1, we first need to rewrite the time-varying linear system of the first-order sensitivity

Input: Reachability problem for (5): $t_0, t_f, X_0 = [\underline{x}, \bar{x}]$
Data: Jacobian bounds $[J^x, \bar{J}^x], [J^{xx}, \bar{J}^{xx}]$
Step 1: Apply (4) to (8) and obtain an interval over-approximation of $S^x([t_0, t_f]; t_0, X_0)$
Step 2: Apply (3) to (9) and obtain an interval over-approximation of $S^{xx}(t_f; t_0, X_0)$
Step 3: Obtain an interval over-approximation of $S^x(t_f; t_0, X_0)$ from the bounds on S^{xx} and the evaluation of $S^x(t_f; t_0, x_0)$ on a finite subset of X_0
Step 4: Obtain an interval over-approximation of $R(t_f; t_0, X_0)$ using the bounds on S^x (Meyer et al., 2018)
Output: Interval solving Problem 4

Algorithm 1: Reachability analysis of system (5).

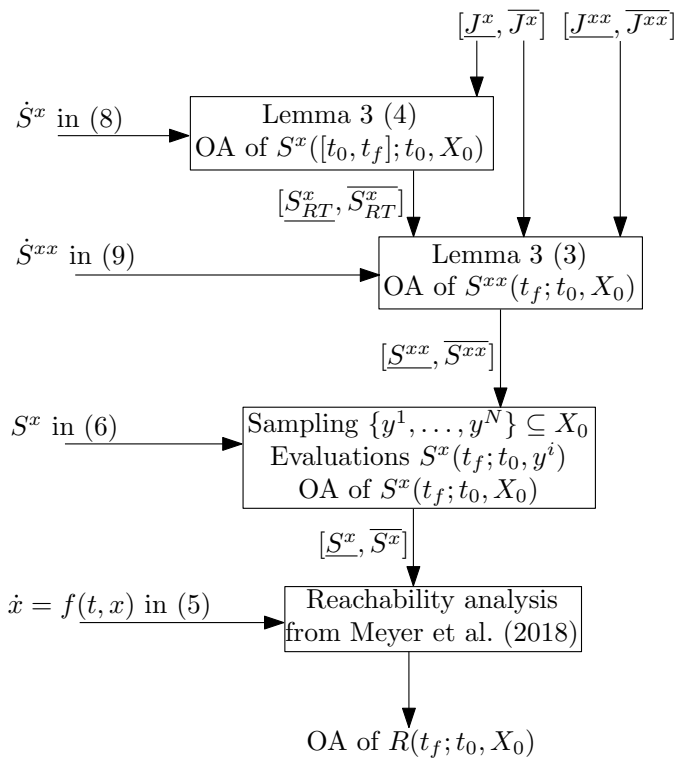


Fig. 1. Sketch of the 4-step reachability procedure in Algorithm 1 where “OA” stands “over-approximation”. For each box, top arrows are the input requirements, side arrows are the equations used and bottom arrows are the output results.

(8) into a linear interval system similarly to (2). This is done using the bounds on J^x from Assumption 5:

$$\dot{S}^x(t; t_0, x_0) \in [J^x, \bar{J}^x] * S^x(t; t_0, x_0). \quad (10)$$

Then, applying (4) in Lemma 3 with $\mathcal{A} = [J^x, \bar{J}^x]$, $\mathcal{B} = \{\mathbf{0}_{n \times n}\}$ and $Z_0 = \{I_n\}$ leads to an over-approximation of the reachable tube $S^x([t_0, t_f]; t_0, [\underline{x}, \bar{x}]) \subseteq [S^x_{RT}, \bar{S}^x_{RT}] \in \mathcal{I}^{n \times n}$ defined as:

$$[S^x_{RT}, \bar{S}^x_{RT}] = H(\{I_n\}, D(t_f - t_0)) + F(t_f - t_0).$$

For the second step in Algorithm 1, we use the bounds $[S^x_{RT}, \bar{S}^x_{RT}]$ obtained in the previous step alongside the Jacobian bounds from Assumption 5 to rewrite the time-varying affine system of the second-order sensitivity (9) into an affine interval system as in (2) with $\mathcal{A} = [J^x, \bar{J}^x]$, $\mathcal{B} = [J^{xx}, \bar{J}^{xx}] * ([S^x_{RT}, \bar{S}^x_{RT}] \otimes [S^x_{RT}, \bar{S}^x_{RT}])$ and

the initial condition $Z_0 = \{\mathbf{0}_{n \times n^2}\}$ from Proposition 6. This leads to an over-approximation of the reachable set $S^{xx}(t_f; t_0, [\underline{x}, \bar{x}]) \subseteq [S^{xx}, \bar{S}^{xx}] \in \mathcal{I}^{n \times n^2}$ defined as:

$$[S^{xx}, \bar{S}^{xx}] = E(t_f - t_0)\mathcal{B}.$$

5.2 Sampling for the first-order sensitivity

Step 3 of Algorithm 1 relies on the evaluation of the first-order sensitivity for some sampled initial states. Let $\{y^1, \dots, y^N\} = Y \subseteq [\underline{x}, \bar{x}]$ be a finite set of N samples in the interval of initial states $[\underline{x}, \bar{x}]$. Similarly to (Tempo et al., 2012, Section 7.4.4), we define below the dispersion of this set of samples, where the infinity norm of a state $x \in \mathbb{R}^n$ is defined as $\|x\|_\infty = \max_{i \in \{1, \dots, n\}} |x_i|$.

Definition 7. Given a finite set $Y \subseteq [\underline{x}, \bar{x}]$, the dispersion of Y in $[\underline{x}, \bar{x}]$ is defined as: $d(Y) = \sup_{x \in [\underline{x}, \bar{x}]} \min_{y \in Y} \|x - y\|_\infty \in \mathbb{R}$.

Smaller values of $d(Y)$ imply that the sample states in Y are well scattered in the interval $[\underline{x}, \bar{x}]$. After evaluating the first-order sensitivity $S^x(t_f; t_0, y^i)$ at time t_f for each of these sampled states through numerical integration of (6) or (8), we can derive guaranteed bounds on the set $S^x(t_f; t_0, [\underline{x}, \bar{x}])$ as follows.

Theorem 8. Given bounds on the second-order sensitivity $S^{xx}(t_f; t_0, [\underline{x}, \bar{x}]) \subseteq [S^{xx}, \bar{S}^{xx}] \in \mathcal{I}^{n \times n^2}$ and a finite set $Y \subseteq [\underline{x}, \bar{x}]$ of sampled initial states, define $M \in \mathbb{R}^{n \times n}$ as

$$M = \max(|S^{xx}|, |\bar{S}^{xx}|) * (I_n \otimes (\mathbf{1}_n * d(Y))),$$

using componentwise absolute value and max operators. Then the set of first-order sensitivity values at time t_f is over-approximated as $S^x(t_f; t_0, [\underline{x}, \bar{x}]) \subseteq [S^x, \bar{S}^x] \in \mathcal{I}^{n \times n}$ with, for all $i, j \in \{1, \dots, n\}$:

$$\bar{S}^x_{ij} = \max_{y \in Y} (S^x_{ij}(t_f; t_0, y)) + M_{ij},$$

$$S^x_{-ij} = \min_{y \in Y} (S^x_{ij}(t_f; t_0, y)) - M_{ij}.$$

The over-approximation interval $[S^x, \bar{S}^x]$ in Theorem 8 thus corresponds to the interval hull of the sampled sensitivity evaluations $\{S^x(t_f; t_0, y) | y \in Y\}$ dilated by M .

From the definition of M in Theorem 8, we can see that the size of the obtained bounds on the first-order sensitivity S^x grows with the dispersion of the sampling set Y . As a consequence, the set Y can be used to tune the tradeoff between reducing the conservativeness of the sensitivity bounds $[S^x, \bar{S}^x]$ and limiting the computation time (related to the number of samples). If computation capabilities were unlimited, Theorem 8 could then provide interval bounds of the first-order sensitivity values with arbitrary precision, as formulated below.

Proposition 9. If the sample number grows to infinity $N \rightarrow \infty$, we can design the sampling set Y such that $[S^x, \bar{S}^x]$ from Theorem 8 converges to the unique tight interval over-approximation of the set $S^x(t_f; t_0, [\underline{x}, \bar{x}])$, i.e. the smallest (in terms of inclusion) interval over-approximation.

5.3 Reachability analysis of the initial system

This section corresponds to step 4 of Algorithm 1 in which we apply the method for reachability analysis introduced

in Meyer et al. (2018). This reachability result is summarized below for self-containment of this paper.

Let $S^{x*} \in \mathbb{R}^{n \times n}$ denote the center of $[\underline{S}^x, \overline{S}^x]$ and define the decomposition function $g : \mathbb{R} \times \mathbb{R}^n \times \mathbb{R}^n \rightarrow \mathbb{R}^n$ whose i^{th} component with $i \in \{1, \dots, n\}$ is

$$g_i(t_0, x, y) = \Phi_i(t_f; t_0, z^i) + \alpha^i(x - y), \quad (11)$$

where the state $z^i = [z_1^i; \dots; z_n^i] \in \mathbb{R}^n$ and row vector $\alpha^i = [\alpha_1^i, \dots, \alpha_n^i] \in \mathbb{R}^{1 \times n}$ are such that for all $j \in \{1, \dots, n\}$,

$$(z_j^i, \alpha_j^i) = \begin{cases} (x_j, \max(0, -S_{ij}^{x*})) & \text{if } S_{ij}^{x*} \geq 0, \\ (y_j, \max(0, \overline{S}_{ij}^x)) & \text{if } S_{ij}^{x*} < 0. \end{cases} \quad (12)$$

Then an over-approximation of the reachable set of (5) is obtained by computing only two evaluations of the decomposition function g .

Lemma 10. (Meyer et al. (2018)). Given bounds on the first-order sensitivity $S^x(t_f; t_0, [\underline{x}, \overline{x}]) \subseteq [\underline{S}^x, \overline{S}^x] \in \mathcal{I}^{n \times n}$ and the definitions in (11)-(12), an over-approximation of the reachable set of (5) is given by:

$$R(t_f; t_0, [\underline{x}, \overline{x}]) \subseteq [g(t_0, \underline{x}, \overline{x}), g(t_0, \overline{x}, \underline{x})].$$

6. NUMERICAL ILLUSTRATION

In this section, we illustrate the approach in Algorithm 1 and compare it to the alternative methods from Meyer et al. (2018) on a numerical example. We consider the continuous-time uncertain unicycle model described as:

$$\dot{x} = [v \cos(x_3) + x_4; v \sin(x_3) + x_5; \omega + x_6; 0; 0; 0], \quad (13)$$

where $[x_1; x_2]$ is the 2D position of the unicycle, x_3 is its orientation, $[x_4; x_5; x_6]$ are constant uncertain parameters in the dynamics of the first three states, $v = 0.25$ is the controlled forward velocity and $\omega = 0.3$ is the controlled angular velocity. Using the conservative bounds $\cos(x_3), \sin(x_3) \in [-1, 1]$, global Jacobian bounds of (13) satisfying Assumption 5 are obtained by taking $[\underline{J}_{1,4}^x, \overline{J}_{1,4}^x] = [\underline{J}_{2,5}^x, \overline{J}_{2,5}^x] = [\underline{J}_{3,6}^x, \overline{J}_{3,6}^x] = \{1\}$, $[\underline{J}_{-1,3}^x, \overline{J}_{1,3}^x] = [\underline{J}_{-2,3}^x, \overline{J}_{2,3}^x] = [\underline{J}_{-1,15}^{xx}, \overline{J}_{1,15}^{xx}] = [\underline{J}_{-2,15}^{xx}, \overline{J}_{2,15}^{xx}] = [-v, v]$ and $[\underline{J}_{-ij}^x, \overline{J}_{ij}^x] = [\underline{J}_{-ij}^{xx}, \overline{J}_{ij}^{xx}] = \{0\}$ for all other elements.

Taking the initial time $t_0 = 0$, we want to evaluate the reachable set of (13) at time $t_f = 10$ for the following interval of initial conditions: $X_0 = [0, 1] \times [0, 1] \times [\frac{\pi}{8}, \frac{2\pi}{8}] \times [-0.05, 0.05] \times [-0.05, 0.05] \times [-0.03, 0.03]$. This reachability problem is solved in five ways described below.

- We first apply Algorithm 1 three times using a uniform grid sampling with an increasing number of samples per dimension of the state space $a \in \{1, 2, 3\}$ (leading to a total number of sample points of $N = a^6 \in \{1, 64, 729\}$). In Figures 2 and 3, these results are plotted in dashed red, dot-dashed blue and plain green, respectively.
- Next we use the one-step interval arithmetics (“IA” in Table 1) approach from Meyer et al. (2018), plotted in dotted purple.
- Finally we apply the sampling and falsification (“SF” in Table 1) approach from Meyer et al. (2018) using $N = 64$ samples, plotted in dashed orange.

The computation times for each of the four steps in Algorithm 1 (or alternatively, for obtaining bounds on

$S^x(t_f; t_0, X_0)$ in both methods from Meyer et al. (2018)) are reported in Table 1. The obtained bounds on $S_{1,3}^x$ and $S_{2,3}^x$ for step 3 are plotted in Figure 2 and the final reachability analysis (step 4) on states x_1 and x_2 is shown in Figure 3. In both figures, the cloud of black dots represents the numerical integration of (6) and (13), respectively, for 500 random samples in X_0 .

From Table 1, we first note that the computation of the final reachable set (step 4) is very fast and identical for all methods since this step is oblivious to the way the sensitivity bounds $[\underline{S}^x, \overline{S}^x]$ are obtained. As expected, the three steps relying on the interval arithmetics results from Lemma 3 (steps 1 and 2 in Algorithm 1 and step 3 in method “IA”) are also achieved quickly. The sampling computations in step 3 of Algorithm 1 naturally grows with the number of samples. For the sampling and falsification approach from Meyer et al. (2018), the sampling time is identical to the one in the second call of Algorithm 1 (due to having the same number of samples $N = 64$), but then the total computation time is increased by the 2 iterations of the falsification procedure used to improve the estimated bounds on S^x . Such expansion of the bounds is not required in Algorithm 1 since from Theorem 8, step 3 is already guaranteed to over-approximate $S^x(t_f; t_0, X_0)$.

In Figure 2, we can first note that, as hinted in Proposition 9, the bounds on the first-order sensitivity obtained in Algorithm 1 shrink as we increase the number of samples. We can also see that the one-step interval arithmetics method from Meyer et al. (2018) gives very conservative bounds on S^x (similar in size to Algorithm 1 with a single sample point). While the sampling and falsification method from Meyer et al. (2018) gives the closest approximation of $S^x(t_f; t_0, X_0)$, the obtained bounds are not actually an over-approximation of this set (despite the 2 iterations of falsification), which means that applying step 4 with such bounds is not sound for the reachability analysis of (13).

Finally, we can combine Figure 3 and Table 1 to conclude on the ability of Algorithm 1 to tune to our needs the tradeoff between computation time and conservativeness. The sampling and falsification approach from Meyer et al. (2018) is discarded from this discussion as we already showed above that it is unreliable when we want guaranteed over-approximations. When computation time is our main concern, we can take $N = 1$ in Algorithm 1 to obtain results comparable to the one-step “IA” method from Meyer et al. (2018), in terms of both conservativeness and low computation time. In particular, although the computation time of the interval arithmetics steps 1-2 would slightly increase with higher state dimension n , the computational complexity of steps 3-4 is constant (i.e. independent of the state dimension) when we take $N = 1$. On the other hand, if more computational power is available, increasing the number of samples tightens the over-approximation and in this example, we can see in Figure 3 that both $N = 64$ and $N = 729$ give tighter bounds than the method from Meyer et al. (2018).

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Samples N	Algorithm 1			IA	SF
	1	64	729		
$[S_{RT}^x, \overline{S}_{RT}^x]$	0.72			-	-
$[S^{xx}, \overline{S}^{xx}]$	0.87			-	-
$[\underline{S}^x, \overline{S}^x]$	0.35	3.2	36	0.44	$3.1 + 4.2$
OA of $R(t_f; t_0, X_0)$	0.07				

Table 1. Time comparison (in seconds) of the steps for reachability analysis in Algorithm 1 with three different sampling grids, and in both methods from Meyer et al. (2018) using a single step interval arithmetics (IA) or sampling and falsification (SF).

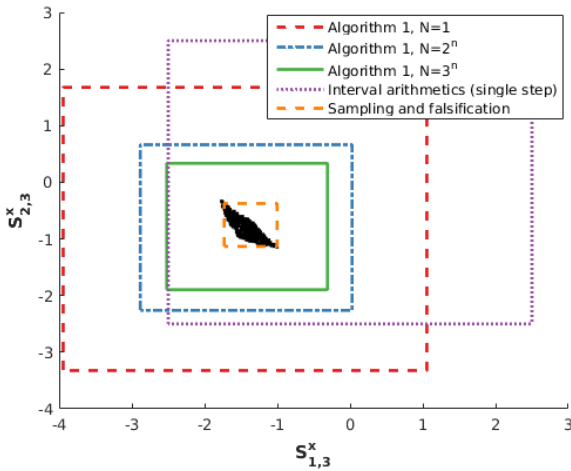


Fig. 2. Over-approximations of the first-order sensitivity components $S_{1,3}^x$ and $S_{2,3}^x$ at time t_f .

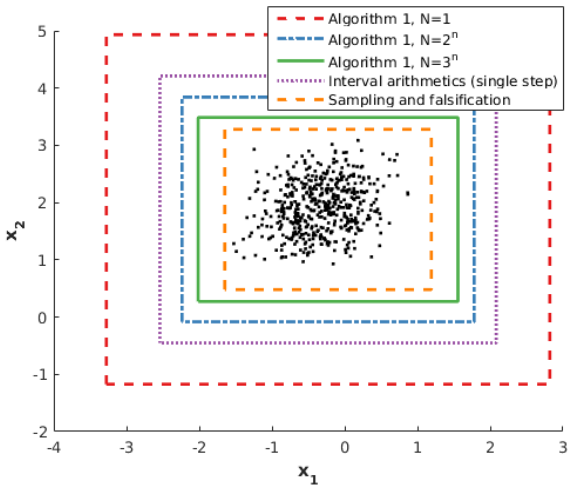


Fig. 3. Comparison of over-approximations of the reachable set of (13) at time t_f for states x_1 and x_2 .

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