

Model Documentation of the 'Chemical reactor model by'

1 Nomenclature

1.1 Nomenclature for Model Equations

- x state vector
- u control input vector
- w noise vector
- z regulated output vector
- y measurement vector

2 Model Equations

State Vector and Input Vector:

$$x \in \mathbb{R}^8 u \in \mathbb{R}^1 w \in \mathbb{R}^1 z \in \mathbb{R}^1 y \in \mathbb{R}^1$$

System Equations:

$$\dot{x}(t) = Ax(t) + B_1w(t) + Bu(t) \quad (1a)$$

$$z(t) = C_1x(t) + D_{11}w(t) + D_{12}u(t) \quad (1b)$$

$$y(t) = Cx(t) + D_{21}w(t) \quad (1c)$$

Outputs: z

2.1 Exemplary parameter values

Symbol	Value							
A	0.5623	-0.01642	0.01287	-0.0161	0.02094	-0.02988	0.0183	0.008743
	0.102	0.6114	-0.02468	0.02468	-0.03005	0.04195	-0.02559	0.03889
	0.1361	0.2523	0.641	-0.03404	0.03292	-0.04296	0.02588	0.08467
	0.09951	0.2859	0.3476	0.6457	-0.03249	0.03316	-0.01913	0.1103
	-0.04794	0.08708	0.3297	0.3102	0.6201	-0.03015	0.01547	0.08457
	-0.1373	-0.1224	0.1705	0.3106	0.191	0.5815	-0.01274	0.05394
	-0.1497	-0.1692	0.1165	0.2962	0.1979	0.07631	0.5242	0.04702
	0	0	0	0	0	0	0	0.6065
B	-0.1774							
	-0.2156							
	-0.2194							
	-0.09543							
	0.0579							
	0.09303							
	0.08962							
	0							
B_1	-0.1774							
	-0.2156							
	-0.2194							
	-0.09543							
	0.0579							
	0.09303							
	0.08962							
	0							
C_1	-0.0465	-0.1135	-0.1909	-0.2619	-0.2634	-0.1422	-0.0002	0.1856
C	-0.0049	0.0049	-0.006	0.01	0.0263	0.3416	0.6759	0
D_{11}	[0]							
D_{12}	[0.1001]							
D_{21}	[1.0]							

3 Derivation and Explanation

This model is part of the "COMPleib" - library and was automatically imported into ACKREP.

The original description was:

REA4 Chemical reactor model by P. M. Maekilae, "Parametric LQ Control", IJOC, Vol. 41, Nr. 6, pp. 1413-1428, 1985 discrete modell

4 Simulation

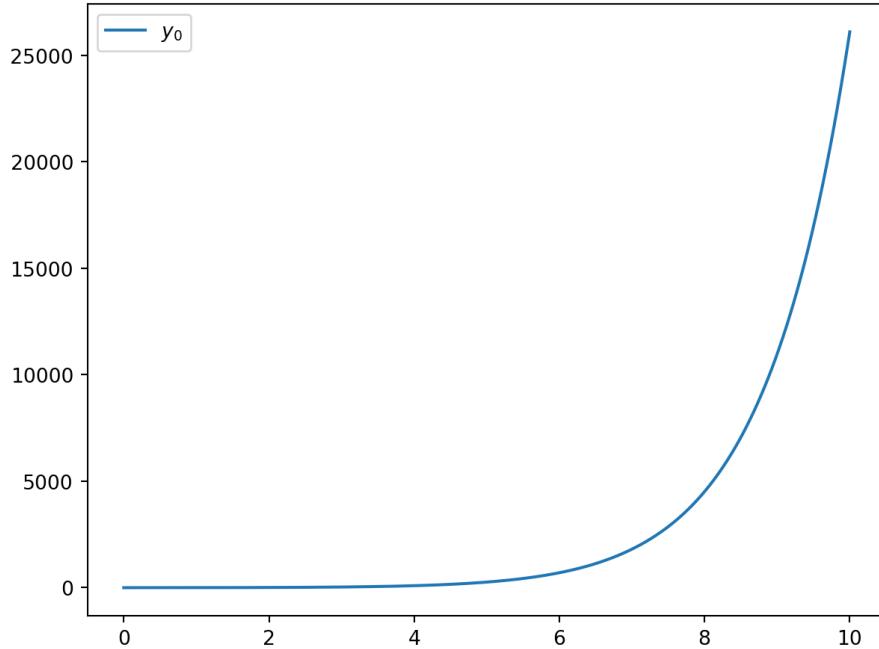


Figure 1: Simulation of the Chemical reactor model by.

References

- [1] . M. Maekilae, "Parametric LQ Control", IJOC, Vol. 41, Nr. 6, pp. 1413-1428, 1985 discrete modell