
Learning Against Uncertainties

Mazen Alamir

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Disclaimer

- The following material **is not supposed to be fully understood!**
- Only its **diversity and wide scope** is to be taken home!
- The major part of the concepts and tools will be **shortly outdated!**
- ... Unless they already are!
- **Conjecture:** the corresponding state of mind **will prevail!**

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- 1 Data-driven partial observability assessment under uncertainties
- 2 Learning-Based Approximate Stochastic NMPC Design
- 3 Tractable stochastic NMPC by supervised clustering
- 4 Data-Driven NMPC by cost function identification
- 5 Learning-Based Monitoring updating period in Real-time NMPC

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What is it about?

(Parametric estimation example)

Article

Mathematical Model of COVID-19 Transmission Dynamics in South Korea: The Impacts of Travel Restrictions, Social Distancing, and Early Detection

Byul Nim Kim ^{1,†}, Eunjung Kim ^{2,†}, Sunmi Lee ^{3,*} and Chunyoung Oh ^{4,*}

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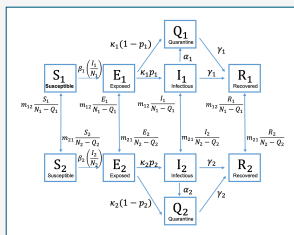
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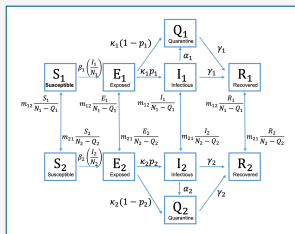
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$$\begin{aligned} \frac{dS_i}{dt} &= -\beta_i \left(\frac{I_i}{N_i} \right) S_i + m_{ji} \frac{S_j}{N_j - Q_j} - m_{ij} \frac{S_i}{N_i - Q_i}, \\ \frac{dE_i}{dt} &= \beta_i \left(\frac{I_i}{N_i} \right) S_i - \kappa_i E_i + m_{ji} \frac{E_j}{N_j - Q_j} - m_{ij} \frac{E_i}{N_i - Q_i}, \\ \frac{dI_i}{dt} &= \kappa_i p E_i - \alpha_i I_i - \gamma_i I_i + m_{ji} \frac{I_j}{N_j - Q_j} - m_{ij} \frac{I_i}{N_i - Q_i}, \\ \frac{dQ_i}{dt} &= \kappa_i (1 - p) E_i + \alpha_i I_i - \gamma_i Q_i, \\ \frac{dR_i}{dt} &= \gamma_i I_i + \gamma_i Q_i + m_{ji} \frac{R_j}{N_j - Q_j} - m_{ij} \frac{R_i}{N_i - Q_i}, \end{aligned}$$

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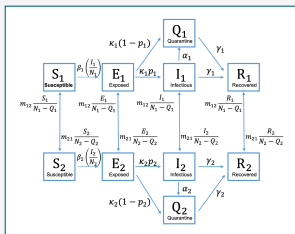
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Model's parameters

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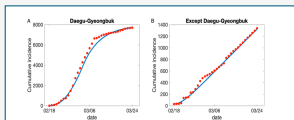
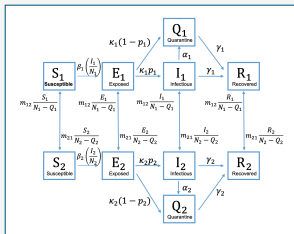


Figure 3. (A) Model calibration to Daegu-Gyeongbuk incidence data (blue solid line: model prediction, red dots data). (B) Model calibration to South Korea incidence data except Daegu-Gyeongbuk area.

Model calibration

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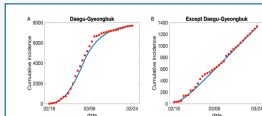
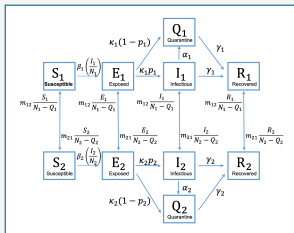


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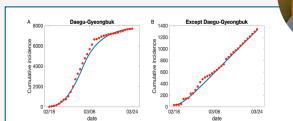
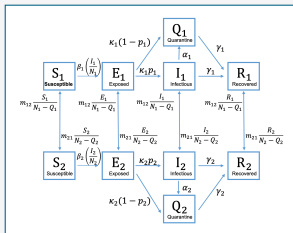


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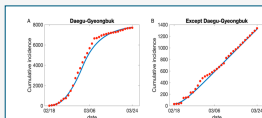
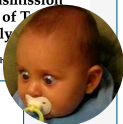
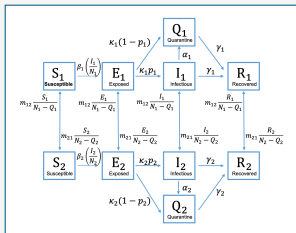


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Annual Reviews in Control

journal homepage: www.elsevier.com/locate/arcontrol

Full Length Article

The Ockham's razor applied to COVID-19 model fitting French data

Mirko Fiacchini ^{*}, Mazen Alami

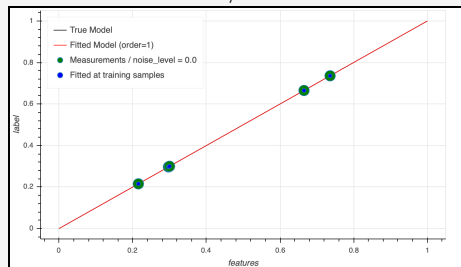
Univ. Grenoble Alpes, CNRS, Grenoble INP, GIPSA-lab, 38000 Grenoble, France

What is it about?

(A simple example)

- True model $y(t) = t$
- # of measurement $n_{obs} = 5$
- Fit polynomials of \neq degrees
- Training measurements
- Fitted values at training samples
- various noise levels

Noise=0 / Order=1

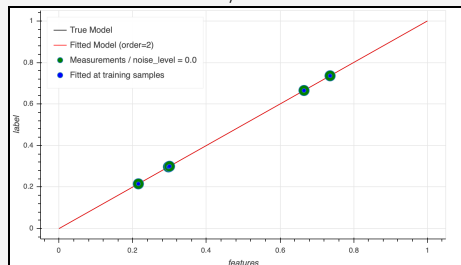


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Noise=0 / Order=2

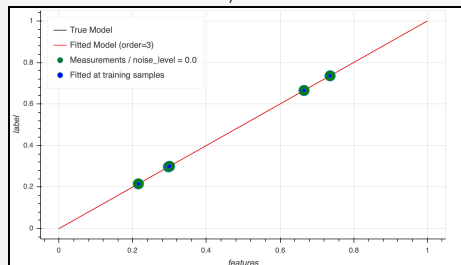


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Noise=0 / Order=3

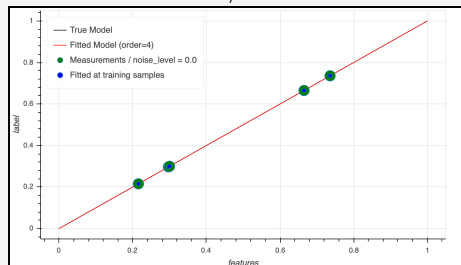


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Noise=0 / Order=4

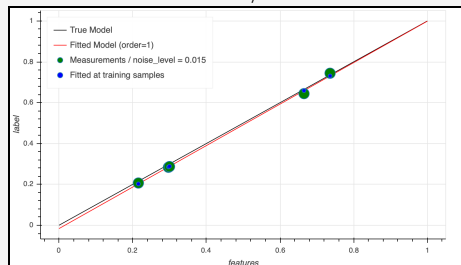


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Noise=0.015 / Order=1

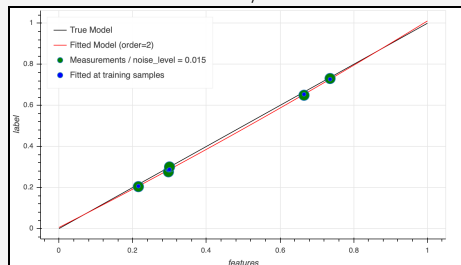


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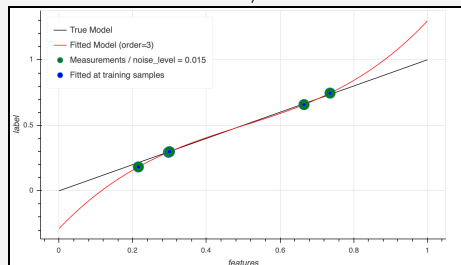


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- # of measurement $n_{obs} = 5$
- Fit polynomials of \neq degrees
- Training measurements
- Fitted values at training samples
- various noise levels

Noise=0.015 / Order=3

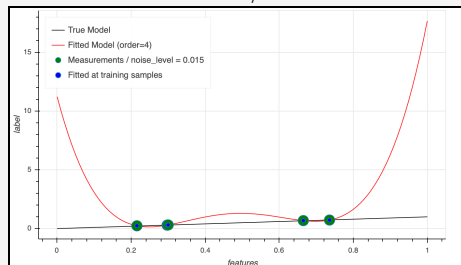


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Noise=0.015 / Order=4

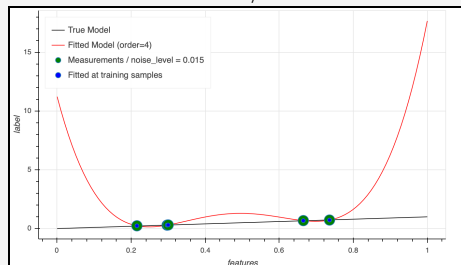


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This phenomenon is called **Over-Fitting**

An over fitted model is dangerous because it can lead to very high prediction errors on unseen realizations.

(This can be avoided by assessing the model by cross-validation techniques)

What is it about?

(Away from identification textbooks)

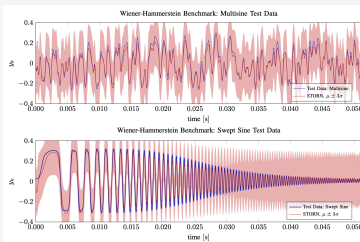
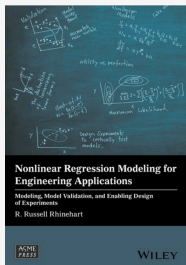


Fig. 12: Wiener Hammerstein benchmark: Time evolution for multisine and swept sine test data set of best results from Tab III, i.e. STORN with $n_{data} = 40$, $n_{fit} = 3$, $n_{noise} = 3$.

- In identification text-books, the previous situation does not prevail.
- The ratio between the samples and the number of parameters is generally quite high.
- Mechatronic systems, electrical systems, . . . etc.
- Identification is done once for all
- Initial state importance is marginal

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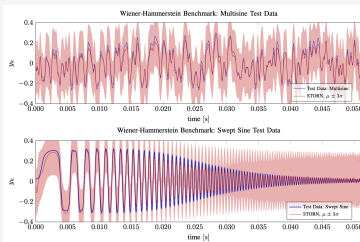
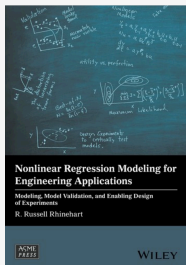


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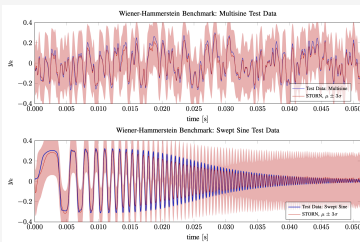
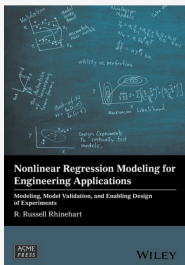


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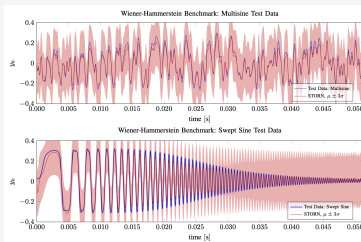
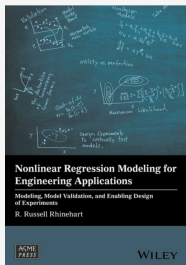


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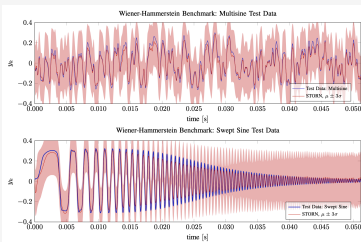
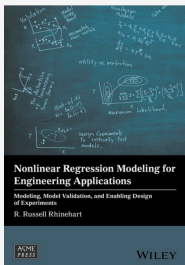


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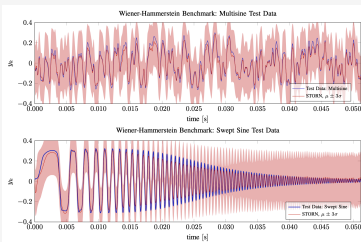
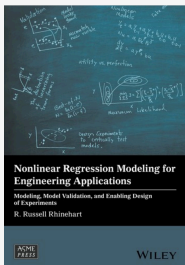


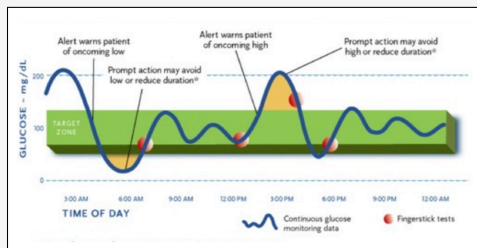
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(The diabetes example)

<https://quizlet.com>

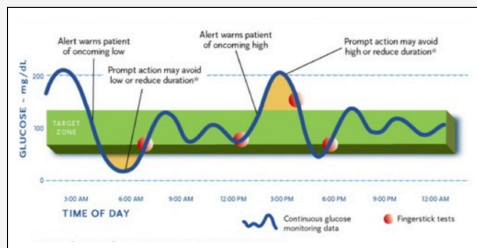


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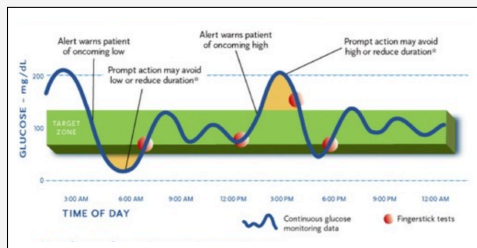


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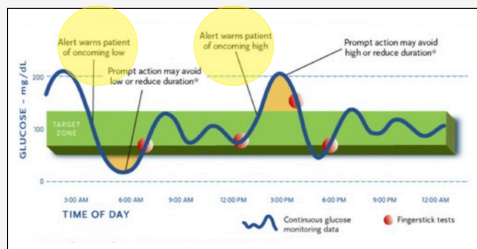


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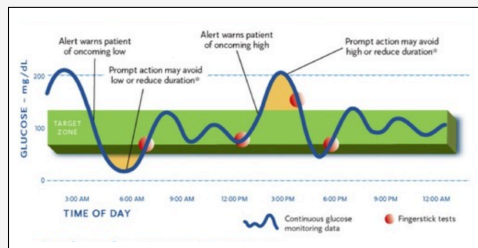


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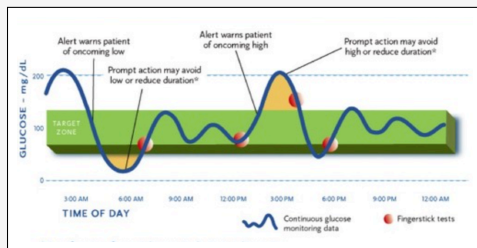


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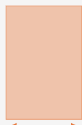
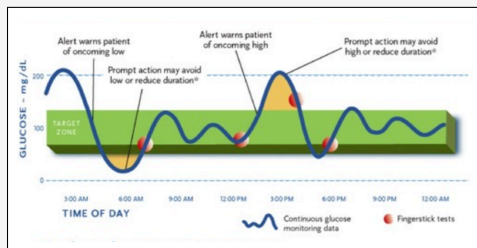


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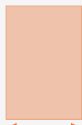
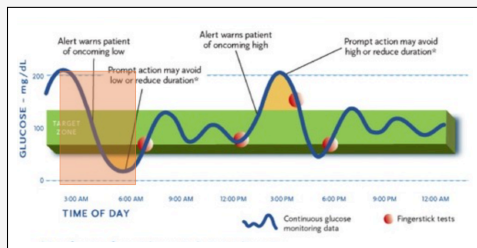
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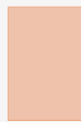
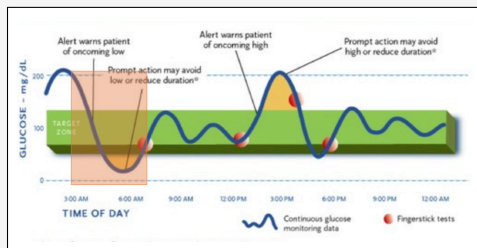
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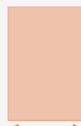
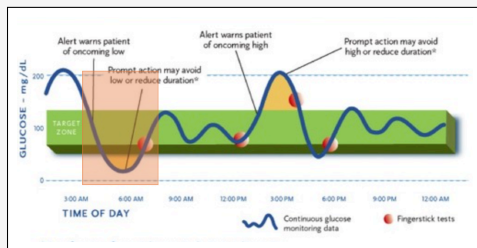
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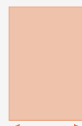
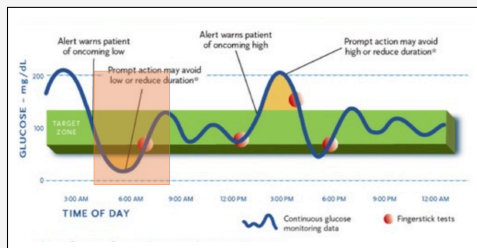
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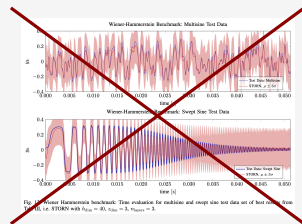


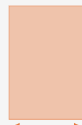
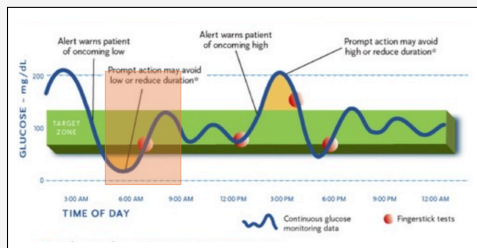
Fig. 10. Wison-Harmonica benchmark: Time evolution for multiple and single size test data set of test cases. (from [10], i.e. STOKO with $\lambda_{\text{min}} = 0$, $\lambda_{\text{max}} = 3$, $\mu_{\text{max}} = 3$)

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(The diabetes example)

https://quizlet.com



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$$\dot{G}(t) = -(S_g + X(t)) \cdot G(t) + S_g \cdot G_b + \frac{R_a(t)}{V_g} + \mathbf{w}(t)$$

$$\dot{X}(t) = -p_2 \cdot X(t) + p_2 \cdot S_1 [I(t) - I_b]$$

$$\dot{Q}_1(t) = -k_\tau \cdot Q_1(t) + m(t)$$

$$\dot{Q}_2(t) = -k_{abs} \cdot Q_2(t) + k_\tau \cdot Q_1(t)$$

$$\dot{I}_{sc1}(t) = -k_d \cdot I_{sc1}(t) + J_{ctrl}(t)$$

$$\dot{I}_{sc2}(t) = -k_d \cdot I_{sc2}(t) + k_d \cdot I_{sc1}(t)$$

$$\dot{I}_p(t) = -k_{cl} \cdot I_p(t) + k_d \cdot I_{sc2}(t)$$

$$R_a(t) = \frac{k_{abs} \cdot f}{BW} \cdot Q_2(t)$$

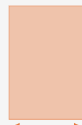
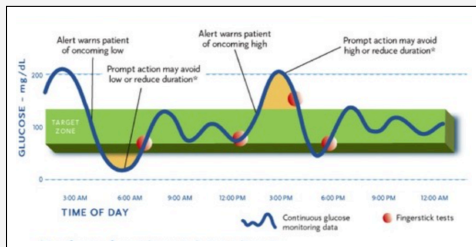
$$I(t) = \frac{I_p}{V_1 \cdot BW}$$

Symbol	Meaning	Value
S_g	Fractional glucose effectiveness	0.01
V_g	Distribution volume of glucose	1.6
k_{abs}	Rate constant—oral glucose consumption	0.01193
k_τ	Time constant related with oral glucose absorption	0.08930
p_2	Rate constant of the remote insulin compartment	0.02
f	Fraction of intestinal absorption	0.9
V_1	Distribution volume of insulin	0.06005
k_d	Rate constant of subcutaneous insulin transport	0.16
k_τ	Rate constant of subcutaneous insulin transport	0.02
S_1	Insulin sensitivity	0.0006
BW	Body weight	Known
G_b	Basal glucose concentration	Eq. (2)
I_b	Reference value for $I(t)$, associated with the fasting plasma glucose concentration.	Steady-state

What is it about?

(The diabetes example)

https://quizlet.com



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Few measurement
Many parameters

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x_1 tumor cell population;
 x_2 circulating lymphocytes population;
 x_3 chemotherapy drug concentration;
 x_4 effector immune cell population;
 u_1 rate of introduction of immunotherapy drug;
 u_2 rate of introduction of chemotherapy drug.

and the dynamics is given by

$$\dot{x}_1 = ax_1(1 - bx_1) - c_1x_4x_1 - k_3x_3x_1$$

$$\dot{x}_2 = -\delta x_2 - k_2x_3x_2 + s_2$$

$$\dot{x}_3 = -\gamma_0x_3 + u_2$$

$$\dot{x}_4 = g \frac{x_1}{h + x_1} x_4 - rx_4 - p_0x_4x_1 - k_1x_4x_3 + s_1u_1.$$

A constraint such as $x_4 \geq \text{Threshold}$ cannot be addressed via data-driven solutions that do not involve measurement of x_4 unless a model and associated state reconstruction algorithm are used.

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Question: Am I in the right room?

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 x_4 effector immune cell population;
 u_1 rate of introduction of immunotherapy drug;
 u_2 rate of introduction of chemotherapy drug.

and the dynamics is given by

$$\dot{x}_1 = ax_1(1 - bx_1) - c_1x_4x_1 - k_3x_3x_1$$

$$\dot{x}_2 = -\delta x_2 - k_2x_3x_2 + s_2$$

$$\dot{x}_3 = -\gamma_0x_3 + u_2$$

$$\dot{x}_4 = g \frac{x_1}{h + x_1} x_4 - rx_4 - p_0x_4x_1 - k_1x_4x_3 + s_1u_1.$$

A constraint such as $x_4 \geq \text{Threshold}$ cannot be addressed via data-driven solutions that do not involve measurement of x_4 unless a model and associated state reconstruction algorithm are used.

Question: Am I in the right room?

Answer: **Definitively!**



What is precisely a data-driven solution?

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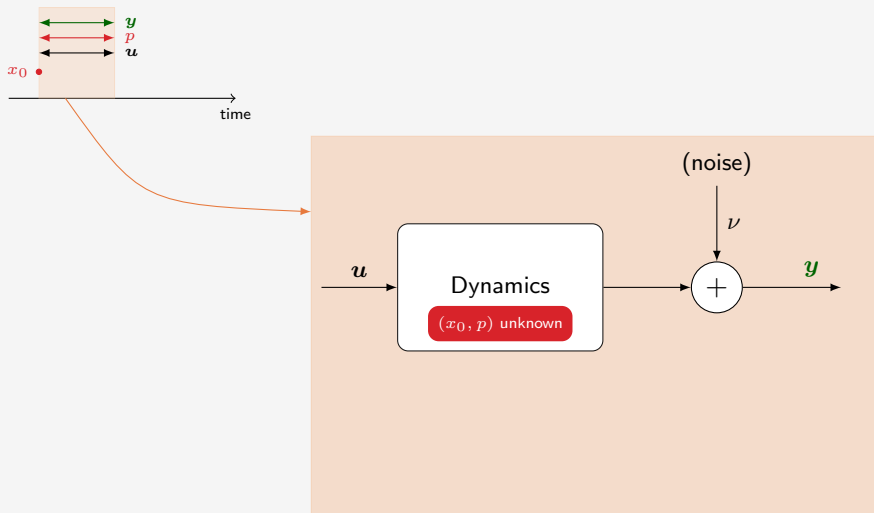


The mathematical model alone and all the related derivations, assumptions and computations **are insufficient** to come out with quantifiable results and or appropriate choices.

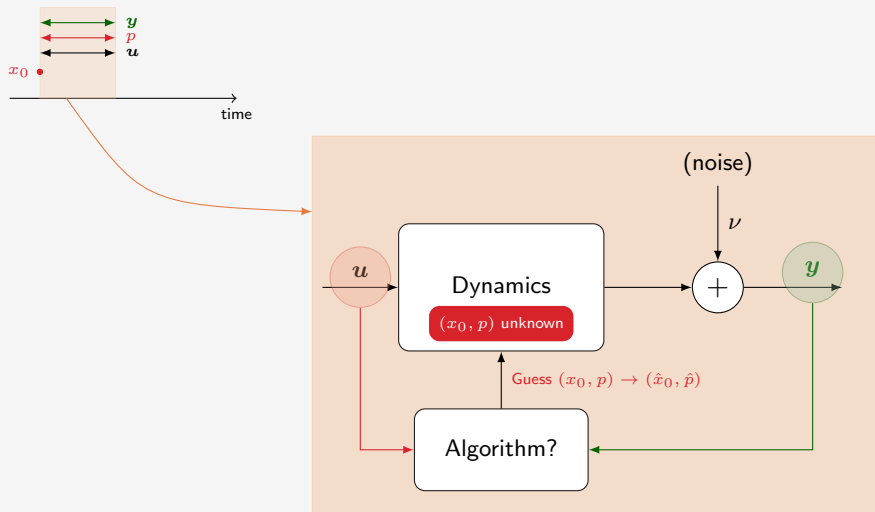
The reality should be partially discovered by a rich set of realizations of the involved uncertain quantities!

A step where a **data-set** is generated/collected for later use in the design/certification process.

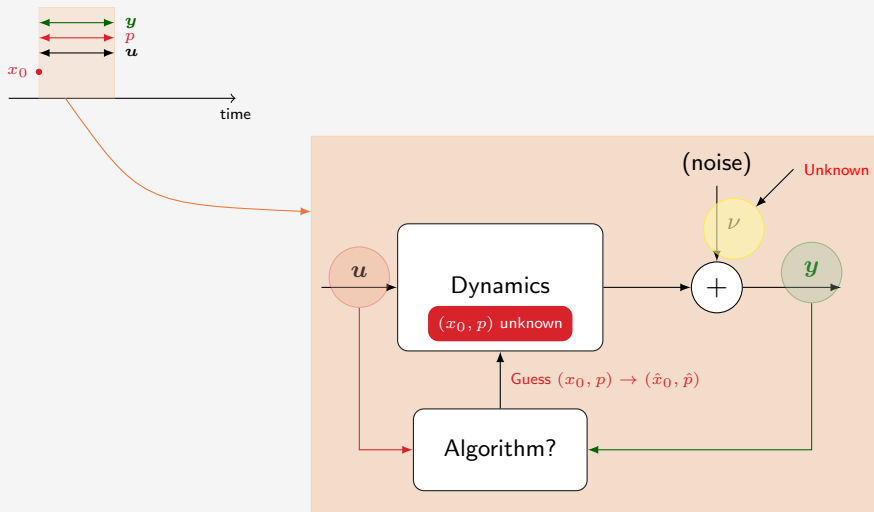
The dynamic extended estimation problem



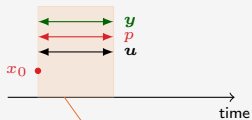
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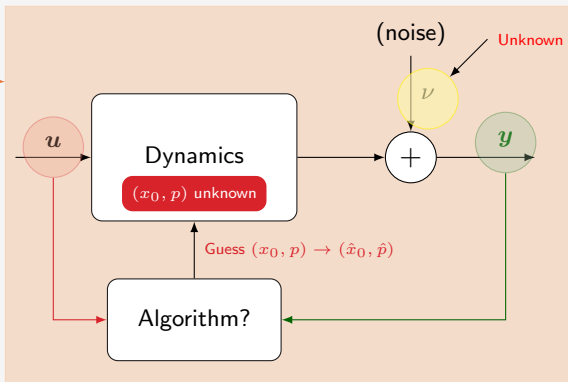
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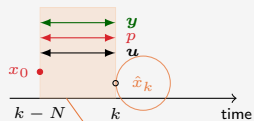
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This has to be repeatedly done on-line at each sampling period



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Given the estimation at instant k
(\hat{x}_0, \hat{p})

The current state is estimated by

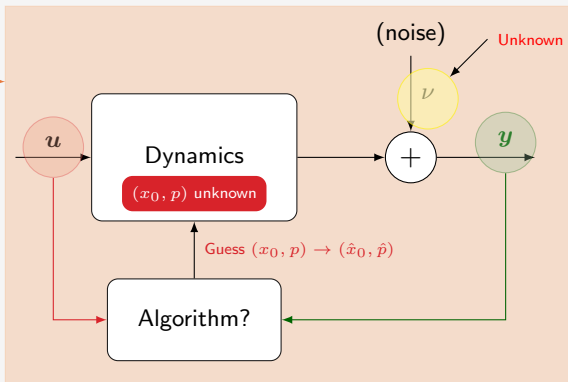
$$\hat{x}_k = X_N(\hat{x}_0, \hat{p}, \mathbf{u})$$

More generally, the value of any quantity:

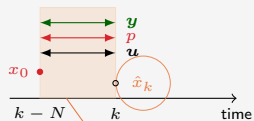
$$z = T(x, p)$$

Can be estimated by:

$$\hat{z}_k = T\left(X_N(\hat{x}_0, \hat{p}, \mathbf{u}), \hat{p}\right)$$



The dynamic extended estimation problem



$z = T(x, p)$ is called hereafter, the **estimation target**.

Example. A feedback law $K(x, p)$ can be viewed as an estimation target in an observer-based control design.

Given the estimation at instant k
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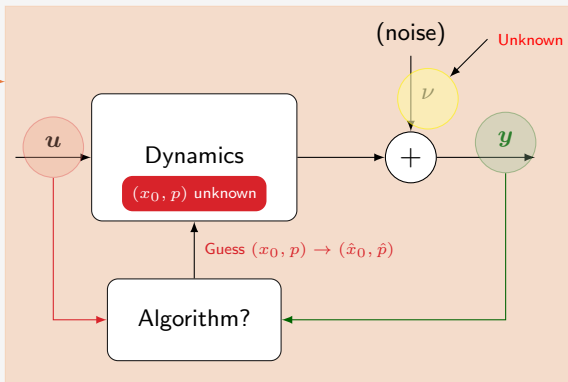
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Problem statement

Given

1. An uncertain model $\dot{x} = f(x, u, p)$
2. Statistics of uncertainties \mathcal{P} on \mathbb{P}
3. Statistics of states \mathcal{X} on \mathbb{X}
4. Statistics of measurement noise \mathcal{V} on \mathbb{V}
5. A given observation target $z = T(x, p)$

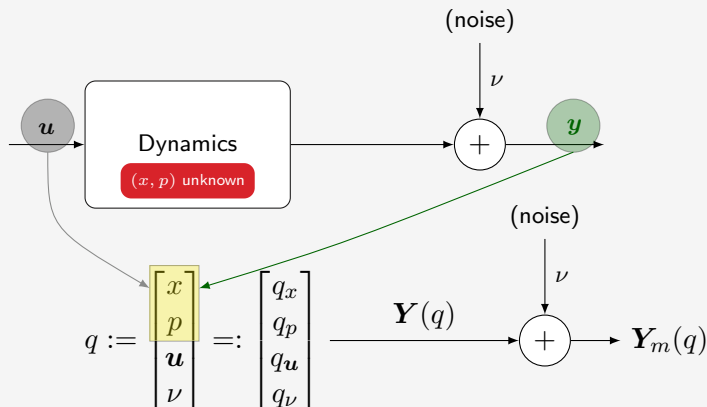
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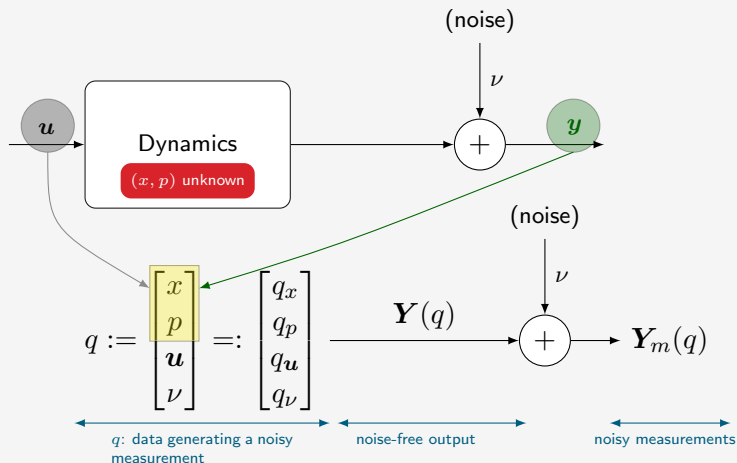
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Derive a certified bound on the excursion of the estimation error on the observation target z together with guidelines towards its minimization.

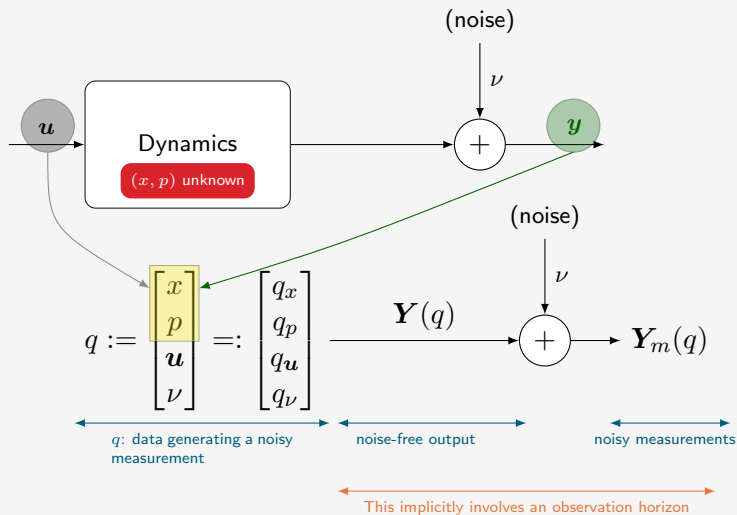
Notation: The scenarios vector q



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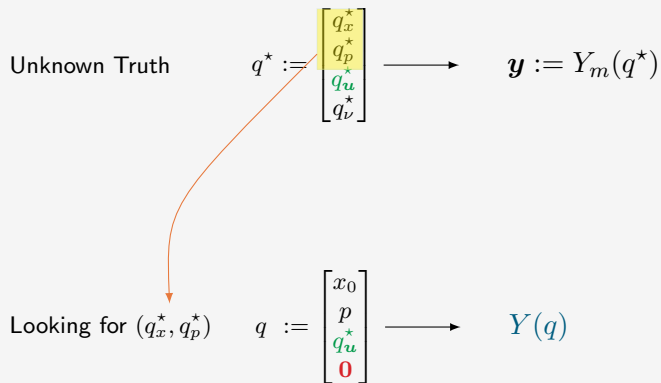


The cost function (Distance between scenarios)

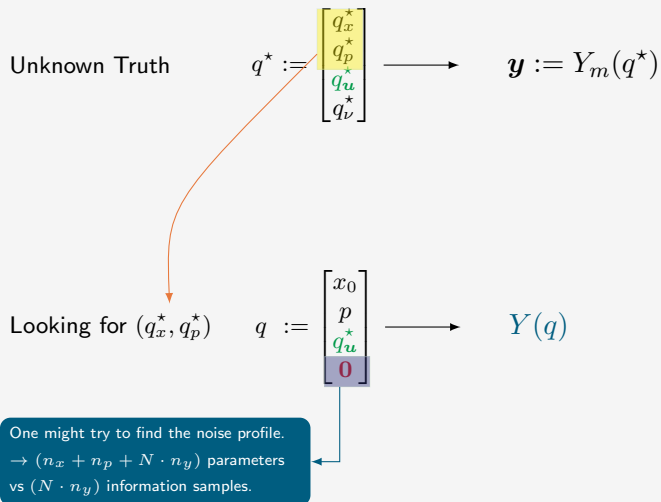
Unknown Truth

$$q^* := \begin{bmatrix} q_x^* \\ q_p^* \\ q_u^* \\ q_v^* \end{bmatrix} \longrightarrow \mathbf{y} := Y_m(q^*)$$

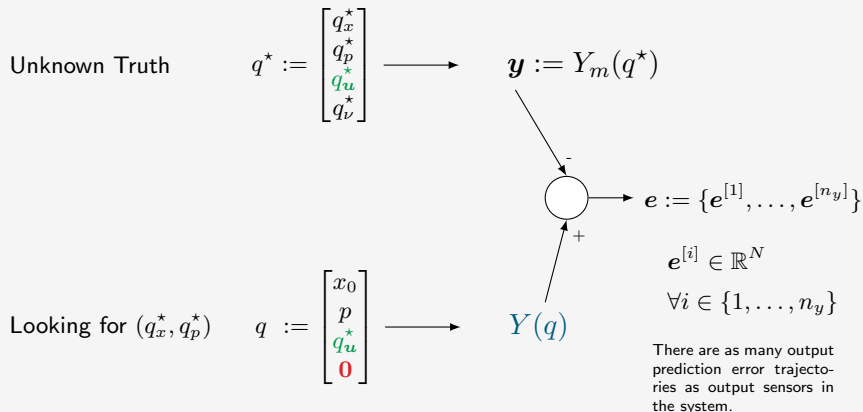
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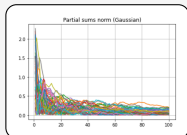
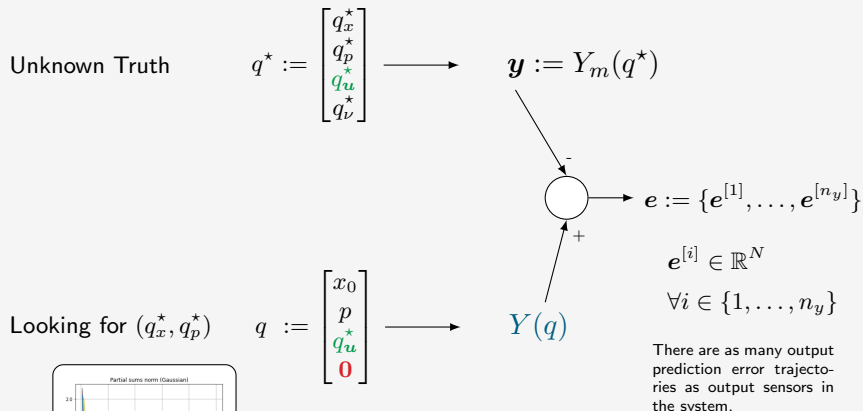
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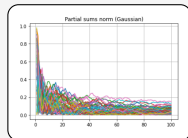
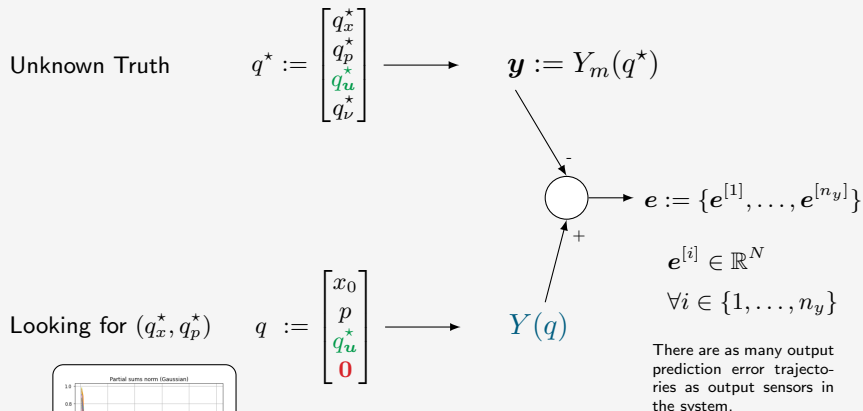


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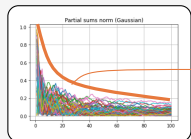
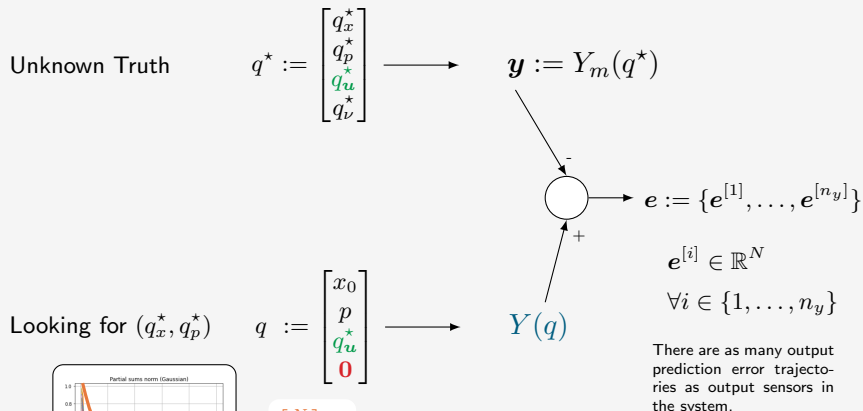
abs(Partial average) of a unitary gaussian noise

The cost function (Distance between scenarios)



abs(Partial average) of a unitary uniform noise

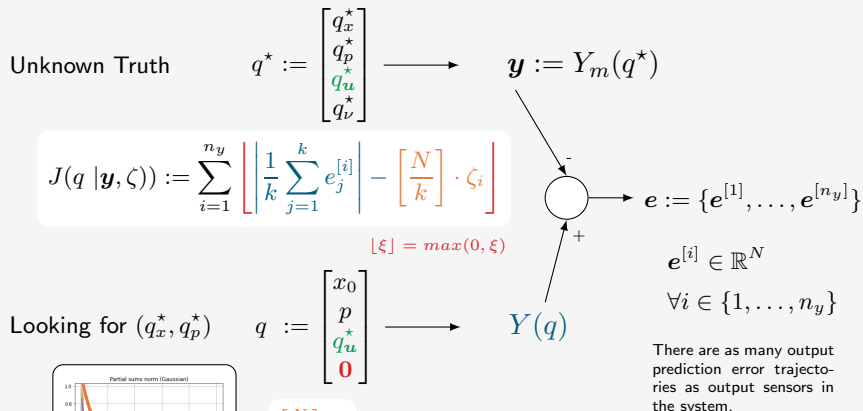
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$$\left[\frac{N}{k} \right] \cdot \zeta$$

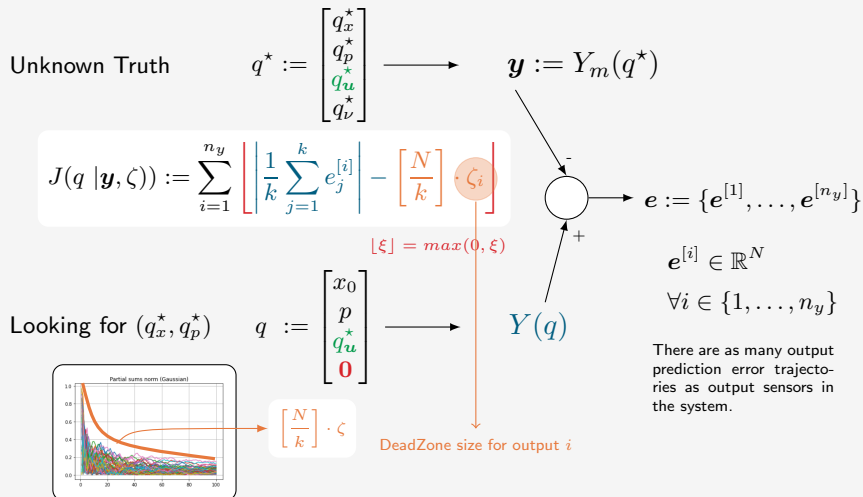
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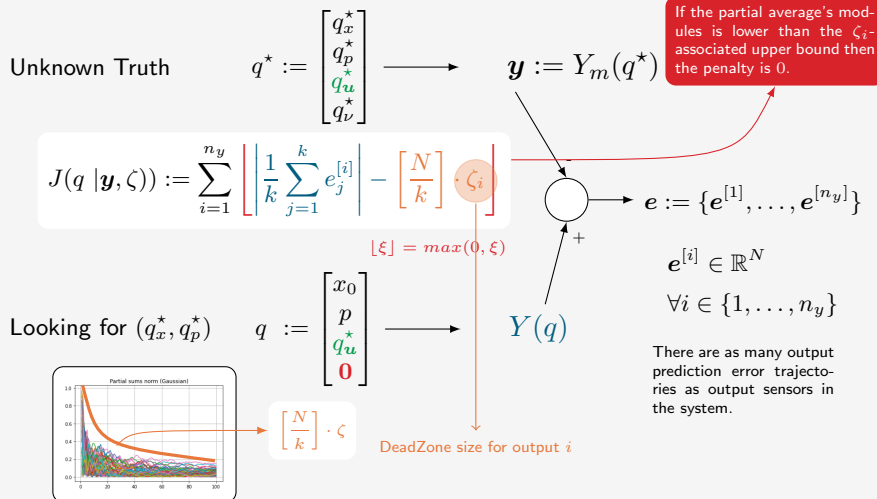
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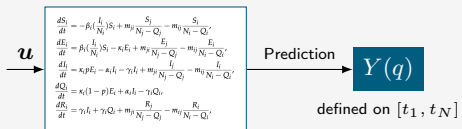
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Estimation as optimization problem (MHE: Moving-Horizon Estimator)

Mathematical model

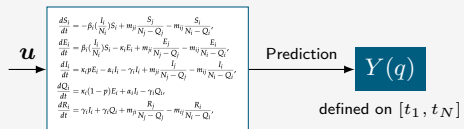


The prediction of the expected measurement should the initial state be x_0 and the true parameter be p under the control profile \mathbf{u} .

Estimation as optimization problem (MHE: Moving-Horizon Estimator)

The prediction of the expected measurement should the initial state be x_0 and the true parameter be p under the control profile u .

Mathematical model



The prediction error associated to q

$$\mathcal{E}(q | \mathbf{y}) := d(Y(q), \mathbf{y})$$

$$\begin{bmatrix} y(t_1) \\ y(t_2) \\ \vdots \\ y(t_N) \end{bmatrix}$$

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$$\begin{aligned} \frac{dS_i}{dt} &= -\beta_i \left(\frac{I_i}{N_i} \right) S_i + m_{ij} \frac{S_j}{N_j - Q_j} - m_{ji} \frac{S_i}{N_i - Q_i}, \\ \frac{dE_i}{dt} &= \beta_i \left(\frac{I_i}{N_i} \right) S_i - \kappa_i E_i + m_{ij} \frac{E_j}{N_j - Q_j} - m_{ji} \frac{E_i}{N_i - Q_i}, \\ \frac{dI_i}{dt} &= \kappa_i p E_i - a_i I_i - \gamma_i I_i + m_{ij} \frac{I_j}{N_j - Q_j} - m_{ji} \frac{I_i}{N_i - Q_i}, \\ \frac{dQ_i}{dt} &= \kappa_i (1-p) E_i + a_i I_i - \gamma_i Q_i, \\ \frac{dR_i}{dt} &= \gamma_i I_i + \gamma_i Q_i + m_{ij} \frac{R_j}{N_j - Q_j} - m_{ji} \frac{R_i}{N_i - Q_i}. \end{aligned}$$

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$$\text{ex. } d(v_1, v_2) = \|v_1 - v_2\|^2$$

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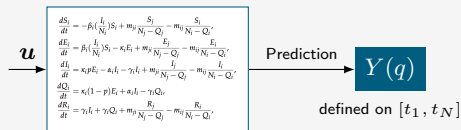
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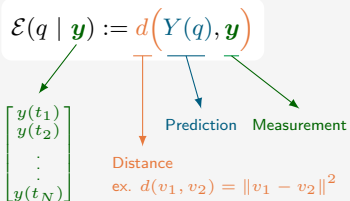
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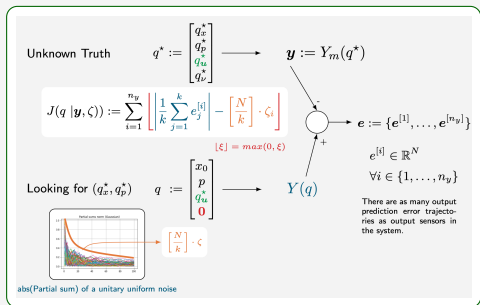
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How to tune the dead-zone size vector $\zeta \in \mathbb{R}_+^{n_y}$?

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Dead-Zone consistency condition

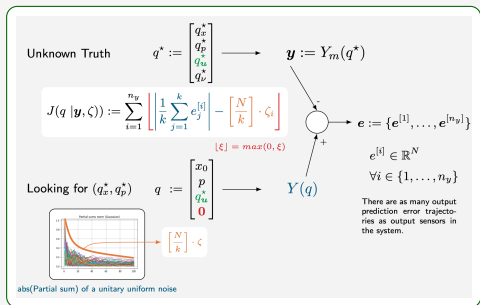


Consistency

A dead-zone size vector ζ is **consistent** (w.r.t the noise) if:

Any candidate scenario q that meets the true value of the triplet (q_x^*, q_p^*, q_u^*) should lead to **zero cost** for any possible realization of the noise.

Dead-Zone consistency condition



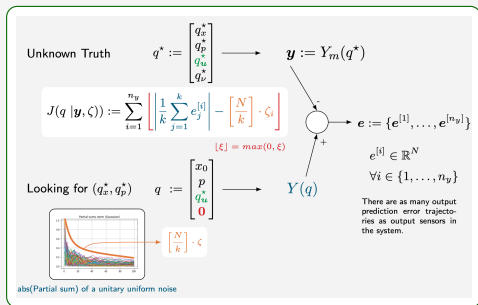
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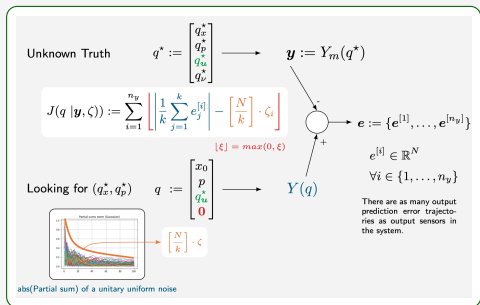
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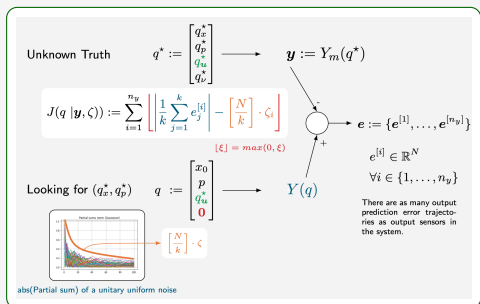
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Note that this can always be made true by taking sufficiently high values of ζ !

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Note that this can always be made true by taking sufficiently high values of ζ !

But this reduces the quality of the estimation!

ϵ -observability of an observation target $q^{(1)}$ $q^{(2)}$

$$\begin{bmatrix} x_0^{(1)} \\ p^{(1)} \\ \mathbf{u} \\ \mathbf{0} \end{bmatrix} \quad \begin{bmatrix} x_0^{(2)} \\ p^{(2)} \\ \mathbf{u} \\ \boldsymbol{\nu} \end{bmatrix}$$

Notation: $q^{(1)} \bowtie q^{(2)}$ (comparable pair)

Two scenarios that share the same input profile while one is free of noise.

ϵ -observability of an observation target

$$\begin{array}{cc}
 q^{(1)} & q^{(2)} \\
 \left[\begin{array}{c} x_0^{(1)} \\ p^{(1)} \\ \mathbf{u} \\ \mathbf{0} \end{array} \right] & \left[\begin{array}{c} x_0^{(2)} \\ p^{(2)} \\ \mathbf{u} \\ \nu \end{array} \right]
 \end{array}
 \xrightarrow{\text{If}}
 J(q^{(1)} \mid Y_m(q^{(2)}), \zeta) = 0$$

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Then: $q^{(1)} \equiv q^{(2)}$ (indistinguishable pair)

Two scenarios that share the same input profile while one is free of noise.

ϵ -observability of an observation target

$$\begin{array}{cc}
 q^{(1)} & q^{(2)} \\
 \left[\begin{array}{c} x_0^{(1)} \\ p^{(1)} \\ \mathbf{u} \\ \mathbf{0} \end{array} \right] & \left[\begin{array}{c} x_0^{(2)} \\ p^{(2)} \\ \mathbf{u} \\ \nu \end{array} \right]
 \end{array}
 \xrightarrow{\text{If}}
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Given an observation target $z = T(x, p)$, z is said to be ϵ -observable on \mathcal{Q} iff:

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- 2 $\forall (q^{(1)}, q^{(2)}) \in \mathcal{Q}^2$:

$$(q^{(1)} \bowtie q^{(2)}) \text{ AND } (q^{(1)} \equiv q^{(2)}) \Rightarrow \|q_z^{(1)} - q_z^{(2)}\| \leq \epsilon$$

where $q_z := T(q_x, q_p)$. In other words, only pairs with ϵ -distant observable targets can be both comparable and indistinguishable.

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 \text{Consistency on } \mathbb{Q} \\
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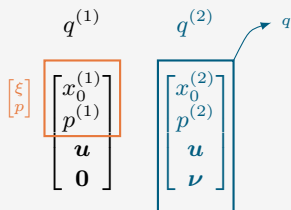
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If

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Then: $q^{(1)} \equiv q^{(2)}$ (indistinguishable pair)

\Rightarrow The ϵ -observability can be investigated by inspecting realizations of the form:

$$w = \begin{bmatrix} q \\ \xi \\ p \end{bmatrix} \in \mathbb{Q} \times \mathbb{X} \times \mathbb{P}$$

for which a constraint of the form:

$$g(\theta, w) = 0 \quad \text{where } \theta := (\epsilon, \zeta)$$

should be satisfied

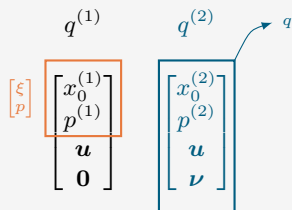
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$$A(w) \Rightarrow B(w) \Leftrightarrow 0 = f(w) = \begin{cases} 0 & \text{if } (\bar{A}) \cup (A \cap B) \\ 1 & \text{otherwise} \end{cases}$$

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Partial ϵ -observability: a robust constraint satisfaction

Given:

- A dynamic model with uncertain parameter & noisy measurements
- An observation target $z = T(x, p)$
- A dead-zone size vector $\zeta \in \mathbb{R}_+^{n_y}$
- A desired precision $\epsilon > 0$
- A set of interest

$$\mathbb{W} := \underbrace{\mathbb{X} \times \mathbb{P} \times \mathbb{U} \times \mathbb{V}}_{\mathbb{Q}} \times \mathbb{X} \times \mathbb{P}$$

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$$\forall w \in \mathbb{W}, \quad g(\theta, w) = 0$$

then, solving:

$$(x_0^*, p^*) \leftarrow \min_{x_0, p} J \left(\begin{pmatrix} x_0 \\ p \\ \mathbf{u} \\ 0 \end{pmatrix} \mid Y_m, \zeta \right)$$

leads to an estimation error on the observation target z that never exceeds ϵ .

and this holds true for any possible realizations inside the sets \mathbb{X} , \mathbb{P} , \mathbb{U} and \mathbb{V}

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Optimal design

$$\min_{\theta=(\epsilon, \zeta)} \theta_1 \mid g(\theta, w) = 0 \quad \forall w \in \mathbb{W}$$

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Optimal design

Intractable
over-pessimistic

$$\min_{\theta=(\epsilon, \zeta)} \theta_1 \mid g(\theta, w) = 0 \quad \forall w \in \mathbb{W}$$

Recalls on probabilistic certification

Probabilistic Certification

A general and **tractable** framework to face high uncertainties and still have something to **guarantee** with an **EXPLICIT** (hopefully high) **probability**.

Probabilistic certification as a relaxed formulation

Optimal Robust constraints satisfaction problem

$$\min_{\theta \in \Theta} J(\theta) \quad \text{s.t.} \quad (\forall w \in \mathbb{W}) \quad g(\theta, w) \leq 0$$

Probabilistic certification as a relaxed formulation

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- 1 Extremely **hard** to solve
- 2 Uselessly **conservative**
 - worst case analysis
 - even extremely rare bad configurations are accommodated for

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Relaxed formulations \Rightarrow **Probabilistic certification**

Probabilistic certification as a relaxed formulation

Define the constraints violation indicator:

$$I(\theta, w) := \begin{cases} 0 & \text{if } g(\theta, w) \leq 0 \\ 1 & \text{otherwise (constraints are violated)} \end{cases}$$

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Robust constraint

$$(\forall w \in \mathbb{W}) \quad g(\theta, w) \leq 0$$

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First Relaxation (R_1)

$$\Pr\{I(\theta, w) = 1\} \leq \eta$$

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Second Relaxation ($R_2(N)$)

$$\frac{1}{N} \sum_{i=1}^N [I(\theta, w^{(i)})] \leq \frac{m}{N} \leq \eta$$

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For any given m , N must be sufficiently high to get

- ✓ $N \geq \frac{m}{\eta}$
- ✓ $\Pr\{R_2(N) \cap (!R_1)\} \leq \delta$

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What value for N ?

Probabilistic certification as a relaxed formulation

When the admissible set Θ is discrete with cardinality n_{Θ}

$$N \geq \frac{1}{\eta} \left(m + \ln\left(\frac{n_{\Theta}}{\delta}\right) + \left(2m \ln\left(\frac{n_{\Theta}}{\delta}\right)\right)^{1/2} \right)$$

more formulas are available (Alamo et al. IEEE-TAC 2009)

- m positive integer ($m = 1, 5, 10$ etc.)
- n_{Θ} $\text{card}(\Theta)$
- η Precision
- δ Confidence parameter

Values of N for $\delta = 0.01$, $m = 1$

n_{Θ}	$\eta = 0.1$	$\eta = 0.05$	$\eta = 0.01$	$\eta = 0.001$
1	132	264	1317	13164
5	154	308	1536	15354
10	163	326	1628	16280
100	193	386	1930	19299
1000	223	445	2225	22249
10000	252	503	2515	25148

NOTA

- ✓ $\dim(w)$ does not matter !!!!
- ✓ $(n_{\Theta}, \delta) \rightarrow$ logarithmically !!

Alamo et al. Randomized strategies for probabilistic solutions of uncertain feasibility and ... IEEE TAC, 2009.

Algorithm & complexity analysis

Optimal Design

$$\min_{\theta := (\epsilon, \zeta)} \epsilon \text{ under } \sum_{i=1}^N I(\theta, w^{(i)}) \leq m$$

where

- $N = N(\delta, \eta, m)$
- I involves the definition of g
- g involves the cost function J

$$J(q | \mathbf{y}, \zeta) := \sum_{i=1}^{n_y} \left[\left| \frac{1}{k} \sum_{j=1}^k e_j^{[i]} \right| - \left[\frac{N}{k} \right] \cdot \zeta_i \right]$$

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Algorithm 1 The certification algorithm

1: **Given:**

- The maps f and h ▷ [see (1) and (2)]
- η, δ, m ▷ Certification parameters [see (41)]
- $\mathbb{X}, \mathbb{P}, \mathbb{U}, \mathbb{V}$ ▷ Working sets
- $\underline{\sigma}_\epsilon, \bar{\sigma}_\epsilon, \underline{\sigma}_\zeta, \bar{\sigma}_\zeta, n_\epsilon, n_\zeta$ ▷ Design Set parameters (46)
- Sampling rule \mathcal{W} on $\mathbb{W} = \mathbb{Q} \times \mathbb{X} \times \mathbb{P}$

2: $n_\Theta \leftarrow n_\epsilon \cdot n_\zeta$ ▷ Cardinality of the design set Θ

3: Compute $\Theta := \{\theta^{(j)}\}_{j=1}^{n_\Theta}$ by (46) in alphabetic order

4: $N_s \leftarrow N_s(\delta, \eta, m, n_\Theta)$ by (41) ▷ Number of scenarios

5: Generate N_s scenarios inputs $\{w^{(i)}\}_{i=1}^{N_s}$ using \mathcal{W}

6: **for** $i \in \{1, \dots, N_s\}$ **do**

7: Compute $e^{(i)}$ ▷ non clipped error for $w^{(i)}$

8: **end for**

9: success \leftarrow False ▷ No solution θ^* found yet

10: **for** $j \in \{1, \dots, n_\Theta\}$ **do**

11: nb_failures $\leftarrow \sum_{i=1}^{N_s} g(w^{(i)}, \theta^{(j)})$

12: **if** nb_failures $\leq m$ **then**

13: success \leftarrow True ▷ Optimal solution found

14: $\theta^* \leftarrow \theta^{(j)}$. **Break.** ▷ Stop the loop

15: **end if**

16: **end for**

17: **Output:** **If** success **return** θ^*

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$$\Theta := \mathbb{L}(\underline{\sigma}_\epsilon, \bar{\sigma}_\epsilon, n_\epsilon) \times \mathbb{L}(\underline{\sigma}_\zeta, \bar{\sigma}_\zeta, n_\zeta) \quad (42)$$

where $\mathbb{L}(\underline{\sigma}, \bar{\sigma}, n)$ is the set of n logarithmically uniformly spaced numbers, namely¹:

$$\mathbb{L}(\underline{\sigma}, \bar{\sigma}, n) = \text{logspace}(\underline{\sigma}, \bar{\sigma}, n)$$

$$:= \left\{ 10^{r_i} \mid r_i = \underline{\sigma} + \frac{(\bar{\sigma} - \underline{\sigma})i}{n-1} \quad i \in \{0, \dots, n-1\} \right\}$$

This obviously leads to a cardinality $n_\Theta = n_\epsilon n_\zeta$

Algorithm & complexity analysis

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11: nb_failures $\leftarrow \sum_{i=1}^{N_s} g(w^{(i)}, \theta^{(j)})$

12: **if** nb_failures $\leq m$ **then**

13: success \leftarrow True ▷ Optimal solution found

14: $\theta^* \leftarrow \theta^{(j)}$. **Break.** ▷ Stop the loop

15: **end if**

16: **end for**

17: **Output:** **If** success **return** θ^*

$$\Theta := \mathbb{L}(\underline{\sigma}_\epsilon, \bar{\sigma}_\epsilon, n_\epsilon) \times \mathbb{L}(\underline{\sigma}_\zeta, \bar{\sigma}_\zeta, n_\zeta) \quad (42)$$

where $\mathbb{L}(\underline{\sigma}, \bar{\sigma}, n)$ is the set of n logarithmically uniformly spaced numbers, namely¹:

$$\mathbb{L}(\underline{\sigma}, \bar{\sigma}, n) = \text{logspace}(\underline{\sigma}, \bar{\sigma}, n)$$

$$:= \left\{ 10^{r_i} \mid r_i = \underline{\sigma} + \frac{(\bar{\sigma} - \underline{\sigma})i}{n-1} \quad i \in \{0, \dots, n-1\} \right\}$$

This obviously leads to a cardinality $n_\Theta = n_\epsilon n_\zeta$

Application example

Parallel reactor $R \rightarrow (P_1, P_2)$

$$\dot{x}_1 = 1 - w_1 x_1^2 e^{-1/x_3} - w_2 e^{-w_3/x_3} - x_1$$

$$\dot{x}_2 = w_1 x_1 - 1^2 e^{-1/x_3} - x_2$$

$$\dot{x}_3 = u - x_3$$

- x_1, x_2 concentrations of R and P_1
- x_3 temperature of the mixture
- u is the jacket temperature
- x_2 and u are measured
- Three possible definitions of z :
 - $z = x$
 - $z = x_1$
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Investigated set/statistics options

⊗ A uniform distribution over a hyperbox \mathbb{X}

ℙ Two possibilities are investigated, namely:

- 1 A uniform distribution over \mathbb{P}
- 2 A Gaussian distribution around a nominal value.

\mathbb{U}^N Two possibilities are investigated, namely:

- 1 A uniform distribution over \mathbb{U}^N
- 2 A random choice consisting in sequences (elements of \mathbb{U}) of the form $(i \in \{1, \dots, N\})$:

$$u_i = \text{Sat}_{\mathbb{U}} \left[\sum_{j=1}^{nf} \beta_j \sin \left(\frac{2j\pi(i\tau)}{N\tau} + \varphi_j \right) \right]$$

\mathbb{V}^N Gaussian white noise (with different variances) are used to represent measurement noise.

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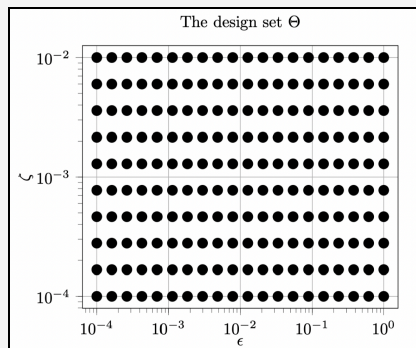
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The design set



The impact of players on partial observability

- Impact of the **observation horizon**
 - Impact of the **measurement noise**
 - Impact of the **uncertainty level**
 - Impact of the **input type**
 - impact of the **certification parameters**
-
- **Computation time**



Impact of the observation horizon

eps1	eps2	eps3	zeta1	zeta2	zeta3	N	noise	rho	std_p	u_mod	p_mod	eta	delta	m
0.379269	0.379269	0.054556	0.000774	0.000774	0.000774	5	0.001000	0.050000	-	Fourier	uniform	0.010000	0.001000	10
0.233572	0.143845	0.054556	0.000774	0.000774	0.000774	10	0.001000	0.050000	-	Fourier	uniform	0.010000	0.001000	10
0.000100	0.000100	0.000100	0.000464	0.000464	0.000464	20	0.001000	0.050000	-	Fourier	uniform	0.010000	0.001000	10

$$z_1 = x, z_2 = x_1, z_3 = x_3$$

Note that $N = 20$ is needed in order to achieve the certification with the lowest values $\epsilon = 10^{-4}$ considered in the design set Θ over the three observation-target variables z_i , $i = 1, 2, 3$.

Otherwise, **indistinguishability might occur** with quite high error values on the targeted indicators.

Note that only three values of N are studied here, lower values of $N \in [10, 20]$ would have probably be sufficient to achieve the high precision certification results.

Impact of the measurement noise

eps1	eps2	eps3	zeta1	zeta2	zeta3	N	noise	rho	std_p	u_mod	p_mod	eta	delta	m
0.000100	0.000100	0.000100	0.000464	0.000464	0.000464	20	0.001000	0.050000	-	Fourier	uniform	0.010000	0.001000	10
0.379269	0.379269	0.088587	0.001292	0.001292	0.001292	20	0.003000	0.050000	-	Fourier	uniform	0.010000	0.001000	10
0.233572	0.143845	0.054556	0.000774	0.000774	0.000774	100	0.003000	0.050000	-	Fourier	uniform	0.010000	0.001000	10

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Note that Increasing the noise from 0.001 to 0.003 leads to a **sensitive degradation in the certifiable reconstruction precision**.

By increasing the observation horizon up to $N = 100$ it is possible to recover the levels of precision of the previous setting which was achievable with $N = 10$ and the previous level (0.001) of the noise.

The higher the noise is the longer the observation horizon should be to achieve the same level of certifiable reconstruction precision.

Impact of the uncertainty level

eps1	eps2	eps3	zeta1	zeta2	zeta3	N	noise	rho	std_p	u_mod	p_mod	eta	delta	m
0.233572	0.088587	0.033598	0.000464	0.000464	0.000464	20	0.001000	0.050000	-	Fourier	uniform	0.010000	0.001000	10
0.000100	0.000100	0.000100	0.000464	0.000464	0.000464	20	0.001000	-	0.200000	Fourier	gaussian	0.010000	0.001000	10
0.088587	0.033598	0.020691	0.000464	0.000464	0.000464	20	0.001000	-	0.300000	Fourier	gaussian	0.010000	0.001000	10

$$z_1 = x, z_2 = x_1, z_3 = x_3$$

Note that The configuration with **uniform distribution** of the parameter vector with $\rho = 0.05$ is **more harmful** to certifiable reconstruction precision than the Gaussian distribution with standard deviation $s_{td} = 0.2$.

Impact of the input type

eps1	eps2	eps3	zeta1	zeta2	zeta3	N	noise	rho	std_p	u_mod	p_mod	eta	delta	m
0.000100	0.000100	0.000100	0.000464	0.000464	0.000464	20	0.001000	0.050000	-	rand	uniform	0.010000	0.001000	10
0.143845	0.033598	0.012743	0.000464	0.000464	0.000464	20	0.001000	0.050000	0.200000	rand	gaussian	0.010000	0.001000	10
0.233572	0.088587	0.054556	0.000464	0.000464	0.000464	20	0.001000	0.050000	0.300000	rand	gaussian	0.010000	0.001000	10
0.000100	0.000100	0.000100	0.000278	0.000278	0.000278	50	0.001000	0.050000	0.300000	rand	gaussian	0.010000	0.001000	10
0.000100	0.000100	0.000100	0.000278	0.000278	0.000278	50	0.001000	0.050000	0.300000	Fourier	gaussian	0.010000	0.001000	10

$$z_1 = x, z_2 = x_1, z_3 = x_3$$

Note that Comparing the first line with the first line of the previous figure suggests that **random input profiles enhance the observability** at least for the setting that is common to these two lines.

Impact of the certification parameters

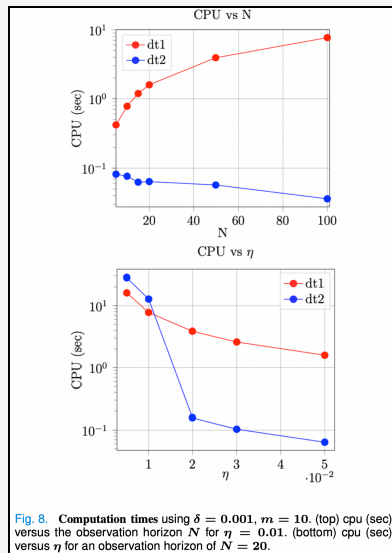
eps1	eps2	eps3	zeta1	zeta2	zeta3	N	noise	rho	std_p	u_mod	p_mod	eta	delta	m
0.615848	0.615848	0.143845	0.000464	0.000464	0.000464	20	0.001000	0.050000	0.300000	rand	gaussian	0.001000	0.001000	10
0.379269	0.379269	0.088587	0.000464	0.000464	0.000464	20	0.001000	0.050000	0.300000	rand	gaussian	0.005000	0.001000	10
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$$z_1 = x, z_2 = x_1, z_3 = x_3$$

Note that A certification precision of $\eta = 10^{-3}$ leads to a certifiable upper bounds on the estimation errors on the observation-target variables which **useless** given the definition of the set \mathbb{X} . This is obviously due to the high level of parameter dispersion $s_{td} = 0.3$.

the last line indicates that up to 5% of the samples correspond to the presence of indistinguishable pairs. For the remaining 95% of the cases, an almost zero reconstruction error can be certified (**provided that the optimization problem is correctly solved**).

CPU



$$\text{CPU} = dt1 + dt2$$

- **dt1** generating the scenarios
- **dt2** finding the optimal design

Recall that the whole scheme is to be executed **off-line** in order to:

- 1 Evaluate the partial observability
- 2 Optimize the observation parameter

The GitHub repository

<https://github.com/mazenalamir/eps-observability>

The screenshot shows a GitHub repository page for the user 'mazenalamir'. The repository name is 'Update README.md', with a commit hash 'c179465' and a date '24 days ago'. There are 9 commits. The main content is the README file, which has the following text:

Python module for checking partial observability of general nonlinear uncertain dynamical systems

This module implements the observability check proposed in the paper:

M. Alamir: Observability Certification and Optimal Design of Nonlinear Observation Parameters in the Presence of Measurement Noise and Model Mismatches. arXiv:2006.11112. June 2021

Code citation:
[DOI: 10.5281/zenodo.4998816](https://doi.org/10.5281/zenodo.4998816)

Problem statement

We consider general uncertain nonlinear dynamical system that can be described by a set of Ordinary Differential System (ODE) of the form:

```
def syst(x,t,u,p):
    ...
    xdot=...
    return xdot
```

where:

- x stands for the state
- u stands for the exogenous input
- p stands for a vector of uncertain parameters

Assume that sensors are available that provides outputs through a dedicated map such as:

```
def output(x, u, p):
    # The equation that defines the measured output other than the input u
    output = ...
    return output
```

Note that we are mainly interested in the case where measurement noise is present that is to be added to the above defined output map.

MA. Partial Extended Observability Certification and Optimal Design of Moving Horizon Estimators Under Uncertainties.

To appear in IEEE Transactions on Automatic control, July, 2022

<https://arxiv.org/abs/2006.11112>

Conclusion

- An example of **Data-driven** assessment of an important control-related problem: *dynamic partial extended estimation*.
- The extensive **Data Generation** step is unavoidable in order to derive non pessimistic results.
- **The choice of the sets/statistics remains a non totally solved issue and is certainly a problem-dependent step.**
- Data-driven methods \neq ML or AI (at least not always)

- 1 Data-driven partial observability assessment under uncertainties
- 2 Learning-Based Approximate Stochastic NMPC Design**
- 3 Tractable stochastic NMPC by supervised clustering
- 4 Data-Driven NMPC by cost function identification
- 5 Learning-Based Monitoring updating period in Real-time NMPC

RECALLS ON STANDARD MPC

Direct single shooting formulation of MPC:

$$\mathbf{u}_k^* \leftarrow \min_{\mathbf{u}_k \in \mathbb{U}^N} [J(\mathbf{u}_k | x_k)] \quad \mathbf{u}_k := \begin{bmatrix} u_k \\ \vdots \\ u_{k+N-1} \end{bmatrix} \in [\mathbb{R}^{n_u}]^N$$



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MPC Feedback Control

$$K_{\text{MPC}}(x_k) = \mathbf{u}_k^*$$

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MPC Feedback Control

$$K_{\text{MPC}}(x_k) = \mathbf{u}_k^*$$

Standard settings

$$J(\mathbf{u}_k \mid x_k) := \sum_{i=1}^N \ell(x_{k+i}, u_{k+i-1})$$

$$x_{k+i+1} = f(x_{k+i}, u_{k+i})$$

PROBLEM STATEMENT

Dynamics

$$x_{k+1} = f(x_k, u_k, w)$$

$$y_k = h(x_k, u_k, w)$$

x the state

u the control

(f, h) known maps

w **constant unknown parameters**

but with known statistics over \mathbb{W}

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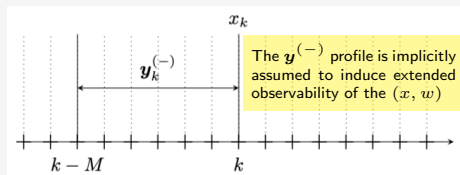
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or gathering all the measurements in $\mathbf{y}^{(-)}$

$$\mathcal{P}(\mathbf{y}^{(-)}) : \min_u \mathbb{E} [J(u \mid \mathbf{y}^{(-)}, \cdot)]$$



More generally

$$\mathcal{P}(\mathbf{y}^{(-)}) : \min_u \mathbb{E}[J] + \alpha \sigma^{\frac{1}{2}} [J]$$

Possible solutions

Option1 (Extended Observer)

Build an extended observer to get

$$(\hat{x}_k, \hat{w}_k)$$

No more need for stochastic setting:

$$\mathbf{u}^* \leftarrow \mathcal{P}(\hat{x}_k, \hat{w}_k)$$

- (+) If possible \rightarrow The best performance
- (-) Needs to build the observer!
- (-) Computation burden adds to NMPC.

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Option2 (Approximate statistics via clustering)

Standard NMPC with the cost given by

$$\frac{1}{n_c} \sum_{i=1}^{n_c} \mu_j J(\mathbf{u} | x_k, w_c^{(j)})$$

- $w_c^{(j)}$ centres of computed clusters
[obtained via supervised clustering]
- μ_j fraction of populations in cluster j
- n_c moderate thanks to clustering

(-) Might still induce heavy computation!

Multiple shooting optimisation → $N(n_c n_x + n_u)$ d.o.f
 $(n_c, N, n_x, n_u) = (10, 100, 3, 1) \rightarrow 3100$ instead of

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POSSIBLE SOLUTIONS

Option3 (Approximate Dynamic Programming)

Approximate the optimal value

$$Q(z) \quad \text{where} \quad z := (x, u)$$

based on the Bellman equation

$$Q(z) = \ell(z) + \gamma \min_v [\hat{\mu} + \alpha \hat{\sigma}^{\frac{1}{2}}][Q(f(z), v)]$$

↑
Fixed-point

- ℓ stage cost
- γ discounting rate
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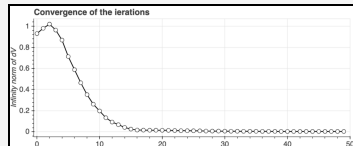
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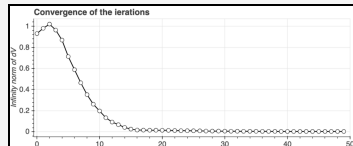
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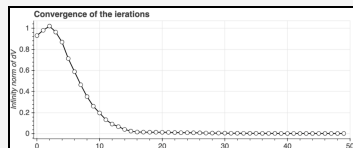
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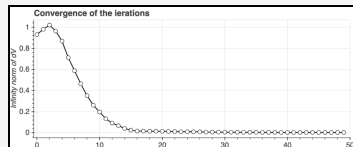
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DEFINITION OF THE ASSOCIATED ML PROBLEM (1)

We would like to have a feedback of the form

$$u_k = \mathbf{ML}(\mathbf{y}_k^{(-)}) \quad \text{where} \quad \mathbf{y}_k^{(-)} = \begin{bmatrix} y_k \\ y_{k-1} \\ \vdots \\ y_{k-M} \end{bmatrix} \in [\mathbb{R}^{n_y}]^{M+1}$$

- $\mathbf{y}_k^{(-)}$ The features
- u_k The label

Where does the label u_k comes from?

DEFINITION OF THE ASSOCIATED ML PROBLEM (2)

Ideally $u_k = \mathbf{u}_0^*(x_k, w)$ where

$$\mathcal{P}(x_k, w) : \quad \mathbf{u}^*(x_k, w) = \arg \min_{\mathbf{u} \in \mathbb{U}^N} [J(\mathbf{u} \mid x_k, w)]$$

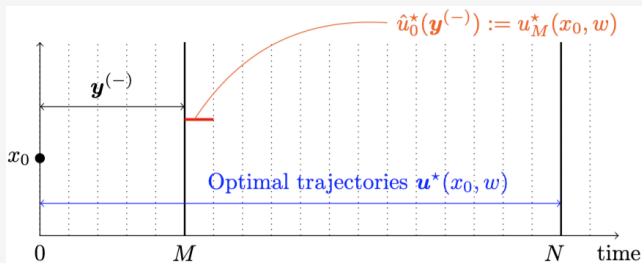
or in terms of $\mathbf{y}_k^{(-)}$ (extended observability required):

$$\begin{aligned} \mathcal{P}(\mathbf{y}_k^{(-)}) : \quad \mathbf{u}^*(\mathbf{y}_k^{(-)}) &= \arg \min_{\mathbf{u} \in \mathbb{U}^N} [J(\mathbf{u} \mid \hat{x}_k, \hat{w})] \\ \hat{x}_k &= \hat{x}(\mathbf{y}_k^{(-)}) \\ \hat{w} &= \hat{w}(\mathbf{y}_k^{(-)}) \end{aligned}$$

But we do not want to build observers! (difficult if not impossible)

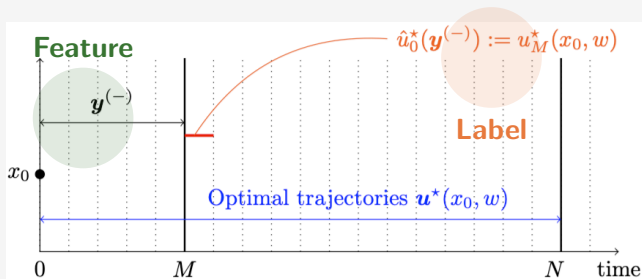
BUILDING THE LEARNING DATA (1)

- 1 Choose arbitrary pair $q := (x_0, w)$
- 2 Solve the problem $\mathcal{P}(x_0, w)$ for known (x_0, w)



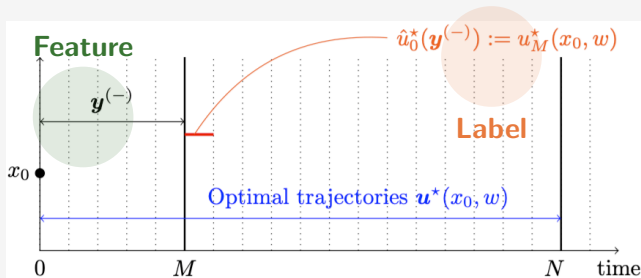
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NOTA

$\hat{u}_0^*(\mathbf{y}^{(-)})$ is only an approximation of $u_0^*(\mathbf{y}_k^{(-)})$ since the **Bellman** equality

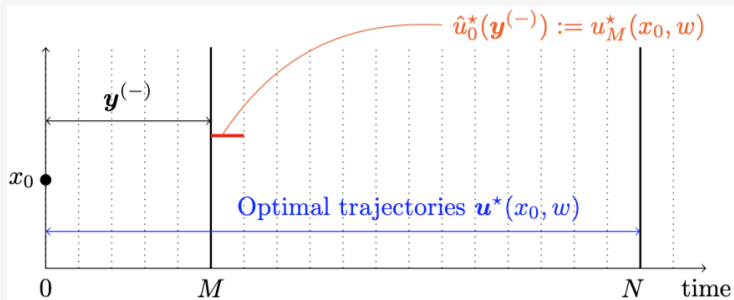
$$u_0^*(x_M, w) = u_M^*(x_0, w)$$

holds only for **infinite horizon**.

BUILDING THE LEARNING DATA (2)

Therefore

$$q = (x_0, w) \xrightarrow{\text{Solve } \mathcal{P}(q)} (\mathbf{y}^{(-)}(q), u_M^*(q))$$

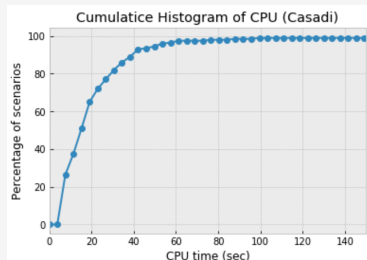


BUILDING THE LEARNING DATA (2)

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$$q = (x_0, w) \xrightarrow{\text{Solve } \mathcal{P}(q)} (\mathbf{y}^{(-)}(q), u_M^*(q))$$

This is far too expensive to get a single sample in the learning data!



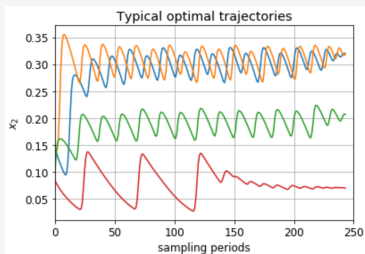
More than **8 days** to get a learning data set containing 50000 samples (Casadi-MS).

BUILDING THE LEARNING DATA (2)

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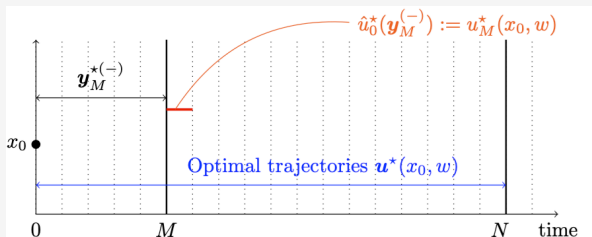
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BUILDING THE LEARNING DATA (3)



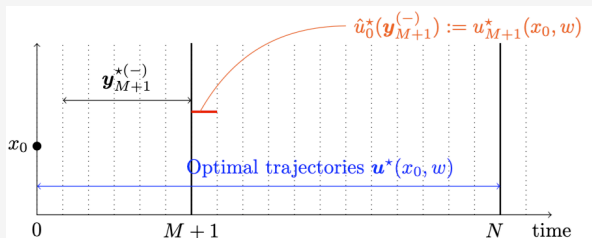
$$\left\{ \mathbf{y}_{M+i}^{*(-)}(q), u_{M+i}^*(q) \right\}_{i=0}^{\mathbf{0}}$$

$$\mathcal{D} := \left\{ [\mathbf{y}_{M+i}^*(q^{(j)})], u_{M+i}^*(q^{(j)}) \right\}_{(i,j) \in \{0, \dots, m\} \times \{1, \dots, n_r\}}$$

m window's moves + n_r different pairs $q \rightarrow \text{card}(\mathcal{D}) = m \cdot n_r$

$m = 250, n_r = 200 \rightarrow 50000$ samples for the price of 200 NLP solution $\rightarrow 50$ minutes!

BUILDING THE LEARNING DATA (3)



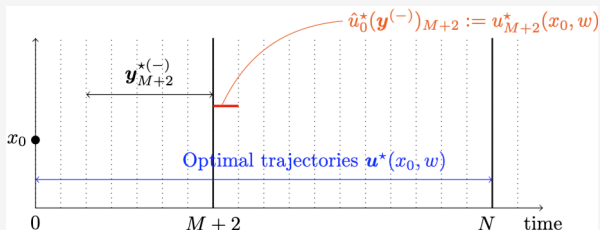
$$\left\{ \mathbf{y}_{M+i}^{*(-)}(q), u_{M+i}^*(q) \right\}_{i=0}^{\mathbf{1}}$$

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BUILDING THE LEARNING DATA (3)



$$\left\{ \mathbf{y}_{M+i}^{*(-)}(q), u_{M+i}^*(q) \right\}_{i=0}^2$$

$$\mathcal{D} := \left\{ [\mathbf{y}_{M+i}^*(q^{(j)})], u_{M+i}^*(q^{(j)}) \right\}_{(i,j) \in \{0, \dots, m\} \times \{1, \dots, n_r\}}$$

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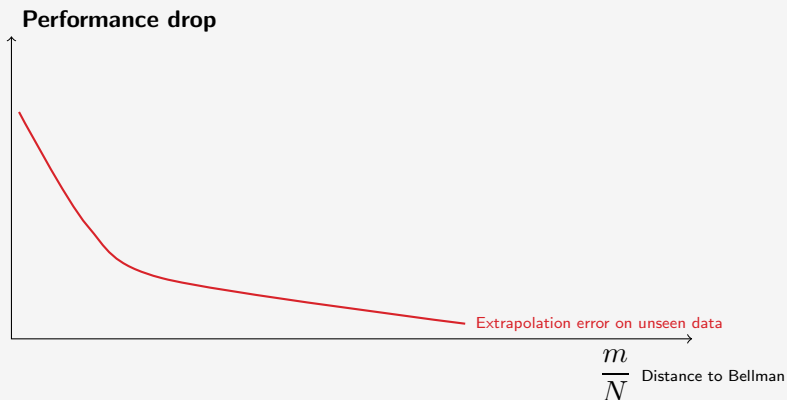
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BUILDING THE LEARNING DATA (4)



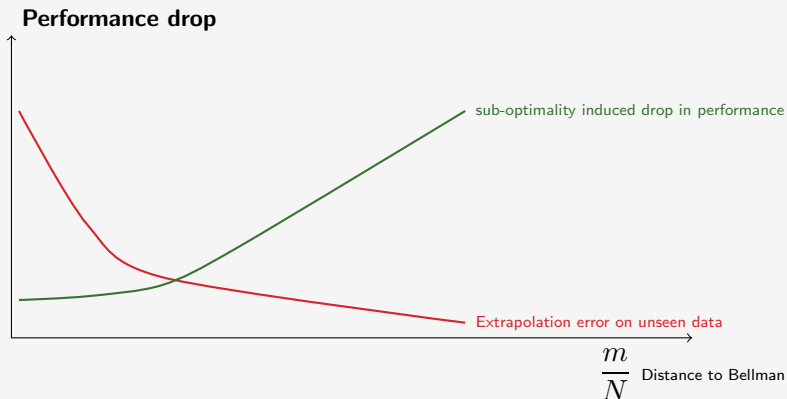
Performance drop for a **Fixed** number of NLP solutions (n_r)

BUILDING THE LEARNING DATA (4)



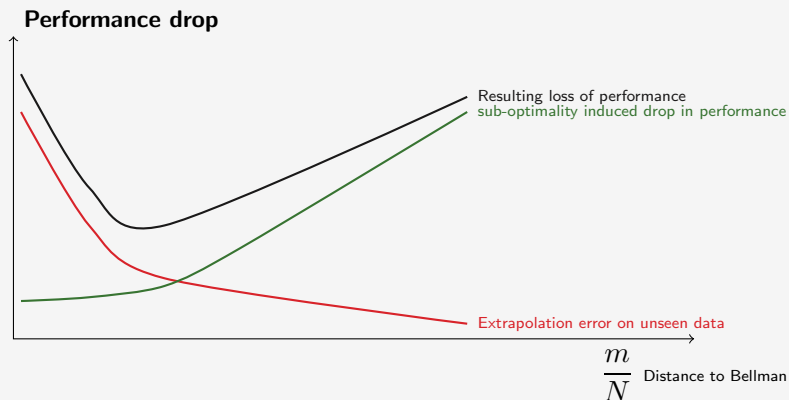
Performance drop for a **Fixed** number of NLP solutions (n_r)

BUILDING THE LEARNING DATA (4)



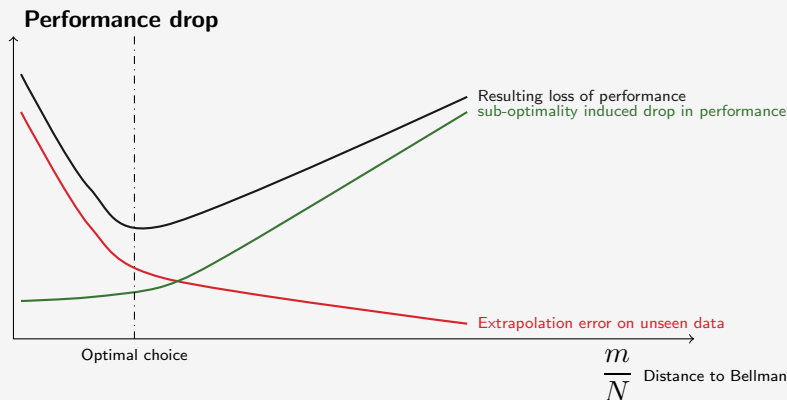
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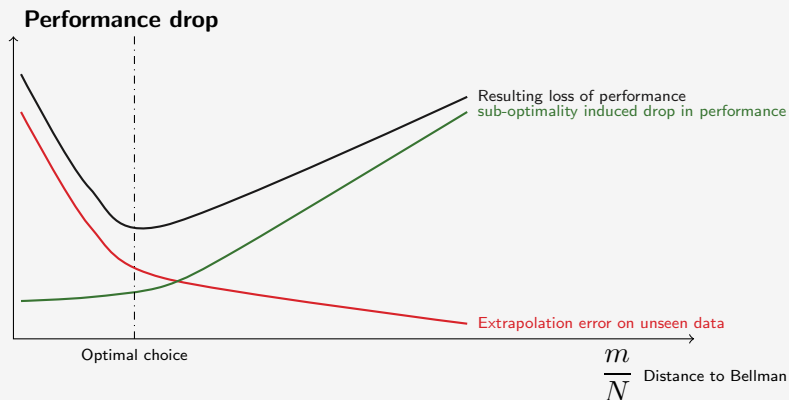
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BUILDING THE LEARNING DATA (4)



Performance drop for a **Fixed** number of NLP solutions (n_r)

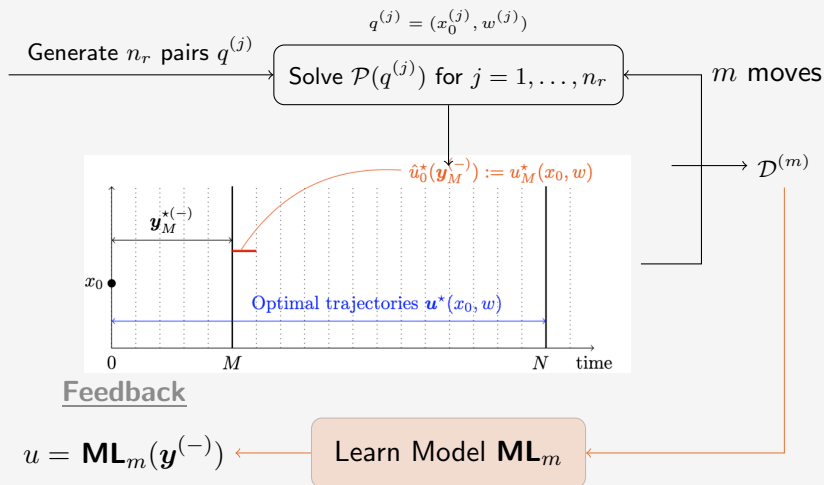
BUILDING THE LEARNING DATA (4)



Performance drop for a **Fixed** number of NLP solutions (n_r)

The optimal choice is problem-dependent.

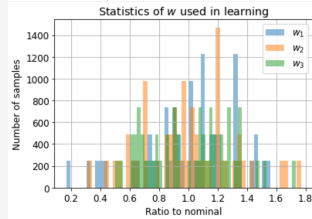
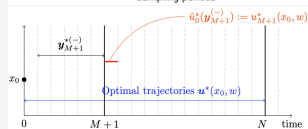
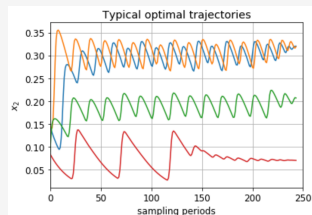
THE WHOLE FLOW IN A GLANCE



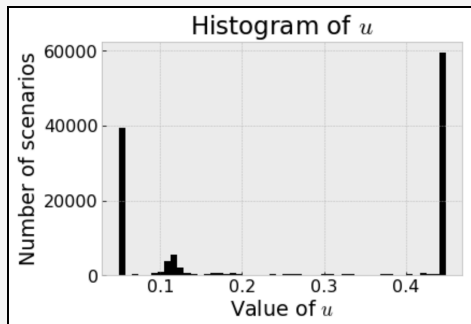
APPLICATION TO THE PARALLEL REACTOR EXAMPLE

Parameters of the method

- $N = 250$
- $M = 10$
- $m \in \{10, 50, 100, 150, 200\}$
- $w^0 = (10^4, 400, 0.55)$
- $w_i = (1 + \nu_i)w_i^0$
- $\nu_i \in \mathcal{N}(0, \sigma_i)$ with $\sigma_i = 0.33$
- $\mathbb{X}_0 := [10^{-4}, 0.5] \times [10^{-4}, 0.2] \times [10^{-4}, 0.25]$
- x_0 sampled uniformly in \mathbb{X}_0
- $n_r = 500$



EXAMINATION OF THE CLOUD OF CONTROL



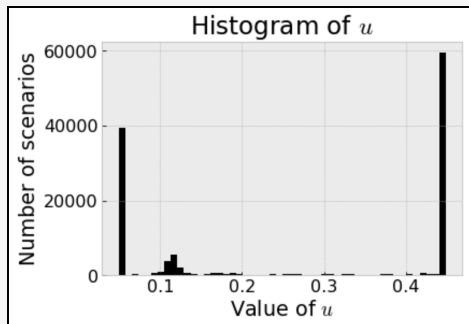
Histogram of the encountered

$$u_{M+i}^*(x_0^{(j)}, w^{(j)})$$

→ `RandomForestClassifier` - `sklearn`

EXAMINATION OF THE CLOUD OF CONTROL

$$\text{label} = \begin{cases} 1 & \text{if } u \leq 0.075 \\ 2 & \text{if } u \in]0.14] \\ 3 & \text{otherwise} \end{cases}$$



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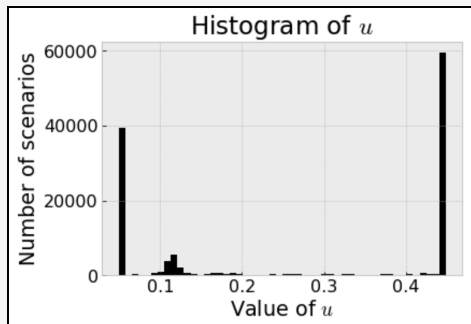
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Based on this observation, a classification approach is adopted for the definition of the **ML**-based feedback

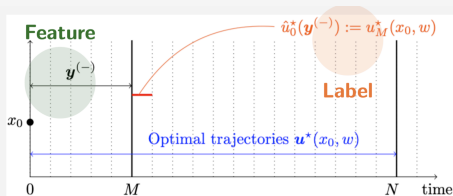


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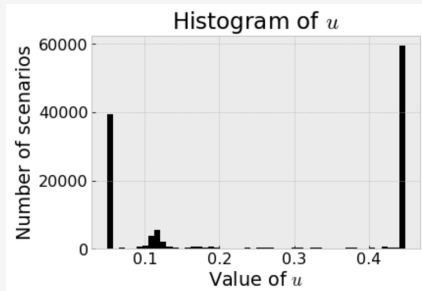
The choice of $\mathbf{y}^{(-)}$



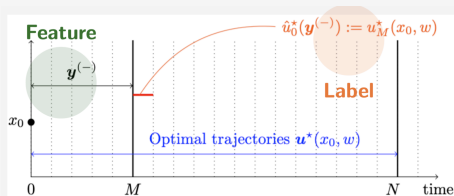
$$\dot{x}_1 = 1 - w_1 x_1^2 e^{-1/x_3} - w_2 x_1 e^{-w_3/x_3} - x_1$$

$$\dot{x}_2 = w_1 x_1^2 e^{-1/x_3} - x_2$$

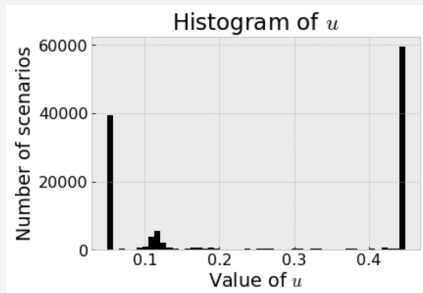
$$\dot{x}_3 = u - x_3$$



The choice of $\mathbf{y}^{(-)}$

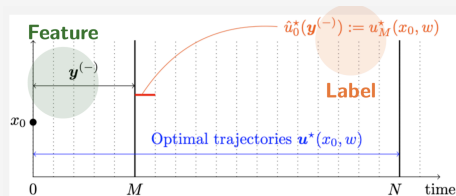


$$\begin{aligned}\dot{x}_1 &= 1 - w_1 x_1^2 e^{-1/x_3} - w_2 x_1 e^{-w_3/x_3} - x_1 \\ \dot{x}_2 &= w_1 x_1^2 e^{-1/x_3} - x_2 \\ \dot{x}_3 &= u - x_3\end{aligned}$$

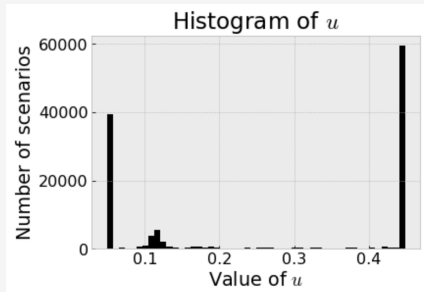


If the control is included in $\mathbf{y}^{(-)}$, then **ML** algorithms would prefer to define the label to be equal to the most recent control as this would correspond to a very high precision.

The choice of $\mathbf{y}^{(-)}$



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By taking $M = 10$, the conjecture is that, for this specific problem, the use of the measurement profile of $\mathbf{y} := x_2$ would be enough to implicitly reveal the value of the state and the parameters and makes the whole framework successful.

LEARNING RESULTS INDICATORS

Training Fitting Results for $m=200$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.864587	0.000876	0.134537
True 2	0.037416	0.910633	0.051952
True 3	0.061044	0.072229	0.866727

Test Fitting Results for $m=200$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.837207	0.001497	0.161296
True 2	0.034890	0.906348	0.058762
True 3	0.085657	0.071689	0.842654

Training Fitting Results for $m=150$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.853440	0.000977	0.145583
True 2	0.032815	0.906244	0.060941
True 3	0.058449	0.069354	0.872197

Test Fitting Results for $m=150$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.802438	0.001393	0.196169
True 2	0.038033	0.885517	0.076450
True 3	0.091729	0.067559	0.840713

LEARNING RESULTS INDICATORS

Training Fitting Results for $m=100$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.827210	0.001355	0.171436
True 2	0.037492	0.893665	0.068843
True 3	0.063109	0.060419	0.876473

Test Fitting Results for $m=100$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.768473	0.002208	0.229319
True 2	0.046135	0.879676	0.074190
True 3	0.107484	0.060515	0.832001

Training Fitting Results for $m=50$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.825142	0.001733	0.173125
True 2	0.079295	0.836123	0.084581
True 3	0.061771	0.026026	0.912203

Test Fitting Results for $m=50$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.701038	0.004540	0.294423
True 2	0.108514	0.824708	0.066778
True 3	0.157318	0.024759	0.817923

LEARNING RESULTS INDICATORS

Training Fitting Results for $m=10$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.995633	0.000000	0.004367
True 2	0.014493	0.956522	0.028986
True 3	0.002632	0.000000	0.997368

Test Fitting Results for $m=10$

	Predicted 1	Predicted 2	Predicted 3
True 1	0.490566	0.001451	0.507983
True 2	0.190476	0.142857	0.666667
True 3	0.211313	0.002134	0.786553

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When test results precision is significantly worse than training precision, this indicate **over-fitting** → bad extrapolation capabilities.

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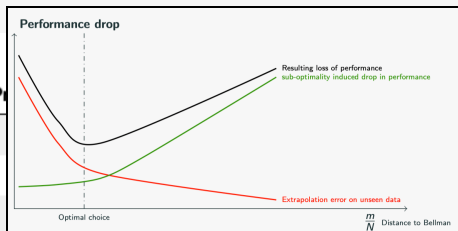
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Validation indicators

Using a new set of 200 pairs of (x_0, w) , closed-loop is simulated over N sampling period. shall compare the following performance indices:

Casadi Optimal

Since the closed-loop simulations last also N sampling periods (The same used for the computation of the open-loop optimal control), the casadi cost is the optimal one since it perfectly knows the parameter values.

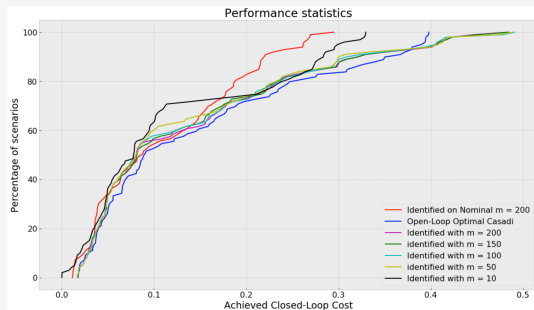
Nominally Optimal

This is the closed loop using learned output feedback from the validation data in which the nominal parameter is used. (Only the initial state is taken from the sampled data).

Proposed for $\neq m$

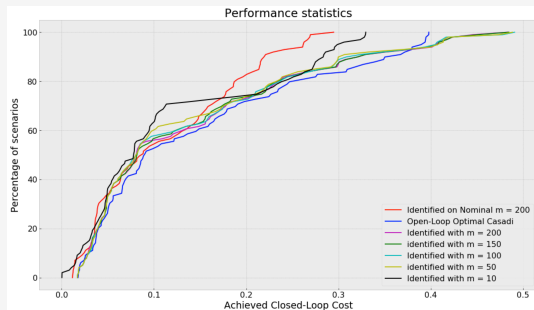
This is the closed loop using learned output feedback from the validation data \mathcal{D}_m obtained by m moves of the window for each x_0 in the learning data.

RESULTS



	mean	relative performance to nominal
Optimal value (Casadi)	0.145409	1.312882
Identified_on_nominal	0.110755	1.000000
Identified with $m = 200$	0.137988	1.245879
Identified with $m = 150$	0.137134	1.238172
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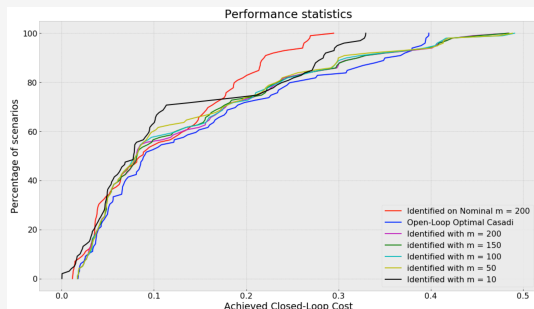
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The proposed method enables to recover 78% of the advantage of knowing perfectly the parameters values when $m \in \{200, 150, 100\}$.

RESULTS

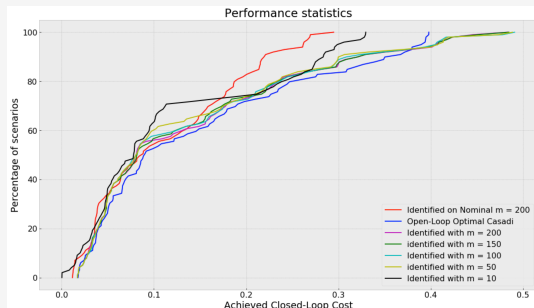


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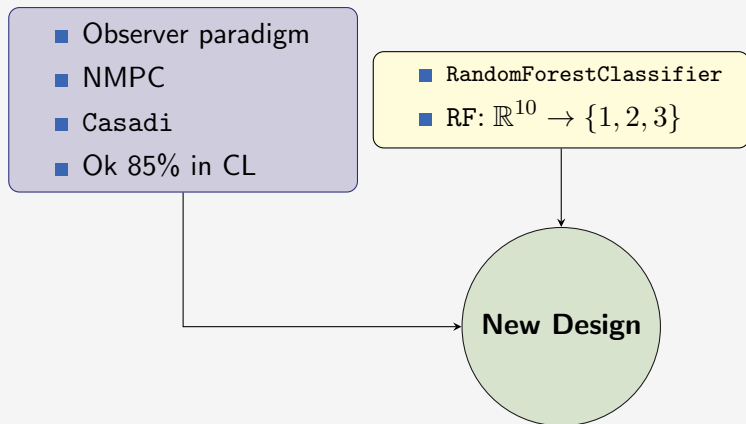
For smaller values of m , data is no more sufficiently representative.

The identified output feedback outperform the casadi output (non achieved optimality) in 10% of the scenarios.

Conclusion

- A heuristic!
- Special care is needed regarding the choice of $\mathbf{y}^{(-)}$
- No need for explicit extended observer.
- Close-to-optimal dynamic output feedback.
- Totally scalable (No exponential explosion with state or parameter dimension).
- Unthinkable without the previous development of NLP (more or less) reliable solvers

TAKE HOME MESSAGE



New Design that would be unconceivable without the mix of two cultures!

- 1 Data-driven partial observability assessment under uncertainties
- 2 Learning-Based Approximate Stochastic NMPC Design
- 3 Tractable stochastic NMPC by supervised clustering**
- 4 Data-Driven NMPC by cost function identification
- 5 Learning-Based Monitoring updating period in Real-time NMPC

STOCHASTIC MPC USING SUPERVISED CLUSTERING

Distributing the data building over the real-life time

Dynamics $\dot{x} = f(x, u, w)$

Cost $J(\mathbf{u}|(x, w))$

Constraints $g(\mathbf{u}|(x, w))$

w-statistics \mathcal{W}

M.A On the use of supervised clustering in stochastic NMPC design. IEEE-TAC. Volume 65, Issue 12, 2020.

STOCHASTIC MPC USING SUPERVISED CLUSTERING

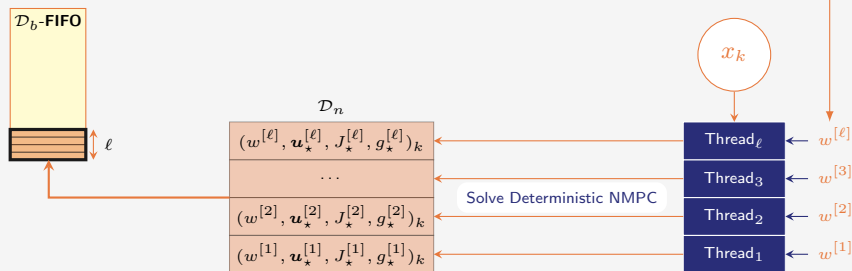
Distributing the data building over the real-life time

Dynamics $\dot{x} = f(x, u, w)$

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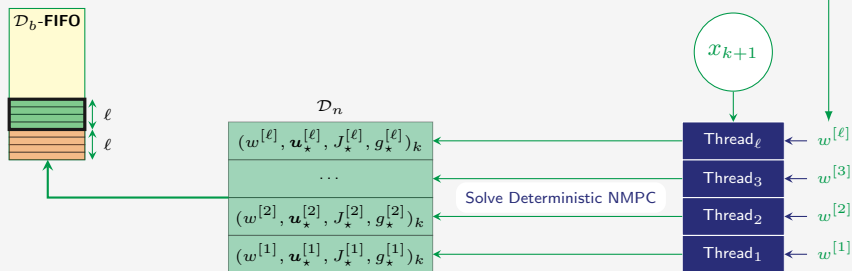
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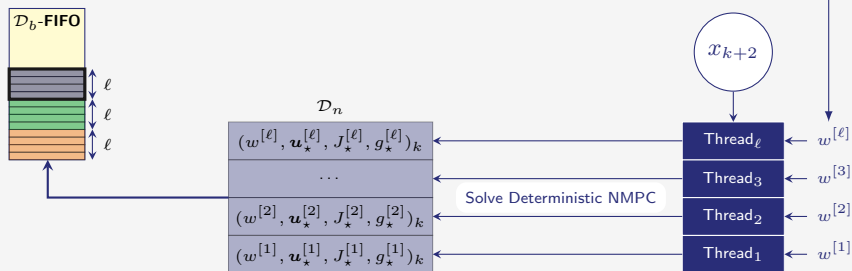
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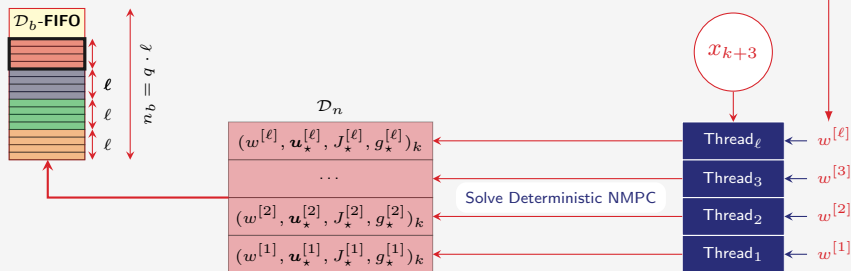
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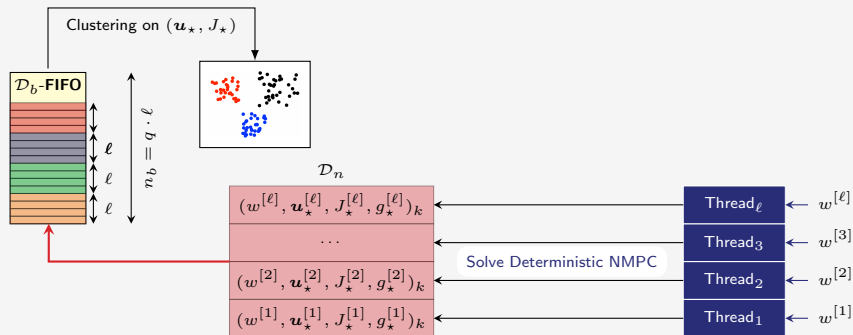
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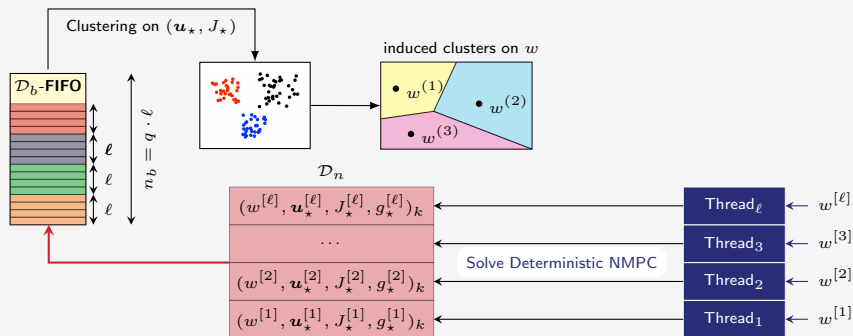
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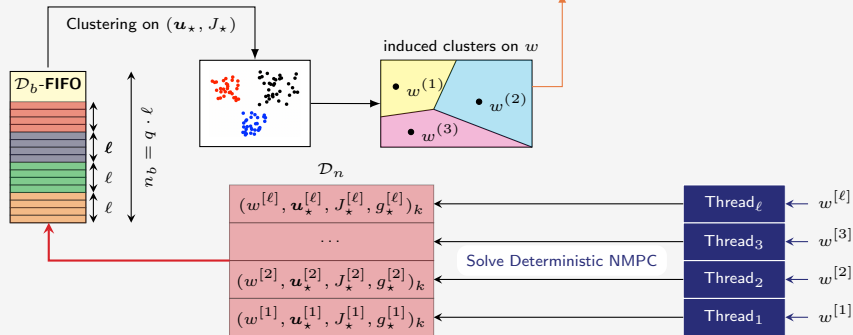
w-statistics \mathcal{W}

$$\min_{\mathbf{u}, \mu \geq 0} \sum_{i=1}^{n_{cl}} p_i \cdot J(\mathbf{u} | (x, w^{(i)}))$$

under

$$g(\mathbf{u} | (x, w^{(i)})) + \frac{1 - \epsilon}{\epsilon} \sigma_g^{(i)} \leq \mu$$

$$i = 1, \dots, n_{cl}$$



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$$i = 1, \dots, n_{cl}$$

x_1 tumor cell population;
 x_2 circulating lymphocytes population;
 x_3 chemotherapy drug concentration;
 x_4 effector immune cell population;
 u_1 rate of introduction of immunotherapy drug;
 u_2 rate of introduction of chemotherapy drug.

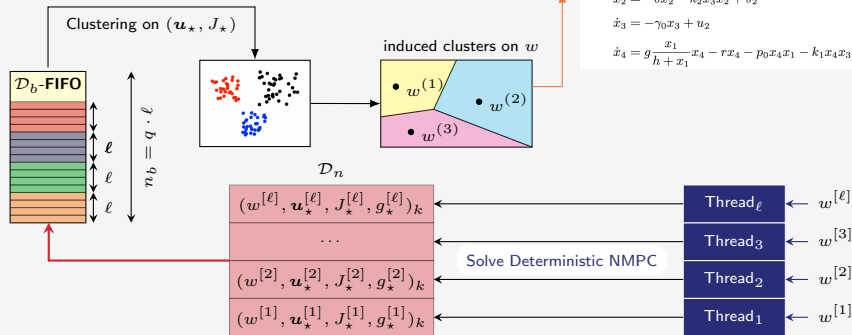
and the dynamics is given by

$$\dot{x}_1 = ax_1(1 - bx_1) - c_1x_4x_1 - k_3x_3x_1$$

$$\dot{x}_2 = -\delta x_2 - k_2x_3x_2 + s_2$$

$$\dot{x}_3 = -\gamma_0x_3 + u_2$$

$$\dot{x}_4 = g \frac{x_1}{h + x_1} x_4 - rx_4 - p_0x_4x_1 - k_1x_4x_3 + s_1u_1.$$



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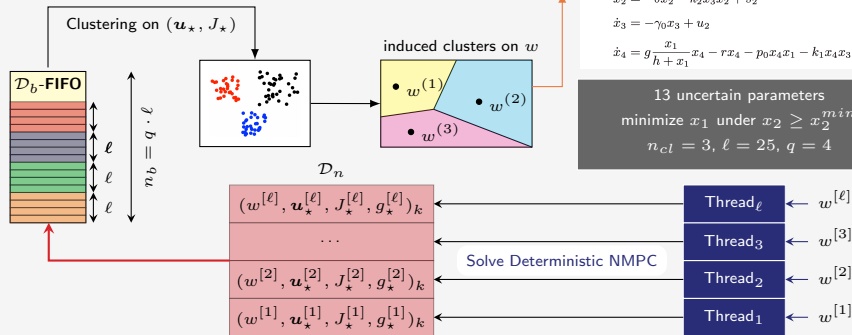
$$\dot{x}_3 = -\gamma_0x_3 + u_2$$

$$\dot{x}_4 = g \frac{x_1}{h + x_1} x_4 - rx_4 - p_0x_4x_1 - k_1x_4x_3 + s_1u_1.$$

13 uncertain parameters

minimize x_1 under $x_2 \geq x_2^{min}$

$n_{cl} = 3, \ell = 25, q = 4$



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Distributing the data building over the real-life time

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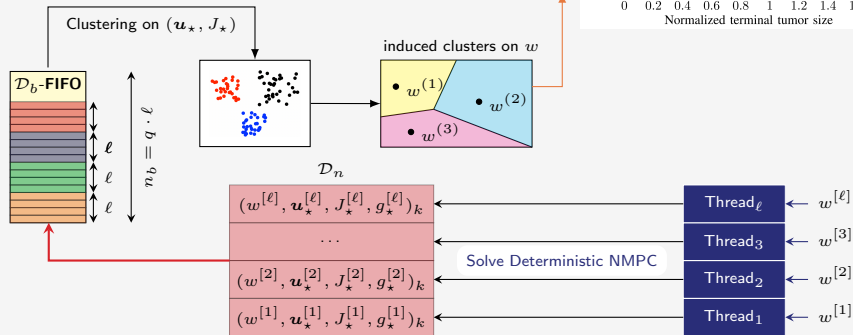
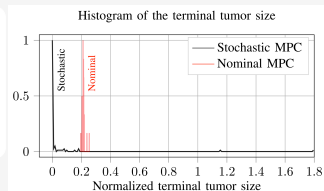
w-statistics \mathcal{W}

$$\min_{u, \mu \geq 0} \sum_{i=1}^{n_{cl}} p_i \cdot J(u|(x, w^{(i)}))$$

under

$$g(u|(x, w^{(i)})) + \frac{1 - \epsilon}{\epsilon} \sigma_g^{(i)} \leq \mu$$

$$i = 1, \dots, n_{cl}$$



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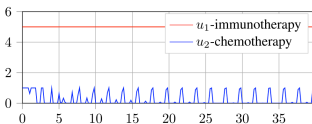
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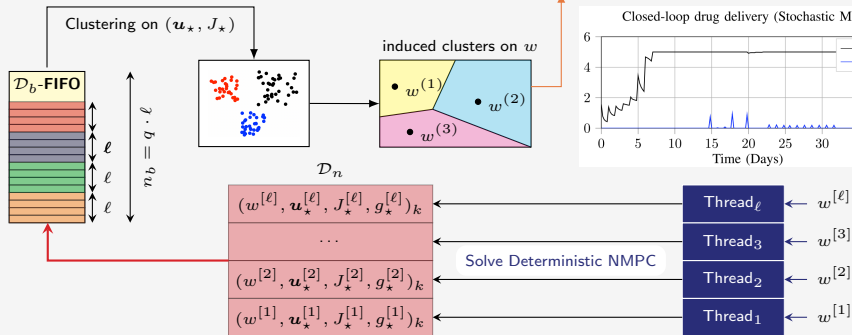
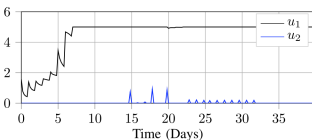
$$g(\mathbf{u}|(x, w^{(i)})) + \frac{1 - \epsilon}{\epsilon} \sigma_g^{(i)} \leq \mu$$

$$i = 1, \dots, n_{cl}$$

Closed-loop drug delivery (Nominal MPC)



Closed-loop drug delivery (Stochastic MPC)



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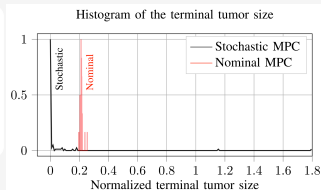
w-statistics \mathcal{W}

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under

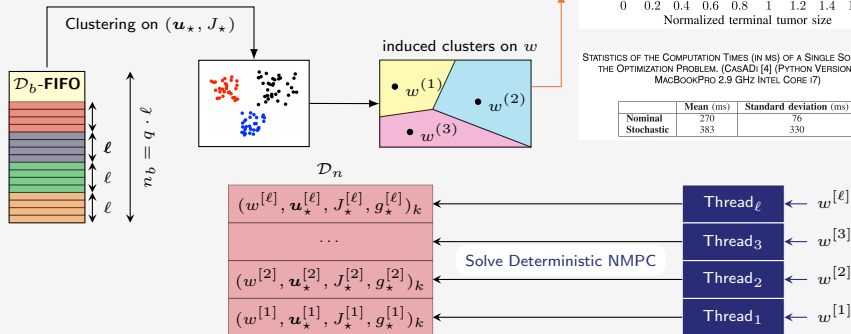
$$g(u|(x, w^{(i)})) + \frac{1 - \epsilon}{\epsilon} \sigma_g^{(i)} \leq \mu$$

$$i = 1, \dots, n_{cl}$$



STATISTICS OF THE COMPUTATION TIMES (IN MS) OF A SINGLE SOLUTION OF THE OPTIMIZATION PROBLEM. (CASADI [4] (PYTHON VERSION) ON A MACBOOKPRO 2.9 GHZ INTEL CORE I7)

	Mean (ms)	Standard deviation (ms)
Nominal	270	76
Stochastic	383	330



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- 1 Data-driven partial observability assessment under uncertainties
- 2 Learning-Based Approximate Stochastic NMPC Design
- 3 Tractable stochastic NMPC by supervised clustering
- 4 Data-Driven NMPC by cost function identification**
- 5 Learning-Based Monitoring updating period in Real-time NMPC

Possible use of ML in MPC design

Direct single shooting formulation of MPC:

$$\mathbf{u}_k^* \leftarrow \min_{\mathbf{u}_k \in \mathbb{U}^N} [J(\mathbf{u}_k | \mathbf{x}_k)]$$

$$\mathbf{u}_k := \begin{bmatrix} u_k \\ \vdots \\ u_{k+N-1} \end{bmatrix} \in [\mathbb{R}^{n_u}]^N$$



MPC Feedback Control

$$K_{\text{MPC}}(\mathbf{x}_k) = \mathbf{u}_k^*$$

Standard settings

$$J(\mathbf{u}_k | \mathbf{x}_k) := \sum_{i=1}^N \ell(\mathbf{x}_{k+i}, \mathbf{u}_{k+i-1})$$

$$\mathbf{x}_{k+i+1} = \mathbf{f}(\mathbf{x}_{k+i}, \mathbf{u}_{k+i})$$

Possible use of ML in MPC design

Option 1 Learn optimal solution u_k^*

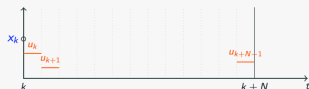
- Known Dynamics
- NMPC solvers (ACADO, GURBOI, CASADI)
- Solve many problems off-line
- Use ML to fit a model with
 - **features:** x_k
 - **Label:** u_k^*

Possible drawbacks:

- (-) Needs the dynamical model
- (-) Expensive computation per samples

Direct single shooting formulation of MPC:

$$u_k^* \leftarrow \min_{u_k \in U^N} [J(u_k | x_k)] \quad u_k := \begin{bmatrix} u_k \\ \vdots \\ u_{k+N-1} \end{bmatrix} \in [\mathbb{R}^{n_u}]^N$$



Standard settings

MPC Feedback Control

$$K_{\text{MPC}}(x_k) = u_k^*$$

$$J(u_k | x_k) := \sum_{i=1}^N \ell(x_{k+i}, u_{k+i-1})$$

$$x_{k+i+1} = f(x_{k+i}, u_{k+i})$$

Possible use of ML in MPC design

Option 2 ML to identify dynamics + on-line solutions

- **Unknown Dynamics**
- Use ML to fit a dynamic model with
 - **Features:** $y_{k-N}, \mathbf{u}_{k-N}, \mathbf{u}_k$
 - **Label:** y_{k+1}
- Use either on-line optimisation or option 1 discussed earlier.

Advantage:

- (-) If successful → Predictive dynamics.

Drawbacks:

- (-) Unnecessarily ambitious for MPC
- (-) Hard (error propagation)

Direct single shooting formulation of MPC:

$$\mathbf{u}_k^* \leftarrow \min_{\mathbf{u}_k \in \mathcal{U}^M} [J(\mathbf{u}_k | x_k)] \quad \mathbf{u}_k := \begin{bmatrix} u_k \\ \vdots \\ u_{k+N-1} \end{bmatrix} \in [\mathbb{R}^{n_u}]^N$$

**Standard settings****MPC Feedback Control**

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$$J(\mathbf{u}_k | x_k) := \sum_{i=1}^N \ell(x_{k+i}, u_{k+i-1})$$

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Possible use of ML in MPC design

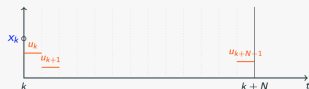
Cool Option Identify cost + on-line solution

- **Unknown Dynamics**
- Gather measurements under
 - open-loop control OR
 - loosely defined feedback
- Use ML to fit a model with
 - **Features:** $\Phi(y_{k-N}, u_{k-N}, u_k)$
 - **Label:** $J(\Phi)$
- Use on-line derivative-free solvers
(ML Models are rarely differentiable)

- (+) fit only what is needed (single model)
- (+) Very light computation per samples
- (+) output MPC feedback (no observer)

Direct single shooting formulation of MPC:

$$u_k^* \leftarrow \min_{u_k \in \mathbb{U}^N} [J(u_k | x_k)] \quad u_k := \begin{bmatrix} u_k \\ \vdots \\ u_{k+N-1} \end{bmatrix} \in [\mathbb{R}^{n_u}]^N$$



Standard settings

MPC Feedback Control

$$K_{\text{MPC}}(x_k) = u_k^*$$

$$J(u_k | x_k) := \sum_{i=1}^N \ell(x_{k+i}, u_{k+i-1})$$

$$x_{k+i+1} = f(x_{k+i}, u_{k+i})$$

The illustrative example

Reactor with two parallel reactions



$$\dot{x}_1 = 1 - 10^4 x_1^2 e^{-1/x_3} - 400 x_1 e^{-0.55/x_3} - x_1$$

$$\dot{x}_2 = 10^4 x_1^2 e^{-1/x_3} - x_2$$

$$\dot{x}_3 = u - x_3$$

- x_1 concentration of R
- x_2 concentration of P_1 (= y)
- x_3 temperature of the mixture
- $u \in [0.049, 0.449]$ heat flow

[Bailey et al. Cyclic operation of reaction systems: effect of heat and mass transfer resistance. AIChE Journal, 17(4):818-825, 1971].

The illustrative example

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Objective

Maximise the production of $y = x_2$

Cost function

$$J := -\left(\frac{1}{N} \mathbf{1}^T \cdot \mathbf{y}_k + \alpha y_{k+N}\right)$$

where

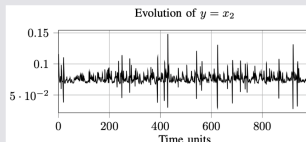
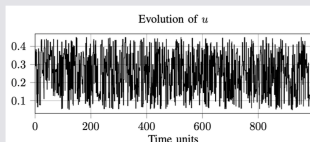
$$\mathbf{y}_k := \begin{bmatrix} y_{k+1} \\ \vdots \\ y_{k+N} \end{bmatrix}$$

Optimal steady state

$$\mathbf{x}_{st} = (0.0832, 0.0846, 0.149)$$

Sketch of the methodology

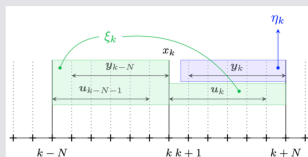
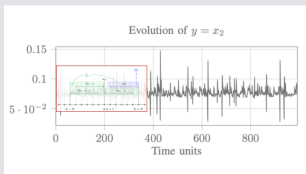
Step 1. Generate Learning Data



- A set of 1000 **interval durations** is randomly sampled uniformly in $[\tau, 100\tau]$.
- A set of 1000 input values are sampled uniformly in $\mathcal{U} := [0.049, 0.149]$.
- → a learning scenario containing **50324 samples**

Sketch of the methodology

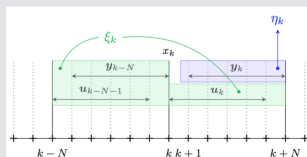
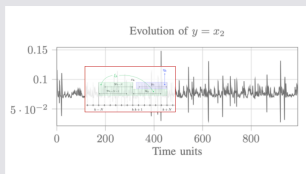
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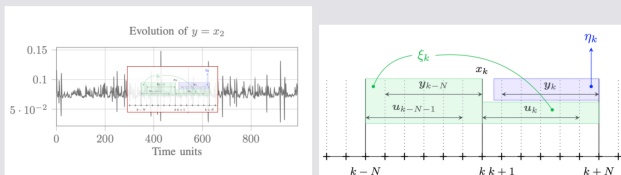
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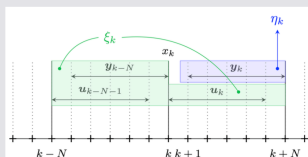
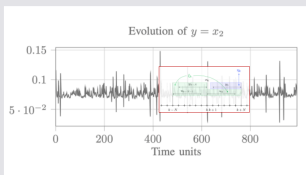
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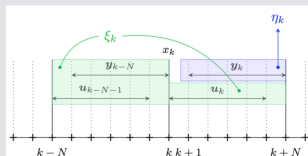
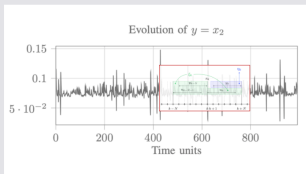
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Sketch of the methodology

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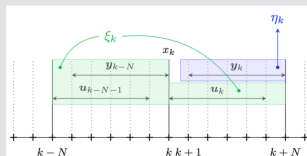
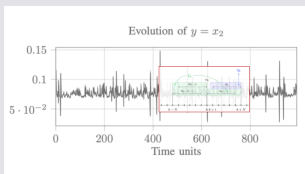
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→ number of samples $n_s = 50324 - 2N$ samples:

$$\left\{ \xi_s, \eta_s \right\}_{s=1}^{n_s} \quad \begin{aligned} \xi_s &= (\mathbf{y}_{k_s-N}, \mathbf{u}_{k_s-N-1}, \mathbf{u}_{k_s}) \in \mathbb{R}^{3N} \\ \eta_s &= J(\mathbf{u}_{k_s} \mid \mathbf{y}_{k_s-N}, \mathbf{u}_{k_s-N-1}) \in \mathbb{R} \\ &= -\left(\frac{1}{N} \mathbf{1}^T \cdot \mathbf{y}_{k_s} + \alpha y_{k_s+N}\right) \end{aligned}$$

Sketch of the methodology

Step 1. Generate Learning Data



- A set of 1000 **interval durations** is randomly sampled uniformly in $[\tau, 100\tau]$.
- A set of 1000 input values are sampled uniformly in $\mathcal{U} := [0.049, 0.149]$.
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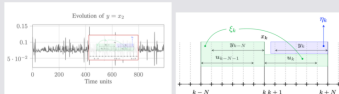
→ number of samples $n_s = 50324 - 2N$ samples:

$$\left\{ \xi_s, \eta_s \right\}_{s=1}^{n_s} \quad \begin{aligned} \xi_s &= (\mathbf{y}_{k_s-N}, \mathbf{u}_{k_s-N-1}, \mathbf{u}_{k_s}) \in \mathbb{R}^{3N} \\ \eta_s &= J(\mathbf{u}_{k_s} \mid \mathbf{y}_{k_s-N}, \mathbf{u}_{k_s-N-1}) \in \mathbb{R} \\ &= -\left(\frac{1}{N} \mathbf{1}^T \cdot \mathbf{y}_{k_s} + \alpha y_{k_s+N}\right) \end{aligned}$$

$N = 100 \rightarrow \xi_s \in \mathbb{R}^{300}$ too high for an appropriate ML design.

Sketch of the methodology

Step 1. Generate Learning Data



- A set of 1000 interval durations is randomly sampled uniformly in $[\tau, 100\tau]$.
- A set of 1000 input values are sampled uniformly in $\mathcal{U} := [0.049, 0.149]$.
- → a learning scenario containing **50324 samples**

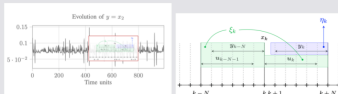
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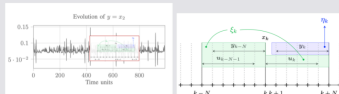
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Step 2. Features selection

$$\Phi(\xi) := \left[\begin{array}{l} \phi_0(\mathbf{y}_{k-N}), \dots, \phi_{n_m-1}(\mathbf{y}_{k-N}), \\ \phi_0(\mathbf{u}_{k-N}), \dots, \phi_{n_m-1}(\mathbf{u}_{k-N}), \\ \phi_0(\mathbf{u}_k), \dots, \phi_{n_m-1}(\mathbf{u}_k) \end{array} \right] \in \mathbb{R}^{3n_m}$$

Sketch of the methodology

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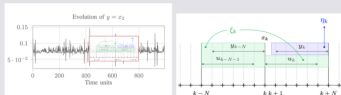
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s : a signal defined on I

$$\phi_j(\mathbf{s}) := \frac{1}{|I|} \sum_{i=1}^{|I|} s^{(j)}(i)$$

Sketch of the methodology

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Machine Learning Problem

$$\eta \approx \text{ML}(\Phi(\xi)) =: \hat{J}(\xi)$$

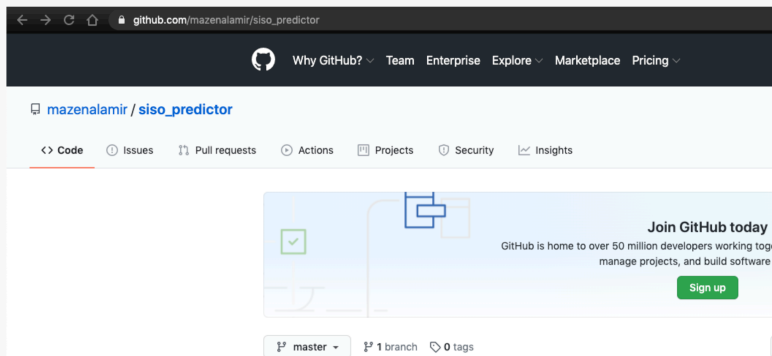
Features $\Phi \in \mathbb{R}^{3n_m} (= \mathbb{R}^9)$

Label $\eta = \text{Cost function}$

s : a signal defined on I


$$\phi_j(\mathbf{s}) := \frac{1}{|I|} \sum_{i=1}^{|I|} s^{(j)}(i)$$








Freely available python implementation on GitHub



The screenshot shows the GitHub interface for the repository 'mazenalamir / siso_predictor'. The browser address bar displays 'github.com/mazenalamir/siso_predictor'. The navigation bar includes the GitHub logo and links for 'Why GitHub?', 'Team', 'Enterprise', 'Explore', 'Marketplace', and 'Pricing'. Below the repository name, there are tabs for 'Code', 'Issues', 'Pull requests', 'Actions', 'Projects', 'Security', and 'Insights'. The 'Code' tab is selected and underlined. A large banner for 'Join GitHub today' is visible, featuring a diagram of a branching strategy and the text 'GitHub is home to over 50 million developers working together to manage projects, and build software faster'. A green 'Sign up' button is present in the banner. At the bottom of the repository view, there are buttons for 'master', '1 branch', and '0 tags'.

Freely available python implementation on GitHub

 **mazenalamir** Update README.md 69dda19 on 4 May 🕒 10 commits

 .ipynb_checkpoints	intial commit	4 months ago
 .vscode	intial commit	4 months ago
 images	add an image example	4 months ago
 README.md	Update README.md	4 months ago
 generate_data.py	intial commit	4 months ago
 main.py	intial commit	4 months ago
 siso_predictor.py	intial commit	4 months ago

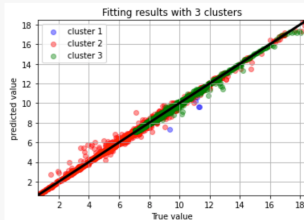
README.md

```
siso_predictor
```

Freely available python implementation on GitHub

Example of use

```
sol = learn_model(
    y=y,
    U=U,
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    N=100,
    n_clusters=3,
    nJump=1,
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    test_size=0.33,
    validation_mode='all',
    plot=True
)
```



Mean abs(Error) on all cluster= 0.022830

abs(Error) - cluster 1 = 0.008946
abs(Error) - cluster 2 = 0.037074
abs(Error) - cluster 3 = 0.022471

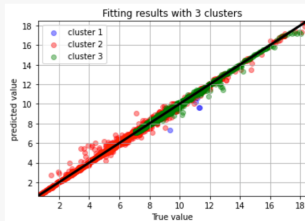
- Random Forest structure (SKLEARN)
- 33% test for train/test data splitting
- Clustering uses KNN (SKLEARN)

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We have a cost function model

$$J = \left[\text{ML} \circ \Phi \right] (\mathbf{y}_{k-N}, \mathbf{u}_{k-N}, \mathbf{u}_k)$$

Solving the optimization problem

The Model $J = [\text{ML} \circ \Phi](\mathbf{y}_{k-N}, \mathbf{u}_{k-N}, \mathbf{u}_k)$

The cost induced by a future control sequence \mathbf{u}_k given the past measurements $(\mathbf{y}_{k-N}, \mathbf{u}_{k-N})$.

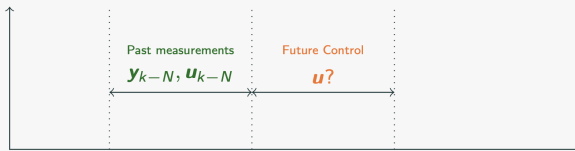
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The Data-Driven MPC Formulation at instant k

$$\mathbf{u}_k^* \leftarrow \arg \min_{\mathbf{u} \in \mathbb{U}^N} [\text{ML} \circ \Phi](\mathbf{y}_{k-N}, \mathbf{u}_{k-N}, \mathbf{u})$$



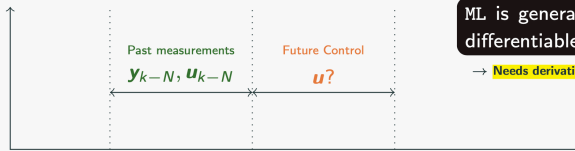
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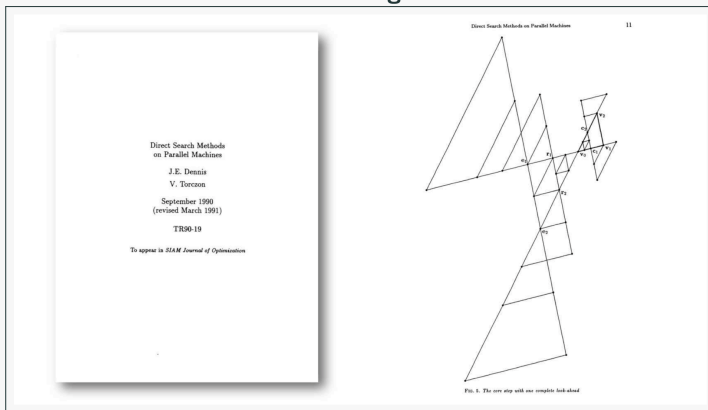
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Solving the optimization problem

Torczon's Algorithm



V. Torczon. *On the convergence of pattern search algorithms.*
SIAM J. on Optimization, 7(1):1-25, 1997

Solving the optimization problem

Suppose n Decision variables

Take $n + 1$ initial nodes
Determine the best node

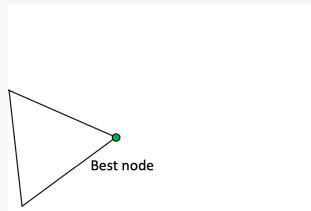


Illustration of Torczon Algorithm for 2D problems

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Try a Reflexion

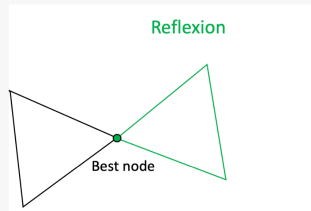


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If successful Try an extension

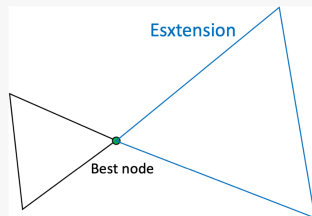


Illustration of Torczon Algorithm for 2D problems

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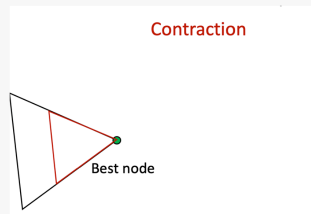


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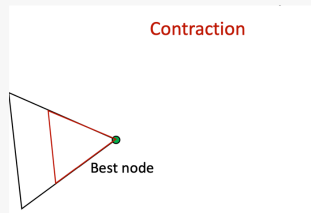


Illustration of Torczon Algorithm for 2D problems

This algorithm is suitable for problems of moderate size!

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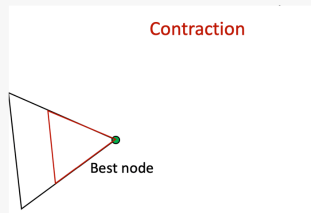


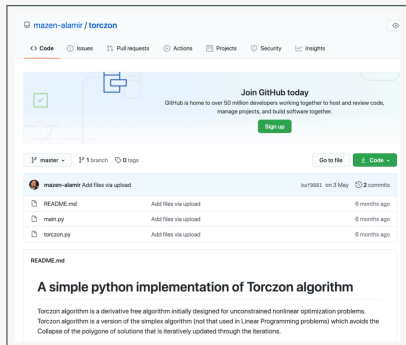
Illustration of Torczon Algorithm for 2D problems

This algorithm is suitable for problems of moderate size!

→ Control parametrisation

[last step!]

Freely available python solver implementation on GitHub



mazen-alamir / torczon

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master 1 branch 0 tags

Go to file Code -

mazen-alamir Add files via upload ba79861 on 3 May 2 commits

README.md	Add files via upload	6 months ago
main.py	Add files via upload	6 months ago
torczon.py	Add files via upload	6 months ago

README.md

A simple python implementation of Torczon algorithm

Torczon algorithm is a derivative free algorithm initially designed for unconstrained nonlinear optimization problems. Torczon algorithm is a version of the simplex algorithm (not that used in Linear Programming problems) which avoids the Collapse of the polygone of solutions that is iteratively updated through the iterations.

Usage

- Define the cost function to be optimized, the cost might embed soft constraint definition via constraint penalty (the penalty can be a part of the variable p that can be any python object or variable).

```
def f(x,p):
    return ..
```

- Call the solver using

```
solve(f_usermf,
      parfp, x0=x0,
      xmin=xmin,
      xmax=xmax,
      Nguess=Nguess,
      Niter=Niter,
      initial_box_width=0.1)
```

where

- $x0$ the initial guess (for the first guess)
- $xmin, xmax$ the box of admissible values
- $Niter$: the number of iterations by single guess
- $Nguess$: the number of initial guesses (randomly sampled using uniform distribution inside the hypercube defined by $xmin$ and $xmax$)
- $initial_box_width$: the amplitude of the initial steps around each initial guess to build the polygone of the torczon algorithm.

Control parametrization

Data-Driven MPC at instant k

$$\mathbf{u}_k^* \leftarrow \arg \min_{\mathbf{u} \in \mathbb{U}^N} [\text{ML} \circ \Phi](\mathbf{y}_{k-N}, \mathbf{u}_{k-N}, \mathbf{u})$$

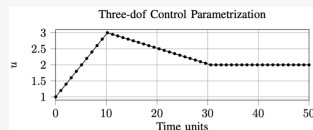
↓

$$\mathbf{p}^* \leftarrow \arg \min_{\mathbf{p} \in \mathcal{P}} [\text{ML} \circ \Phi](\mathbf{y}_{k-N}, \mathbf{u}_{k-N}, \mathcal{U}(\mathbf{p}))$$

↓

MPC Feedback

Apply the first control input in $\mathcal{U}(\mathbf{p}^*)$



$\mathbf{u} = \mathcal{U}(\mathbf{p})$ where (here)

$$\mathbf{p} := [u(0), u(10), u(30)]$$

Results

Two settings are tested ($\tau = 0.02, N = 100$), single cluster

Setting 1

- $p := (u(0), u(10)) \in \mathbb{U}^2$
- $N_{guess} = 1, n_{iter}^{max} = 30$ (Torczon)
- $\rightarrow 1 \times 30 \times (2 + 4) = 180$ functions evals

Setting 2

- $p := (u(0), u(5), u(10), u(20)) \in \mathbb{U}^4$
- $N_{guess} = 3, n_{iter}^{max} = 50$ (Torczon)
- $\rightarrow 3 \times 50 \times (4 + 4) = 1200$ functions evals

Single evaluation of the random-forest cost model = 1.7 msec

[Python 3 jupyter-notebook on a MacBook-Pro 2017, Mojave 10.14.5 with a 2.9 GHz Core i7 processor]

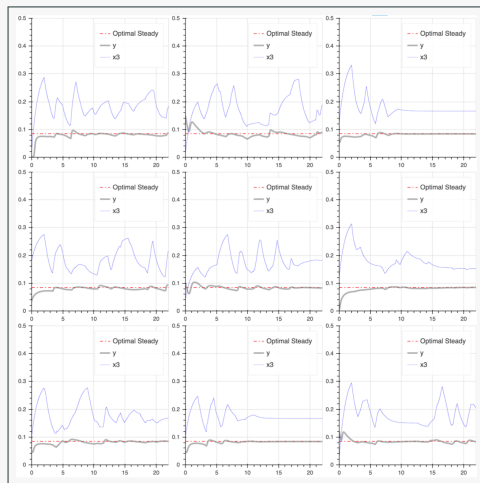
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NOTA

The first two seconds ($N \times \tau$) under open-loop control waiting for the buffer $\mathbf{y}_{k-N}, \mathbf{u}_{k-N}$ to be available.



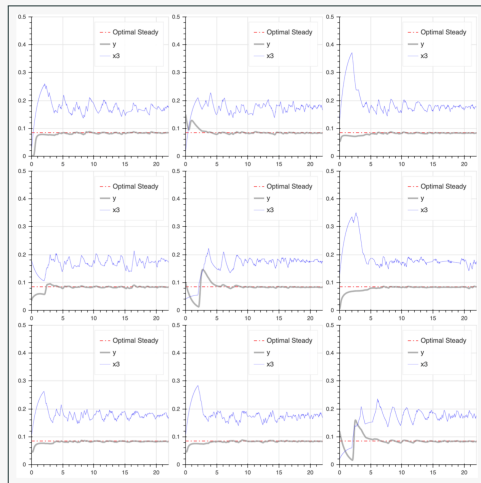
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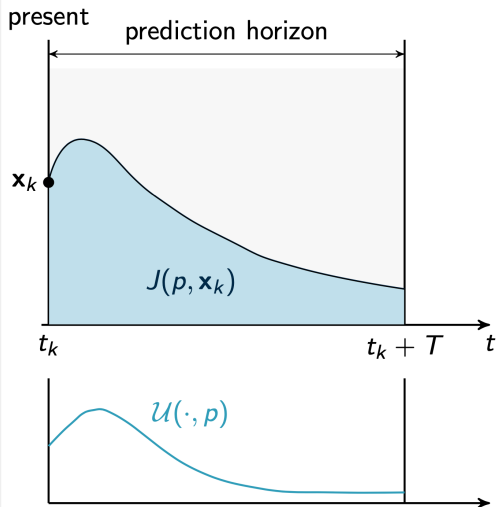
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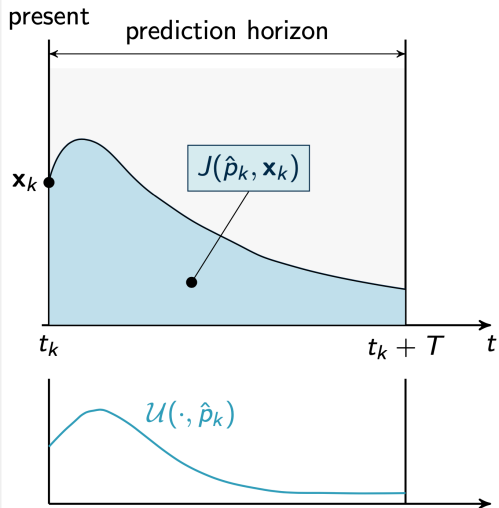


- 1 Data-driven partial observability assessment under uncertainties
- 2 Learning-Based Approximate Stochastic NMPC Design
- 3 Tractable stochastic NMPC by supervised clustering
- 4 Data-Driven NMPC by cost function identification
- 5 Learning-Based Monitoring updating period in Real-time NMPC**

Ideal Framework: Recalls & Basic Notations



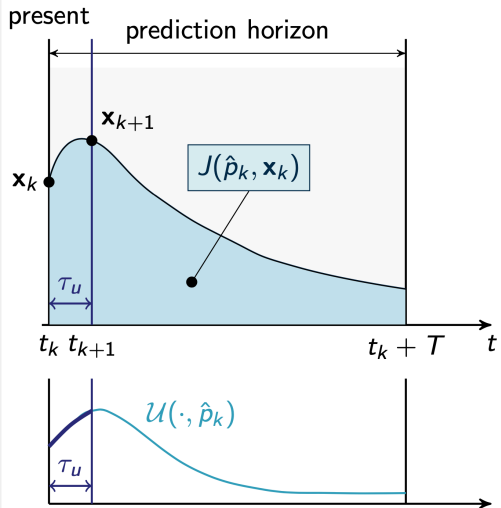
Ideal Framework: Recalls & Basic Notations



$$\min_p J(p, \mathbf{x}_k) \text{ s.t. } C(p, \mathbf{x}_k) \leq 0$$

$$\hat{p}_k = \hat{p}(\mathbf{x}_k)$$

Ideal Framework: Recalls & Basic Notations

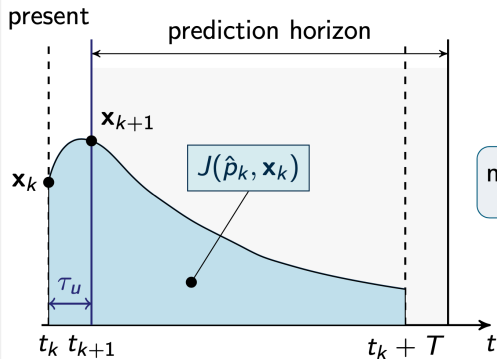


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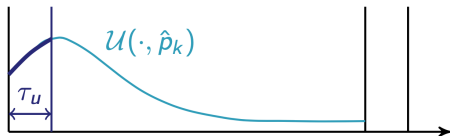
$$\hat{p}_k = \hat{p}(\mathbf{x}_k)$$

Apply $U(\cdot, \hat{p}(\mathbf{x}_k))$ during τ_u

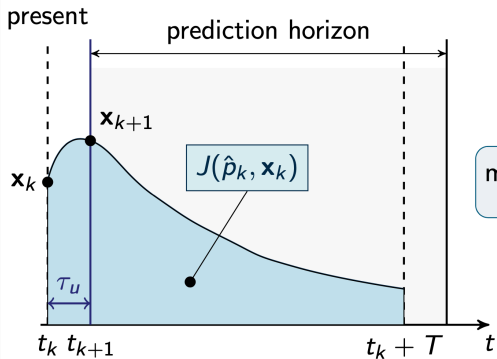
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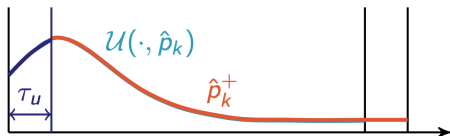
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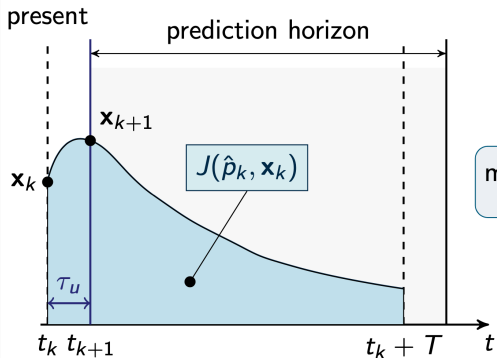
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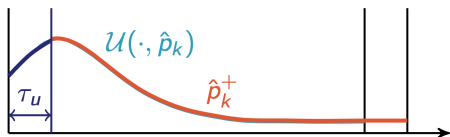
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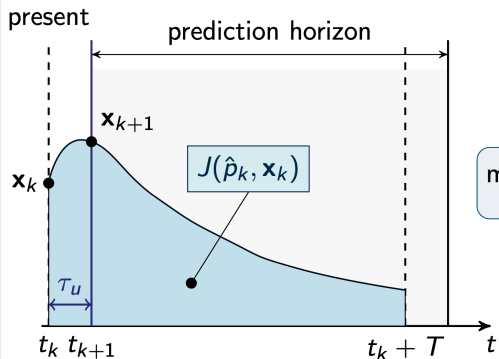
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Initialize at \hat{p}_k^+ (hot start)

$$\hat{p}_{k+1} = \hat{p}(\mathbf{x}_{k+1})$$



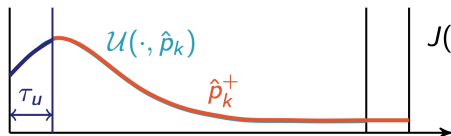
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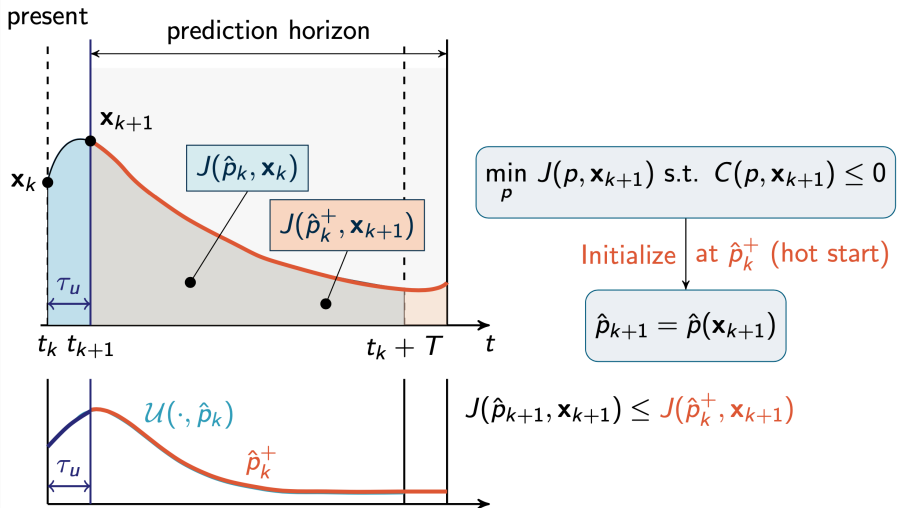
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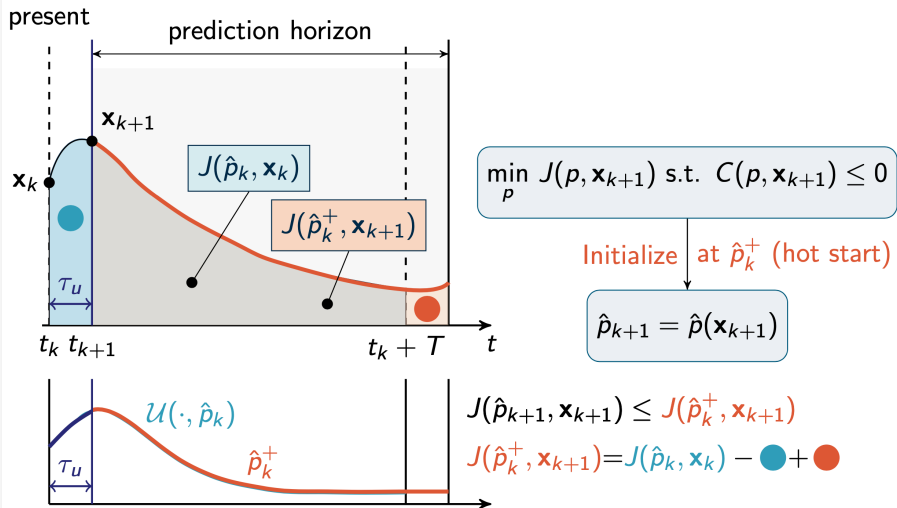


$$J(\hat{p}_{k+1}, \mathbf{x}_{k+1}) \leq J(\hat{p}_k^+, \mathbf{x}_{k+1})$$

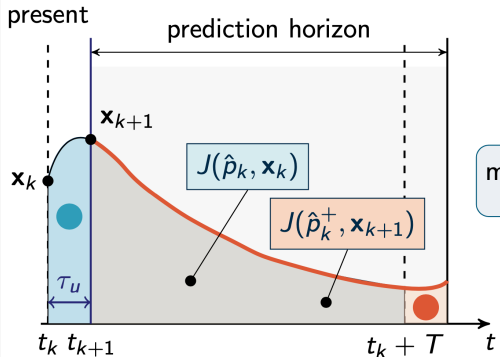
Ideal Framework: Recalls & Basic Notations



Ideal Framework: Recalls & Basic Notations



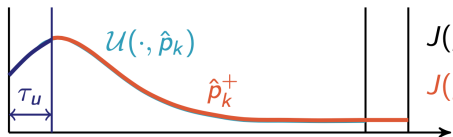
Ideal Framework: Recalls & Basic Notations



$$\min_p J(p, \mathbf{x}_{k+1}) \text{ s.t. } C(p, \mathbf{x}_{k+1}) \leq 0$$

Initialize at \hat{p}_k^+ (hot start)

$$\hat{p}_{k+1} = \hat{p}(\mathbf{x}_{k+1})$$

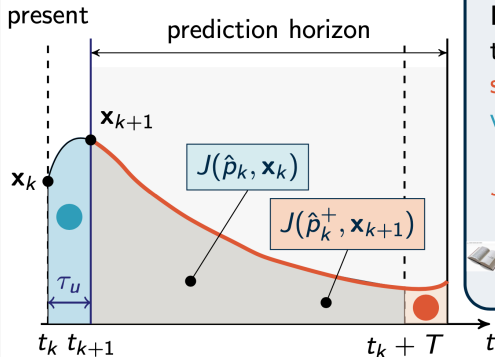


$$J(\hat{p}_{k+1}, \mathbf{x}_{k+1}) \leq J(\hat{p}_k^+, \mathbf{x}_{k+1})$$

$$J(\hat{p}_k^+, \mathbf{x}_{k+1}) = J(\hat{p}_k, \mathbf{x}_k) - \text{blue circle} + \text{red circle}$$

$\leq 0 ?$

Ideal Framework: Recalls & Basic Notations



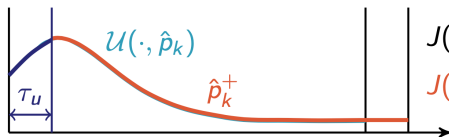
Keep in mind

In the **ideal framework**, when the horizon moves, the **hot start** \hat{p}_k^+ computed from the **previous optimal solution** \hat{p}_k satisfies:

$$J(\hat{p}_k^+, \mathbf{x}_{k+1}) \leq J(\hat{p}_k, \mathbf{x}_k)$$



Mayne et al. Automatica (2000)



$$J(\hat{p}_{k+1}, \mathbf{x}_{k+1}) \leq J(\hat{p}_k^+, \mathbf{x}_{k+1})$$

$$J(\hat{p}_k^+, \mathbf{x}_{k+1}) = J(\hat{p}_k, \mathbf{x}_k) - \text{blue circle} + \text{red circle} \leq 0 ?$$

Ideal MPC: The key assumptions

Keep in mind

In the **ideal framework**, when the horizon moves, the **hot start** \hat{p}_k^+ computed from the **previous optimal solution** \hat{p}_k satisfies:

$$J(\hat{p}_k^+, \mathbf{x}_{k+1}) \leq J(\hat{p}_k, \mathbf{x}_k)$$



Mayne et al. Automatica (2000).

- ▶ Formulation involving Final constraints
- ▶ (hot start)-compatible parametrization
- ▶ \hat{p}_k sufficiently good

Ideal MPC: The key assumptions

Keep in mind

In a **realistic framework**, when the horizon moves, the **hot start** p_k^+ computed from the **previous solution** p_k satisfies:

$$J(p_k^+, \mathbf{x}_{k+1}) = J(p_k, \mathbf{x}_k) + D(\tau_u)$$

$$D(0) = 0$$

$$D(\tau_u) \text{ is **not necessarily** } \leq 0.$$

- ▶ Formulation involving Final constraints
- ▶ (hot start)-compatible parametrization
- ▶ \hat{p}_k sufficiently good

Ideal MPC: The key assumptions

Keep in mind

In a **realistic framework**, when the horizon moves, the **hot start** p_k^+ computed from the **previous solution** p_k satisfies:

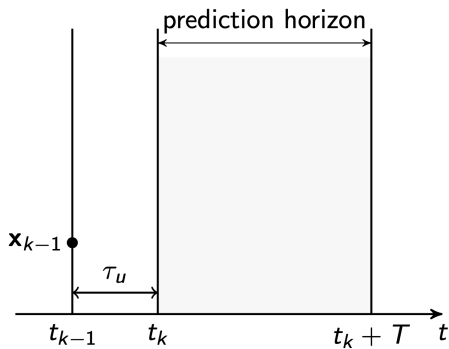
$$J(p_k^+, \mathbf{x}_{k+1}) = J(p_k, \mathbf{x}_k) + D(\tau_u)$$

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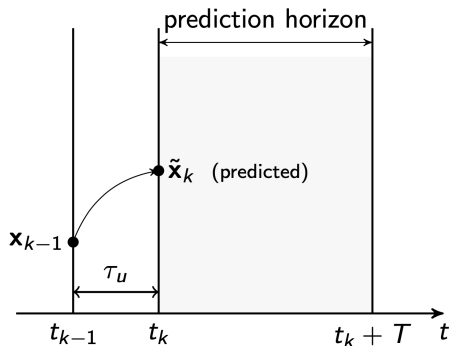
$D(\tau_u)$ is **not necessarily** ≤ 0 .

Even with **perfect undisturbed** model

Preparation & Feedback Steps

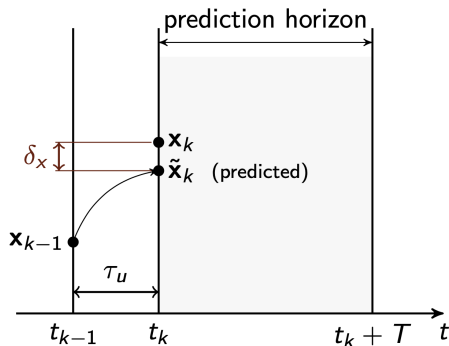


Preparation & Feedback Steps



1. Predict $\tilde{\mathbf{x}}_k$
2. During $[t_{k-1}, t_k]$
 Compute $\hat{p}(\tilde{\mathbf{x}}_k)$ [and $\frac{\partial \hat{p}_k}{\partial \mathbf{x}}$]

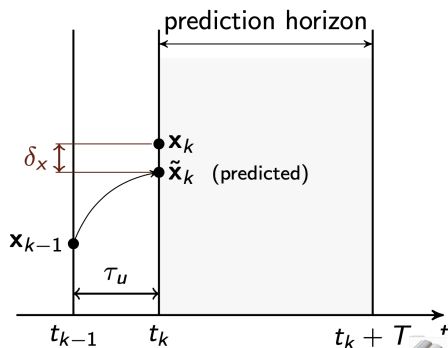
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2. During $[t_{k-1}, t_k]$
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3. Once \mathbf{x}_k is available:

$$\hat{p}_k \leftarrow \hat{p}(\tilde{\mathbf{x}}_k) + \left[\frac{\partial \hat{p}_k}{\partial \mathbf{x}} \right] \cdot \delta_x$$

Preparation & Feedback Steps



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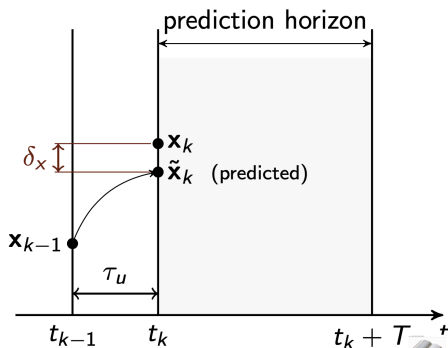
$$\hat{p}_k \leftarrow \hat{p}(\tilde{\mathbf{x}}_k) + \left[\frac{\partial \hat{p}_k}{\partial \mathbf{x}} \right] \cdot \delta_{\mathbf{x}}$$

(*feedback step*)

 Diehl et al. SIAM J. Ctrl and Opt. (2005)

 Zavala and Biegler. Automatica (2009)

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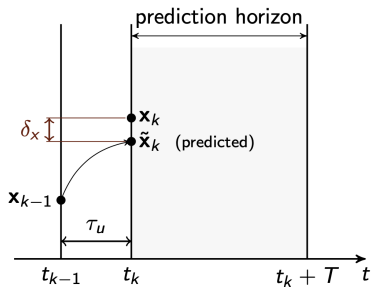
(*feedback step*)

 Diehl et al. SIAM J. Ctrl and Opt. (2005)

 Zavala and Biegler. Automatica (2009)

NOTA: Even with feedback, uncertainty potentially increases $D(\tau_u)$

Definition of Fast NMPC Problems

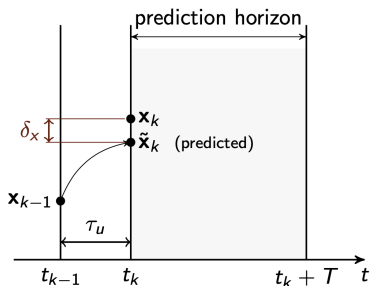


τ_u is the time between two control updating

τ_u is the time during which there is **no feedback**

$$\Rightarrow \tau_u \leq \tau_u^{max}$$

Definition of Fast NMPC Problems



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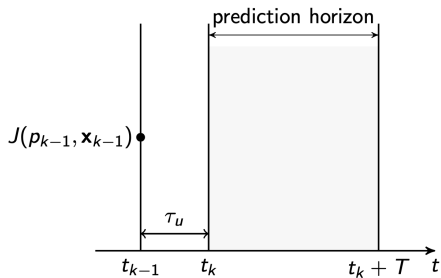
$$\Rightarrow \tau_u \leq \tau_u^{max}$$

Fast NMPC Problems

Fast NMPC problems are those for which

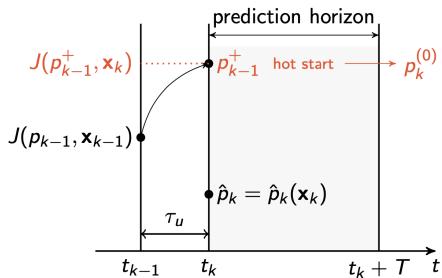
$$\tau_{solve}(NLP(\tilde{\mathbf{x}}_k)) \geq \tau_u^{max}$$

The Iterative Process



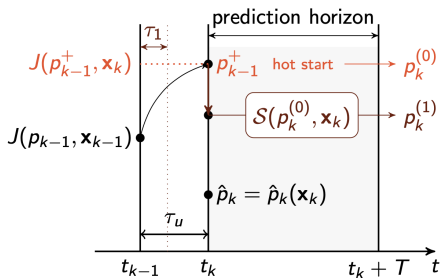
$$p^{(i+1)} \leftarrow \mathcal{S}(p^{(i)}, \mathbf{x})$$

The Iterative Process



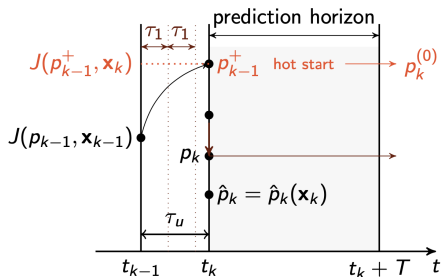
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The Iterative Process



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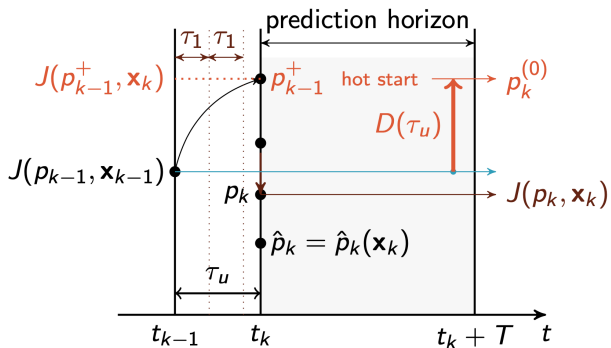
$$p^{(i+1)} \leftarrow \mathcal{S}(p^{(i)}, \mathbf{x})$$

$$q = \text{int}\left(\frac{\tau_u}{\tau_1}\right)$$

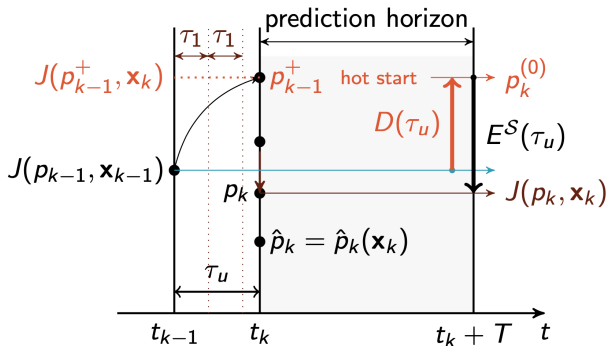
dynamic equation for p

$$p_k := \mathcal{S}^{(q)}(p_{k-1}^+, \mathbf{x}_{k-1})$$

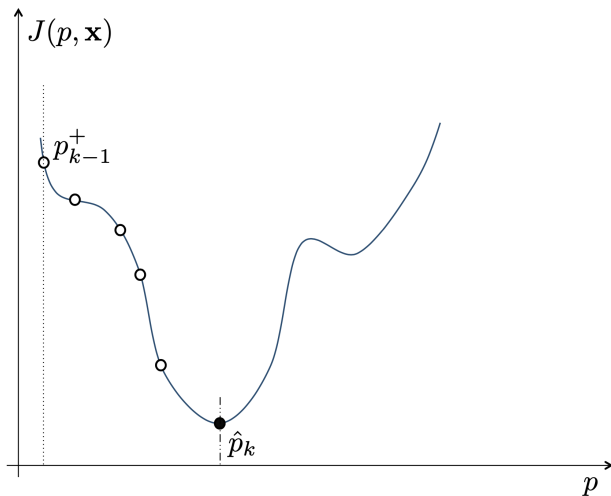
Sufficient Conditions of Success



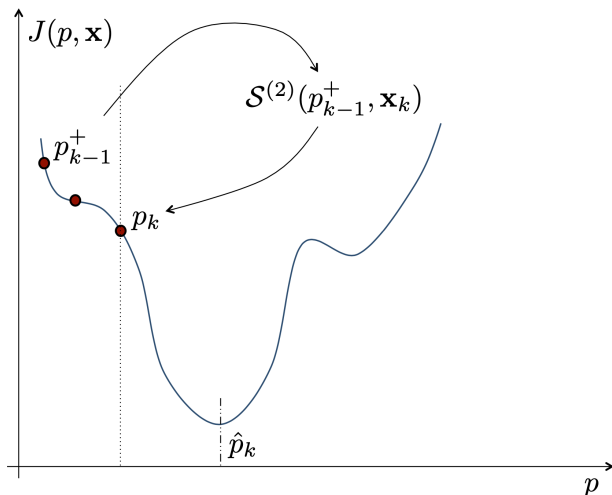
Sufficient Conditions of Success



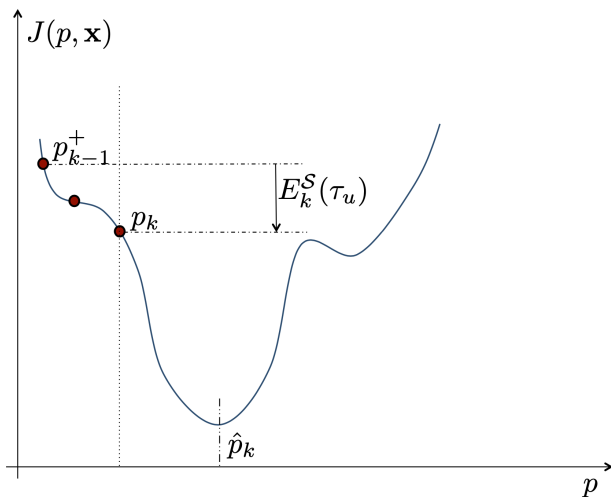
Closed-Loop Evolution of the Cost Function



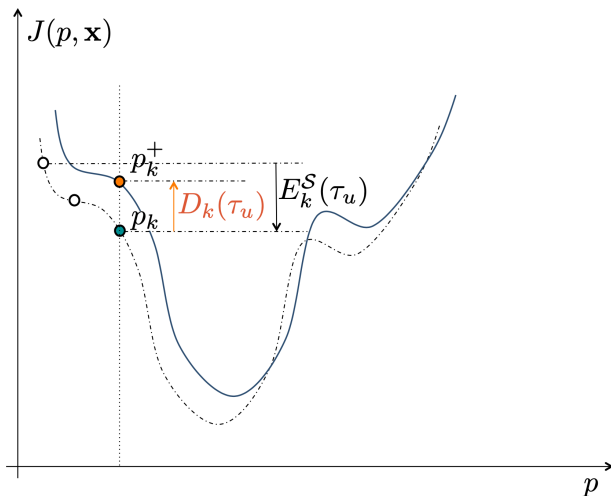
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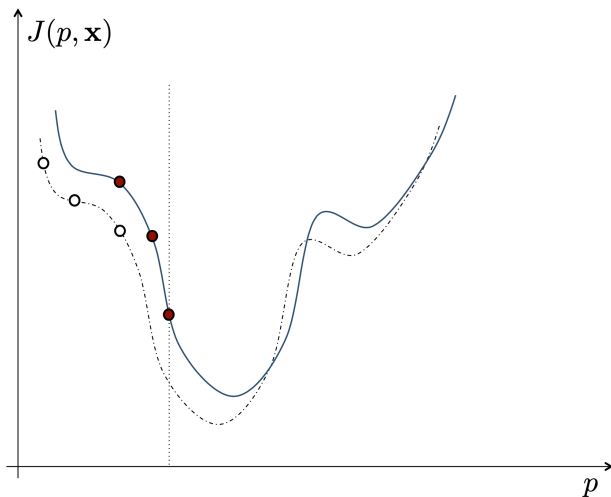
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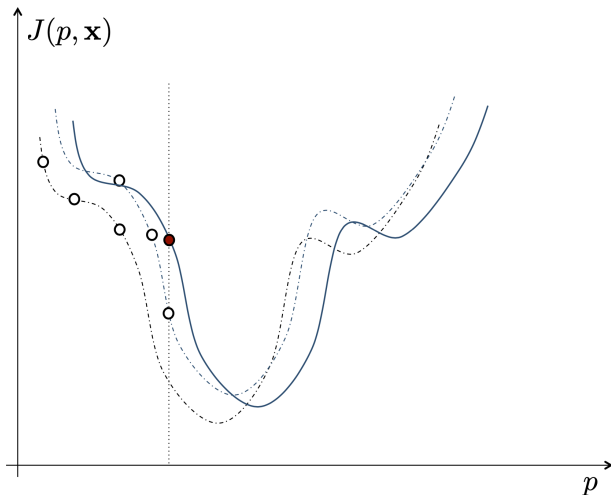
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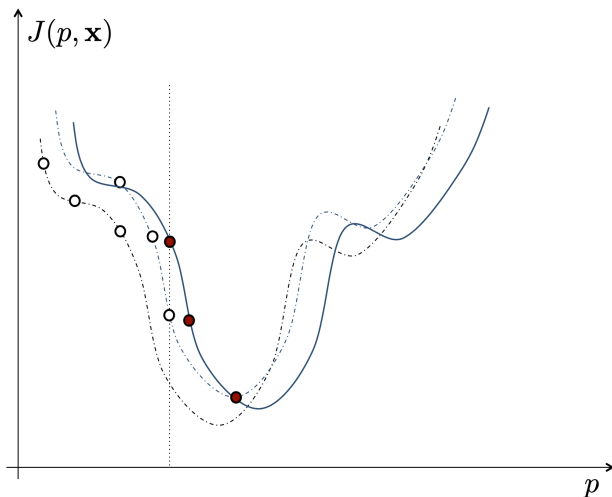
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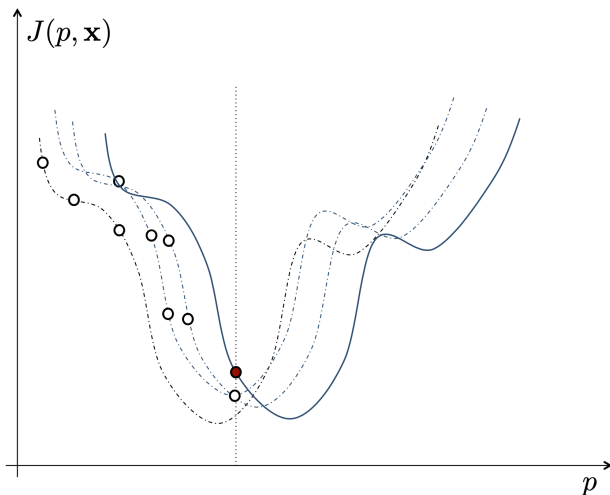
Closed-Loop Evolution of the Cost Function



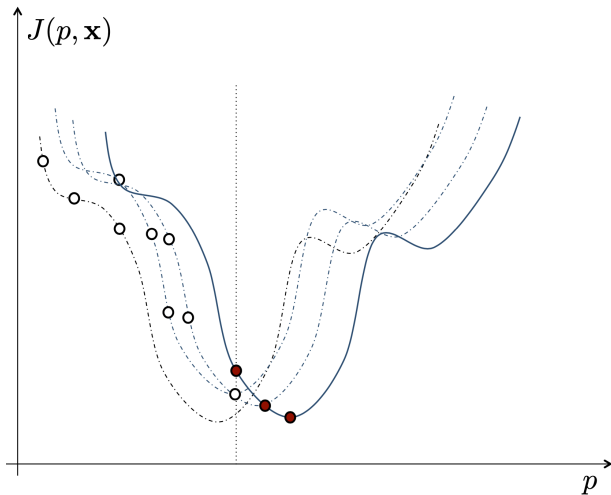
Closed-Loop Evolution of the Cost Function



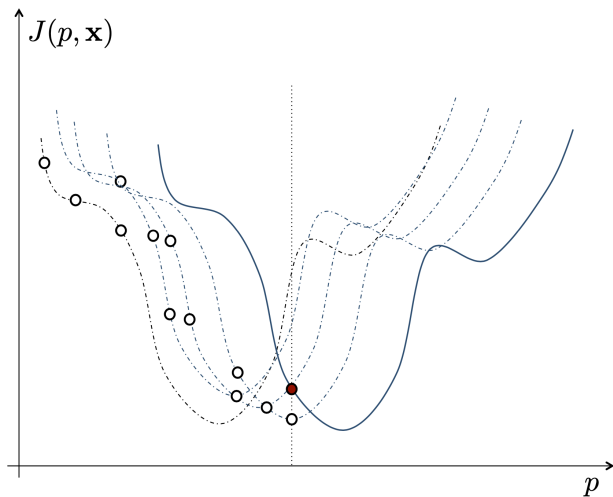
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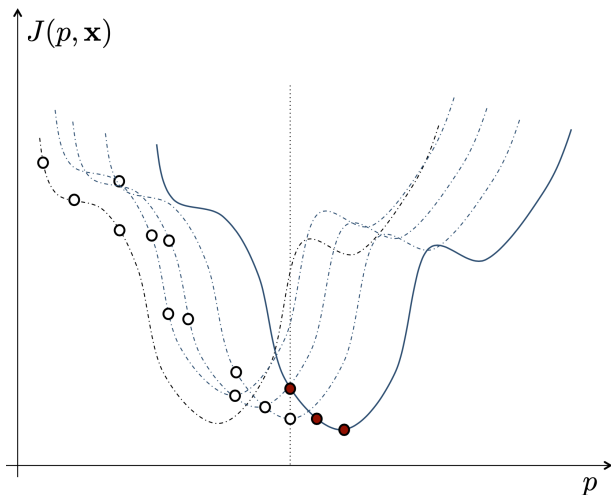
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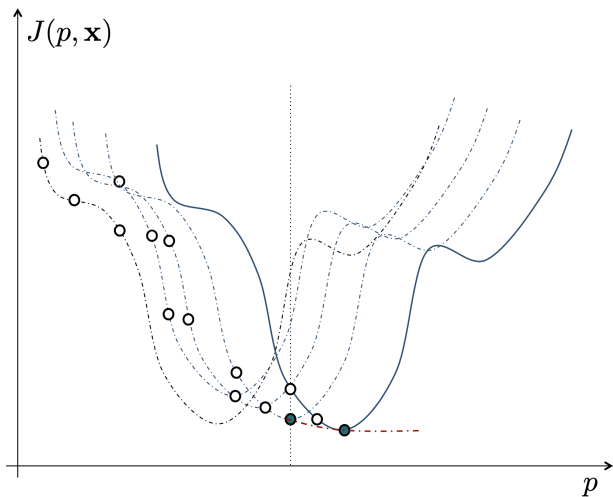
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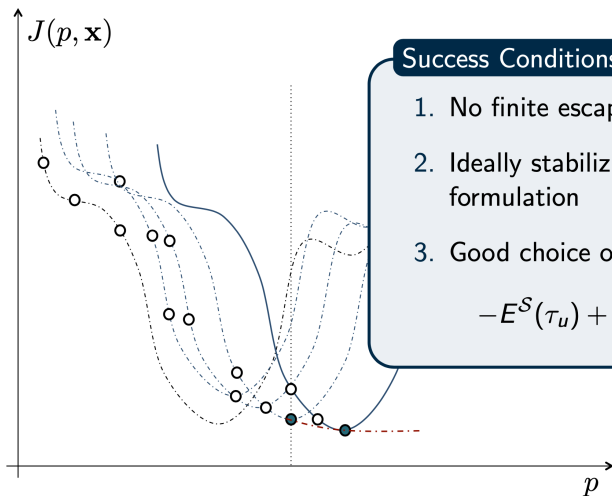
Closed-Loop Evolution of the Cost Function



Closed-Loop Evolution of the Cost Function



Closed-Loop Evolution of the Cost Function



Success Conditions

1. No finite escape time
2. Ideally stabilizing NMPC formulation
3. Good choice of (\mathcal{S}, τ_u) :

$$-E^{\mathcal{S}}(\tau_u) + D(\tau_u) < 0$$

Reminder

$$D(\tau_u) - E^S(\tau_u) < 0$$

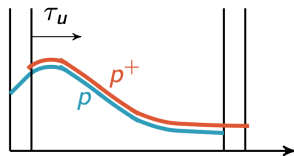
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 $D(\tau_u)$ $-E^S(\tau_u)$

Reminder

$$D(\tau_u) - E^{\mathcal{S}}(\tau_u) < 0$$

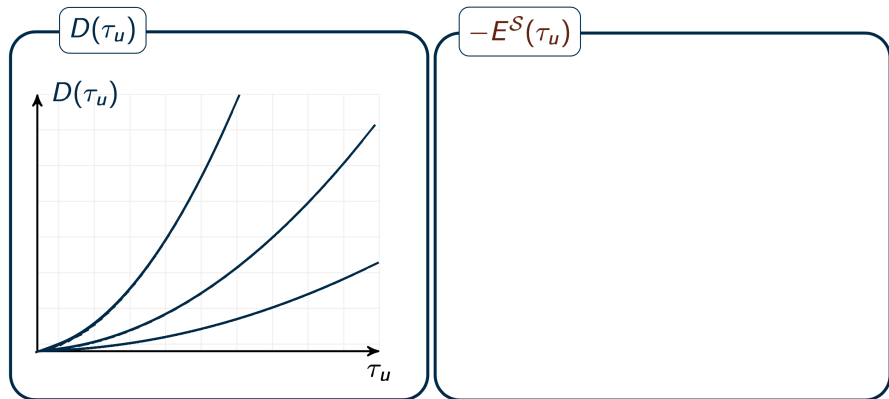
 $D(\tau_u)$


- ▶ $D(\tau_u) := J(p^+, x^+) - J(p, x)$
- ▶ $D(0) = 0$, $D(\tau_u)$ can be ≥ 0
- ▶ $\tau_u \in [0, \tau_u^{max}]$
- ▶ Independent of the solver \mathcal{S}

 $-E^{\mathcal{S}}(\tau_u)$

Reminder

$$D(\tau_u) - E^S(\tau_u) < 0$$



Reminder

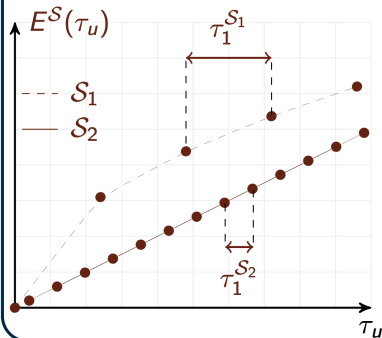
$$D(\tau_u) - E^S(\tau_u) < 0$$

 $D(\tau_u)$
 $-E^S(\tau_u)$

- ▶ $E^S(\tau_u) := J(p^{(0)}, \mathbf{x}) - J(p^{(q^S)}, \mathbf{x})$
- ▶ $q^S = \text{int}(\tau_u / \tau_1^S)$
- ▶ τ_1^S time for a single iteration
- ▶ $\tau_u \in \{1, 2, \dots\} \times \tau_1^S$
- ▶ $E^S(0) = 0$

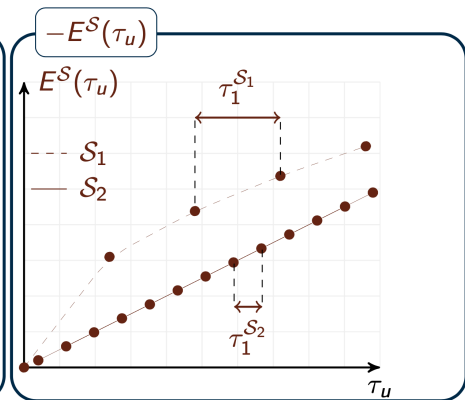
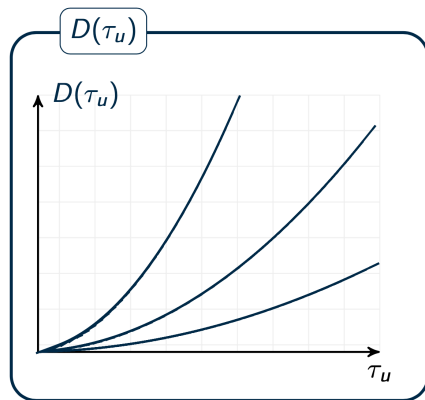
Reminder

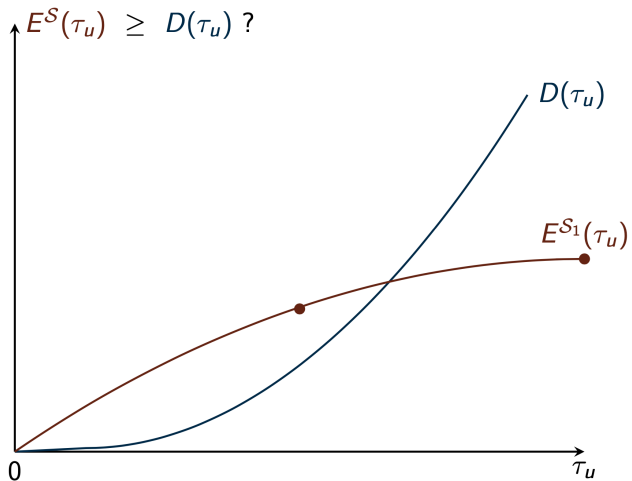
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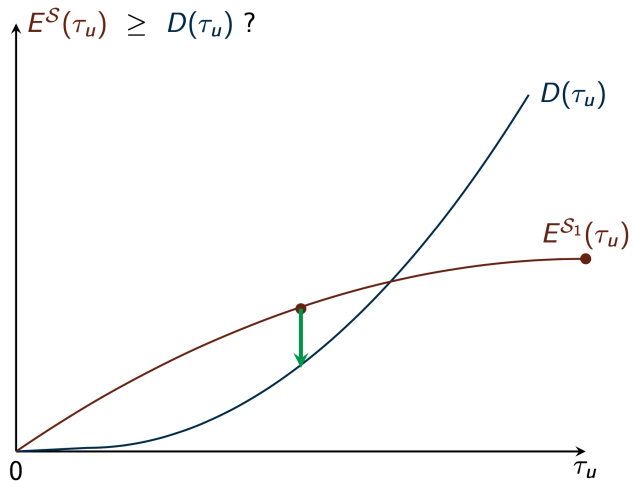
 $D(\tau_u)$
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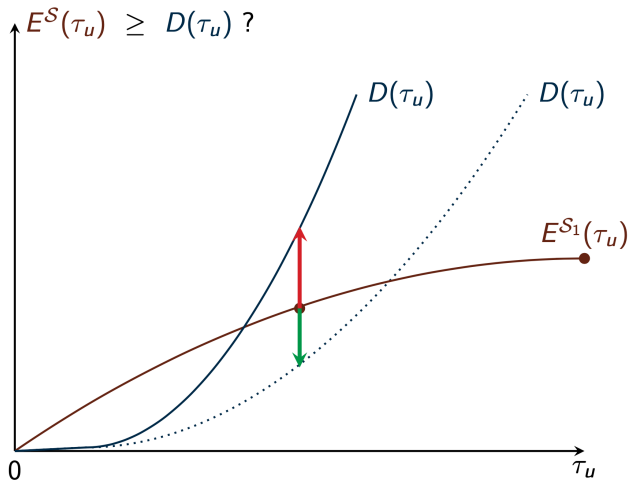
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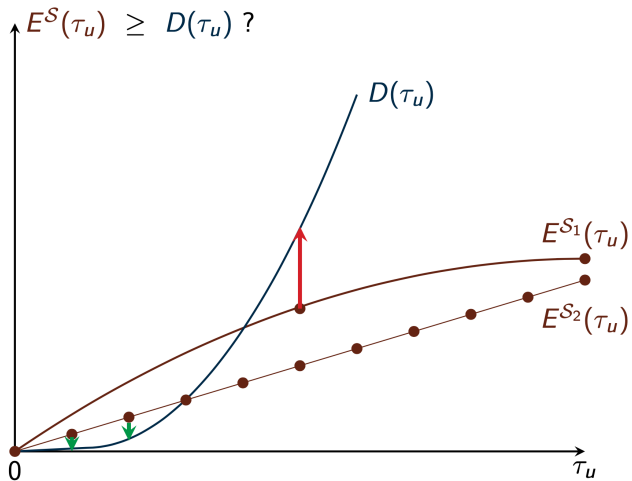
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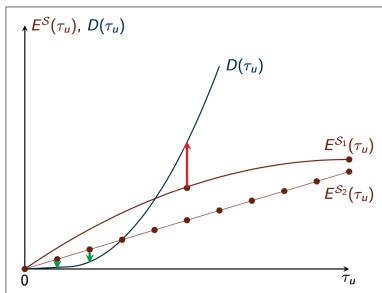








Key properties of a solver

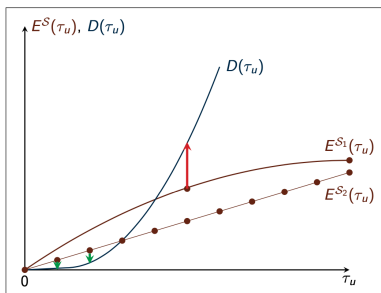


Keep in Mind

It is sometimes **better** to choose a **less efficient^a** solver with **shorter preparation step** duration τ_1 .

^aper iteration

Key properties of a solver



Keep in Mind

It is sometimes **better** to choose a **less efficient**^a solver with **shorter preparation step** duration τ_1 .

^aper iteration

Gradient-based studies



Bemporad and Patrino (2012), Jones et al. (2012), MA (2013).

Heuristics for second order methods



Bock et al. SIAM (2007)

Comparing Algorithms: An Example



Houska et al. Automatica (2011)

Model

$$\dot{x}_1 = x_2$$

$$\dot{x}_2 = u \quad u \in [-1, +1]$$

$$\dot{x}_3 = x_4$$

$$\dot{x}_4 = -g \sin x_3 - u \cos x_3 - bx_4$$

Comparing Algorithms: An Example



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Cost

$$\sum_{i=1}^{N_p} \|\mathbf{x}_{k+i}^T\|_Q^2 + \|\mathbf{u}_{k+i-1}^T\|_R^2$$

Comparing Algorithms: An Example



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Cost

$$\sum_{i=1}^{N_p} \|\mathbf{x}_{k+i}^T\|_Q^2 + \|\mathbf{u}_{k+i-1}^T\|_R^2$$

▶ $N_p = 10$ and $T_p = 3$ sec

▶ Two solvers:

\mathcal{S}_1 Full gradient

\mathcal{S}_2 Partial gradient

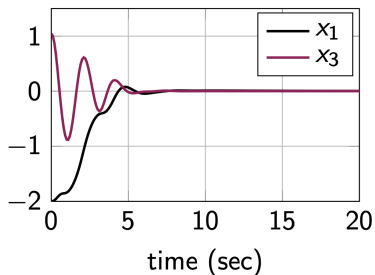
▶ $\tau_1^{\mathcal{S}_2} = \left[\frac{4}{13}\right] \times \tau_1^{\mathcal{S}_1}$

▶ $\tau_u(\mathcal{S}_2) = \left[\frac{4}{13}\right] \times \tau_u(\mathcal{S}_1)$

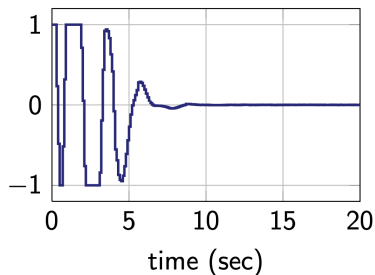
▶ $\tau_u(\mathcal{S}_1) \in \{5, 50, 100\}$ ms

Typical Closed-Loop Behavior

Closed-Loop State Trajectory

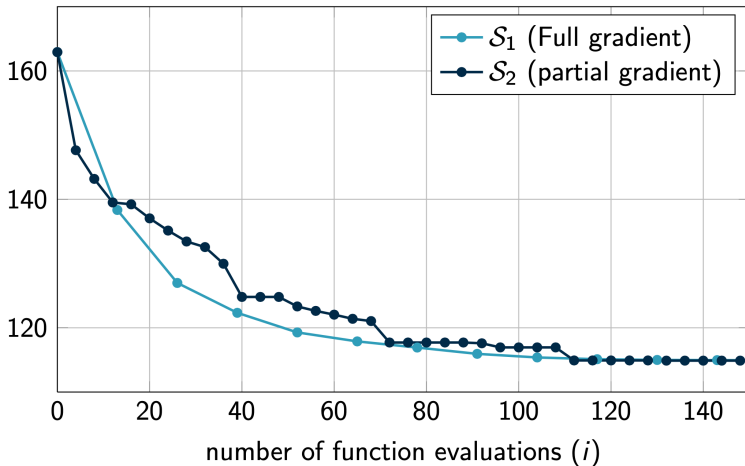


Closed-Loop Control Trajectory



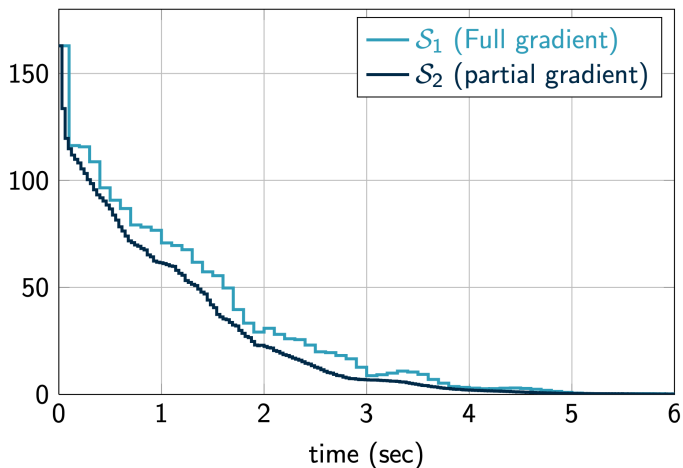
\mathcal{S}_1 is more efficient than \mathcal{S}_2 ... **per iteration** !

$$J(p^{(i)}, \mathbf{x}_0)$$



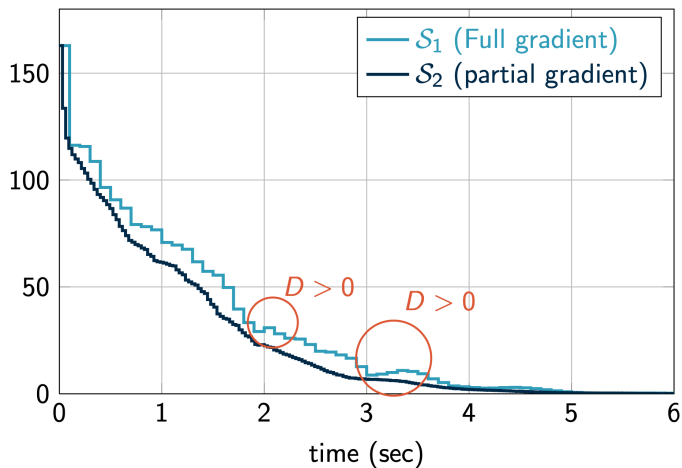
Closed-Loop Evolutions of the Predicted Cost

$$J(p_k, \mathbf{x}_k)$$



Closed-Loop Evolutions of the Predicted Cost

$$J(p_k, \mathbf{x}_k)$$



Comparison of Effective Closed-Loop Cost

$$J_{cl} := \frac{1}{N_{sim}} \sum_{k=1}^{N_{sim}} \left[\|\mathbf{x}_k\|_Q^2 + \|u_{k-1}\|_R^2 \right]$$

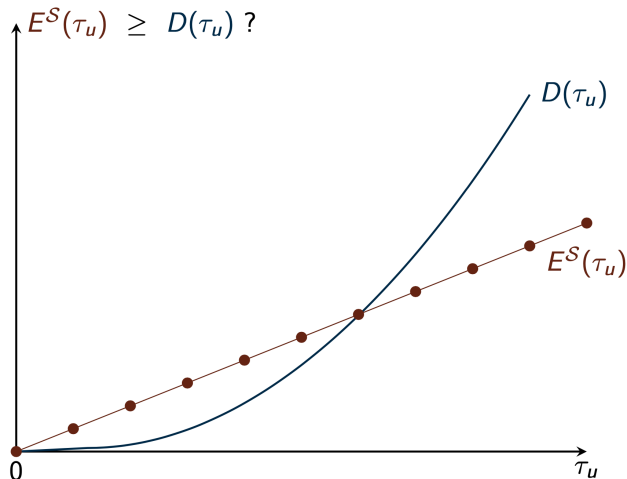
$\tau_u(\mathcal{S}_1)$	100 ms	50 ms	5 ms
Full Gradient	0.728	0.430	0.027
Partial Gradient	0.300	0.110	0.008
Gain %	59%	74%	70%

Updating Scheme For a Given Solver

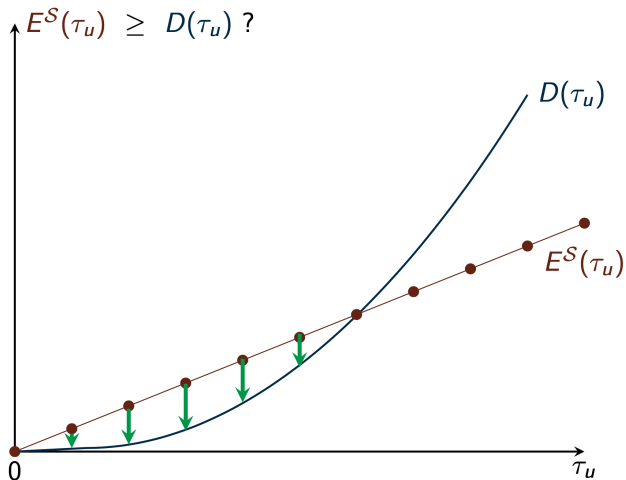
Assume that a solver \mathcal{S} has been chosen . . .

Is there any remaining choice ?

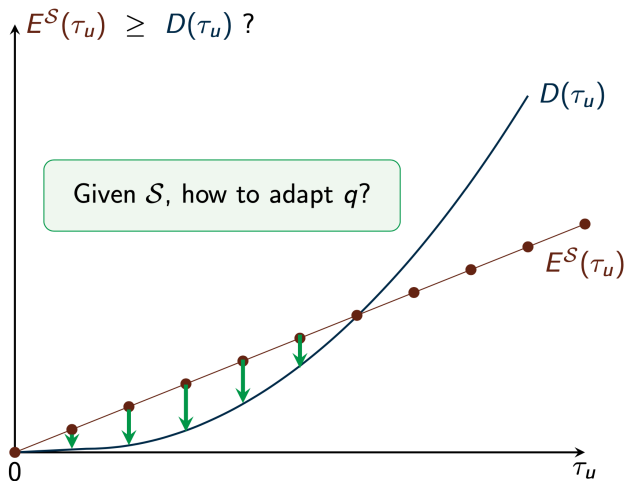
What is the optimal τ_u for a given solver?



What is the optimal τ_u for a given solver?



What is the optimal τ_u for a given solver?



Updating τ_u is a control problem ...!

$$p_{k+1} = \mathcal{S}^{(q(\tau_u))}(p_k^+, x_k)$$

Updating τ_u is a control problem ...!

$$\begin{aligned} p_{k+1} &= \mathcal{S}^{(q(\tau_u))}(p_k^+, \mathbf{x}_k) \\ \mathbf{x}_{k+1} &= f(\mathbf{x}_k, \mathcal{U}(0, p_k)) \end{aligned}$$

Updating τ_u is a control problem ...!

$$\begin{pmatrix} p_{k+1} \\ \mathbf{x}_{k+1} \end{pmatrix} = F\left(\begin{pmatrix} p_k \\ \mathbf{x}_k \end{pmatrix}, \tau_u\right)$$

Updating τ_u is a control problem ...!

$$\begin{aligned} \begin{pmatrix} p_{k+1} \\ \mathbf{x}_{k+1} \end{pmatrix} &= F\left(\begin{pmatrix} p_k \\ \mathbf{x}_k \end{pmatrix}, \tau_u\right) \\ y &= J(p_k, \mathbf{x}_k) \end{aligned}$$

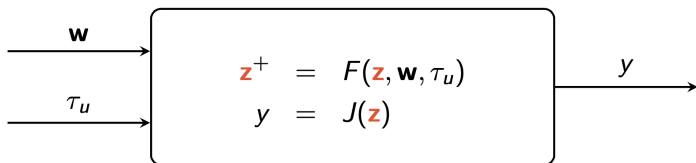
Updating τ_u is a control problem ...!

$$\begin{aligned} \mathbf{z}^+ &= F(\mathbf{z}, \tau_u) \\ y &= J(\mathbf{z}) \end{aligned}$$

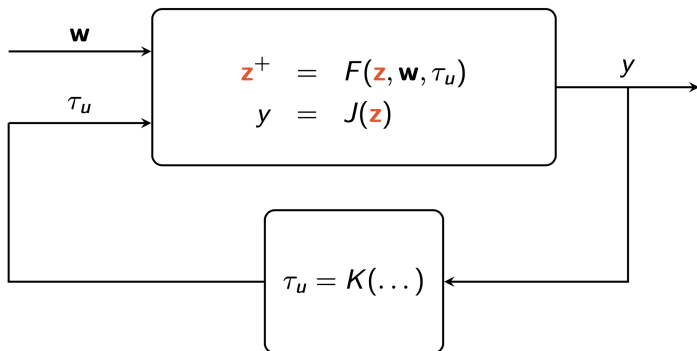
Updating τ_u is a control problem ...!

$$\begin{aligned} \mathbf{z}^+ &= F(\mathbf{z}, \mathbf{w}, \tau_u) \\ y &= J(\mathbf{z}) \end{aligned}$$

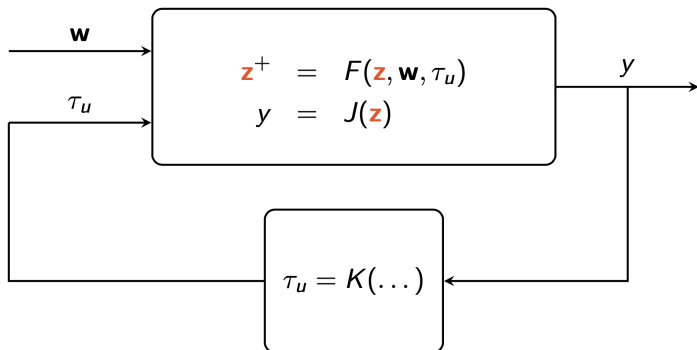
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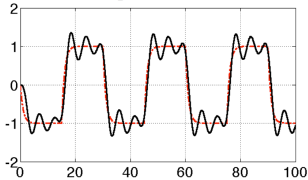
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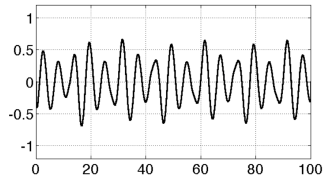
M.A. ECC (2013)

$q = 2$ without adaptation

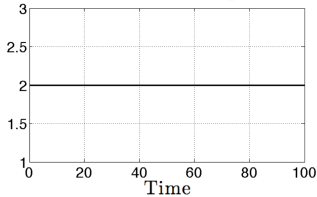
Output Evolution



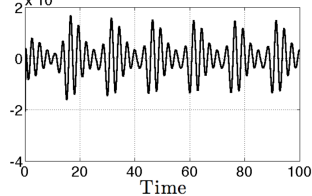
Control Evolution



Evolution of q



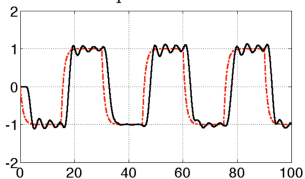
Evolution of α_D



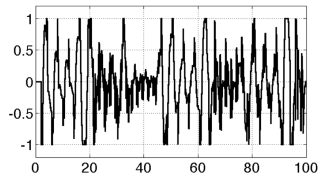
MA, ECC (2013)

$q = 100$ without adaptation

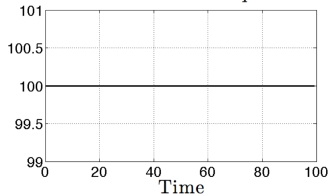
Output Evolution



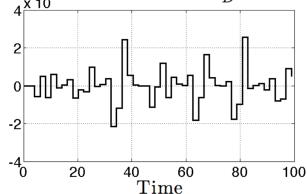
Control Evolution



Evolution of q



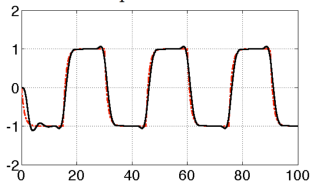
Evolution of α_D



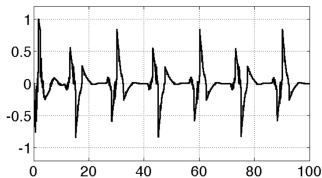
MA, ECC (2013)

$q^{(0)} = 2$ with adaptation

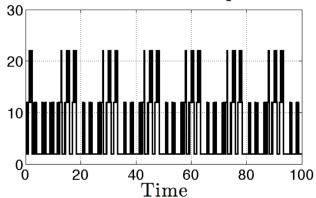
Output Evolution



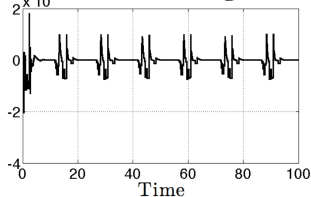
Control Evolution



Evolution of q



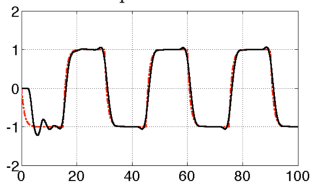
Evolution of α_D



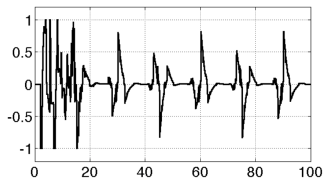
MA, ECC (2013)

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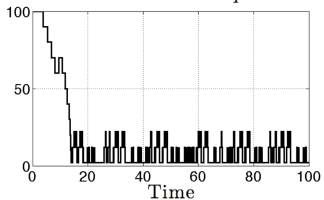
Output Evolution



Control Evolution



Evolution of q



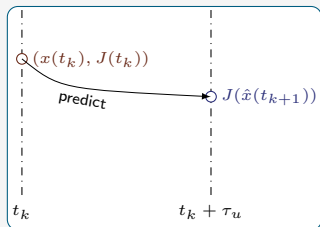
Evolution of α_D



MA, ECC (2013)

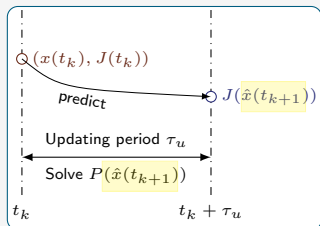
MONITORING CONTROL UPDATING PERIOD

New ML-induced opportunities (preliminary)



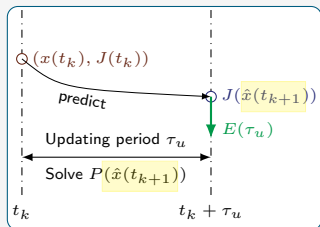
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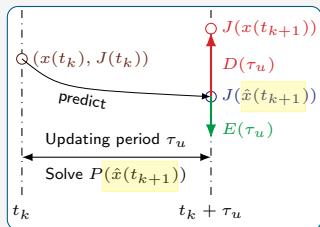
MONITORING CONTROL UPDATING PERIOD

New ML-induced opportunities (preliminary)



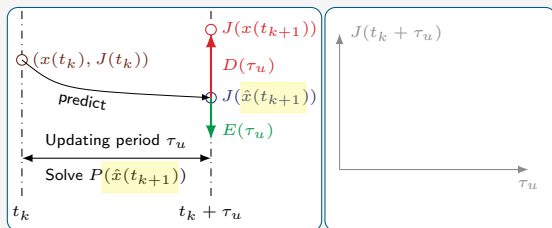
MONITORING CONTROL UPDATING PERIOD

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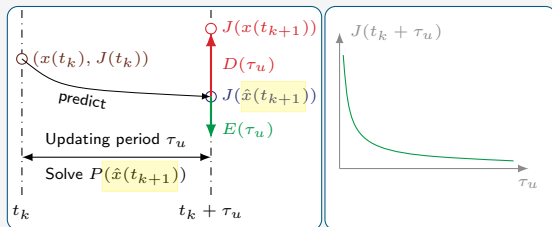
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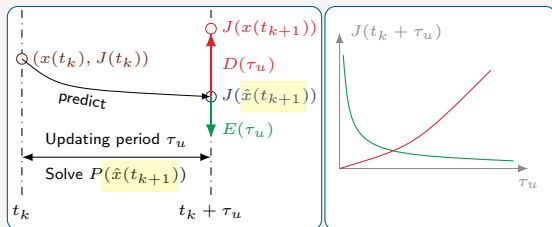
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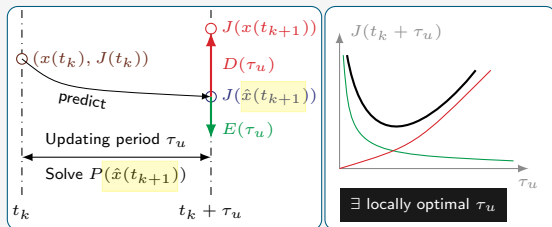
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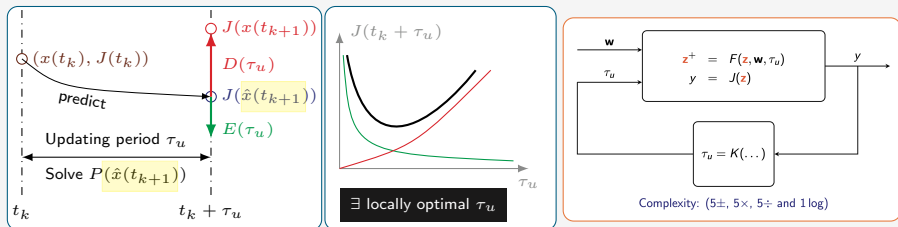
MONITORING CONTROL UPDATING PERIOD

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MONITORING CONTROL UPDATING PERIOD

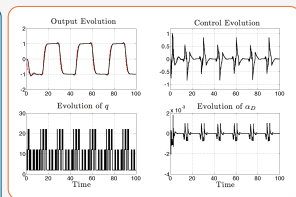
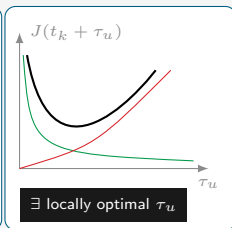
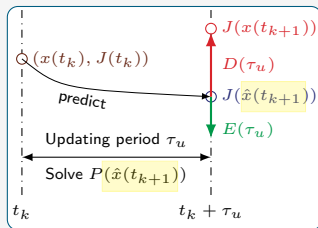
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M.A. Monitoring Control Updating Period In Fast Gradient-Based NMPC. ECC2013, 2013

MONITORING CONTROL UPDATING PERIOD

New ML-induced opportunities (preliminary)

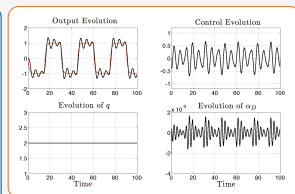
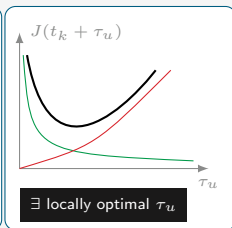
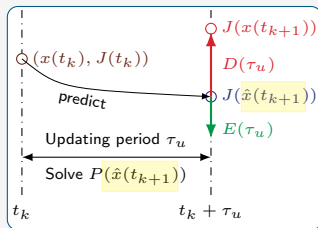


$\tau_u := q \times \text{cpu of a single iteration}$

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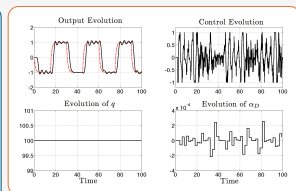
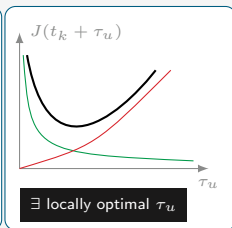
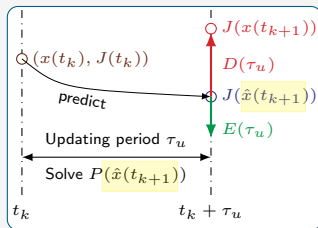


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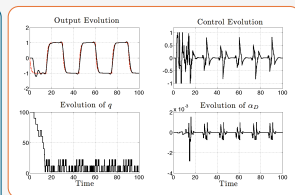
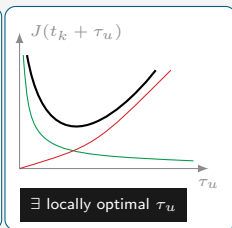
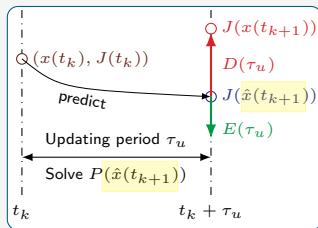


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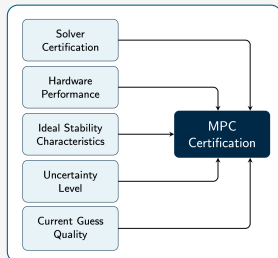
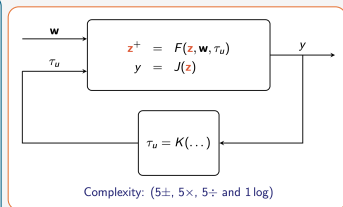
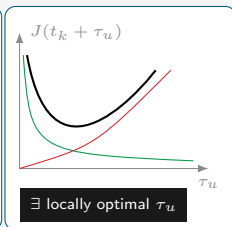
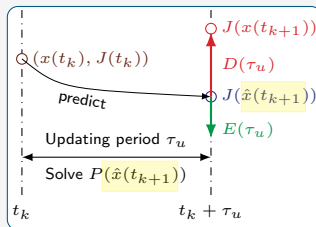


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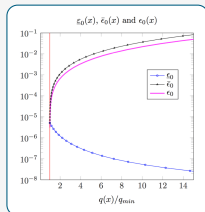
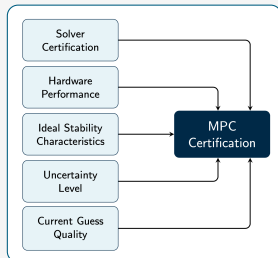
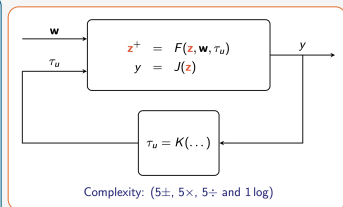
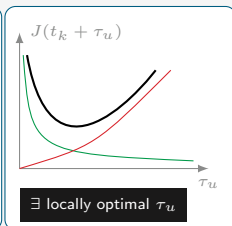
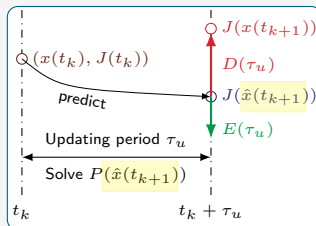
New ML-induced opportunities (preliminary)



M. Alamir. A State-Dependent Updating Period For Certified Real-Time MPC. IEEE TAC, 2017.

MONITORING CONTROL UPDATING PERIOD

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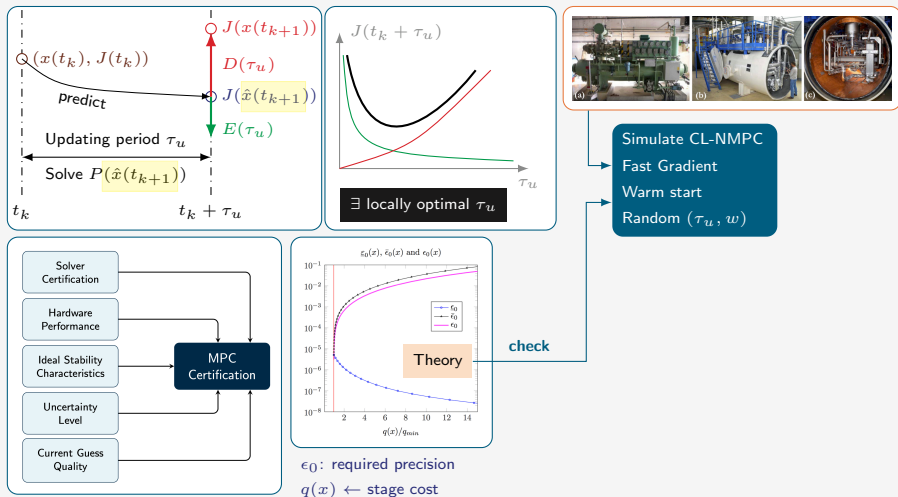


ϵ_0 : required precision
 $q(x) \leftarrow$ stage cost

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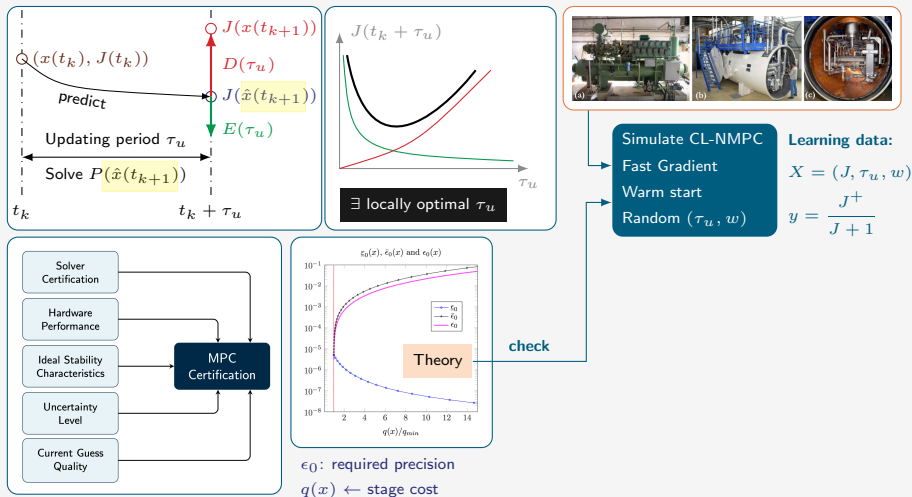
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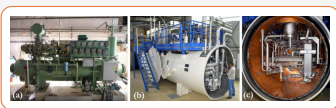
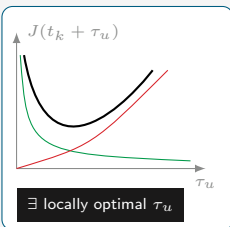
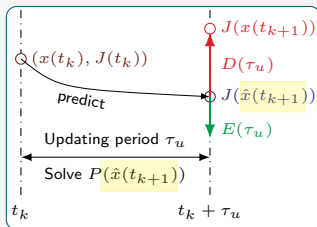
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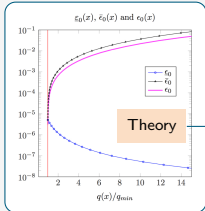
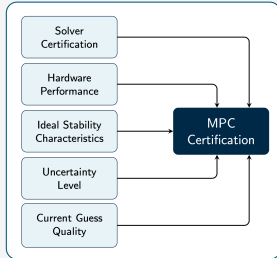
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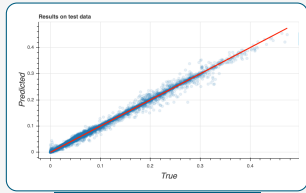


Simulate CL-NMPC
Fast Gradient
Warm start
Random (τ_u, w)

Learning data:
 $X = (J, \tau_u, w)$
 $y = \frac{J^+}{J+1}$



ϵ_0 : required precision
 $q(x) \leftarrow$ stage cost

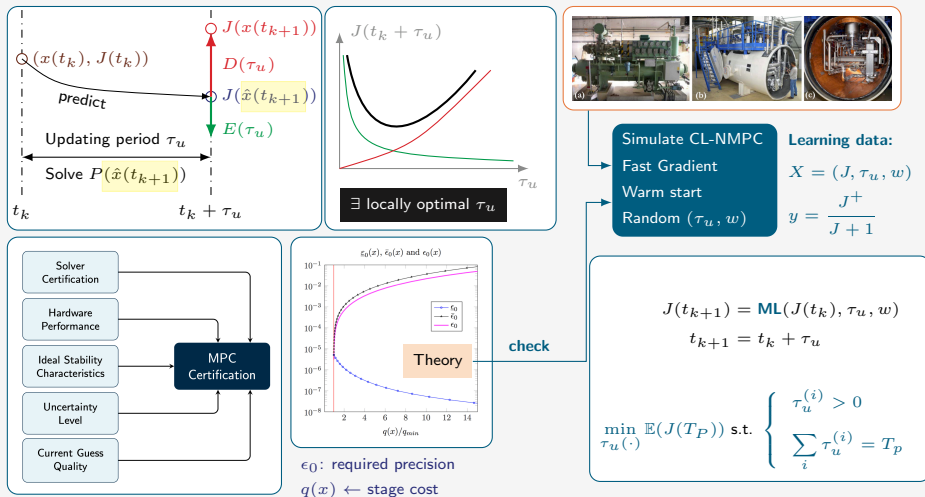


$$J^+ = \text{ML}(J(x), \tau_u, w)$$

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MONITORING CONTROL UPDATING PERIOD

New ML-induced opportunities (preliminary)



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MONITORING CONTROL UPDATING PERIOD

New ML-induced opportunities (preliminary)

$$J(t_{k+1}) = \mathbf{ML}(J(t_k), \tau_u, w)$$

$$t_{k+1} = t_k + \tau_u$$

$$\min_{\tau_u(\cdot)} \mathbb{E}(J(T_P)) \text{ s.t. } \begin{cases} \tau_u^{(i)} > 0 \\ \sum_i \tau_u^{(i)} = T_P \end{cases}$$

Approximate Stochastic
Dynamic Programming

$$\tau_u(t_k) := \tau^*(J(t_k) | \mathcal{W})$$

Option3 (Approximate Dynamic Programming)

Approximate the optimal value

$$Q(z) \text{ where } z := (x, u)$$

based on the Bellman equation

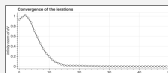
$$Q(z) = \ell(z) + \gamma \min_v [\hat{\mu} + \alpha \hat{\sigma}^{\frac{1}{2}}] [Q(f(z), v)]$$

Fixed-point

- ℓ stage cost
- γ discounting rat
- $\hat{\mu}$ approximated statistics over realisations of w



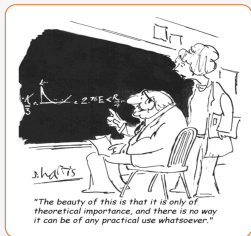
- Choose a ML-based structure for Q
- Fixed-point iteration on Q



[MA]. Explicit approximation of stochastic optimal control for combined therapy of cancer. arXiv:1905.04937, 2019.

- (-) Curse of dimensionality
- (-) Fragile choice of the support state domain

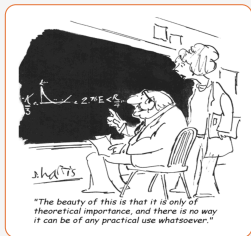
CONCLUSION



US



CONCLUSION

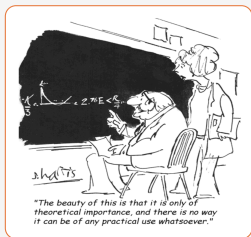


US

Machine Learning



CONCLUSION



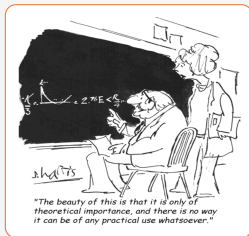
US

Computational technologies
GPU, FPGA, Containers, ...

Machine Learning



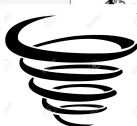
CONCLUSION



US

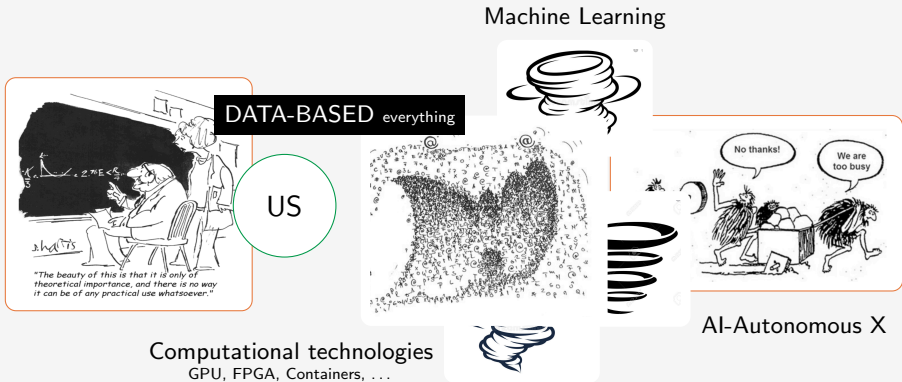
Computational technologies
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Machine Learning



AI-Autonomous X

CONCLUSION



CONCLUSION

Business as usual is not an option!

