Basic Numerical Calculus

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1 DISCRETE FUNCTIONS AND DATA

1.1 Discrete Functions

A discrete function is a function that only has defined values at discrete points in its domain. While a continuous function may be defined as y = f(x), a discrete function is written as y vs. x. Essentially, the vector x is a collection of x values, while the vector y is a collection of the corresponding y values. By convention, we say that there are N + 1 points in x (see Section 1.2 for why this is the case).

$$\mathbf{x} = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_{N+1} \end{bmatrix}, \qquad \mathbf{y} = \begin{bmatrix} y_1 \\ y_2 \\ \vdots \\ y_{N+1} \end{bmatrix}$$

1.2 Continuous Functions and Their Discretization

When analyzing a *continuous* function numerically, the first thing we do is discretize the interval we are analyzing. We refer to this discretized interval as the **computational domain**. Essentially, we consider a function f(x) not as a continuous function, but rather as values corresponding to discrete locations, called **nodes**, in space. To discretize the domain, we first need to specify three quantities:

- 1. a: the left endpoint of the domain (i.e. the minimum value of x)
- 2. b: the right endpoint of the domain (i.e. the maximum value of x)
- 3. N: the number of subintervals (if we specify N subintervals, we will have N+1 nodes)

The length L of the domain is then

$$L = b - a$$

The discrete values of x (i.e. $x_1, ..., x_{N+1}$) are the nodes. Collectively, the set of nodes is referred to as the **mesh**. There are many different ways to create a mesh. In our case, we use a uniform mesh; this means that the nodes are equally spaced [2]. Thus, for a uniform mesh, the **grid spacing** is given by

$$\Delta x = \frac{L}{N}$$

The vector \mathbf{x} storing the nodes can be defined as

$$\mathbf{x} = \begin{bmatrix} x_1 \\ x_2 \\ x_3 \\ \vdots \\ x_i \\ \vdots \\ x_N \\ x_{N+1} \end{bmatrix} = \begin{bmatrix} a \\ a + \Delta x \\ a + 2\Delta x \\ \vdots \\ a + (i-1)\Delta x \\ \vdots \\ a + (i-1)\Delta x \\ \vdots \\ a + (N-1)\Delta x \\ a + N\Delta x \end{bmatrix}$$
(1)

We can then also define a vector y to store all the y_i 's, where y_i is the evaluation of f(x) at $x = x_i$.

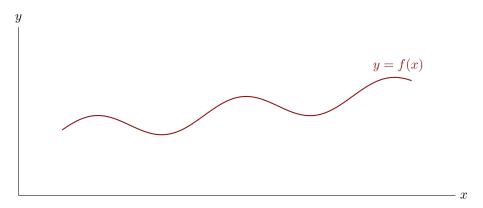
$$\mathbf{y} = \begin{bmatrix} y_1 \\ \vdots \\ y_{N+1} \end{bmatrix} = \begin{bmatrix} f(x_1) \\ \vdots \\ f(x_{N+1}) \end{bmatrix}$$

Thus, the vectors y and x essentially form a discrete function (essentially a discretized representation of y = f(x)) as defining a discrete form of f(x).

$$y = f(x)$$
 discretization \mathbf{y} vs. \mathbf{x}

f vs. **x**
$$\equiv \{(x_i, f(x_i))\}_{i=1}^{N+1}$$

An example of the discretization of a univariate function onto a uniform 1D mesh is shown in Fig. 1 below.



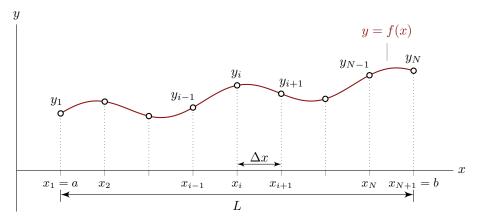


Figure 1: Discretization of a continuous function.

When discretizing a continuous function it is easiest to discretize it onto a uniform mesh. However, unevenly spaced nodes may be used as well. The numerical calculus techniques introduced in Sections 2 and 3 make no assumptions with regards to how the nodes are spaced.

1.3 Interpreting Data as a Discrete Function

When we sample data from the physical world, we built a data set

$$\{(x_i, y_i)\}_{i=1}^{N+1}$$

If we store all the x_i 's in a vector \mathbf{x} , and all the y_i 's in a vector \mathbf{y} , then we have the discrete function

Thus, a data set is essentially a discrete function.

$$\mathbf{y}$$
 vs. $\mathbf{x} \equiv \{(x_i, y_i)\}_{i=1}^{N+1}$

In reality, the data may be explained by some underlying, continuous function y = f(x). Let's say for some reason we needed either the derivative,

$$\frac{dy}{dx} = f'(x)$$

the definite integral,

$$\int_{a}^{b} f(x) \, dx$$

or a function defined by an integral,

$$\int_{a}^{b} f(x) dx$$
$$g(x) = \int_{a}^{x} f(x) dx$$

If we knew f(x), we could (in almost all cases) obtain its derivative, f'(x). However, even with knowledge of f(x), it is likely that we would not be able to evaluate either of the integrals shown above. The numerical differentiation and integration methods presented in Sections 2 and 3 will allow us to approximate these derivatives/integrals using the discrete function y vs. x (i.e. the data set).

2 NUMERICAL DIFFERENTIATION

2.1 Finite Difference Approximations [2, 5]

2.1.1 | Forward Difference Approximation

$$\left| \frac{dy}{dx} \right|_{x=x_i} \approx \frac{y_{i+1} - y_i}{x_{i+1} - x_i} \tag{2}$$

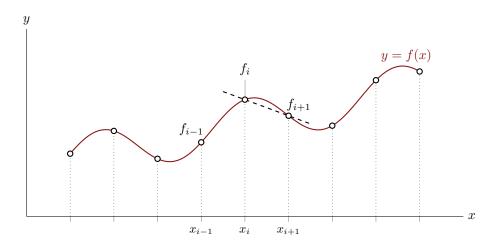


Figure 2: Forward approximation.

2.1.2 | Backward Difference Approximation

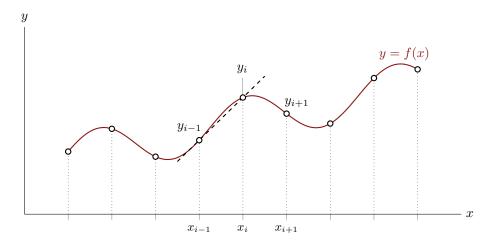


Figure 3: Backward approximation.

2.1.3 | Central Difference Approximation

$$\frac{dy}{dx}\Big|_{x=x_i} \approx \frac{y_{i+1} - y_{i-1}}{x_{i+1} - x_{i-1}}$$
 (4)

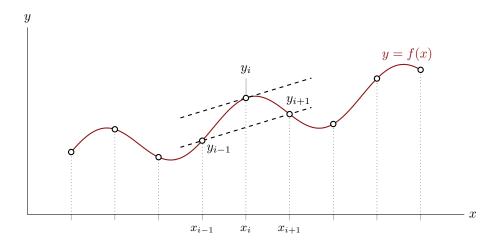


Figure 4: Central approximation.

The central approximation is of higher accuracy than the forward and backward approximations.

2.2 Differentiation Over an Interval (Cumulative Differentiation)

Consider the vectors y and x storing sampled points from an underlying function y = f(x). We can consider these vectors as a set of points or a data set:

$$\mathbf{y}$$
 vs. $\mathbf{x} \equiv \{(x_i, y_i)\}_{i=1}^{N+1}$

$$\therefore \mathbf{x} = \begin{bmatrix} x_1 \\ \vdots \\ x_{N+1} \end{bmatrix}, \qquad \mathbf{y} = \begin{bmatrix} y_1 \\ \vdots \\ y_{N+1} \end{bmatrix}$$

Note that $y_i = f(x_i)$.

Our goal is to find the derivative dy/dx = f'(x), but without knowledge of f(x), we cannot use simple algebraic differentiation rules. Instead, since we know the discrete y vs. x that essentially stores sampled values of y = f(x), we can numerically estimate the value of the derivative at all the points stored in the vector x. The result of this numerical differentiation is a vector dy that stores the numerical evaluation of dy/dx = f'(x) at all the points in x.

$$\mathbf{dy} = \begin{bmatrix} dy_1 \\ \vdots \\ dy_{N+1} \end{bmatrix} = \begin{bmatrix} \frac{dy}{dx} \Big|_{x=x_1} \\ \vdots \\ \frac{dy}{dx} \Big|_{x=x_{N+1}} \end{bmatrix}$$

I refer to this numerical differentiation process as **cumulative differentiation**, analogous to cumulative integration in the case of numerical integration (see Section 3.3). The algorithm I use to perform cumulative differentiation is rather

simple. At all interior nodes, I use a central approximation to approximate the derivative. At the left endpoint, I use a forward approximation, since there is no x_1 or y_1 . At the right endpoint, I use a backward approximation, since there is no x_{N+2} or y_{N+2} .

Algorithm 1:

Cumulative differentiation.

Given:

- $\begin{array}{ll} \bullet \ \ \mathbf{x} \in \mathbb{R}^{N+1} & \text{- vector of } x \text{ values} \\ \bullet \ \mathbf{y} \in \mathbb{R}^{N+1} & \text{- vector storing evaluations of } y = f(x) \text{ at every point in } \mathbf{x} \end{array}$

Procedure:

- 1. Determine the number of subintervals, N, given that $\mathbf{x}, \mathbf{y} \in \mathbb{R}^{N+1}$.
- 2. Preallocate $\mathbf{dy} \in \mathbb{R}^{N+1}$ to store the cumulative derivative.
- 3. Calculate derivative at left endpoint using forward difference approximation.

$$dy_1 = \frac{y_2 - y_1}{x_2 - x_1}$$

4. Calculate derivative at right endpoint using backward difference approximation.

$$dy_{N+1} = \frac{y_{N+1} - y_N}{x_{N+1} - x_N}$$

5. Calculate derivatives at all other points using central difference approximation.

for
$$i=2$$
 to N
$$dy_i=\frac{y_{i+1}-y_{i-1}}{x_{i+1}-x_{i-1}}$$
 end

• $\mathbf{dy} \in \mathbb{R}^{N+1}$ - vector storing the evaluation of dy/dx = f'(x) at every point in \mathbf{x}

Differentiation at a Point (Point Differentiation)

Previously, we introduced an algorithm (Algorithm 1) for approximating the derivative dy/dx = f'(x) at every node x_i . In this section, we only want to approximate the derivative at a specific point (or at a specific set of points). Note that these points do not have to be the nodes we used to discretize y = f(x) (or if using a data set, we can approximate derivatives at points not included in the data set). We refer to this as **point differentiation**. To perform point differentiation, we first find the derivative at every node using cumulative differentiation (i.e. Algorithm 1). Then, we use linear interpolation to linearly interpolate a value for $f'(x_i^*)$ at every x_i^* .

Consider the case where there are p points x_i^* (where j = 1, ..., p) at which we wish to evaluate the derivative of y = f(x). We then define a vector \mathbf{x}^* as

$$\mathbf{x}^* = \begin{bmatrix} x_1^* \\ \vdots \\ x_p^* \end{bmatrix}$$

Let dy^* be the vector in which we store the evaluations of $f'(x_i^*)$. Then

$$\mathbf{dy}^* = \begin{bmatrix} dy_1 \\ \vdots \\ dy_p \end{bmatrix} = \begin{bmatrix} \frac{dy}{dx} \Big|_{x=x_1^*} \\ \vdots \\ \frac{dy}{dx} \Big|_{x=x_p^*} \end{bmatrix}$$

Algorithm 2:

Point differentiation.

Given:

• $\mathbf{x} \in \mathbb{R}^{N+1}$ - vector of x values

• $\mathbf{y} \in \mathbb{R}^{N+1}$ - vector storing evaluations of y = f(x) at every point in \mathbf{x}

 $\mathbf{x}^* \in \mathbb{R}^p$ - point(s) at which to differentiate

Procedure:

1. Find dy (i.e. the cumulative derivative of f vs. x) using Algorithm 1.

2. Find dy* (i.e. the point derivatives at x*) by linearly interpolating/extrapolating dy at every point in x* (can be done using MATLAB's interp1 function with the `linear` and `extrap` options specified).

Return:

• $dy^* \in \mathbb{R}^p$ - vector storing the evaluation of dy/dx = f'(x) at every point in \mathbf{x}^*

3 NUMERICAL INTEGRATION

3.1 Types of Integration

In most math courses, we encounter two main types of integration: **definite integration** and **indefinite integration**. At its core, the definite integral of a function f(x) over the interval [a, b] computes the area bounded by f, the x-axis, x = a, and x = b. The actual calculation is performed as

$$\int_{a}^{b} f(x) dx = F(b) - F(a)$$

where F(x) is the *anti*derivative of f(x). On the other hand, the indefinite integral of f(x) is just its antiderivative plus a constant C.

$$\int f(x) \, dx = F(x) + C$$

Whereas the definite integral gives us a number, an indefinite integral gives us a *family* of functions (since there are infinite possibilities for *C*). Definite integration is discussed in Section 3.2, while the numerical analog of indefinite integration, called **cumulative integration**, is discussed in Section 3.3.

3.1.1 | Motivation for Numerical Integration

Numerical integration methods are use to approximate integration. There are three cases in which numerical integration is most often used [3]:

- 1. The integrand y = f(x) may only be known at certain points (i.e. you have the data set/discrete function y vs. x).
- 2. It may be impossible to find the antiderivative of f(x) in closed form.
- 3. It may be possible to find the antiderivative of f(x) in closed form, but it is easier to just integrate numerically.

3.2 Definite Integration

The definite integral of a univariate function f(x) on the interval [a,b] computes the area bounded by f, the x axis, and the vertical lines x=a and x=b. To approximate a definite integral, we can split the interval [a,b] into smaller intervals, approximate the area bounded by each of these smaller intervals, and then sum the areas of the smaller intervals. If the function y=f(x) is unknown and we know only the sampled values y_i at some x values x_i , then approximating the integral is in fact all we can do. The **trapezoidal rule** approximates a definite integral by splitting up the area under the curve into multiple trapezoids, as shown in Fig. 5. The area A of a trapezoid with base b and

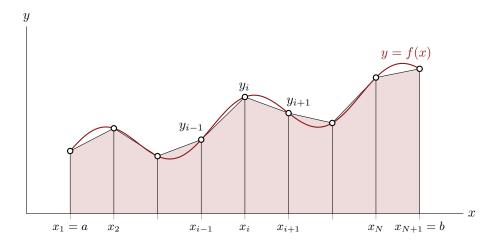


Figure 5: Trapezoidal rule.

heights h_1 (at left boundary) and h_2 (at right boundary) is

$$A = \frac{b\left(h_1 + h_2\right)}{2}$$

The area of the trapezoid defined by (x_i, y_i) and (x_{i+1}, y_{i+1}) is then

$$A = \underbrace{\frac{\sum_{i=1}^{b} \frac{h_1 + h_2}{(y_i + y_{i+1})}}_{2}}_{2}$$

Thus, to approximate the integral using the trapezoidal rule, we simply add up the areas of all the individual trapezoids [4].

$$\int_{a}^{b} f(x) dx \approx \sum_{i=1}^{N} \left[\frac{(y_{i+1} + y_i)(x_{i+1} - x_i)}{2} \right]$$
 (5)

Algorithm 3: definite integral

Definite integration.

Given:

- $\mathbf{x} \in \mathbb{R}^{N+1}$ vector of x values
- $\mathbf{y} \in \mathbb{R}^{N+1}$ vector storing evaluations of y = f(x) at every point in \mathbf{x}

Procedure:

- 1. Determine the number of subintervals, N, given that $\mathbf{x}, \mathbf{y} \in \mathbb{R}^{N+1}$.
- 2. Initialize a variable, I, to store the definite integral.

$$I = 0$$

3. Evaluate the definite integral using the trapezoidal rule.

for
$$i=1$$
 to N
$$I=I+\frac{(y_i+y_{i-1})(x_i-x_{i-1})}{2}$$
 and

Return:

• $I \in \mathbb{R}$ - definite integral of f(x) over the interval defined by \mathbf{x} (and where y = f(x) is given in discrete form as \mathbf{y} vs. \mathbf{x})

3.3 Cumulative Integration

Definite integration calculates a single area over an interval. What if we want to know the definite integral from the lower bound to *every single* point x in the interval [a, b]? Instead of specifying the upper bound of integration, we can leave it parameterized as x.

$$\int_{a}^{x} f(x) dx = F(x) - F(a)$$

We can note that F(x) is the antiderivative of f(x), while F(a) is this antiderivative evaluated at x=a (and therefore a constant). Thus, a cumulative integral produces a *function*.

We know that using a numerical approach, we can never return a continuous function; we can only return values of a function at specified, discrete points (recall that we refer to this set of points as a discrete function). Therefore, from a discrete standpoint, we can consider a data set \mathbf{f} vs. \mathbf{x} , or a function f sampled at discrete values of x. Either way, we have a set of N+1 points (x_i,y_i) (there are N+1 nodes since we define a computational domain as having N subintervals – see Section 1.2). Performing cumulative integration numerically should thus return an array of N+1 values, where the value at the i^{th} index represents the definite integral of elements 0 through i of the original data set [1].

Let CI(x) represent the cumulative integral of f(x) at x. In a continuous form, we can write

$$CI(x) = \int_{a}^{x} f(x) dx \tag{6}$$

At an arbitrary node x_i , we can approximate $CI(x_i) = CI_i$ using the trapezoidal rule defined by Eq. (5).

$$CI_{j} = \int_{a}^{x_{j}} f(x) dx \approx \sum_{i=1}^{j-1} \left[\frac{(y_{i+1} + y_{i})(x_{i+1} - x_{i})}{2} \right]$$
 (7)

Similarly, to find CI_{j+1} , we *could* calculate

$$CI_{j+1} = \int_{a}^{x_j} f(x) dx \approx \sum_{i=1}^{j} \left[\frac{(y_{i+1} + y_i)(x_{i+1} - x_i)}{2} \right]$$

However, we can notice that when calculating CI_{j+1} in the form above, we are essentially repeating almost every single calculation performed when calculating CI_j , which would make this a very inefficient way to program the cumulative integral. Instead, we can split up the summation.

$$CI_{j+1} \approx \sum_{i=1}^{j-1} \left[\frac{(y_{i+1} + y_i)(x_{i+1} - x_i)}{2} \right] + \frac{(y_{j+1} + y_j)(x_{j+1} - x_j)}{2}$$
(8)

Substituting Eq. (7) into Eq. (8),

$$CI_{j+1} = CI_j + \frac{(y_{j+1} + y_j)(x_{j+1} - x_j)}{2}$$
(9)

With Eq. (9), it becomes simple to write an algorithm to compute the cumulative integral of a function.

Algorithm 4: cumulative_integral

Cumulative integration.

Given:

- $\mathbf{x} \in \mathbb{R}^{N+1}$ vector of x values
- $\mathbf{y} \in \mathbb{R}^{N+1}$ vector storing evaluations of y = f(x) at every point in \mathbf{x}

Procedure:

- 1. Determine the number of subintervals, N, given that $\mathbf{x}, \mathbf{y} \in \mathbb{R}^{N+1}$.
- 2. Preallocate $CI \in \mathbb{R}^{N+1}$ to store the cumulative integral.
- 3. Set the first element of CI equal to 0 (since the integral from x = a to x = a is 0). Note: This step is only necessary if CI was NOT initialized as a vector of zeros.

$$CI_1 = 0$$

4. Evaluate the cumulative integral.

for
$$i=1$$
 to N
$$\qquad \qquad \mathrm{CI}_{i+1} = \mathrm{CI}_i + \frac{(y_i+y_{i-1})\,(x_i-x_{i-1})}{2}$$
 end

Return:

• $\mathbf{CI} \in \mathbb{R}^{N+1}$ - cumulative integral of f(x) over the interval defined by \mathbf{x} (and where y = f(x) is given in discrete form as \mathbf{y} vs. \mathbf{x})

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