

# Computational complexity of solving polynomial differential equations over unbounded domains

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Joint work with Daniel Graça

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# Ordinary Differential Equations (ODEs)

System of ODEs:

$$\begin{cases} y_1(0) = y_{0,1} \\ \vdots \\ y_n(0) = y_{0,n} \end{cases} \quad \begin{cases} y_1'(t) = f_1(y_1(t), \dots, y_n(t), t) \\ \vdots \\ y_n'(t) = f_n(y_1(t), \dots, y_n(t), t) \end{cases}$$

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In this talk: autonomous first order explicit system of ODEs

$$y(0) = y_0 \quad y' = f(y) \quad y : (a, b) \rightarrow \mathbb{R}^n$$

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In this talk (unless specified)

We use Computable Analysis.



## Computability of solutions: the theory

Let  $I = (a, b)$  and  $f \in C^0(\mathbb{R}^d)$ . Assume  $y \in C^1(I, \mathbb{R}^d)$  satisfies  $\forall t \in I$ :

$$y(0) = 0, \quad y'(t) = f(y(t)). \quad (1)$$

Given  $t \in I$  and  $n \in \mathbb{N}$ , can we compute  $q \in \mathbb{Q}^d$  s.t.  $\|q - y(t)\| \leq 2^{-n}$  ?

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**Theorem (Collins and Graça)**

The map  $f \mapsto y(\cdot)$  for those  $f$  where  $y$  is unique, is computable.

## Complexity of solutions: typical textbook result

Assume  $f$  **Lipschitz** and computable, and  $y : I \rightarrow \mathbb{R}^d$  satisfies  $\forall t \in I$ :

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## Problems with this approach:

- Accuracy of the result?  $\mathcal{O}(h^4) \leq Ah^4$  but  $A$  is **unknown**
- Same problem with complexity
- $f$  is **Lipschitz**: typically only holds over **compact** domains

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- you know  $K$ :  $f$  must be Lipschitz on “ $\{y(u) : u \in I\}$ ” or globally
- you know  $M$ : **but it depends on  $y$  !!**

**Chicken-and-egg problem**: the constant in the accuracy bound depends on computing the solution.



## Complexity of solutions: the rescaling “myth”

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**To compute  $y(T)$  we could:**

- 1 Define  $z(u) = y(Tu)$ , then

$$y(T) = z(1)$$

- 2 Observe that

$$z'(u) = Tf(z) =: f_T(z)$$

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**Accuracy:**  $A_{K_T, M_z} h$  where

$K_T =$  Lipschitz constant of  $f_T$

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## Conclusion

This tells us **nothing** about the complexity of the problem.

## Side note on practical methods

Assume  $y : [0, 1] \rightarrow \mathbb{R}^d$  satisfies  $\forall t \in [0, 1]$ :

$$y(0) = 0, \quad y'(t) = f(y(t)).$$

There exists methods of the form:

given  $h$  and  $t$ , compute  $q \in \mathbb{Q}^d$  and  $\varepsilon > 0$  such that  $\|y(t) - q\| \leq \varepsilon$  with the guarantee that  $\varepsilon \rightarrow 0$  as  $h \rightarrow 0$ .

These methods have **no upper bound** on complexity.

They usually rely on interval arithmetic.

# Nonuniform complexity-theoretic approach

Assume  $y : [0, 1] \rightarrow \mathbb{R}^d$  satisfies  $\forall t \in [0, 1]$ :

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| <b>Assumption on <math>f</math></b> | <b>Lower bound on <math>y</math></b> | <b>Upper bound on <math>y</math></b> |
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| PTIME + analytic        | —                  | PTIME                |

But those results can be **deceiving**...

$$\begin{cases} y_1(0) = 1 \\ y_2(0) = 1 \\ \vdots \\ y_d(0) = 1 \end{cases} \quad \begin{cases} y'_1 = y_1 \\ y'_2 = y_1 y_2 \\ \vdots \\ y'_n = y_{d-1} y_n \end{cases} \quad \rightarrow \quad \begin{matrix} y(t) = \mathcal{O} \left( e^{e^{\dots e^t}} \right) \\ y \text{ is PTIME over } [0, 1] \end{matrix}$$

# Nonuniform complexity: limitation

Example:

$f$  PTIME analytic  $\Rightarrow y$  PTIME  $\Rightarrow y(t) \pm 2^{-n}$  in time  $An^k$

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## Conclusion

This only **slightly** better than the previous approach.

# Uniform (operator) complexity approach

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$$T(t, n, K_d, K_f)$$

where

- $K_d$ : depends on the dimension  $d$
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# Uniform (operator) complexity approach

Assume  $y : I \rightarrow \mathbb{R}^d$  satisfies  $\forall t \in I$ :

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**Problem:** we cannot predict the behaviour of  $y$  based on  $f$  only.

# Are you confused?

You should be!

- practical methods: “no complexity”
- nonuniform complexity: misleading
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**Question:** are we looking at the problem the wrong way?

# Parametrized complexity approach

**Goal:** Assume  $y : I \rightarrow \mathbb{R}^d$  satisfies  $\forall t \in I$ :

$$y(0) = 0, \quad y'(t) = f(y(t)),$$

where  $f : \mathbb{R}^d \rightarrow \mathbb{R}^d$  is **nice**. Then  $y(t) \pm 2^{-n}$  can be computed in time

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**Important** differences with “textbook” approach:

- Result is always correct
- $K_y$  not assumed to be known (e.g.  $K$  and  $M$  of previous slides)

# Parametrized complexity result

Assume  $y : I \rightarrow \mathbb{R}^d$  satisfies  $\forall t \in I$ :

$$y(0) = 0, \quad y'(t) = p(y(t)),$$

where  $p : \mathbb{R}^d \rightarrow \mathbb{R}^d$  is **vector of multivariate polynomials**.

## Theorem (TCS 2016)

Assuming  $t \in I$ , computing  $y(t) \pm 2^{-n}$  takes time:

$$\text{poly}(\text{deg } p, \log \Sigma p, n, \ell_y(t))^d$$

where:

- $\Sigma p$ : sum of absolute value of coefficients of  $p$

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where:

- $\Sigma p$ : sum of absolute value of coefficients of  $p$
- $\ell_y(t)$ : “length” of  $y$  over  $[0, t]$

$$\ell_y(t) = \int_0^t \max(1, \|y'(u)\|) du$$

**Note:** the algorithm find  $\ell(t)$  automatically, **it is not part of the input**



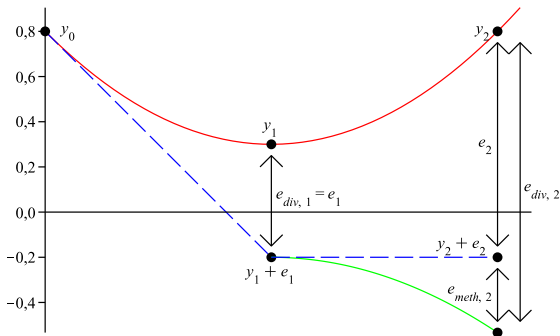
# Euler method

$$y(0) = 0 \quad y'(t) = p(y(t))$$

Time step  $h$ , discretize and compute  $\tilde{y}^i \approx y(ih)$ :

$$y(t+h) \approx y(t) + hy'(t) \quad \leadsto \quad \tilde{y}^{i+1} = \tilde{y}^i + hp(\tilde{y}^i)$$

Linear approximation at each step.



$$\begin{array}{l} \text{---} y(t) = \phi(y_0, 0, t) \\ \text{---} \phi(y_1 + e_1, 1, t) \end{array}$$

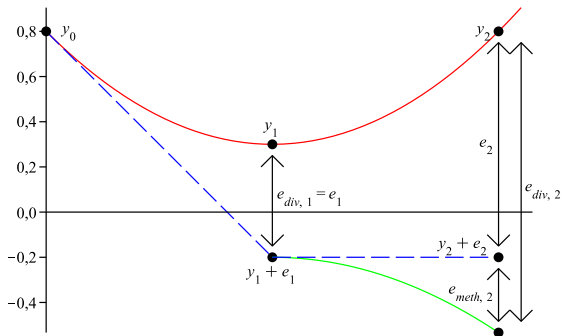
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Linear approximation at each step. Does **not work well** in practice.



$$\text{— red — } y(t) = \phi(y_0, 0, t) \quad \text{— green — } \phi(y_1 + e_1, 1, t)$$

# Taylor method

$$y(0) = 0 \quad y'(t) = p(y(t))$$

Time step  $h$ , discretize and compute  $\tilde{y}^i \approx y(ih)$ :

$$y(t+h) \approx y(t) + \sum_{i=1}^{\omega} h^i y^{(i)}(t) \quad \text{using } y^{(i)}(t) = \text{poly}_i(y(t))$$

Do a  $\omega$ -th order Taylor approximation at each step.

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Works well for  $\omega \geq 3$  but

- How to choose  $h$  and  $\omega$ ? **One more parameter to choose!**
- Error analysis is less obvious
- Complexity increases with  $\omega$

# Adaptive Taylor method

Adapt  $h$  and  $\omega$  at each step.

$$y(0) = 0 \quad y'(t) = p(y(t))$$

Time step  $h_i$ , discretize and compute  $\tilde{y}^i \approx y(\sum_{j \leq i} h_j)$ :

$$y(t + h_i) \approx y(t) + \sum_{i=1}^{\omega_i} h_i^i y^{(i)}(t) \quad \text{using } y^{(i)}(t) = \text{poly}_i(y(t))$$

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Do a  $\omega_i$ -th order Taylor approximation at each step.

Adapt the amount of computation to the hardness of the problem but

- Many more parameters to choose
- Error analysis is challenging
- Complexity analysis usually not done

# Adaptive Taylor method: parameter choice

How to choose the time steps  $h_i$  and orders  $\omega_j$ :

- $h_i$ : estimate the radius of convergence
- $\omega_j$ : try to guess the accuracy loss

Use voodoo magic and interval arithmetic to ensure correctness.

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**Our idea:** we need to choose  $h_i, \omega_i$  based on some high-level geometrical feature.

**Our algorithm in one sentence:** choose  $h_i, \omega_i$  so that

at each step, we increase the length of the solution by 1

# Interesting (practical ?) consequences

Compute  $y(t) \pm \varepsilon$

| Method         | Max. Order   | Number of steps   |
|----------------|--------------|---|
| Fixed $\omega$ | $\omega - 1$ | $\mathcal{O}\left(L^{\frac{\omega+1}{\omega-1}} \varepsilon^{-\frac{1}{\omega-1}}\right)$ |

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where 
$$L \approx \int_0^t \max(1, \|y'(u)\|) du$$

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| Smart ( $\omega = 1 + \log \frac{L}{\varepsilon}$ ) | $\log \frac{L}{\varepsilon}$ | $\mathcal{O}\left(L^{\sim 1}\right)$  |

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| Variable  | $\mathcal{O}\left(\log \frac{L}{\varepsilon}\right)$ | $\mathcal{O}(L)$  |

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# Conclusion

Solving Ordinary Differential Equations numerically:

- vastly different algorithms/results for vastly different expectations
- practical methods: no complexity
- nonuniform complexity: imprecise/misleading
- uniform worst-case complexity: everything is hard
- uniform parametrized complexity: encouraging

Questions:

- how far can we push parametrized complexity?
- can theory bring insight to practice?
- geometric complexity?

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$$y(t+h) \approx \sum_{j=0}^K \frac{h^j}{j!} y^{(j)}(t) \quad \rightsquigarrow \quad \tilde{y}^{i+1} = \sum_{j=0}^K \frac{h^j}{j!} P_k(\tilde{y}^i)$$

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What about  $h$  ?

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What about  $h$  ?

- **Fixed**  $h$ : wasteful
- **Adaptive**  $h$ : choose  $h$  depending on  $i, p, n$  and  $\tilde{y}^i$



# Choice of the parameters

Choice of  $h$  based on an effective lower bound on radius of convergence of the Taylor series:

**Lemma:** If  $y' = p(y)$ ,  $\alpha = \max(1, \|y_0\|)$ ,  $k = \deg(p)$ ,  $M = (k-1)\sum p\alpha^{k-1}$  then:

$$\|y^{(k)}(t) - P_k(y(t))\| \leq \frac{\alpha(Mt)^k}{1 - Mt}$$

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- local error  $\approx (Mt)^k \approx 2^{-k}$ : order gives the number of correct bits

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I spare you the analysis of the global error !

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This is impossible, right ?!

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## Example

$$\begin{cases} x(t) = t^{u(t)} \\ u(t) = e^{-t} - (1 - e^{-t}) \frac{1}{v(t)} \\ v(t) = v_0 \end{cases} \rightsquigarrow \begin{cases} x(t) \sim t^{\frac{1}{v_0}} \\ u(t) \rightarrow \frac{1}{v_0} \\ v(t) = v_0 \end{cases}$$

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## Remark

- All parameters are fixed except  $y_0 = (1, 1, v_0)$

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This is impossible, right ?!

## Example

$$\begin{cases} x(t) = t^{u(t)} \\ u(t) = e^{-t} - (1 - e^{-t}) \frac{1}{v(t)} \\ v(t) = v_0 \end{cases} \rightsquigarrow \begin{cases} x(t) \sim t^{\frac{1}{v_0}} \\ u(t) \rightarrow \frac{1}{v_0} \\ v(t) = v_0 \end{cases}$$

## Remark

- All parameters are fixed except  $y_0 = (1, 1, v_0)$
- Value are time  $t = 2$  can be arbitrary large for arbitrary small  $v_0$

# But wait...

This is impossible, right ?!

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## Theorem

There is no universal bound in  $\rho$ ,  $y_0$ ,  $t_0$ ,  $t$  and  $\mu$ .