

Wiki: Linear multistep method

"Adams method" redirects here. For the electoral apportionment method, see [Method of smallest divisors](#).

Linear multistep methods are used for the [numerical solution of ordinary differential equations](#). Conceptually, a numerical method starts from an initial point and then takes a short **step** forward in time to find the next solution point. The process continues with subsequent steps to map out the solution. Single-step methods (such as [Euler's method](#)) refer to only one previous point and its derivative to determine the current value. Methods such as [Runge-Kutta](#) take some intermediate steps (for example, a half-step) to obtain a higher order method, but then discard all previous information before taking a second step. Multistep methods attempt to gain efficiency by keeping and using the information from previous steps rather than discarding it. Consequently, multistep methods refer to several previous points and derivative values. In the case of *linear* multistep methods, a [linear combination](#) of the previous points and derivative values is used.

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1. Definitions

Numerical methods for ordinary differential equations approximate solutions to initial value problems of the form

$$y' = f(t, y), \quad y(t_0) = y_0.$$

The result is approximations for the value of $y(t)$ at discrete times t_i :

$$\begin{aligned}
 t_i &= t_0 + ih \\
 y_i &= y(t_i) = y(t_0 + ih) \\
 f_i &= f(t_i, y_i)
 \end{aligned}$$

where h is the time step (sometimes referred to as Δt).

A linear multistep method uses a linear combination of y_i and y'_i to calculate the value of y for the desired current step.

Multistep method will use the previous s steps to calculate the next value. Consequently, the desired value at the current processing stage is y_{n+s} .

A linear multistep method is a method of the form

where h denotes the step size and f the right-hand side of the differential equation. The coefficients a_0, \dots, a_{s-1} and b_0, \dots, b_s determine the method. The designer of the method chooses the coefficients; often, many coefficients are zero.

Typically, the designer chooses the coefficients so they will exactly interpolate $y(t)$ when it is an n th order polynomial.

If the value of b_s is nonzero, then the value of y_{n+s} depends on the value of $f(t_{n+s}, y_{n+s})$. Consequently, the method is implicit if $b_s \neq 0$. In that case, the formula can directly compute y_{n+s} . If $b_s = 0$ then the method is explicit and the equation for y_{n+s} must be solved. [Iterative methods](#) such as [Newton's method](#) are often used to solve the implicit formula.

Sometimes an explicit multistep method is used to "predict" the value of y_{n+s} . That value is then used in an implicit formula to "correct" the value. The result is a [Predictor-corrector method](#).

2. Examples

Consider for an example the problem

$$y' = y, \quad y(0) = 1.$$

The exact solution is $y(t) = e^t$.

2.1. One-Step Euler

A simple numerical method is Euler's method:

$$y_{n+1} = y_n + hf(t_n, y_n).$$

Euler's method can be viewed as an explicit multistep method for the degenerate case of one step.

This method, applied with step size $h = \frac{1}{2}$ on the problem $y' = y$, gives the following results:

2.2. Two-Step Adams Bashforth

Euler's method is a one-step method. A simple multistep method is the two-step Adams-Bashforth method

$$y_{n+2} = y_{n+1} + \frac{3}{2}hf(t_{n+1}, y_{n+1}) - \frac{1}{2}hf(t_n, y_n).$$

This method needs two values, y_{n+1} and y_n , to compute the next value, y_{n+2} . However, the initial value problem provides only one value, $y_0 = 1$. One possibility to resolve this issue is to use the y_1 computed by Euler's method as the second value. With this choice, the Adams-Bashforth method yields (rounded to four digits):

The exact solution at $t = t_4 = 2$ is $e^2 = 7.3891\dots$, so the two-step Adams-Bashforth method is more accurate than Euler's method. This is always the case if the step size is small enough.

3. Multistep Method Families

Three families of linear multistep methods are commonly used: Adams-Bashforth methods, Adams-Moulton methods, and the [backward differentiation formulas](#) (BDFs).

3.1. Adams-Bashforth methods

The Adams-Bashforth methods are explicit methods. The coefficients are $a_{s-1} = -1$ and $a_{s-2} = \dots = a_0 = 0$, while the b_j are chosen such that the method has order s (this determines the method uniquely).

The Adams-Bashforth methods with $s = 1, 2, 3, 4, 5$ are ([Hairer, Nørsett & Wanner 1993](#), §III.1; [Butcher 2003](#), p. 103):

- $y_{n+1} = y_n + hf(t_n, y_n)$ —this is simply the Euler method;
- $y_{n+2} = y_{n+1} + h \left(\frac{3}{2}f(t_{n+1}, y_{n+1}) - \frac{1}{2}f(t_n, y_n) \right)$;

- $y_{n+3} = y_{n+2} + h \left(\frac{23}{12} f(t_{n+2}, y_{n+2}) - \frac{4}{3} f(t_{n+1}, y_{n+1}) + \frac{5}{12} f(t_n, y_n) \right);$
- $y_{n+4} = y_{n+3} + h \left(\frac{55}{24} f(t_{n+3}, y_{n+3}) - \frac{59}{24} f(t_{n+2}, y_{n+2}) + \frac{37}{24} f(t_{n+1}, y_{n+1}) - \frac{3}{8} f(t_n, y_n) \right);$
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The coefficients b_j can be determined as follows. Use [polynomial interpolation](#) to find the polynomial p of degree $s - 1$ such that

$$p(t_{n+i}) = f(t_{n+i}, y_{n+i}), \quad \text{for } i = 0, \dots, s - 1.$$

The [Lagrange formula](#) for polynomial interpolation yields

$$p(t) = \sum_{j=0}^{s-1} \frac{(-1)^{s-j-1} f(t_{n+j}, y_{n+j})}{j!(s-j-1)!h^{s-1}} \prod_{\substack{i=0 \\ i \neq j}}^{s-1} (t - t_{n+i}).$$

The polynomial p is locally a good approximation of the right-hand side of the differential equation $y' = f(t, y)$ that is to be solved, so consider the equation $y' = p(t)$ instead. This equation can be solved exactly; the solution is simply the integral of p . This suggests taking

$$y_{n+s} = y_{n+s-1} + \int_{t_{n+s-1}}^{t_{n+s}} p(t) dt.$$

The Adams-Bashforth method arises when the formula for p is substituted. The coefficients b_j turn out to be given by

$$b_{s-j-1} = \frac{(-1)^j}{j!(s-j-1)!} \int_0^1 \prod_{\substack{i=0 \\ i \neq j}}^{s-1} (u+i) du, \quad \text{for } j = 0, \dots, s-1.$$

Replacing $f(t, y)$ by its interpolant p incurs an error of order h^s , and it follows that the s -step Adams-Bashforth method has indeed order s ([Iserles 1996](#), §2.1)

The Adams-Bashforth methods were designed by [John Couch Adams](#) to solve a differential equation modelling capillary action due to [Francis Bashforth](#). [Bashforth \(1883\)](#) published his theory and Adams' numerical method ([Goldstine 1977](#)).

3.2. Adams-Moulton methods

The Adams-Moulton methods are similar to the Adams-Bashforth methods in that they also have $a_{s-1} = -1$ and $a_{s-2} = \dots = a_0 = 0$. Again the b coefficients are chosen to obtain the highest order possible. However, the Adams-Moulton methods are implicit methods. By removing the restriction that $b_s = 0$, an s -step Adams-Moulton method can reach order $s+1$, while an s -step Adams-Bashforth method has only order s .

The Adams-Moulton methods with $s = 0, 1, 2, 3, 4$ are ([Hairer, Nørsett & Wanner 1993](#), §III.1; [Quarteroni, Sacco & Saleri 2000](#)):

- $y_n = y_{n-1} + hf(t_n, y_n)$ – this is the [backward Euler method](#);
- $y_{n+1} = y_n + \frac{1}{2}h(f(t_{n+1}, y_{n+1}) + f(t_n, y_n))$ – this is the [trapezoidal rule](#);
- $y_{n+2} = y_{n+1} + h\left(\frac{5}{12}f(t_{n+2}, y_{n+2}) + \frac{2}{3}f(t_{n+1}, y_{n+1}) - \frac{1}{12}f(t_n, y_n)\right)$;

$$\begin{aligned} & \cdot y_{n+3} = y_{n+2} + h \left(\frac{3}{8} f(t_{n+3}, y_{n+3}) + \frac{19}{24} f(t_{n+2}, y_{n+2}) - \frac{5}{24} f(t_{n+1}, y_{n+1}) + \frac{1}{24} f(t_n, y_n) \right). \\ & \cdot \end{aligned}$$

The derivation of the Adams-Moulton methods is similar to that of the Adams-Bashforth method; however, the interpolating polynomial uses not only the points t_{n-1}, \dots, t_{n-s} , as above, but also t_n . The coefficients are given by

$$b_{s-j} = \frac{(-1)^j}{j!(s-j)!} \int_0^1 \prod_{\substack{i=0 \\ i \neq j}}^s (u+i-1) du, \quad \text{for } j = 0, \dots, s.$$

The Adams-Moulton methods are solely due to [John Couch Adams](#), like the Adams-Bashforth methods. The name of [Forest Ray Moulton](#) became associated with these methods because he realized that they could be used in tandem with the Adams-Bashforth methods as a [predictor-corrector](#) pair ([Moulton 1926](#)); [Milne \(1926\)](#) had the same idea. Adams used [Newton's method](#) to solve the implicit equation ([Hairer, Nørsett & Wanner 1993](#), §III.1).

4. Analysis

The central concepts in the analysis of linear multistep methods, and indeed any numerical method for differential equations, are [convergence, order, and stability](#).

The first question is whether the method is consistent: is the difference equation

a good approximation of the differential equation $y' = f(t, y)$? More precisely, a multistep method is *consistent* if the local error goes to zero as the step size h goes to zero, where the *local error* is defined to be the difference between the result y_{n+s} of the method, assuming that all the previous

values y_{n+s-1}, \dots, y_n are exact, and the exact solution of the equation at time t_{n+s} , divided by h . A computation using [Taylor series](#) shows out that a linear multistep method is consistent if and only if

$$\sum_{k=0}^{s-1} a_k = -1 \quad \text{and} \quad \sum_{k=0}^s b_k = s + \sum_{k=0}^{s-1} k a_k.$$

All the methods mentioned above are consistent ([Hairer, Nørsett & Wanner 1993](#), §III.2).

If the method is consistent, then the next question is how well the difference equation defining the numerical method approximates the differential equation. A multistep method is said to have *order* p if the local error is of order $O(h^{p+1})$ as h goes to zero. This is equivalent to the following condition on the coefficients of the methods:

$$\sum_{k=0}^{s-1} a_k = -1 \quad \text{and} \quad q \sum_{k=0}^s k^{q-1} b_k = s^q + \sum_{k=0}^{s-1} k^q a_k \quad \text{for } q = 1, \dots, p.$$

The s -step Adams-Bashforth method has order s , while the s -step Adams-Moulton method has order $s + 1$ ([Hairer, Nørsett & Wanner 1993](#), §III.2).

These conditions are often formulated using the *characteristic polynomials*

$$\rho(z) = z^s + \sum_{k=0}^{s-1} a_k z^k \quad \text{and} \quad \sigma(z) = \sum_{k=0}^s b_k z^k.$$

In terms of these polynomials, the above condition for the method to have order p becomes

$$\rho(e^h) - h\sigma(e^h) = O(h^{p+1}) \quad \text{as } h \rightarrow 0.$$

In particular, the method is consistent if it has order one, which is the case

$$\text{if } \rho(1) = 0 \quad \text{and} \quad \rho'(1) = \sigma(1).$$

If the roots of the characteristic polynomial ρ all have modulus less than or equal to 1 and the roots of modulus 1 are of multiplicity 1, we say that the [root condition](#) is satisfied. The method is convergent [if and only if](#) it is consistent and the root condition is satisfied. Consequently, a consistent method is stable if and only if this condition is satisfied, and thus the method is convergent if and only if it is stable.

Furthermore, if the method is stable, the method is said to be *strongly stable* if $z = 1$ is the only root of modulus 1. If it is stable and all roots of modulus 1 are not repeated, but there is more than one such root, it is said to be *relatively stable*. Note that 1 must be a root; thus stable methods are always one of these two.

4.1. Example

Consider the Adams-Bashforth three-step method

$$y_{n+1} = y_n + h \left(\frac{23}{12} f(t_n, y_n) - \frac{16}{12} f(t_{n-1}, y_{n-1}) + \frac{5}{12} f(t_{n-2}, y_{n-2}) \right).$$

The characteristic equation is thus

$$q(z) = z^3 - z^2 = z^2(z - 1)$$

which has roots $z = 0, 1$, and the conditions above are satisfied. As $z = 1$ is the only root of modulus 1, the method is strongly stable.

5. First and second Dahlquist barriers

These two results were proved by [Germund Dahlquist](#) and represent an important bound for the order of convergence and for the [A-stability](#) of a linear multistep method.

5.1. First Dahlquist barrier

A [zero-stable](#) and linear q -step multistep method cannot attain an order of convergence greater than $q + 1$ if q is odd and greater than $q + 2$ if q is even. If the method is also explicit, then it cannot attain an order greater than q ([Hairer, Nørsett & Wanner 1993](#), Thm III.3.5).

5.2. Second Dahlquist barrier

There are no explicit [A-stable](#) and linear multistep methods. The implicit ones have order of convergence most 2 ([Hairer & Wanner 1996](#), Thm V.1.4).

6. See also

- [Digital energy gain](#)

7. References

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