

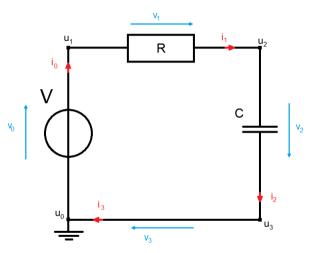
Circuit Modelling



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Introducing example: Charging of a capacitor:





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Formulating a Mathematical Model



Network Topology



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Define the incidence matrix $A = (a_{ij}) \in \mathbb{R}^{k \times l}$:

$$\tilde{\alpha}_{ij} = \begin{cases} 1 & \text{edge } j \text{ starts at node } i, \\ -1 & \text{edge } j \text{ ends at node } i, \\ 0 & \text{else.} \end{cases}$$

With

$$\begin{split} \mathcal{N} &= (n_0, n_1, n_2, ..., n_k) & \cdots & \text{nodes,} \\ \mathcal{E} &= \{e_j: j=1, ..., l\} & \cdots & \text{edges,} \end{split}$$

furthermore

$$\mathfrak{u}=(\mathfrak{u}_0,\mathfrak{u}_1,\mathfrak{u}_2,...)$$
 ... corresponding electrical potentials at the nodes.

By grounding node 0, i.e. $u_0=0$ we obtain the reduced incidence matrix.



Formulating a Mathematical Model



Energy Conservation Laws



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• Kirchhoff's voltage law (KVL):

The sum of voltages along each loop of the network must equal to zero.

$$\to A^{\top} u = v. \tag{1}$$

• Kirchhoff's current law (KCL):

For any node, the sum of currents flowing into the node is equal to the sum of currents flowing out of the node.

$$\rightarrow Ai = 0. (2)$$



Formulating a Mathematical Model



Electrical Components and their Relations



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Resistor

$$v = R i$$
 or $i = G u$. (3)

Figure: resistor symbol

Capacitor

$$Q = C v$$
 and by derivation in t $I = C \frac{d}{dt} v = C v'$. (4)



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Inductor (Coil)

$$\Phi = L \ i \quad \text{and by derivation in } t \quad \nu = L \ i'. \tag{5}$$



Figure: inductor symbol

Voltage Source

$$v = v_{\rm src} \tag{6}$$



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• Current Source

$$i = i_{src}$$
 (7

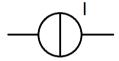


Figure: current source symbol

Formulating a Mathematical Model



Modified Nodal Analysis - MNA



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Rearrange the columns of the reduced incidence matrix A into

$$A = (A_R A_C A_L A_V A_I)$$

 A_R , A_C , A_L , A_V and A_I ... columns related to components Represent voltages:

$$v = A^{\mathsf{T}} u$$

 \rightarrow rearrange ν into $\nu = (\nu_R, \nu_C, \nu_L, \nu_{src}, \nu_I)$ and i into $i = (i_R, i_C, i_L, i_V, i_src)$. Rewrite component relations:

$$i_R = G v_R = G A_R^\top u,$$

 $i_C = C v_C' = C A_C^\top u'.$

Kirchhoffs current law:

$$A_C i_C + A_R i_R + A_L i_L + A_V i_V = -A_I i_{src}.$$

Combine:

$$A_{c}CA_{c}^{\dagger}II' + A_{b}GA_{b}^{\dagger}II + A_{t}I_{t} + A_{t}I_{t} = -A_{t}I_{c}$$



Together with component law for inductors (5) and potential-voltage relation for voltage sources (6):

$$\begin{split} A_{C}CA_{C}^{\top}u' + A_{R}GA_{R}^{\top}u + A_{L}i_{L} + A_{V}i_{V} &= -A_{I}i_{src}, \\ Li'_{L} - A_{L}^{\top}u &= 0, \\ -A_{V}^{\top}u &= -\nu_{src}. \end{split}$$

In matrix form:

$$\begin{pmatrix} A_{C}CA_{C}^{\top} & 0 & 0 \\ 0 & L & 0 \\ 0 & 0 & 0 \end{pmatrix} * \begin{pmatrix} u' \\ i'_{L} \\ i'_{V} \end{pmatrix} + \begin{pmatrix} A_{R}GA_{R}^{\top} & A_{L} & A_{V} \\ -A_{L}^{\top} & 0 & 0 \\ -A_{V}^{\top} & 0 & 0 \end{pmatrix} * \begin{pmatrix} u \\ i_{L} \\ i_{V} \end{pmatrix} = \begin{pmatrix} -A_{I}i_{src} \\ 0 \\ -\nu_{src} \end{pmatrix}.$$
(8)



Differential Algebraic Equations



Types of DAEs



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In the most general form a DAE can be written as: Find $y: \mathbb{R} \to \mathbb{R}^n$ such that

$$F(t, y(t), y'(t)) = 0, \qquad \forall t \in I$$
(9)

with $F: \mathbb{R} \times \mathbb{R}^n \times \mathbb{R}^n \to \mathbb{R}^n$ sufficiently smooth and I the time-interval.

• Linear systems with constant coefficcients

are systems of the form: find y such that

$$Ay'(t) + By(t) = f(t),$$
(10)

with $A, B \in \mathbb{R}^{n \times n}$, A singular, B regular and $f : \mathbb{R} \to \mathbb{R}^n$ a function in time.

• Linear time dependent systems are systems of the form: find y such that

$$A(t)y'(t) + B(t)y(t) = f(t),$$

with $A, B : \mathbb{R} \to \mathbb{R}^{n \times n}$, $f : \mathbb{R} \to \mathbb{R}^n$ functions, such that for every $t \in \mathbb{R}$ the matrix A(t) is singular and the matrix B(t) regular.

Structured (non-linear) systems
 are semi-explicit systems of the form: find (y, z) such that

$$y'(t) = f(t, y(t), z(t)),$$
 (11)

$$0 = g(t, y(t), z(t)),$$
 (12)

with $f: \mathbb{R} \to \mathbb{R}^n$ and $g: \mathbb{R} \to \mathbb{R}^d$ functions.



Differential Algebraic Equations



Weierstrass-Kronecker Normalform



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prerequisites:

Definition

The matrix pencil $\{A,B\}$ is called *regular* if there exists some $c \in \mathbb{R}$, such that (cA+B) is regular (i.e. $det(cA+B) \neq 0$), otherwise it is called singular.

Theorem

For every matrix $Q \in \mathbb{R}^{n \times n}$ there exists a regular matrix $T \in \mathbb{C}^{n \times n}$, such that

$$T^{-1}QT = J = \text{diag}(J_1,...,J_r) \quad \textit{with} \quad J_i = \begin{pmatrix} \lambda_i & 1 & & 0 \\ 0 & \lambda_i & \ddots & \vdots \\ & \ddots & \ddots & 1 \\ 0 & \dots & 0 & \lambda_i \end{pmatrix} \in \mathbb{C}^{m_i \times m_i}$$

and $n = m_1 + ... + m_r$.



Theorem

Let $\{A, B\}$ be a regular matrix pencil. There exist $P, Q \in \mathbb{C}^{n \times n}$ such that

$$PAQ = \begin{pmatrix} I_d & 0 \\ 0 & N \end{pmatrix}, PBQ = \begin{pmatrix} R & 0 \\ 0 & I_{n-d} \end{pmatrix}$$

where

$$N = diag(N_1, ..., N_r)$$
 with $N_i = \begin{pmatrix} 0 & 1 & & 0 \\ & \ddots & \ddots & \\ & & & 0 & 1 \\ 0 & & & 0 \end{pmatrix} \in \mathbb{R}^{n_i \times n_i}$

and R has Jordan Normalform. By I_k we denote the identity matrix of size $k \times k$.

using these findings: Using the findings above we are able to transform the initial DAE (10) using the matrix P from Theorem 3. By multiplying with P from the left, we obtain

$$PAy'(t) + PBy(t) = Pf(t).$$

Setting

$$y(t) = Q\begin{pmatrix} u(t) \\ v(t) \end{pmatrix}, \quad Pf(t) = \begin{pmatrix} s(t) \\ q(t) \end{pmatrix},$$

with $u(t), s(t) : \mathbb{R} \to \mathbb{R}^d$ and $q(t), v(t) : \mathbb{R} \to \mathbb{R}^{n-d}$.

We get a system of the form

$$u'(t) + Ru(t) = s(t),$$

 $Nv'(t) + v(t) = q(t),$
(13)

where
$$PAQ = \begin{pmatrix} I & \\ & N \end{pmatrix}$$
 and $PBQ = \begin{pmatrix} R & \\ & I \end{pmatrix}$.



$$v(t) = q(t) - Nv'(t) = q(t) - N(\underline{q(t) - Nv'(t)})' = q - Nq' + N^{2}v''$$

$$= q - Nq' + N^{2}(q - Nv')'' = q - Nq' + N^{2}q'' - N^{3}v'''$$

$$\vdots$$

$$= q - Nq' + ... + (-1)^{k-1}N^{k-1} \underbrace{\frac{d^{k}}{dt^{k}}q}_{:=q^{(k-1)}} + (-1)\underbrace{N^{k}v^{(k)}}_{=0}$$

$$= \sum_{i=0}^{k-1} (-1)^{i}N^{i}q^{(i)}(t)$$
(14)

where k is the nilpotency index of N.

Definition



Differential Algebraic Equations



Weierstrass-Kronecker Normalform Index of a Differential Algebraic Equation



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Definition

Consider the differential algebraic equation (9) to be uniquely locally solvable and F sufficiently smooth. For a given $m \in \mathbb{N}$ consider the equations

$$F(t, y, y') = 0,$$

$$\frac{dF(t, y, y')}{dt} = 0,$$

$$\vdots$$

$$\frac{d^{m}F(t, y, y')}{dt^{m}} = 0.$$

The smallest natural number m for which the above system results in an explicit system of ordinary differential equations (ODEs), i.e. it has the form

$$y' = \phi(t, y),$$

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Definition

Let y(t) be the exact solution of *Abstract-DAE!!!!!!!!* and $\tilde{y}(t)$ be the solution of the perturbed system $F(t, \tilde{y}, \tilde{y}') = \delta(t)$. The smallest number $k \in \mathbb{N}$ such that

$$\|y(t) - \tilde{y}(t)\| \leq C \left(\|y(t_0) - \tilde{y}(t_0)\| + \sum_{j=0}^k \max_{t_0 \leq \xi \leq T} \left\| \int_{t_0}^{\xi} \frac{\mathrm{d}^j \delta}{\mathrm{d} \tau^j}(\tau) d\tau \right\| \right)$$

for all $\tilde{y}(t)$, is called the **perturbation index** of this system.



Differential Algebraic Equations



Weierstrass-Kronecker Normalform Consistent Initial Values



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index v = 0.

Case: Index v = 1.

By rewriting our system into the form

$$y'(t) = f(t, y(t), z(t)),$$

 $0 = g(t, y(t), z(t)).$

we are able to give conditions for consistent initial values. Namely y_0 and z_0 are consistent initial values for this system, if $g(t_0, y_0, z_0) = 0$ holds.

Case: Index v=2

For index-2 systems we rewrite our system into

$$y' = f(t, y(t), z(t)),$$

$$0 = g(t, y(t)).$$

Consistent initial values y_0 , z_0 for this case not only have to fulfill $g(t_0, y_0) = 0$ but also the hidden constraint $g_t(t_0, y_0) + g_u(t_0, y_0) f(t_0, y_0, z_0)$. By g_t and g_u we denote the derivative of q with respect to t or y, respectively.

Index Analysis of the Modified Nodal Analysis



General Index Analysis



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content...



Index Analysis of the Modified Nodal Analysis



Topological Conditions



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Theorem (Index conditions [shashkov_tuprints27452])

Let the matrices of the capacitances, inductances and resistances be positive definite.

 $\ker([A_R, A_C, A_V]^\top) = \{0\} \text{ and } \ker([A_C, A_V]) = \{0\}$

If

$$ker([A_R, A_C, A_V, A_L]^\top) = \{0\} \text{ and } ker(A_V) = \{0\}$$
 (15)

holds, then the MNA (8) leads to a system with index $v \le 2$.

If additionally

holds, then the system is of index $v \le 1$

If further

$$\ker(A_{\mathbf{C}}^{\top}) = \{0\}$$
 and $\dim(v_{\mathtt{src}}) = 0$

holds, then the system has index v = 0.

(16)

(17)

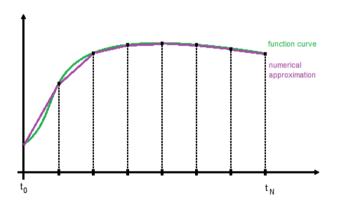
- Condition (15) can be interpreted, as the circuit neither containing loops of voltage sources nor cutsets of current sources.
- Condition (16) can be interpreted, as the circuit containing neither loops of capacitors and/or voltage sources nor cutsets of inductors and/or current sources.
- Condition (17) can be interpreted, as every node in the circuit being connected to the reference node (ground) through a path containing only the capacitors.



general initial value problem Find y, such that

$$y'(t) = f(t,y), \quad t \in [t_0, t_1],$$
 (18)

$$y(t_0) = y_0.$$
 (19)



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Numerical Solutions



Single-Step Methods



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Definition

A numerical method to approximate a differential equation 18 on a time-grid $t_0, ..., t_l$ with the intermediate values $y_0, ..., y_l$ is called a single-step method if it is of the form

$$y_{j+1} = y_j + h_j \phi(t_j, y_j, y_{j+1}, h_j).$$
 (20)

We call ϕ the *procedural function*. If ϕ does not depend on y_{j+1} , then the method is called *explicit*, otherwise it is called *implicit*.



Numerical Solutions



Single-Step Methods
Consistency, Stability and Convergence



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Let y_{m+1} be the result of one step of a single step method (20) with the exact start-vector $y_m = y(t_m)$ then

$$\delta_{m+1} = \delta(t_m + h) = y(t_{m+1}) - \tilde{y}_{m+1}, \quad m = 0, ..., N-1$$
 (21)

is called the *local discretization error* of the single step method at the point t_{m+1} .



A single-step method is called *consistent* if for all initial value problems (18)

$$\lim_{h\to 0}\frac{\|\delta(t+h)\|}{h}=0\quad \text{for}\quad t_0\leq t\leq t_l \tag{22}$$

holds.

It is called *consistent of order p*, if for a sufficiently smooth function f

$$\|\delta(t+h)\| \le Ch^{p+1} \quad \text{for all} \quad h \in (0,H] \quad \text{and} \quad t_0 \le t \le t_1 - h \tag{23}$$

holds with C independent of h.

A single-step method is called *convergent*, if for all initial value problems (18) for the *global discretization error*

$$e_{\mathfrak{m}} = \mathfrak{y}(\mathfrak{t}_{\mathfrak{m}}) - \mathfrak{y}_{\mathfrak{m}}$$

holds that

$$\max_{m} \|e_m\| \to 0$$
 for $h_{max} \to 0$.

The single-step method is called to have the *convergence order* p, if

$$\max_{m}\|e_{m}\| \leq Ch_{m\alpha x}^{p} \quad \text{for} \quad h_{m\alpha x} \in (0,H] \quad \text{with} \quad t_{0} \leq t_{m} \leq t_{l}$$

with the constant C not dependent on the step size h.

A single-step method is called (discretely) stable if for grid-functions y_h and \tilde{y}_h with

$$y_{i+1} = y_i + h\phi(t_i, y_i),$$
 (24)

$$\tilde{y}_{i+1} = \tilde{y}_i + h[\phi(t_i, \tilde{y}_i) + \theta_i], \tag{25}$$

and perturbations $\theta_i=\theta_h(t_i)$ of the right side as well as a bounded perturbation in the initial-values $y_0-\tilde{y}_0$ the error is bounded by

$$\|y_h - \tilde{y}_h\|_{\infty,h} \le C(\|y_0 - \tilde{y}_0\|_{L^2} + \|\theta_h\|_{\infty,h})$$

with a constant C which is not dependent on h. The norm $\|.\|_{\infty,h}$ denotes the maximum norm over the time-grid, i.e. for a function $b:T=t_0,...,t_N\to\mathbb{R}^d$ we have $\|b\|_{\infty,h}=\max_{t\in T}\|b(t)\|,\|b\|$ is the euclidean norm.



Single-Step Methods further stability properties



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Dahlquist equation, i.e. find y such that

$$y' = \lambda y, \quad t > 0$$
 (26)
 $y(0) = y_0$ (27)

with $\lambda \in \mathbb{C}$ and \mathfrak{u}_0 fixed.

1. If a single-step method can be written in the form

$$y_{i+1} = R(z) y_i, \quad z := h\lambda$$
 (28)

then we call $R: \mathbb{C} \to \mathbb{C}$ the *stability function* of the single-step method.

2. The set

$$S := \{ z \in \mathbb{C} : |R(z)| \le 1 \} \tag{29}$$

is called the region of stability of the method.

- 3. A single-step method is called
 - \circ *0-stable*, if $0 \in S$.
 - \circ *A-stable*, if $\mathbb{C}^- \subset S$.
 - \circ *L-stable*, if $R(z) \to 0$ for $Re(z) \to -\infty$.





Multistep Methods



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For given $\alpha_0, ..., \alpha_k$ and $\beta_0, ..., \beta_k$ the iteration rule

$$\sum_{l=0}^{k} \alpha_{l} y_{m+l} = h \sum_{l=0}^{k} \beta_{l} f(t_{m+l}, y_{m+l}), \quad m = 0, 1, ..., N - k$$
 (30)

is called a *linear multistep method* (linear k-step method). It is always assumed that $\alpha_k \neq 0$ and $|\alpha_0| + |\beta_k| > 0$. If $\beta_k = 0$ holds, then the method is called explicit, otherwise implicit.





Multistep Methods
Consistency, Convergence and Stability

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Let y_{m+k} be the result of one step of the multi-step method (30) with the start-values given as the evaluations of the exact solution $y_{m+l} = y(t_{m+l})$ at $0 \le l < k$. This means

$$\alpha_k \tilde{u}_{m+k} = \sum_{l=0}^{k-1} (h\beta_l f(t_{m+l}, y(t_{m+l})) - \alpha_l y(t_{m+l})) + h\beta_k f(t_{m+k}, y_{m+k}).$$

Then

$$\delta_{m+k} = \delta(t_{m+k}) = y(t_{m+k}) - y_{m+k}, \quad m = 0, 1, ..., N - k$$

is called the *local discretization error* (local error) of the linear multi-step method, see Def. 30 at the point t_{m+k} .

A linear multi-step method is called *consistent*, if for all functions $y(t) \in C^2([t_0, t_l])$

$$\lim_{h\to 0}\frac{1}{h}L[y(t),h]=0$$

holds. It has the *consistency order* p, if for all functions $y(t) \in C^{p+1}[t_0, t_1]$

$$L[y(t), h] = \mathcal{O}(h^{p+1})$$
 for $h \to 0$

holds.



We say that a linear multi-step method is convergent if for a solution y of the problem a solution vector created by an LMSM y_j for $j \in 0, ..., k$ we have that

$$\lim_{h\to\infty}\max_{0\leq j\leq k}\|y(t_j)-y_j\|=0.$$



A linear multi-step method is called (discretely) stable, if for solutions y_h and \tilde{y}_h of

$$\sum_{l=0}^{k} \alpha_{l} y_{m+l} = h \sum_{l=0}^{k} \beta_{l} f(t_{m+l}, y_{m+l}),$$
 (31)

$$\sum_{l=0}^{k} \alpha_{l} \tilde{y}_{m+l} = h \sum_{l=0}^{k} \beta_{l} f(t_{m+l}, \tilde{y}_{m+l}) + h \theta_{n}$$
 (32)

and bounded initial values $y_j - \tilde{y}_j$ for $j \in 0, ..., k$ we have that

$$\max_{t_0 \le t_n \le T} \|y_n - \tilde{y}_n\| \le C \sum_{i=0}^{k-1} \|y_i - \tilde{y}_i\| + \max_{t_0 \le t_n \le T} \|\theta_n\|.$$





Multistep Methods further stability properties



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Dahlquist test problem as a model problem, find y such that

$$y' = \lambda y, \quad t > 0 \tag{33}$$

$$y(0) = y_0 \tag{34}$$

with $\lambda \in \mathbb{C}$ and y_0 fixed.

Thus the resulting linear multistep method is of the form

$$\sum_{l=0}^{k} \alpha_{l} y_{n+l} = h \sum_{l=0}^{k} \beta_{l} \lambda y_{n+l}$$

$$\iff \sum_{l=0}^{k} [\alpha_{l} - h \beta_{l} \lambda] y_{n+l}$$

1. The set

$$S := \{ z \in \mathbb{C} : \rho(\xi) - z\sigma(\xi) = 0 \implies \xi \in \mathbb{C} \text{ and } |\xi| \le 1.$$
If ξ has multiplicity greater than 1, then $|\xi| < 1 \}$ (35)

is called the region of stability of the method.

- 2. A linear multistep method is called
 - \circ *0-stable*, if $0 \in S$.
 - \circ stable in the point $z \in \mathbb{C}$, if $z \in S$.
 - o $A(\alpha)$ -stable, if it is stable in all z that lie within the set $\{z \in \mathbb{C}^- : |arg(z) \pi| \le \alpha\}$ for $\alpha \in (0, \frac{\pi}{2})$.

Theorem

Let f(t, y) be sufficiently smooth and the linear multi-step method be zero-stable and consistent of order p, then it is also convergent of order p.





Implicit Linear Multistep Formulas



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Implicit Linear Multistep Formulas BDF-k Methods



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The backward differentiation formula (BDF) is a family of implicit linear multistep methods. They have the general form

$$\sum_{k=0}^{s} \alpha_k y_{n+k} = h\beta f(t_{n+s}, y_{n+s})$$
 (36)

The BDF or BDF-k formulas for k = 1, ..., 3 have the following form

$$k = 1 : hf_{m+1} = y_{m+1} - y_m$$

$$k = 2 : hf_{m+2} = \frac{1}{2}(3y_{m+2} - 4y_{m+1} + y_m)$$

$$k = 3 : hf_{m+3} = \frac{1}{6}(11y_{m+3} - 18y_{m+2} + 9y_{m+1} - 2y_m)$$



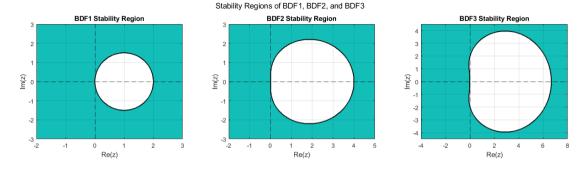


Figure: stability regions of BDF-schemes

Theorem (NumerikGewÄühnlicherDifferentialgleichungen))

The BDF-k methods have consistency order p = k.





Implicit Linear Multistep Formulas Trapezoidal rule



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This procedure is repeated for small subsections of the interval [a,b]. Thus we obtain the iteration formula

$$u_h(t+h) = u_h(t) + \frac{h}{2}[f(t,u_h(t)) + f(t+h,u_h(t+h))].$$





Numerical Examples



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Numerical Examples Example1



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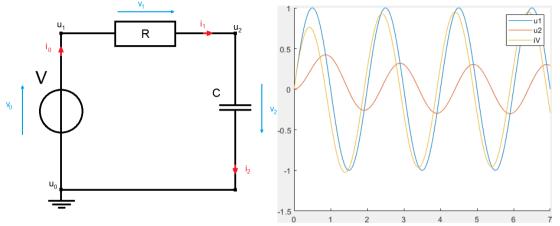


Figure: charging capacitor with series resistor and voltage source

Figure: Exact solution for example 1.



h	k = 1		k = 2		k = 3		trapezoidal	
	u2	iV	u2	iV	u2	iV	u2	iV
0.1	4.620×10^{-2}	4.620×10^{-2}	9.567×10^{-3}	9.567×10^{-3}	3.057×10^{-3}	3.057×10^{-3}	3.344×10^{-3}	3.344×10^{-3}
0.05	2.339×10^{-2}	2.339×10^{-2}	2.454×10^{-3}	2.454×10^{-3}	6.083×10 ⁻⁴	6.083×10^{-4}	8.367×10^{-4}	8.367×10^{-4}
0.025	1.178×10 ⁻²	1.178×10^{-2}	6.264×10^{-4}	6.264×10^{-4}	1.672×10 ⁻⁴	1.672×10^{-4}	2.092×10^{-4}	2.092×10^{-4}

Table: Resulting errors for the BDF-k methods and ther trapezoidal rule.





Numerical Examples Example 2



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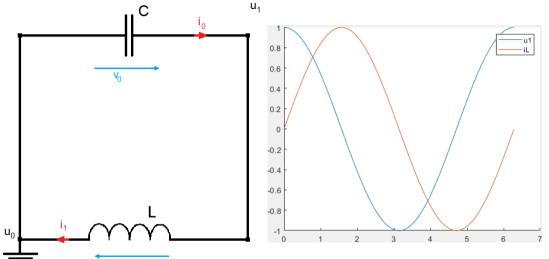


Figure: Exact solution for example 2.



h	k = 1		k = 2		k = 3		trapezoidal	
	u1	iL	u1	iL	u1	iL	u1	iL
0.1	7.145×10^{-1}	6.905×10^{-1}	7.763×10^{-2}	8.060×10^{-2}	5.395×10^{-3}	5.180×10^{-3}	1.963×10^{-2}	2.087×10^{-2}
0.05	4.659×10^{-1}	4.448×10^{-1}	1.964×10^{-2}	2.066×10^{-2}	5.938×10^{-4}	5.579×10^{-4}	4.912×10^{-3}	5.224×10^{-3}
0.025	2.695×10^{-1}	2.551×10^{-1}	4.924×10^{-3}	5.216×10^{-3}	5.773×10^{-5}	4.740×10^{-5}	1.228×10^{-3}	1.308×10^{-3}

Table: Resulting errors for the BDF-k methods and ther trapezoidal rule.





Numerical Examples Example 3



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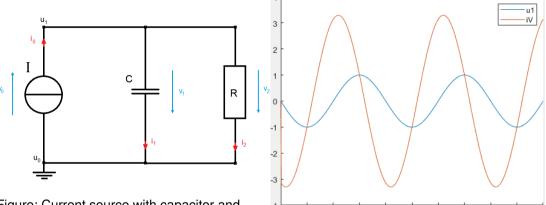


Figure: Current source with capacitor and resistor.

Figure: Exact solution for example 3.



h	k = 1	k = 2	k = 3	trapezoidal	
	iV	iV	iV	iV	
0.1	4.894×10^{-1}	1.023×10^{-1}	2.530×10^{-2}	5.219×10^{-2}	
0.05	2.462×10^{-1}	2.577×10^{-2}	6.426×10^{-3}	1.295×10^{-2}	
0.025	1.233×10^{-1}	6.456×10^{-3}	1.613×10^{-3}	3.232×10^{-3}	

Table: Resulting errors for the BDF-k methods and the trapezoidal rule.

