

Analysis  
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# 1 Set Theory

... Thus, we see that  $\{0,1\}^X$  denotes the power set of  $X$ ; it is standard to abbreviate this to  $2^X$ , with the interpretation of 2 as a set with 2 elements and not the number 2, but with the potential confusion this may cause with notation such as “ $\mathbb{R}^2$ ”, we instead choose to use the (fancier) notation of  $\mathcal{P}(X)$  to denote the power set of  $X$ .

# 2 Real Numbers

“The great discovery of the late nineteenth century was that numbers can be understood abstractly via axioms, without necessarily needing a concrete model; of course a mathematician can use any of these models when it is convenient, to aid his or her intuition and understanding, but they can also be just as easily discarded when they begin to get in the way.”

Tao [3, 19]

## 2.1 Axiomatisation of the Reals

The existence of a model (a structure satisfying the axioms) will be given in the next section. The standard axiomatisation of  $\mathbb{R}$  is as follows:

1.  $\mathbb{R}$  is a field with respect to  $+$  and  $*$ .
2.  $\mathbb{R}$  is totally ordered with respect to  $\leq$ . <sup>(1)</sup>
3. The two operations preserve order;  $0 \leq x \wedge 0 \leq y \Rightarrow 0 \leq x * y$  and  $x \leq y \Rightarrow x + z \leq y + z$ .
4. The relation  $\leq$  is complete; any non-empty subset of  $\mathbb{R}$  bounded above has a least upper bound.

We note that Axiom 4 is nonfirstorderisable. There is an alternative, more concise axiomatisation due to Tarski.

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<sup>(1)</sup>For  $\leq$  to be totally ordered means that:

- (a)  $\forall x, x \leq x$  (Reflexivity)
- (b)  $\forall x, y, x \leq y \wedge y \leq x \Rightarrow x = y$  (Antisymmetry)
- (c)  $\forall x, y, z, x \leq y \wedge y \leq z \Rightarrow x \leq z$  (Transitivity)
- (d)  $\forall x, y, x \leq y \vee y \leq x$  (Totality)

## 2.2 Cauchy Sequences

### 2.2.1 Definitions

While  $\mathbb{Q}$  is dense (by its property that  $\forall x, y \exists z x < y \Rightarrow x < z < y$ ), it is not “complete” in the sense that certain definable concepts (e.g.  $\sqrt{2}$ ) are not in  $\mathbb{Q}$ . From the following definitions, equivalence classes of Cauchy sequences will form a model for the axioms from 2.1.

**Definition 2.2.1.1.** A sequence  $(a_q)_{q=m}^{\infty}$  is a mapping from  $\{n \in \mathbb{Z} \mid n \geq m\}$  to  $\mathbb{Q}$ . (Should the mapping be computable, definable, or not necessarily either?) (What is the definition of mapping?)

**Definition 2.2.1.2.** A *Cauchy sequence* is a sequence satisfying the property that for any rational number  $\varepsilon > 0$ , there is a finite  $N \in \mathbb{N}$  such that  $\forall i, j > N, d(a_i, a_j) < \varepsilon$ . (Is it always possible to determine whether a (computable) sequence is Cauchy?)

**Definition 2.2.1.3.** Two Cauchy sequences  $(a_i)$  and  $(b_i)$  are equivalent (written  $(a_i) \sim (b_i)$ ) if for any rational number  $\varepsilon > 0$ , there is a finite  $N \in \mathbb{N}$  such that  $\forall i > N, d(a_i, b_i) < \varepsilon$ . (Is equivalence computable?)

**Remark 2.2.1.1.** The notion of equivalence defined above forms an equivalence relation.

*Proof.* Reflexivity and symmetry are obvious. Given  $(a_i) \sim (b_i)$  and  $(b_i) \sim (c_i)$ , we can use  $N_1$  and  $N_2$  to write:

$$\forall i > \max(N_1, N_2), d(a_i, c_i) \leq d(a_i, b_i) + d(b_i, c_i) < \frac{\varepsilon}{2} + \frac{\varepsilon}{2} = \varepsilon.$$

Note that we have used the triangle inequality axiom for metric spaces. □

### 2.2.2 Arithmetic Operations on $\mathbb{R}$

We now define the arithmetic operations on the equivalence classes.

**Definition 2.2.2.1.**  $[(a_n)] + [(b_n)] := [(a_n + b_n)]$ , where  $(a_n + b_n)$  is the sequence where the  $n^{\text{th}}$  term is  $a_n + b_n$ .

To make sure this definition of addition is well-defined, we need to make sure that the addition of two Cauchy equivalence classes is a unique Cauchy equivalence class.

“You see, there’s a catch when you define things using equivalence relations. If you ever wish to define some operation on equivalence classes, you can’t just willy-nilly define it for one member of the class and expect it to make sense for all members of the class.”

Kemp [1, 6]

Firstly, we show that the addition of two Cauchy equivalence classes is a Cauchy equivalence class. Noting that:

$$\begin{aligned} d(a_i + b_i, a_j + b_j) &= |(a_i + b_i) - (a_j + b_j)| \\ &= |(a_i - a_j) + (b_i + b_j)| \\ &\leq |(a_i - a_j)| + |(b_i + b_j)| \\ &= d(a_i, a_j) + d(b_i + b_j), \end{aligned}$$

(is this true for general metrics?) we can write:

$$\forall i > \max(N_1, N_2), d(a_i + b_i, a_j + b_j) \leq d(a_i, a_j) + d(b_i + b_j) < \frac{\varepsilon}{2} + \frac{\varepsilon}{2} = \varepsilon.$$

Also, additions of equivalent sequences are equivalent; given  $(a_n) = (a'_n)$ , we will show that  $(a_n + b_n) = (a'_n + b_n)$ .

$$\begin{aligned} \forall i > N, d(a_i + b_i, a'_i + b_i) &= |(a_i + b_i) - (a'_i + b_i)| \\ &= |a_i - a'_i| < \varepsilon. \end{aligned}$$

Thus,  $(a_n + b_n) \sim (a'_n + b_n) \sim (a'_n + b'_n)$ . □

Multiplication is defined similarly, but with caution with 0 and reciprocation.

### 2.2.3 $\leq$ on $\mathbb{R}$

**Definition 2.2.3.1.**  $x \leq y$  iff  $y - x$  is 0 or positive.

### 2.2.4 Least Upper-bound Axiom

From the structure we have built, it is not too difficult to show that it models the first three axioms from 2.1. We now show that Axiom 4 is satisfied by our construction.

#### Definition 2.2.4.1.

**Theorem 2.2.4.1.** Any non-empty set  $S \subseteq \mathbb{R}$  with an upper bound (a real number  $M$  satisfying  $\forall x \in S, x \leq M$ ) has exactly one minimal upper bound  $u$  such that  $u \leq M$  for all upper bounds  $M$  of  $S$ .

*Proof.* Let  $u_0 = M$  for some upper bound  $M$  of  $S$  and  $\ell_0 = x$  for some  $x \in S$ . If  $a_n = \frac{\ell_n + s_n}{2}$  is an upper bound for  $S$ , assign  $\ell_{n+1} = \ell_n$  and  $u_{n+1} = a_n$ . Otherwise, assign  $\ell_{n+1} = a_n$  and  $u_{n+1} = u_n$ . This assignment preserves the upper-bound property of  $u_n$ , and so by induction,  $\forall n \in \mathbb{N}$ ,  $u_n$  is an upper bound for  $S$ . Similarly,  $\forall n \in \mathbb{N}$ ,  $\exists x \in S, \ell_n \leq x$ .

We want to show that  $(u_n)$  and  $(\ell_n)$  are Cauchy sequences, but a caveat is that they are sequences of real numbers, while our definition only allowed sequences of rationals. We may easily extend our definition of a sequence in 2.2.1.1 as a mapping to  $\mathbb{R}$  instead of  $\mathbb{Q}$ . This is not circular as we have already built  $\mathbb{R}$  independently from  $\mathbb{Q}$  (i.e. it is not circular as we are not re-defining the reals, but effectively defining a new construct which happens to have the same name “sequence” as our previous construct). A crucial nontrivial property is that equivalence classes of Cauchy sequences of reals are isomorphic to the reals. That is to say, the equivalence class of a real Cauchy sequence is equivalent to the equivalence class of some rational Cauchy sequence. (Equivalently, Cauchy sequences of reals converge to a real number.)

The proof is as follows. First we note that rationals can get arbitrarily close to any real number. Now, given a real sequence  $(r_n)$ , we construct a rational sequence  $(q_n)$  such that  $|u_n - q_n| < \frac{1}{n}$ . Then, by construction,  $[(r_n)] = [(q_n)]$ . All we need to show is that  $(q_n)$  is Cauchy. To do so, we note, given  $n, m > \max(N_1, N_2)$  for appropriate  $N_1, N_2$ :

$$\begin{aligned} |q_n - q_m| &= |(q_n - u_n) + (u_n - u_m) + (u_m - q_m)| \\ &\leq |q_n - u_n| + |u_n - u_m| + |u_m - q_m| \\ &< \frac{\varepsilon}{3} + \frac{\varepsilon}{3} + \frac{\varepsilon}{3} = \varepsilon. \end{aligned}$$

For  $N_1$  we need a natural number satisfying  $\frac{1}{N_1} < \frac{\varepsilon}{3}$ , which is guaranteed by the Archimedean property (proof omitted).

$[(u_n)]$  is an upper bound for  $S$ . If not, then for some  $s \in S$ ,  $[(u_n)] < s \Rightarrow [(u_n)] + \varepsilon < s$  for some small enough  $\varepsilon$ . But as  $(u_n)$  is non-increasing, we can find a  $n$  such that  $u_n < [(u_n)] + \varepsilon$ .<sup>(2)</sup> But then,  $u_n < [(u_n)] + \varepsilon < s$ , which is a contradiction as  $u_n$  is always an upper bound for  $S$ .

Using a similar argument, no real number smaller than  $[(\ell_n)]$  is an upper bound for  $S$ . Suppose otherwise. Let  $s$  be an upper bound of  $S$  such that  $s < [(\ell_n)] \Rightarrow s < [(\ell_n)] - \varepsilon$  for some small enough  $\varepsilon$ . But as  $(\ell_n)$  is non-decreasing, we can find a  $n$  such that  $\ell_n > [(\ell_n)] - \varepsilon$ . But then,

$$\ell_n > [(\ell_n)] - \varepsilon > s \Rightarrow \ell_n \text{ is greater than all elements in } S,$$

which is a contradiction by construction of  $\ell_n$ . □

---

<sup>(2)</sup>Choose  $n$  such that  $d(u_n, [(u_n)]) < \varepsilon$ . By the non-increasing property we have  $u_n \geq [(u_n)] \Rightarrow |u_n - [(u_n)]| = (u_n - [(u_n)])$ . Then,

$$\begin{aligned} u_n - d(u_n, [(u_n)]) &= u_n - (u_n - [(u_n)]) \\ &= [(u_n)] \\ \Rightarrow u_n - d(u_n, [(u_n)]) + \varepsilon &= [(u_n)] + \varepsilon \\ \Rightarrow u_n &< [(u_n)] + \varepsilon \end{aligned}$$

**Remark 2.2.4.1.** A potential point of confusion is that the property that  $u_n$  is an upper bound extends to  $\lim_{n \rightarrow \infty} u_n$ <sup>(3)</sup>, while the property  $\forall n \in \mathbb{N}, \exists x \in S, \ell_n \leq x$  of  $\ell_n$  does not extend to  $\lim_{n \rightarrow \infty} \ell_n$ . As a counterexample consider the set  $[0, 1)$ .

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<sup>(3)</sup>Where  $(a_n)$  is a Cauchy sequence,  $\lim_{n \rightarrow \infty} a_n$  is defined to be the real number corresponding to  $[(a_n)]$ .

## 3 Metric Spaces

### 3.1 Definitions

**Definition 3.1.0.1.** A metric space is a tuple  $(X, d)$  of a set  $X$  and a function  $d : X \times X \rightarrow [0, \infty)$  satisfying:

1.  $d(x, x) = 0$ .
2.  $x \neq y \Rightarrow d(x, y) > 0$ .
3.  $d(x, y) = d(y, x)$ .
4.  $d(x, z) \leq d(x, y) + d(y, z)$ .

**Example 3.1.0.1** (Euclidean Metric). asd

**Example 3.1.0.2** (Discrete Metric). The metric

$$d(x, y) := 1 - \delta(x, y) = \begin{cases} 0, & x = y \\ 1, & x \neq y \end{cases}$$

is called the *discrete metric*. The discrete metric counterintuitively only has clopen sets. There are two ways to explain this:

1. A singleton (which is always (and usually exclusively) closed) is open under the discrete metric. To see this, note that  $U_x(0.556) = \{x\} \in \{x\}$ . This in turn shows that any set is open under the discrete metric, as for every point in it one can find an open ball contained in the set. The set of complements of all subsets of the space  $X$  however, is equal to the set of all subsets of  $X$ , and so any set in  $X$  is also closed.
2. From the boundary-point definition of open sets, we see that any set in the discrete metric cannot have a boundary point.

#### 3.1.1 Open and Closed Sets

**Definition 3.1.1.1.** The open ball in a metric space  $(X, d)$  is defined as

$B(x_0, r) := \{x \in X \mid d(x, x_0) < r\}$ . Munkres writes this as the  $r$ -neighbourhood at  $x_0$ , or  $U(x_0, r)$ .

**Definition 3.1.1.2** (Interior, Exterior, Boundary Points).

- An *interior* point of a set  $E$  is a point such that there is an open ball (of positive radius) centred at it which is completely contained in  $E$ .
- If there is such a ball completely disjoint from  $E$ , then it is an *exterior* point.
- If neither, the point is a *boundary* point.

The set of boundary points is denoted by  $\partial E$ .

**Definition 3.1.1.3** (Open, Closed, Clopen Sets).

A set is

- *closed* if it contains all its boundary points;
- *open* if it contains none;
- *neither* if neither;
- *clopen*, if it has no boundary.

(Note that a clopen set is vacuously open and closed.)

**Remark 3.1.1.1.** Open balls are open. Consider  $B = U(x, r)$  and  $x' \in B$ . Let  $r' = r - d(x', x)$ , and  $B' = U(x', r')$ . Then, for any  $y \in B$ ,  $d(y, x) \leq d(y, x') + d(x', x) < r' + d(x', x) = r \Rightarrow y \in B$ .

**Definition 3.1.1.4.** A cluster point, or a point of accumulation, of a non-empty set  $S$  is a point such that every punctured ball centred on it contains points in  $S$ .

**Remark 3.1.1.2.** In order to deal with annoying exam questions whose proofs may depend on definitions, we demonstrate here the equivalence of different definitions of closed and open sets.

1. A set is open iff it contains none of its boundary points, and closed if it contains all.
2. A set is open iff for every point in it, there is some open ball centred there contained in the set.
3. A set is closed/open iff its complement is open/closed.
4. A set is closed iff it contains all its cluster points.

*Proof.* From its definition, we can see that a set can never contain its exterior points. So, 1 implies that a set is open iff it only contains interior points. But then,  $1 \Rightarrow 2$  follows from the definition of interior points. Also,  $2 \Rightarrow 1$  as 2 implies every point in an open set is an interior point.

To show 3, we note that a set and its complement share the same boundary;  $\partial S = \partial S'$ . This follows from the definition; the boundary points of  $S'$  are the points such that every open ball centred on it contains both points in  $S$  and points not in  $S$ ; this is precisely the boundary for  $S$ . Thus, a set contains all its boundary points iff its complement contains none, and so  $1 \iff 3$ .

Finally, cluster points by definition cannot be exterior. So if a set is closed, it must contain all cluster points (as closed sets contain all non-exterior points). Conversely, we note that every boundary point is a cluster point, as per their definitions. Thus, a set containing all its cluster points contains all its boundary points, and we are done.  $\square$

**Proposition 3.1.1.1.** Closed sets are closed under (finite applications of) the set union operation.

*Proof.* Suppose otherwise; that for some closed sets  $A$  and  $B$ ,  $A \cup B$  is not closed. Then there is some boundary point  $x$  of  $A \cup B$ , such that  $x \notin A \cup B$ . However,  $x$  cannot be an exterior point for both  $A$  and  $B$ , as every open ball centred at it contains points either always in  $A$  or always in  $B$ . WLOG,  $x$  is a boundary point for  $A$ , which contradicts the assumption that  $A$  is closed.  $\square$

**Proposition 3.1.1.2.** Closed sets are closed under (possibly infinite applications of) the set union operation.

*Proof.* Suppose otherwise; that for some set of closed sets  $A$ ,  $\bigcap A$  is not closed. Then there is some boundary point  $x$  of  $\bigcap A$ , such that  $x \notin B$ ,  $\forall B \in \bigcap A$ . However,  $x$  cannot be an exterior point for any  $B \in \bigcap A$ , as every open ball centred at it contains points common to  $B$ ,  $\forall B \in \bigcap A$ . Thus,  $x \in B$ ,  $\forall B \in \bigcap A$  (as closed sets contain all non-boundary points), a contradiction.  $\square$

**Remark 3.1.1.3.** For analogous results for open sets, use De Morgan's laws;  $(A \cup B)' = A' \cap B'$  and  $(A \cap B)' = A' \cup B'$ .

**Definition 3.1.1.5.** A set is bounded if there is a ball containing it.

**Theorem 3.1.1.1** (Bolzano-Weierstrass). Every bounded set  $S \in \mathbb{R}^n$  with infinitely many elements has a cluster point.

*Proof.* Take a hypercube bounding  $S$  of length  $L$ , and slice it into  $2^n$  smaller hypercubes, each with side length  $\frac{L}{2}$ . The union of these pieces contains infinitely many points of  $S$ , so at least one of these slices must contain infinitely many points in  $S$ . Choose one and repeat; the sequence of centre-points for the chosen hypercubes is a Cauchy sequence for a cluster point of  $S$ . To see why, create a punctured ball on the point represented by the Cauchy sequence. Now, we note that the hypercubes corresponding to the entries in the Cauchy sequence get arbitrarily small, and their centres arbitrarily close to the limit, and so a certain hypercube in the sequence is contained within the punctured ball. Also, all hypercubes in the sequence contain infinitely many points in  $S$ , and we are done.  $\square$

**Remark 3.1.1.4.** The theorem does not hold for general metric spaces.

**Corollary 3.1.1.1** (Bolzano-Weierstrass for subsequences). Every bounded sequence (in  $\mathbb{R}^n$ ) has a convergent subsequence.



*Proof.* Take the Cauchy sequence from before, and replace each entry with any point in the corresponding hypercube, whose index in the sequence is greater than the last entry's. We now have some subsequence of the original, which is Cauchy.  $\square$

### 3.1.2 Continuity

**Definition 3.1.2.1.** Where  $(X, d_x)$  and  $(Y, d_y)$  are metric spaces, a function  $f : X \rightarrow Y$  is continuous at  $x \in X$  iff  $f(x)$  is equal to the mapping of  $f$  to any Cauchy sequence of  $x$  (under the metric  $d_y$ ) i.e.  $f(x) = (f(x_0), f(x_1), \dots)$ .

**Theorem 3.1.2.1.** This is equivalent to the usual epsilon-delta definition:  $f$  is continuous at  $x$  if

$$\forall \varepsilon > 0 \exists \delta > 0, d_x(x_i, x) < \delta \Rightarrow d_y(f(x_i), f(x)) < \varepsilon,$$

or

$$\forall \varepsilon > 0 \exists \delta > 0, f(U_X(x, \delta)) \subseteq U_Y(f(x), \varepsilon).$$

*Proof.* As  $x_i$  gets arbitrarily close to  $x$  (under  $d_x$ ),  $f(x_i)$  gets arbitrarily close to  $f(x)$  (under  $d_y$ ). For a given  $\varepsilon > 0$ , find the index  $i$  s.t.  $\forall i, d_y(f(x_i), f(x)) < \varepsilon$  (which we are guaranteed as the sequence is Cauchy). Then, choose  $d_x(x_i, x)$  as our  $\delta$  – then, the index  $j$  such that  $\forall j, d_x(x_j, x) < \delta$  satisfies  $j > i$ . The case of  $\delta = 0$  can be avoided by choosing a Cauchy sequence of  $x$  whose entries are never equal to  $x$ . If no such Cauchy sequence can be found, choose a  $\delta$  such that the antecedent is always false; then  $f$  is vacuously continuous at  $x$ .  $\square$

**Remark 3.1.2.1.** Continuity is preserved under composition, and continuity preserves connectedness (proof omitted).

**Theorem 3.1.2.2.**  $f$  is continuous  $\iff f^-$  preserves openness/closedness of subsets of the codomain.

*Proof.*

$L \Rightarrow R$ :

Preservation of openness: Let  $V$  be an open subset of the codomain. Choose  $x$  in  $f^-(V)$ . By openness,  $\exists \varepsilon, U_Y(f(x), \varepsilon) \subseteq V$ . But by continuity, this implies that  $\exists \delta, f(U_X(x, \delta)) \subseteq U_Y(f(x), \varepsilon)$ . But then, by definition,  $U_X(x, \delta) \subseteq f^-(U_Y(f(x), \varepsilon)) \subseteq f^-(V)$ .

Preservation of closedness: Let  $V$  be a closed subset of the codomain. Consider any Cauchy  $(x_i)$  with entries in  $f^-(V)$  which converges to  $x$ . Then, by closedness and continuity,  $(f(x_i)) \in V \Rightarrow f(x) \in V \Rightarrow x \in f^-(V)$ .

$R \Rightarrow L$ :

Preservation of openness: Where  $f : X \rightarrow Y$ , choose  $\varepsilon > 0$  and consider  $f(x)$  for any  $x \in X$ . Then,  $U_Y(f(x), \varepsilon)$  is open in  $Y$  (neighbourhoods are always open). Thus,  $f^-(U_Y(f(x), \varepsilon))$  is open, and so we can find a neighbourhood centred at  $x$  contained in it:

$\exists \delta, U_X(x, \delta) \subseteq f^-(U_Y(f(x), \varepsilon)) \Rightarrow f(U_X(x, \delta)) \subseteq U_Y(f(x), \varepsilon)$  and we are done.

Unlike a direct method like before, this time we will use the complementary relationship between open and closed sets. Consider an open set  $U$  in  $Y$ . Then  $Y \setminus U$  is closed in  $Y$  (because  $U \setminus Y$  is open). Then, by assumption,  $f^-(Y \setminus U)$  is closed, but  $f^-(Y \setminus U) = X \setminus f^-(U)$  (from some set algebra) so  $f^-(U)$  is open, and we are done.  $\square$

**Remark 3.1.2.2.** Maps preserving openness are called open maps.  $f$  is an open map  $\niff f$  is continuous. The respective counterexample are:

$$L \niff R: f : \mathbb{R} \rightarrow \mathbb{Z}, f(x) = \lfloor x \rfloor$$

$$R \niff L: f : \mathbb{R} \rightarrow \mathbb{R}, f(x) = 0.$$

### 3.1.3 Connectedness

Before defining connectedness, it will be convenient to discuss openness and closedness of subspaces (subspace implying that the measure of the metric space is kept) of metric spaces. In particular, we note how  $[0, 1)$  is neither closed or open in  $(\mathbb{R}, d)$  (where  $d$  is Euclidean) but open in the subspace  $X = ([0, 2], d)$ ;  $[0, 1) = U_X(0, 1)$ . We may characterise open/closed sets in subspaces as follows:

**Theorem 3.1.3.1.** Consider a subspace  $X$  of the metric space  $(M, d)$ .  $A \subseteq X$  is open/closed in  $X$  iff there is an open/closed  $\mathcal{O} \in M$  such that  $A = X \cap \mathcal{O}$ .

*Proof.* We consider the open case first.

$L \Rightarrow R$ : Construct an open set  $\mathcal{O}$  in  $M$  as:

$$\mathcal{O} = \bigcup_{a \in A} U_M(a, \varepsilon_a)$$

where  $\varepsilon_a$  is chosen so that  $U_X(a, \varepsilon_a) \in A$ . Then,

$$\begin{aligned} X \cap \mathcal{O} &= X \cap \bigcup_{a \in A} U_M(a, \varepsilon_a) \\ &= \bigcup_{a \in A} (X \cap U_M(a, \varepsilon_a)) \\ &= \bigcup_{a \in A} (U_X(a, \varepsilon_a)) \\ &= A \end{aligned} \tag{1}$$

$$\tag{2}$$

$R \Rightarrow L$ : We want to show that, given a set  $\mathcal{O}$  open in  $M$ ,  $A = X \cap \mathcal{O}$  is open in  $X$ . Choose any  $a \in A$ . We know that  $a \in \mathcal{O}$ , so by openness of  $\mathcal{O}$  we can choose a  $\varepsilon$  such that  $U_M(a, \varepsilon) \subseteq \mathcal{O}$ . But this gives  $U_X(a, \varepsilon) = X \cap U_M(a, \varepsilon) \subseteq X \cap \mathcal{O} = A$ .

The case for closed sets follow easily. Let  $B$  be a subset of  $X$ .

$$\begin{aligned} B \text{ is closed in } X &\iff X \setminus B \text{ is open in } X \\ &\iff X \setminus B = X \cap \mathcal{O}, \text{ for some open } \mathcal{O} \subseteq M \\ &\iff X \cap (X \setminus B)' = X \cap (X \cap \mathcal{O})' \\ &\iff X \cap (X \cap B')' = X \cap (X' \cup \mathcal{O}') \\ &\iff X \cap (X' \cup B) = X \cap \mathcal{O}' \\ &\iff X \cup B = X \cap \mathcal{C} \text{ for some closed } \mathcal{C} \subseteq M \\ &\iff B = X \cap \mathcal{C}. \end{aligned}$$

□

**Definition 3.1.3.1.** A set  $S$  is disconnected iff there exists two non-empty disjoint sets covering it, but not individually. This is equivalent to saying that  $S$  has a non-trivial proper clopen set in  $S$ .

*Proof.*

$L \Rightarrow R$ : Suppose we are working in the metric space  $(X, d)$ , and that the open sets  $A$  and  $B$  demonstrate the disconnectedness of  $S \subseteq X$ , i.e.  $S \subseteq A \cup B$ ,  $A \cap B = \emptyset$ ,  $A \cap S \neq \emptyset \neq B \cap S$ . Now,  $A \cap S$  is open in  $S$  as  $A$  is open in  $X$ . It follows that  $S \setminus (A \cap S) = B \cap S$  <sup>(1)</sup> is closed in  $S$ , but by symmetry of  $A$  and  $B$ ,  $B \cap S$  is also open and  $A \cap S$  is also closed.

<sup>(1)</sup>An open ball in  $X$  is equivalent to the same open ball in  $M$ , but consisting only of the points belonging to  $X$ ;  $U_X(x, r) = X \cap U_M(x, r)$ .

<sup>(2)</sup>Let  $B = \bigcup_{a \in A} (U_X(a, \varepsilon_a))$ .  $A \subseteq B$  as  $B$  runs over all points in  $A$ .  $B \subseteq A$  as all neighbourhoods in the union are contained in  $A$ .

$R \Rightarrow L$ : Suppose that a set  $S$  in a metric space  $(X, d)$  has no non-trivial proper subsets which are clopen in  $S$ . Assume for the sake of contradiction that  $S$  is disconnected. Then by the  $L \Rightarrow R$  proof above we can construct non-trivial proper clopen subsets in  $S$ , a contradiction.  $\square$

**Theorem 3.1.3.2.** Continuity preserves connectedness.

*Proof.* Where  $f : X \rightarrow Y$  is continuous and a non-empty  $S \subseteq X$  is connected, let  $\mathcal{O}_1$  and  $\mathcal{O}_2$  satisfy the conditions to show that a  $f(S)$  is disconnected, for the sake of contradiction. We will show that  $f^-(\mathcal{O}_1)$  and  $f^-(\mathcal{O}_2)$  satisfy the conditions to show that  $S$  is disconnected. Both contain points in  $S$  as  $\mathcal{O}_1$  and  $\mathcal{O}_2$  contain points in  $f(S)$ . This also shows that they are non-empty. They are disjoint as otherwise, if  $x$  was common to both,  $f(x) \in \mathcal{O}_1 \cap \mathcal{O}_2$ , a contradiction. Finally, they collectively cover  $S$  as  $f(S) \subseteq \mathcal{O}_1 \cup \mathcal{O}_2 \Rightarrow S \subseteq f^-(\mathcal{O}_1 \cup \mathcal{O}_2) = f^-(\mathcal{O}_1) \cup f^-(\mathcal{O}_2)$ .  $\square$

**Definition 3.1.3.2.** An *interval* is a set in the extended reals  $\mathbb{R}^*$  denoted by  $[a, b] := \{x \mid a \leq x \leq b\}$ , with either “ $\leq$ ” replaceable by a “ $<$ ”, along with a change from a square bracket to a paranthesis.

**Theorem 3.1.3.3.**  $X$  is an interval  $\iff a, b \in X \wedge a < b \Rightarrow \forall x \in (a, b), x \in X$

*Proof.*

$L \Rightarrow R$ : Follows trivially by definition of an interval.

$R \Rightarrow L$ : Let  $u = \sup(X)$  and  $l = \inf(X)$ . We want to show that  $X$  contains all elements “in between”  $u$  and  $l$ . For any  $x$  such that  $l < x < u$ , by definition of supremum and infimum, we can find  $a, b \in X$  such that  $a < x < b$ , and thus  $x \in X$ .  $\square$

**Theorem 3.1.3.4.** Where  $S \subseteq \mathbb{R}$ ,  $S$  is connected  $\iff S$  is an interval.

*Proof.*

$L \Rightarrow R$ : Consider any two points  $a, b$  in  $S$  and suppose that  $x \in (a, b)$ ,  $x \notin S$ . But then  $(-\infty, x)$  and  $(x, \infty)$  demonstrate the disconnectedness of  $S$ , contradicting the hypothesis.

$R \Rightarrow L$ : We first show the results for closed, bounded intervals. For the sake of contradiction, let  $U, V \subseteq [a, b]$  be clopen subsets (in  $[a, b]$ ) demonstrating disconnectedness of the closed interval. Assume wlog that  $b \in V$  and let  $u$  be the supremum of  $U$ .  $u \in U$  by closedness of  $U$ . Noting that  $(b - \epsilon, b] \in V$  by openness, it follows that  $u < b$  by disjointness, but then,  $[u, u + \epsilon) \in U$  as  $U$  is open in  $[a, b]$ , contradicting that  $u$  is a supremum.

For general intervals, the results follows from the fact that connectedness is preserved under certain unions, and all intervals can be generated from a union of closed bounded intervals. See “4-connected.pdf”.  $\square$

**Corollary 3.1.3.1** (Intermediate Value Theorem). Where  $f : S \rightarrow \mathbb{R}$  is continuous,  $S$  is connected, and  $a < b$ ,  $a, b \in f(S) \Rightarrow [a, b] \subseteq f(S)$ .

**Remark 3.1.3.1.** Connected open sets are polygonally connected.

*Proof.* Let  $x$  be a point in the open connected set  $S$ , and  $X$  be the smallest subset of  $S$  containing  $x$  closed under polygonal connection; that is,  $y$  is polygonally connected to  $x \iff y \in X$ . We note that  $X$  must be open, as if it were to have a boundary point  $b$ , by openness of  $S$  there is a neighbourhood of  $b$  in  $S$ , then in turn, this neighbourhood would be contained in  $X$  (as every point in a neighbourhood is polygonally connected to its centre), contradicting the assumption that  $b$  is a boundary of  $X$ . Similarly, suppose that  $S \setminus X = S \cap X'$  has a boundary point  $b$ . This means that every open ball centred at  $b$  contains points in  $(S \setminus X)' = S' \cup X$ . In particular, every neighbourhood of  $b$  must contain points in  $X$ , as otherwise this contradicts  $S$  being open. But then, as this open ball contains a point in  $X$ , its centre  $b$  must also be polygonally connected to  $x$ , a contradiction, as  $S \setminus X$  is supposed to be points in  $S$  unreachable from  $x$ . Thus,  $S \setminus X$  is open.

Now, notice that  $S \setminus X$  and  $S$  are open, disjoint, and collectively cover  $S$ . As  $S$  is connected, at least one of them cannot contain points in  $S$  – this must be  $S \setminus X$  as  $X$  contains  $x$ ; thus,  $S \setminus X = \emptyset$ .  $\square$

### 3.1.4 Compactness

**Definition 3.1.4.1.** A compact set is a set which is closed and bounded. This definition is equivalent to saying that all sequences in a compact set  $S$  have a subsequence converging to a point in  $S$ .

*Proof.* A closed set contains all cluster points, so any sequence in it must converge to it. Conversely, the set must be closed as it contains all its cluster points. Also, it must be bounded as otherwise, one can construct an unbounded sequence, which has no convergent subsequences.  $\square$

**Remark 3.1.4.1.** For general topological spaces this definition is somewhat problematic, so a definition using open covers may instead be used.

**Theorem 3.1.4.1** (Heine-Borel).

$K \subseteq \mathbb{R}$  is compact  $\iff (C \text{ covers } K \Rightarrow \text{a finite subset of } C \text{ covers } K)$ . A set  $C$  covers  $B$  iff  $B \subseteq \bigcup C$  and  $\forall A \in C, A$  is open.

*Proof.*

$L \Rightarrow R$ : Given  $K \subseteq \mathbb{R}$  is compact, assume bwoc that coverings of  $K$  have no finite subcover. Using the hypercube argument from before, at least one of the smaller hypercubes must require an infinite number of elements from a given covering  $C$ . Let  $L$  be the limit point of the Cauchy sequence of the centres of the hypercubes.  $C$  must have some open set  $B$  containing  $L$ , but this contradicts the fact that each of our hypercubes required an infinite amount of elements of  $C$  to be covered. (We can find an arbitrarily small hypercube in our sequence that fits inside  $B$ .)

$R \Rightarrow L$ : Given the RHS, assume bwoc that  $K$  is not closed i.e. there exists some point  $c \notin K$  whose Cauchy sequence is in  $K$ . The set of neighbourhoods of all  $x \in K$  with their radii chosen not to include some neighbourhood of  $c$  (say, with half the distance from  $x$  to  $c$ ) covers  $K$ , but has no finite subcovering – given a finite subcovering, each of its elements is disjoint from a certain neighbourhood of  $c$ , although that neighbourhood certainly contains elements in  $K$ . Thus,  $K$  must be closed.

Also, given the RHS, consider the covering  $C = \{B(x, 1) \mid x \in K\}$ . A finite subcovering of this consists of a finite number of balls of radius 1, so  $K$  must be bounded.  $\square$

**Lemma 3.1.4.1.** Continuous maps preserve compactness.

*Proof.* Let  $f$  be a continuous map and  $X$  a closed set.

$$\begin{aligned} (f(x_0), f(x_1), \dots) &= f((x_0, x_1, \dots)) && \text{(by continuity of } f) \\ &= f(x), \quad x \in X && \text{(by closedness of } X) \end{aligned}$$

Thus, any Cauchy sequence in  $f(X)$  is contained in  $f(X)$  – closedness is preserved by continuous maps.

Now, assume  $X$  is bounded (but not necessarily closed). For  $f(X)$  to be unbounded means that there exists a sequence in  $f(X)$  which has no convergent subsequence. For such a sequence  $(f(x_0), f(x_1), \dots)$ , consider the corresponding sequence in  $X$ ,  $(x_0, x_1, \dots)$ . While this sequence may not be Cauchy it is bounded, so there is some convergent subsequence  $(x_{\phi(n)})$ . But then,  $(f(x_{\phi(0)}), f(x_{\phi(1)}), \dots) = f((x_{\phi(n)}))$  by continuity of  $f$ , so  $(f(x_i))$  has a subsequence that converges – a contradiction.  $\square$

**Theorem 3.1.4.2.** Compactness ensures attainment of maxima. ( $f$  attains a maximum on  $c \in \text{dom } f$  if  $f(c) = \sup_{\text{dom } f} f$ .)

*Proof.* Let  $X = \text{dom } f$ .  $\sup_X f = (f(x_0), f(x_1), \dots)$  for some sequence  $(x_0, x_1, \dots)$  in  $X$ . By boundedness, some subsequence  $(x_{\phi(n)})$  converges, and is guaranteed to be in  $X$  by closedness. In particular, we have  $f((x_{\phi(n)})) = (f(x_{\phi(0)}), f(x_{\phi(1)}), \dots) = \sup_X f$  and we are done.  $\square$

**Definition 3.1.4.2.** A function from metric space  $(X, d_x)$  to  $(Y, d_y)$  is uniformly continuous if for every  $\varepsilon > 0$  there is a global parameter  $\delta$  such that  $d_x(x_0, x_1) < \delta \Rightarrow d_y(f(x_0) - f(x_1)) < \varepsilon$ .

**Theorem 3.1.4.3** (Heine-Cantor). Continuity implies uniformness on compact sets.

*Proof.* For a continuous  $f : X \rightarrow Y$  where  $X$  is compact, fix  $\varepsilon > 0$  and assign each  $x \in X$  with a  $\delta_x$  such that  $d_x(x, a) < \delta_x \Rightarrow d_y(f(x), f(a)) < \frac{\varepsilon}{2}$ , which we are guaranteed by continuity. Now, we notice that  $d_x(a, x), d_x(x, b) < \min(\delta_a, \delta_b) \Rightarrow d_y(f(a), f(b)) \leq d_y(f(a), f(x)) + d_y(f(x), f(b)) \leq \frac{\varepsilon}{2} + \frac{\varepsilon}{2}$ . So setting  $\delta = \min_{x \in X}(\delta_x)$  gives uniform continuity, but the problem is that as there are potentially an infinite number of  $\delta_x$  to consider, which may set  $\delta$  to 0 (this is not allowed by definition, where  $\delta > 0$ .)

The trick here is to use the Heine-Borel theorem to reduce the number of  $\delta_x$  needed to a finite amount. The set

$$C = \left\{ B\left(x, \frac{\delta_x}{2}\right) \mid x \in X \right\}$$

covers  $X$  so there is a finite subset  $C' \subseteq C$  also covering  $X$ .

Now take  $\delta$  to be the minimal radius of the balls in  $C'$  (this will be the global parameter we need). For any  $x \in X$ , we can find some ball in  $C'$  containing it, say  $B(x_i, \frac{\delta_{x_i}}{2})$ . Given  $d_x(x, y) < \delta$ ,  $y$  must also be contained in this ball;

$$d_x(x_i, y) \leq d_x(x_i, x) + d_x(x, y) \leq \frac{\delta_{x_i}}{2} + \delta \leq \delta_{x_i}.$$

Then,

$$\begin{aligned} d_y(f(x), f(y)) &\leq d_y(f(x), f(x_i)) + d_y(f(x_i), f(y)) \\ &\leq \frac{\varepsilon}{2} + \frac{\varepsilon}{2} = \varepsilon. \end{aligned}$$

□

**Theorem 3.1.4.4** (Rolle's Theorem). Where  $f : [a, b] \rightarrow \mathbb{R}$  is continuous on its domain, differentiable on  $(a, b)$ , and satisfies  $f(a) = f(b)$ ,  $\exists x \in (a, b)$ ,  $f'(x) = 0$ .

*Proof.* The domain is compact  $f$  attains its maxima somewhere in it. Suppose that they at least one of them is contained in  $(a, b)$  – say (wlog) a maximum on  $x \in (a, b)$ . By definition, we have  $f(x \pm h) \leq f(x)$ . Thus,

$$\begin{aligned} f'(x) &= \lim_{h \rightarrow 0^+} \frac{f(x+h) - f(x)}{h} \\ &\leq 0. \end{aligned}$$

The inequality follows as the term in the limit is always non-positive; if every term is non-positive/non-negative in a Cauchy sequence, the corresponding real number will be non-positive/non-negative. (Why?) But also,

$$\begin{aligned} f'(x) &= \lim_{h \rightarrow 0^-} \frac{f(x+h) - f(x)}{h} \\ &\geq 0. \end{aligned}$$

As  $f$  is differentiable on  $(a, b)$ ,  $f'(x)$  exists and must be 0.

If both extrema are located on  $a$  and  $b$ , then by definition  $\forall x \in [a, b]$ ,  $f(x) = f(a)$  and it follows that  $f'$  is 0 everywhere on  $(a, b)$ . □

**Corollary 3.1.4.1** (Mean-value Theorem). Where  $f : [a, b] \rightarrow \mathbb{R}$  is continuous on its domain and differentiable on  $(a, b)$ ,  $\exists x \in (a, b)$ ,  $f'(x) = \frac{f(b)-f(a)}{b-a}$ .

*Proof.* Consider  $g(x) = f(x) - mx$ , where  $m = \frac{f(b)-f(a)}{b-a}$ . Then  $g$  satisfies the conditions for Rolle's theorem, and we know that:

$$\begin{aligned} \exists x \in [a, b], \quad g'(x) &= 0 \\ \Rightarrow \frac{d}{dx}(f(x) - mx) &= 0 \\ \Rightarrow f'(x) &= m \end{aligned}$$

□

### 3.1.5 Limits

**Definition 3.1.5.1.** Where  $f : X \rightarrow Y$ ,  $\lim_{x \rightarrow x_0} f(x) := [f(x_i)]$  where  $[(x_i)] = x_0$ . This is equivalent to the epsilon-delta definition.

*Proof.* Choose  $f(x_i)$  in the sequence such that it is within  $\varepsilon$  of its limit. Then,  $d_x(x_i, x_0)$  is the desired  $\delta$ .  $\square$

**Proposition 3.1.5.1.** Given  $f : \mathbb{R}^m \rightarrow \mathbb{R}^n$ ,  $f = (f_1, f_2, \dots, f_n)$ ,  
 $\lim_{\vec{x} \rightarrow \vec{a}} f(\vec{x}) = \vec{L} \iff \forall i, \lim_{\vec{x} \rightarrow \vec{a}} f_i(\vec{x}) = L_i$ .

*Proof.*

$$\begin{aligned} \|f(\vec{x}) - \vec{L}\| \rightarrow 0 &\iff \sqrt{\sum_{q=1}^n (f_q(\vec{x}) - L_q)^2} \rightarrow 0 \\ &\iff \sum_{q=1}^n (f_q(\vec{x}) - L_q)^2 \rightarrow 0 \\ &\iff \forall i, (f_i(\vec{x}) - L_i)^2 \rightarrow 0 && \text{(each term is positive)} \\ &\iff \forall i, f_i(\vec{x}) \rightarrow L_i \end{aligned}$$

$\square$

**Corollary 3.1.5.1.** It follows that a vector-valued function is continuous iff all of its components are.

## 3.2 Differentiation

### 3.2.1 Definitions

Recall our definition of discrete derivatives in which we took differences in consecutive values of a sequence  $f$ . Extending this definition to functions on continuous domains, we find the problem that we can't find consecutive values by completeness. However, the natural extension here is to take the Cauchy sequence of the normalised  $h$ -step difference:

$$\begin{aligned} Df(x) &:= \left( [\hat{\Delta}_h f](x) \right)_{h \rightarrow 0} \\ &= \lim_{h \rightarrow 0} \frac{f(x+h) - f(x)}{h} \end{aligned}$$

The generalisation to multi-dimensional inputs is simple, once one accepts the corresponding extension of normalised  $h$ -step differences to vector inputs:

$$\hat{\Delta}_{\vec{h}} f(\vec{x}) := \frac{f(\vec{x} + \vec{h}) - f(\vec{x})}{\|\vec{h}\|}.$$

This generalisation preserves the meaning of the  $h$ -step normalised difference of  $f$  at  $x$  as “the unit change in  $f$  per unit distance travelled along the shortest path from  $x$  to  $x+h$ ”. The Cauchy sequence extension of this is, not (what is defined as) the standard multivariable derivative, but the directional derivative; Then, the derivative becomes:

$$\begin{aligned} D_{\vec{a}} f(\vec{x}) &:= \left( \hat{\Delta}_{h\vec{a}} f(\vec{x}) \right)_{h \rightarrow 0} \\ &= \lim_{h \rightarrow 0} \frac{f(\vec{x} + h\vec{a}) - f(\vec{x})}{\|h\vec{a}\|}. \end{aligned}$$

(wait is this right? so in the limit, the magnitude of the direction vec doesnt matter?) Of course, the problem arises that this limit may depend on the sequence in which  $\vec{h}$  approaches  $\vec{0}$  – in such cases we say that the derivative is not defined.

A (perhaps not too helpful) analogue in the discrete context will now be explored. Consider the function  $f : \mathbb{N}^2 \rightarrow \mathbb{N}^3$ ,  $f(m, n) = (n - m, m^2, mn)$ :

$f(x, y)$	$f(1, *)$	$f(2, *)$	$f(3, *)$	$f(4, *)$	$f(5, *)$
$f(*, 1)$	(0, 1, 1)	(1, 1, 2)	(2, 1, 3)	(3, 1, 4)	(4, 1, 5)
$f(*, 2)$	(-1, 4, 2)	(0, 4, 4)	(1, 4, 6)	(2, 4, 8)	(3, 4, 10)
$f(*, 3)$	(-2, 9, 3)	(-1, 9, 6)	(0, 9, 9)	(1, 9, 12)	(2, 9, 15)
$f(*, 4)$	(-3, 16, 4)	(-2, 16, 8)	(-1, 16, 12)	(0, 16, 16)	(1, 16, 20)
$f(*, 5)$	(-4, 25, 5)	(-3, 25, 10)	(-2, 25, 15)	(-1, 25, 20)	(0, 25, 25)

Our function is no longer a sequence, and so a natural extension of the discrete derivative is not immediately clear. We can see that there are two ways we can proceed; we can make a sequence out of this table by choosing a line (either downwards, rightwards, or diagonally) and calculating differences as if values along that line formed a sequence, or we can try to model how the function changes as its input increases by a distance of 1 in the Euclidean/taxicab metric. The former way is the spirit of taking a directional derivative, while the latter is  $\nabla$ . Rearranging the discrete derivative equation  $((\Delta f(n)) * k = f(n + k) - f(n))$  may make generalisation clearer. Extending that equation to our current example, we get something like:

$$(\Delta f(m, n)) \begin{bmatrix} a \\ b \end{bmatrix} = f(m + a, n + b) - f(m, n).$$

That is,  $\Delta f(m, n)$  is a  $3 \times 2$  matrix. From this equation alone however, we are not sure whether  $\Delta f(m, n)$  is well-defined (that is, it only has one value) – its value may as well change for different values of  $a$  and  $b$ . Let us try calculating  $\Delta f(2, 2)$  for our example. Then, letting the entries of  $\Delta f(2, 2)$  be represented by  $a_i$ , we may calculate the difference of  $f(2, 2)$  with  $f(2, 3)$  and  $f(3, 2)$  (points that distance 1 away in the Euclidean metric):

$$\begin{aligned} \begin{bmatrix} a_1 & a_2 \\ a_3 & a_4 \\ a_5 & a_6 \end{bmatrix} \begin{bmatrix} 1 \\ 0 \end{bmatrix} &= f \left( \begin{bmatrix} 2 \\ 2 \end{bmatrix} + \begin{bmatrix} 1 \\ 0 \end{bmatrix} \right) - f \left( \begin{bmatrix} 2 \\ 2 \end{bmatrix} \right) \\ &= \begin{bmatrix} 1 \\ 4 \\ 6 \end{bmatrix} - \begin{bmatrix} 0 \\ 4 \\ 4 \end{bmatrix} \\ &= \begin{bmatrix} 1 \\ 0 \\ 2 \end{bmatrix} \end{aligned}$$

This gives  $a_1 = 1$ ,  $a_3 = 0$ ,  $a_5 = 2$ , but leaves no information for the other entries. They can be obtained by using the offset (0,1):

$$\begin{aligned} \begin{bmatrix} a_1 & a_2 \\ a_3 & a_4 \\ a_5 & a_6 \end{bmatrix} \begin{bmatrix} 0 \\ 1 \end{bmatrix} &= f \left( \begin{bmatrix} 2 \\ 2 \end{bmatrix} + \begin{bmatrix} 0 \\ 1 \end{bmatrix} \right) - f \left( \begin{bmatrix} 2 \\ 2 \end{bmatrix} \right) \\ &= \begin{bmatrix} -1 \\ 5 \\ 2 \end{bmatrix} \end{aligned}$$

Giving  $a_2 = -1$ ,  $a_4 = 5$ ,  $a_6 = 2$ . So our discrete derivative would be:

$$\Delta f(2, 2) = \begin{bmatrix} 1 & -1 \\ 0 & 5 \\ 2 & 2 \end{bmatrix}.$$



From this example, it is easy to see that discrete derivatives are always well-defined, when differences between coordinates which are only 1 unit ahead are considered.

Now, to generalise to the reals, we have to consider the limits of values as the offset approaches  $\vec{0}$ . A naive generalisation from the discrete equation gives the derivative of  $f$  at  $\vec{x}$  as the unique matrix  $D_f(\vec{x})$  satisfying

$$\lim_{\vec{a} \rightarrow \vec{0}} f(\vec{x} + \vec{a}) - f(\vec{x}) - D_f(\vec{x})\vec{a} = 0.$$

But the immediate problem is that any matrix satisfies this equation as  $\vec{a} \rightarrow \vec{0}$ . To generalise to the reals, we note that we want our derivative matrix to represent the change in the function given a unit offset in the input. The change in the function is given by the Cauchy sequence:

$$\left\{ \frac{f(\vec{x} + \vec{a}) - f(\vec{x})}{\|\vec{a}\|} \right\}_{\vec{a}}$$

Of course, this value depends on the choice of the Cauchy sequence for  $\vec{a}$  (even if they're all equal to  $\vec{0}$ ). Now,  $D_f(\vec{x})\vec{a}$  should approximate the change in the function as the input of the function is offset by  $\vec{a}$ . Normalising this to represent an offset by a unit distance gives  $\frac{D_f(\vec{x})\vec{a}}{\|\vec{a}\|}$ . We want (why?)

$$\left\{ \frac{D_f(\vec{x})\vec{a}}{\|\vec{a}\|} \right\}_{\vec{a}} = \left\{ \frac{f(\vec{x} + \vec{a}) - f(\vec{x})}{\|\vec{a}\|} \right\}_{\vec{a}}$$

, where the sequence of  $\vec{a}$  is the same for both sequences. Rephrasing this gives:

$$\lim_{\vec{a} \rightarrow \vec{0}} \frac{f(\vec{x} + \vec{a}) - f(\vec{x}) - D_f(\vec{x})\vec{a}}{\|\vec{a}\|} = 0.$$

Again, it may as well be that  $D_f(\vec{x})$  depends on the sequence of  $\vec{a}$ . However, the standard definition for derivatives demands (?) that  $D_f(\vec{x})$  should be the same for any such sequence.

To demonstrate the uniqueness of the derivative matrix, suppose that  $D_1$  and  $D_2$  satisfy the conditions to be a derivative matrix. Then:

$$\begin{aligned} \lim_{\vec{a} \rightarrow \vec{0}} \left( \frac{f(\vec{x} + \vec{a}) - f(\vec{x}) - D_1\vec{a}}{\|\vec{a}\|} - \frac{f(\vec{x} + \vec{a}) - f(\vec{x}) - D_2\vec{a}}{\|\vec{a}\|} \right) &= 0 \\ \Rightarrow \lim_{\vec{a} \rightarrow \vec{0}} \frac{D_2\vec{a} - D_1\vec{a}}{\|\vec{a}\|} &= 0 \\ \Rightarrow \lim_{\vec{a} \rightarrow \vec{0}} (D_2 - D_1)\hat{a} &= 0 \end{aligned}$$

$(D_2 - D_1)$  must be the zero matrix, as  $(D_2 - D_1)\hat{a} = 0$  for all unit vectors  $\hat{a}$ .

We note that the existence of a derivative matrix for the reals hinges on the property that the change in the function depends only on how the function changes when one goes along the basiss vectors; that is, we are effectively defining the matrix in the same way for discrete derivatives, by considering the function's behaviour along the axes, but then rejecting the derivative if it fails to predict behaviour in any other direction.

Thus, we see that a differentiable function has defined partial derivatives but not necessarily the converse.

(meta: how do lienar functions fit into this? why not multiplicative?)

### 3.2.2 Theorems

**Theorem 3.2.2.1.** A  $C^1$  function is differentiable.

*Proof.* The key idea is that we can successively apply the single-variable mean-value theorem, to approximate one component at a time. Fixing all components but the first, we can treat  $f$  as a

single-variable function with derivative  $f_{x_1}$ , and the mean-value theorem asserts the existence of some  $b_1 \in (0, a_1)$  such that:

$$\begin{aligned} f_{x_1} \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} + \begin{bmatrix} b_1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} \right) &= \frac{f \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} + \begin{bmatrix} a_1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} \right) - f \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} \right)}{a_1} \\ \Rightarrow f \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} + \begin{bmatrix} a_1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} \right) &= f \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} \right) + a_1 f_{x_1} \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} + \begin{bmatrix} b_1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} \right) \end{aligned}$$

Similarly, fixing all variables but the second and applying the mean-value theorem again gives:

$$f \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} + \begin{bmatrix} a_1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} + \begin{bmatrix} 0 \\ a_2 \\ \vdots \\ 0 \end{bmatrix} \right) = f \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} \right) + a_1 f_{x_1} \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} + \begin{bmatrix} b_1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} \right) + a_2 f_{x_2} \left( \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} + \begin{bmatrix} a_1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} + \begin{bmatrix} 0 \\ b_2 \\ \vdots \\ 0 \end{bmatrix} \right)$$

Continuing in this manner gives:

$$f(\vec{x} + \vec{a}) = f(\vec{x}) + \sum_{i=1}^n a_i f_{x_i}(\vec{c}_i),$$

where

$$\vec{c}_i = \vec{x} + \sum_{j=1}^{i-1} a_j \hat{e}_j + b_i \hat{e}_i.$$

Now, let our derivative matrix  $D = [f_{x_1}(\vec{x}) \quad f_{x_2}(\vec{x}) \quad \cdots \quad f_{x_n}(\vec{x})]$ . Then:

$$\begin{aligned} \lim_{\vec{a} \rightarrow \vec{0}} \frac{f(\vec{x} + \vec{a}) - f(\vec{x}) - D\vec{a}}{\|\vec{a}\|} &= \lim_{\vec{a} \rightarrow \vec{0}} \frac{\sum_{i=1}^n a_i f_{x_i}(\vec{c}_i) - \sum_{i=1}^n a_i f_{x_i}(\vec{x})}{\|\vec{a}\|} \\ &= \lim_{\vec{a} \rightarrow \vec{0}} \sum_{i=1}^n \frac{a_i}{\|\vec{a}\|} (f_{x_i}(\vec{c}_i) - f_{x_i}(\vec{x})) \\ &= 0 \end{aligned}$$

The last equality follows by noting that

$$\begin{aligned} \lim_{\vec{a} \rightarrow \vec{0}} \vec{c}_i &= \lim_{\vec{a} \rightarrow \vec{0}} \left( \vec{x} + \sum_{j=1}^{i-1} a_j \hat{e}_j + b_i \hat{e}_i \right) \\ &= \vec{x}, \end{aligned}$$

as  $b_j \in (0, a_j) \Rightarrow (a_j \rightarrow 0 \Rightarrow b_j \rightarrow 0)$ . It follows that  $\lim_{\vec{a} \rightarrow \vec{0}} (f_{x_i}(\vec{c}_i) - f_{x_i}(\vec{x})) = 0$ . Also  $\frac{a_i}{\|\vec{a}\|} \in [0, 1]$ . Thus, each term in the sum of the limit above tends to 0, and so the entire limit is equal to 0 (assuming finite dimensions).  $\square$

**Theorem 3.2.2.2** (Clairaut's Theorem). Where  $f$  is  $C^2$ ,  $f_{xy} = f_{yx}$ .

*Proof.* First, we provide some intuition, starting with discrete case.

$$\begin{array}{ccc}
 \begin{array}{|c|c|c|} \hline 1 & 3 & 4 \\ \hline 7 & 5 & 6 \\ \hline 2 & 1 & 5 \\ \hline \end{array} & \xrightarrow{\partial_x} & \begin{array}{|c|c|} \hline 2 & 1 \\ \hline -2 & 1 \\ \hline -1 & 4 \\ \hline \end{array} \\
 \downarrow \partial_y & & \downarrow \partial_y \\
 \begin{array}{|c|c|c|} \hline 6 & 2 & 2 \\ \hline -5 & -4 & 1 \\ \hline \end{array} & \xrightarrow{\partial_x} & \begin{array}{|c|c|} \hline -4 & 0 \\ \hline 1 & 3 \\ \hline \end{array}
 \end{array}$$

Clairaut's theorem always holds in the discrete case. To see why this is, consider the case for a general 2-by-2 grid:

$$\begin{array}{ccc}
 \begin{array}{|c|c|} \hline a & b \\ \hline c & d \\ \hline \end{array} & \xrightarrow{\partial_x} & \begin{array}{|c|} \hline b-a \\ \hline d-c \\ \hline \end{array} \\
 \downarrow \partial_y & & \downarrow \partial_y \\
 \begin{array}{|c|c|} \hline c-a & d-b \\ \hline \end{array} & \xrightarrow{\partial_x} & \begin{array}{|c|} \hline a+d-b-c \\ \hline \end{array}
 \end{array}$$

The theorem follows from the fact that  $(d-b) - (c-a) = (d-c) - (b-a)$ .

We use a similar logic to prove the case for reals, using the fact that

$$(f(a+h, b+k) - f(a+h, b)) - (f(a, b+k) - f(a, b)) = (f(a+h, b+k) - f(a, b+k)) - (f(a+h, b+k) - f(a, b)).$$

We work with the LHS first. We note that the LHS is a difference between the function where only the first variable has changed. Thus, we may reduce to an expression with  $\partial_x$  using the mean-value theorem:

$$\begin{aligned}
 (f(a+h, b+k) - f(a+h, b)) - (f(a, b+k) - f(a, b)) &= h\partial_x[f(x, y+k) - f(x, y)]_{(a+h', b)} \\
 &= h\partial_x[kf_y(x, y+k')]_{(a+h', b)} \\
 &= hkf_{yx}(a+h', b+k')
 \end{aligned}$$

And by similar logic on the RHS:

$$(f(a+h, b+k) - f(a, b+k)) - (f(a+h, b+k) - f(a, b)) = khf_{xy}(a+h'', b+k'')$$

As LHS=RHS, we get:

$$\begin{aligned}
 hkf_{yx}(a+h', b+k') &= khf_{xy}(a+h'', b+k'') \\
 \Rightarrow f_{yx}(a+h', b+k') &= f_{xy}(a+h'', b+k'') \quad ((h, k) \neq (0, 0)) \\
 \Rightarrow \lim_{(h, k) \rightarrow (0, 0)} f_{yx}(a+h', b+k') &= \lim_{(h, k) \rightarrow (0, 0)} f_{xy}(a+h'', b+k'') \\
 \Rightarrow f_{yx}(a, b) &= f_{xy}(a, b).
 \end{aligned}$$

The last implication follows as  $f$  is  $C^2$ ; the second partial derivatives are continuous.  $\square$

**Theorem 3.2.2.3** (Chain Rule). Consider  $f: \mathbb{R}^a \rightarrow \mathbb{R}^b$  and  $g: \mathbb{R}^b \rightarrow \mathbb{R}^c$ , with  $\vec{x} \in \text{dom}g$ . Then:

$$D_{g \circ f}(\vec{x}) = D_g(f(\vec{x}))D_f(\vec{x}).$$

*Proof.* We want to say something about  $(g \circ f)(\vec{x} + \vec{a})$  with information about  $(g \circ f)(\vec{x})$ . From the definition of the derivative we can write:

$$f(\vec{x} + \vec{a}) = f(\vec{x}) + D_f(\vec{x})\vec{a} + o(a)$$

with  $o(\vec{a})$  denoting the little-o notation;  $o(\vec{a})$  satisfies

$$\lim_{\vec{a} \rightarrow 0} \frac{o(\vec{a})}{\|\vec{a}\|} = 0.$$

Now,

$$\begin{aligned} (g \circ f)(\vec{x} + \vec{a}) &= g(f(\vec{x} + \vec{a})) \\ &= g(f(\vec{x}) + D_f(\vec{x})\vec{a} + o(\vec{a})) \\ &= g(f(\vec{x})) + D_g(f(\vec{x}))(D_f(\vec{x})\vec{a} + o(\vec{a})) + o(D_f(\vec{x})\vec{a} + o(\vec{a})) \\ &= g(f(\vec{x})) + D_g(f(\vec{x}))D_f(\vec{x})\vec{a} + D_g(f(\vec{x}))o(\vec{a}) + o(D_f(\vec{x})\vec{a} + o(\vec{a})) \end{aligned}$$

Now it suffices to show that  $D_g(f(\vec{x}))o(\vec{a}) + o(D_f(\vec{x})\vec{a} + o(\vec{a})) = o(\vec{a})$ . As  $D_g(f(\vec{x}))$  is independent of  $\vec{a}$ , it follows that  $D_g(f(\vec{x}))o(\vec{a}) = o(\vec{a})$ . For brevity, let  $\vec{w} = D_f(\vec{x})\vec{a} + o(\vec{a})$ . We now just want to show that  $o(\vec{w}) = o(\vec{a})$ , i.e.

$$\vec{a} \rightarrow 0 \Rightarrow \frac{o(\vec{w})}{\|\vec{a}\|} \rightarrow 0$$

By definition, we know that  $\vec{w} \rightarrow 0 \Rightarrow \frac{o(\vec{w})}{\|\vec{w}\|} \rightarrow 0$ . But note that if  $\vec{a} \rightarrow 0$ , then  $\vec{w} = D_f(\vec{x})\vec{a} + o(\vec{a}) \rightarrow 0$ , as  $D_f(\vec{x})$  is independent of  $\vec{a}$ . So we have

$$\vec{a} \rightarrow 0 \Rightarrow \frac{o(\vec{w})}{\|\vec{w}\|} \rightarrow 0.$$

To show that  $\frac{o(\vec{w})}{\|\vec{a}\|} \rightarrow 0$ , we will use the fact that  $\frac{o(\vec{w})}{\|\vec{a}\|} = \frac{o(\vec{w})}{\|\vec{w}\|} \frac{\|\vec{w}\|}{\|\vec{a}\|}$ , and show that  $\frac{\|\vec{w}\|}{\|\vec{a}\|}$  is finite.

$$\begin{aligned} \|D_f(\vec{x})\vec{a} + o(\vec{a})\| &\leq \|D_f(\vec{x})\vec{a}\| + \|o(\vec{a})\| \\ &\leq M \|\vec{a}\| + \frac{\|o(\vec{a})\|}{\|\vec{a}\|} \|\vec{a}\| \quad (4) \end{aligned}$$

So,

$$\lim_{\vec{a} \rightarrow \vec{0}} \frac{\|\vec{w}\|}{\|\vec{a}\|} = \lim_{\vec{a} \rightarrow \vec{0}} M + \frac{\|o(\vec{a})\|}{\|\vec{a}\|} = M.$$

Thus,

$$\vec{a} \rightarrow 0 \Rightarrow \frac{o(\vec{w})}{\|\vec{w}\|} \rightarrow 0 \wedge \frac{\|\vec{w}\|}{\|\vec{a}\|} \rightarrow M \Rightarrow \frac{o(\vec{w})}{\|\vec{w}\|} \frac{\|\vec{w}\|}{\|\vec{a}\|} = \frac{o(\vec{w})}{\|\vec{a}\|} \rightarrow 0.$$

□

### 3.2.3 Inverse Function Theorem

Apparently, this theorem is important. (reflect later)

**Definition 3.2.3.1.** Where  $f$  maps a metric space to itself, it is a *contraction* if  $d(f(x), f(y)) \leq c(x \leftrightarrow y)$  with  $0 < c \leq 1$ , and a strict contraction if  $0 < c < 1$ .

**Theorem 3.2.3.1** (Banach Fixed-Point Theorem). Strict contractions have at most one fixed point, and exactly one if  $X$  is non-empty and complete.

*Proof.* If  $f$  were to have two fixed points  $x$  and  $y$ , then

$$\begin{aligned} f(x) \leftrightarrow f(y) &< x \leftrightarrow y \\ &\Rightarrow x \leftrightarrow y < x \leftrightarrow y, \end{aligned}$$

a contradiction.

---

<sup>(4)</sup>Where  $T$  is a matrix,  $\|T\vec{v}\| = \|\vec{v}\| \|T\hat{v}\|$ . But  $f : \{\vec{x} \mid \|\vec{x}\| = 1\} \rightarrow \mathbb{R}$ ,  $f(\vec{x}) = \|T\hat{v}\|$  has a compact domain, so attains a maximum, say a value of  $M$ . Then,  $\|T\vec{v}\| \leq M \|\vec{v}\|$ .

Now suppose that the metric space  $X$  is non-empty and complete. Choose a  $x_0 \in X$  and consider the sequence  $x_n = [f]^n(x_0)$ . That this sequence is Cauchy can be established by noting that  $x_{n+1} \leftrightarrow x_n \leq c(x_n \leftrightarrow x_{n-1}) \Rightarrow x_{n+1} \leftrightarrow x_n \leq c^n(x_1 \leftrightarrow x_0)$ . Then, one can use the triangle inequality to expand  $x_m \leftrightarrow x_n$  and use the geometric series formula. We propose that the limit of the sequence,  $x^*$  (which exists by completeness), is the fixed point of  $f$ .

$$x^* = \langle x_n \rangle = \langle f(x_{n-1}) \rangle = f(\langle x_{n-1} \rangle) = f(x^*).$$

The penultimate equality is justified by continuity of  $f$ . (Why is  $f$  continuous?) □

**Theorem 3.2.3.2** (Inverse Function Theorem). Where  $f$  is  $C^1$  on an open domain with an invertible derivative at  $x_0$ , there is an open subset  $U$  of the domain containing  $x_0$  on which  $f$  is bijective. Also,  $f^-$  (on this restriction) satisfies

$$[Df^-](f(x_0)) = ([Df^-](x_0))^-.$$

The main part of the proof, as presented by Tao, is not exactly intuitive. It hinges on a corollary of the fixed-point theorem that states that if  $g$  is strictly contractive on  $U(0, r)$ ,  $f = g + I$  is injective on that neighbourhood and  $U(0, (1 - c)r) \subset f(U(0, r)) \subset U(0, (1 + c)r)$ . The proof is as follows.

If  $g(x) = g(y)$ , then as  $g$  is a contraction, this implies that  $x = y$ .

To show that the image of  $U(0, r)$  under  $f$  includes  $U(0, (1 - c)r)$ , we “cleverly” abuse the fixed-point theorem. see wikipedia.

*Proof.* The second part is trivial. Assume that  $f$  has an inverse. Then,

$$D_{f \circ f}(x) = I \Rightarrow D_{f^-}(f(x))D_f(x) = I \Rightarrow D_{f^-}(f(x)) = (D_f(x))^-.$$

I cbf writing this shit just read the stuff on wikipedia the c1 argument is better done in Tao, tho. □

### 3.3 Integration

The integral aims to generalise the notion of sums to continuous domains. In the discrete case, we have seen that sums can accumulate successive differences to return values of the original function; for example, we have

$$\begin{aligned} f(a+3) &= f(a) + (f(a+1) - f(a)) + (f(a+2) - f(a+1)) + (f(a+3) - f(a+2)) \\ &= f(a) + \Delta f(a) + \Delta f(a+1) + \Delta f(a+2) \\ &= f(a) + \sum_{q=a}^{a+3-1} \Delta f(q). \end{aligned}$$

Thus, a characterisation of the sum is as follows:

**Definition 3.3.0.1.** The sum,  $\Sigma$ , is an operator acts on a function's difference, with parameters  $a$  and  $a+j-1$ , to give the  $j$ -step difference of the function at  $f(a)$ ; that is,

$$\sum_{q=a}^{a+j-1} \Delta f(q) = \Delta_j f(a).$$

We want to extend this interpretation of the sum into the reals:

**Definition 3.3.0.2.** The integral,  $\int$ , is an operator acts on a function's derivative, with parameters  $a$  and  $a+j$ , to give the  $j$ -step difference of the function at  $f(a)$ ; that is,

$$\int_{q=a}^{a+j} Df(q) = \Delta_j f(a).$$

We have given a declarative definition; we ought to give an imperative construction. Actually, the following imperative definition is also declarative in a sense. As we did with the difference/derivative, we will reframe the sum where the gap between each term becomes arbitrary, and take a sequence where that gap approaches 0. To make this gap arbitrary, we introduce the notion of partitions. Let  $I$  be the interval we are working with (with endpoints  $a$  and  $b$ ), and  $P$  be a partition operator.  $p_i^*$  is the lower bound of the interval  $p_i$ . (Later on, it will be fine for  $p_i^*$  to be any point in the interval.)

$$\sum_{p_i \in P(I)} \Delta_{|p_i|} f(p_i^*) = \Delta_{|I|} f(a).$$

Noting that the above equality holds for any partition at all, it follows that the sequence of the LHS as  $\max_{p_i \in P(I)} |p_i| \rightarrow 0$  is constant, and so converges to the RHS. Let us define  $\text{mesh } P(I) := \max_{p_i \in P(I)} |p_i|$ . Then,

$$\left\langle \sum_{p_i \in P(I)} \Delta_{|p_i|} f(p_i^*) \right\rangle_{\text{mesh } P(I) \rightarrow 0} = \Delta_{|I|} f(a).$$

The summand looks somewhat like the derivative, which we can make more explicit by:

$$\left\langle \sum_{p_i \in P(I)} \hat{\Delta}_{|p_i|} f(p_i^*) |p_i| \right\rangle_{\text{mesh } P(I) \rightarrow 0} = \Delta_{|I|} f(a).$$

It seems reasonable to replace the normalised difference with the derivative as the gap is approaching is 0, and indeed, I will show that

$$\left\langle \sum_{p_i \in P(I)} \hat{\Delta}_{|p_i|} f(p_i^*) |p_i| \right\rangle_{\text{mesh } P(I) \rightarrow 0} = \left\langle \sum_{p_i \in P(I)} Df(p_i^*) |p_i| \right\rangle_{\text{mesh } P(I) \rightarrow 0}.$$

The equivalence follows exactly due to the intuition; the summands become arbitrarily close as one is the convergent value of the other. To make this explicit, let us subtract two terms from the sequence (where each sequence has the same domain/partition structure):

$$\begin{aligned}
 \sum_{p_i \in P(I)} \hat{\Delta}_{|p_i|} f(p_i^*) |p_i| - \sum_{p_i \in P(I)} Df(p_i^*) |p_i| &= \sum_{p_i \in P(I)} \left[ \hat{\Delta}_{|p_i|} f - Df \right] (p_i^*) |p_i| \\
 &= \sum_{p_i \in P(I)} \frac{\varepsilon}{|I|} |p_i| \\
 &= \frac{\varepsilon}{|I|} \sum_{p_i \in P(I)} |p_i| \\
 &= \frac{\varepsilon}{|I|} |I| \\
 &= \varepsilon.
 \end{aligned}$$

The second line can be justified by choosing an partition small enough so that  $Df$  and  $\hat{\Delta}_{|p_i|} f$  are arbitrarily close. Thus, our imperative definition of the integral is:

$$\int_I Df = \left\langle \sum_{p_i \in P(I)} Df(p_i^*) |p_i| \right\rangle_{\text{mesh } P(I) \rightarrow 0}.$$

Comment; this roughly corresponds to riemann integration. An immediate problem is the function  $Q(x)=1$  if  $x$  is irrational, 0 otherwise. Depending on how one chooses tags/partition, the integral evaluates to 0 or 1. we need the stiefjes integral or lebegsue, maybe even kurzweil henstock. discuss this later.

### 3.3.1 Line Integrals

Everything becomes much simpler when interpreting a curve  $\tilde{\gamma}$  as a coordinate transform/change-of-variables, which is discussed elsewhere (actually a bit different because  $\gamma$  changes dimensions; so we need to show that  $\left\langle \frac{|\tilde{\gamma}(p_i)|}{|p_i|} \right\rangle_{|p_i| \rightarrow 0} = \|D\tilde{\gamma}\|$ ). Then, we have

$$\int_{\tilde{\gamma}(I)} f = \int_I f \circ \tilde{\gamma} \|D\tilde{\gamma}\|$$

We can show that this definition satisfies a fundamental-theorem-of-calculus of sorts; let  $\mathbf{T} = \text{unit}(\tilde{\gamma}' \circ t)$ ;  $\mathbf{T}$  is a function only accepting vectors mapped to by  $\gamma$ , and gives the derivative of  $\gamma$  at that point. This is set up so that  $\mathbf{T} \circ \tilde{\gamma} = \gamma'$ . Then,

$$\int_{\tilde{\gamma}(I)} D_{\mathbf{T}} f = \Delta_{|\tilde{\gamma}|} f(\mathbf{a})$$

(Where we define  $|\tilde{\gamma}| = \text{end point} - \text{start point}$ ) This follows easily:

$$\begin{aligned} \int_{\tilde{\gamma}(I)} D_{\mathbf{T}} f &= \int_{\tilde{\gamma}(I)} (Df) \mathbf{T} \\ &= \int_I [Df \circ \tilde{\gamma}] \mathbf{T} \circ \tilde{\gamma} \|D\tilde{\gamma}\| \\ &= \int_I [Df \circ \tilde{\gamma}] D\tilde{\gamma} \\ &= [f \circ \tilde{\gamma}](\mathbf{b}) - [f \circ \tilde{\gamma}](\mathbf{a}) \\ &= \Delta_{|\tilde{\gamma}|} f(\mathbf{a}). \end{aligned}$$

### 3.3.2 Path Integral on Vector Fields

Mathematicians (or some other people) thought it would be good notation to write

$$\int_{\tilde{\gamma}} \mathbf{F} \cdot ds := \int_{\tilde{\gamma}} \mathbf{F} \cdot \mathbf{T}.$$

This is because it gives a neat form of a pseudo-fundamental-theorem-of-calculus:

$$\begin{aligned} \int_{\tilde{\gamma}} \nabla f \cdot ds &= \int_{\tilde{\gamma}} \nabla f \cdot \mathbf{T} \\ &= \int_{\tilde{\gamma}} D_{\mathbf{T}} f \\ &= \Delta_{|\tilde{\gamma}|} f(\mathbf{a}) \end{aligned}$$

and call this the “line integral of a vector field”. We will just use  $\int_{\tilde{\gamma}} \mathbf{F} \cdot \mathbf{T}$  instead of unnecessarily introducing new notation.



### 3.4 Multivariable Integration

We again consider the discrete 2-dimensional difference grid to illustrate concepts in multivariable integration. We want to extend our notion of the integral as an operator which accumulates differences into a difference of a larger interval to multiple dimensions. While the difference of a one dimensional interval  $[a, b]$  is simply  $b - a$ , it isn't immediately clear what the difference of the interval  $[a, b] \times [c, d]$  should be, let alone for arbitrary 2-dimensional regions. Below is a grid of  $f(x, y) = S_1(x + y, x)$ , with partial differences given in blue. For example,  $f(0, 0) = 1$ ,  $f(2, 3) = 35$ , and  $\Delta_x f(1, 2) = 14$ .

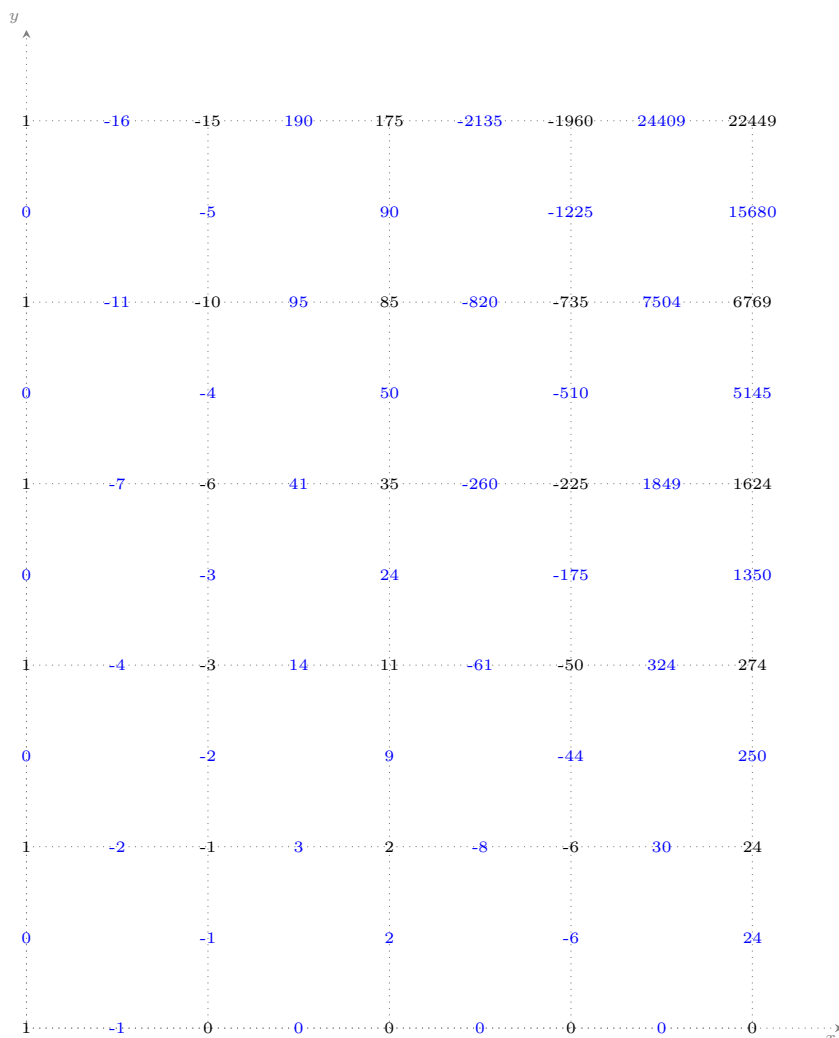
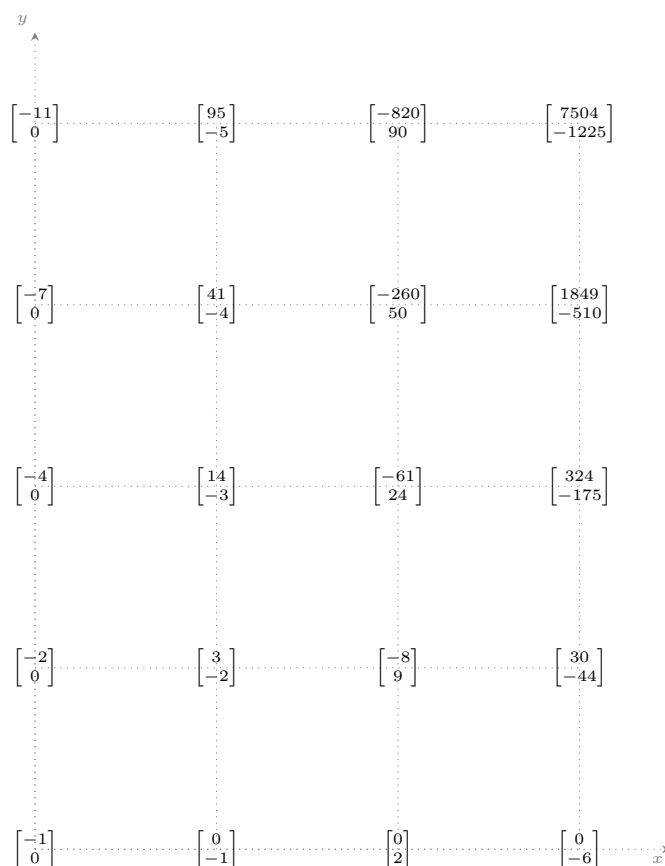


Figure 1: Graph of  $f(x, y) = S_1(x + y, x)$ , with the origin at the bottom-left.

Now, consider the grid of  $\Delta f$ .

Figure 2: Graph of  $\Delta f(x, y)$ , with the origin at the bottom-left.

Let's calculate the sum of  $\Delta f$  over the square  $[0, 1] \times [0, 1]$ , to see what that might represent in [Figure 1](#):

$$\begin{bmatrix} -1 \\ 0 \end{bmatrix} + \begin{bmatrix} 0 \\ -1 \end{bmatrix} + \begin{bmatrix} -2 \\ 0 \end{bmatrix} + \begin{bmatrix} 3 \\ -2 \end{bmatrix} = \begin{bmatrix} 0 \\ -3 \end{bmatrix}$$

It is still not too clear what this value represents. It will help to expand the difference values into their definitions, and mark what this means on the original grid. By summing over  $[0, 2] \times [0, 2]$  we are effectively calculating:

$$\begin{aligned} \Delta f(0, 0) + \Delta f(0, 1) + \Delta f(1, 0) + \Delta f(1, 1) &= \begin{bmatrix} \Delta_x f(0, 0) \\ \Delta_y f(0, 0) \end{bmatrix} + \begin{bmatrix} \Delta_x f(0, 1) \\ \Delta_y f(0, 1) \end{bmatrix} + \begin{bmatrix} \Delta_x f(1, 0) \\ \Delta_y f(1, 0) \end{bmatrix} + \begin{bmatrix} \Delta_x f(1, 1) \\ \Delta_y f(1, 1) \end{bmatrix} \\ &= \begin{bmatrix} f(2, 0) - f(0, 0) + f(2, 1) - f(0, 1) \\ f(0, 2) - f(0, 0) + f(1, 2) - f(1, 0) \end{bmatrix}. \end{aligned}$$

Marking points that are added/subtracted with  $+/-$ , and colouring the ones that count to the first part of the derivative with red, and second part blue, on the original grid:

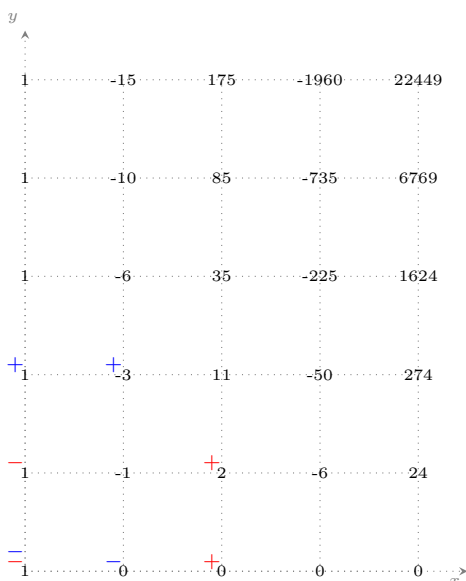


Figure 3: Graph of  $f(x, y) = S_1(x + y, x)$ , with the origin at the bottom-left.

There are a few (equivalent) interpretations of what is going on; that summing the derivative over the square results in

- some sum around the border, where one side is negative and the other positive;
- a collection of  $j$ -step differences in the  $x$  and  $y$  directions;
- a of  $j$ -step difference of a line in the  $x$  and  $y$  directions.

God I hate this subject, I'm just going to write whatever.

$$\begin{aligned} \sum_D \Delta f &= \sum_{(x,y) \in [a,b] \times [c,d]} [\Delta_x f, \Delta_y f] \\ &= \left[ \sum_{y=c}^d (f(b+1, y) - f(a, y)), \sum_{x=a}^b (f(x, d+1) - f(x, c)) \right] \end{aligned}$$

To get the standard orientation, we instead need  $\sum_{x=a}^b (f(x, d+1) - f(x, c))$ ; the sum across the bottom row ( $\sum_{x=a}^b f(x, c)$ ) should be positive, and the top negative. So we have:

$$\sum_D \Delta f \cdot [1 \quad -1] =: \sum_{\partial D} f.$$

We could generalise  $Df$  to just be  $[\Delta_x Q \quad \Delta_y P]$  in which case we get:

$$\sum_D [\Delta_x Q \quad \Delta_y P] \cdot [1 \quad -1] = \sum_{\partial D} \begin{bmatrix} P \\ Q \end{bmatrix} \cdot ds.$$

## 3.5 Shitty notes on topology for gen stokes

### 3.5.1 Definitions

Unlike metric spaces whose foundations centre around the notion of a metric, the more general topological space centres its foundations on the notion of open sets.

**Definition 3.5.1.1.** A *topological space* is a set  $X$  associated with a certain  $\mathcal{T} \subseteq \mathcal{P}(X)$  satisfying:

- $\emptyset, X \in \mathcal{T}$ ;
- closure under the intersection operation;
- closure under the arbitrary union operation.

Any  $\mathcal{T}$  which satisfies the above is called a *topology* on  $X$ , and the elements of the topology are called the *open sets*.

**Example 3.5.1.1.**

- $\mathcal{T} = \{\emptyset, X\}$ , the minimal topology on  $X$ , called the *indiscrete topology*;
- $\mathcal{T} = \mathcal{P}(X)$ , the maximal topology on  $X$ , called the *discrete topology*.
  - Like sets under a discrete metric, every set is open in a discrete topology.

The following definition is basically identical to the one we gave for metric spaces.

**Definition 3.5.1.2** (Interior, Exterior, Boundary Points).

- An *interior* point of a set  $E$  is a point such that there is a neighbourhood containing it, which itself is a subset of  $E$ .
- Otherwise, if there is a neighbourhood containing the point that is completely disjoint from  $E$ , then the point is *exterior*.
- If neither, the point is a *boundary* point.

The set of boundary points is denoted by  $\partial E$ .

**Definition 3.5.1.3** (Topological Convergence). An infinite sequence of points  $(x_i)$  in  $X$  is said to *converge* to  $x \in X$  if for any neighbourhood  $U$  of  $x$ ,  $\exists N \in \mathbb{N}$  s.t.  $i > N \Rightarrow x_i \in U$ .

**Exercise 3.5.1.1.** Give an example of a sequence which converges to more than one point.

*Solution.* A particularly boring solution is to consider the indiscrete topology on  $\mathbb{R}$ , in which any infinite sequences converges to any infinite number. A more insightful solution will be to think more generally; if the point  $y$  is included in every neighbourhood of  $x$ , then we have

$$x \text{ is a limit of the sequence} \iff y \text{ is a limit of the sequence.}$$

So, consider the topology  $\{\emptyset, \mathbb{R}\} \cup \{(a, \infty) \mid a \in \mathbb{R}\}$ . Any neighbourhood of  $a$  contains any number larger than  $a$ . This concept gives rise to the following definition. □

**Definition 3.5.1.4** (Hausdorff Spaces). A topological space is Hausdorff if  $x \neq y \Rightarrow \exists U_x, U_y$  s.t.  $U_x \cap U_y = \emptyset$ .

**Remark 3.5.1.1.** Limit of sequence unique only if topological space is Hausdorff.

**Definition 3.5.1.5** (Open, Closed, Clopen Sets).

A set is

- *closed* if it contains all its boundary points;
- *open* if it contains none;
- *neither* if neither;
- *clopen*, if it has no boundary.

(Note that a clopen set is vacuously open and closed.)

## 4 Applications

### 4.1 Taylor Series

#### 4.1.1 Discrete Analogy

We have previously discussed Taylor series for discrete sequences. We revisit the discrete case but in a slightly different perspective, to allow a clearer generalisation to the continuous case.

Noting that  $f(x+1) = f(x) + \Delta f(x)$ ,  $f(x+2) = f(x) + \Delta f(x) + \Delta f(x+1)$ , and so on, we can write:

$$\begin{aligned} f(x+a) &= f(x) + \Delta f(x) + \Delta f(x+1) + \cdots + \Delta f(x+a-1) \\ &= f(x) + \sum_{q=0}^{a-1} \Delta f(x+q) \end{aligned}$$

But now, we note that in turn, we can write

$$\begin{aligned} \Delta f(x+q) &= \Delta f(x) + \Delta^2 f(x) + \Delta^2 f(x+1) + \cdots + \Delta^2 f(x+q-1) \\ &= \Delta f(x) + \sum_{q_2=0}^{q-1} \Delta^2 f(x+q_2). \end{aligned}$$

Substituting this into ( ) we get:

$$\begin{aligned} f(x+a) &= f(x) + \sum_{q=0}^{a-1} \Delta f(x+q) \\ &= f(x) + \sum_{q=0}^{a-1} \left( \Delta f(x) + \sum_{q_2=0}^{q-1} \Delta^2 f(x+q_2) \right) \\ &= f(x) + \Delta f(x)a + \sum_{q=0}^{a-1} \sum_{q_2=0}^{q-1} \Delta^2 f(x+q_2) \end{aligned}$$

Repeating the process once more:

$$\begin{aligned} f(x+a) &= f(x) + \Delta f(x)a + \sum_{q=0}^{a-1} \sum_{q_2=0}^{q-1} \Delta^2 f(x+q_2) \\ &= f(x) + \Delta f(x)a + \sum_{q=0}^{a-1} \sum_{q_2=0}^{q-1} \left( \Delta^2 f(x) + \sum_{q_3=0}^{q_2-1} \Delta^3 f(x+q_3) \right) \\ &= f(x) + \Delta f(x)a + \Delta^2 f(x) \sum_{q=0}^{a-1} \sum_{q_2=0}^{q-1} (1) + \sum_{q=0}^{a-1} \sum_{q_2=0}^{q-1} \sum_{q_3=0}^{q_2-1} \Delta^3 f(x+q_3) \\ &= f(x) + \Delta f(x)a + \Delta^2 f(x) \frac{a^2}{2} + \sum_{q=0}^{a-1} \sum_{q_2=0}^{q-1} \sum_{q_3=0}^{q_2-1} \Delta^3 f(x+q_3) \end{aligned}$$

Noting that

$$\begin{aligned} \sum_{q_1=0}^{a-1} \sum_{q_2=0}^{q_1-1} \cdots \sum_{q_{k-2}=0}^{q_{k-3}-1} \sum_{q_{k-1}=0}^{q_{k-2}-1} \sum_{q_k=0}^{q_{k-1}-1} (1) &= \sum_{q_1=0}^{a-1} \sum_{q_2=0}^{q_1-1} \cdots \sum_{q_{k-2}=0}^{q_{k-3}-1} \sum_{q_{k-1}=0}^{q_{k-2}-1} (q_{k-1} - 0) \\ &= \sum_{q_1=0}^{a-1} \sum_{q_2=0}^{q_1-1} \cdots \sum_{q_{k-2}=0}^{q_{k-3}-1} \left( \frac{q_{k-2}^2}{2} - 0 \right) \\ &\vdots \end{aligned}$$

$$= \frac{a^k}{k!},$$

we see that the above formulation is equal to the Newton series we derived before, as eventually  $\delta^i f$  becomes 0, and so the iterated sum at the end eventually disappears:

$$f(x+a) = \sum_{q=0}^k \frac{k^q}{q!} [\Delta^q f](x).$$

However when discussing error terms for the continuous Taylor series, it is more convenient to work with this form:

$$f(x+a) = \underbrace{f(x) + \Delta f(x)a + \Delta^2 f(x) \frac{a^2}{2}}_{\text{Taylor Polynomial}} + \underbrace{\sum_{q=0}^{a-1} \sum_{q_2=0}^{q_1-1} \sum_{q_3=0}^{q_2-1} \Delta^3 f(x+q_3)}_{\text{Error}}$$

We can give an upper bound for the error term. For the error term for the  $k-1$ <sup>st</sup> Taylor polynomial, let

$$M = \max_{n \in [0, a]} \{ \Delta^k f(x+q) \}.$$

Then:

$$\begin{aligned} \sum_{q_1=0}^{a-1} \sum_{q_2=0}^{q_1-1} \cdots \sum_{q_k=0}^{q_{k-1}-1} \Delta^k f(x+q_k) &\leq \sum_{q_1=0}^{a-1} \sum_{q_2=0}^{q_1-1} \cdots \sum_{q_k=0}^{q_{k-1}-1} M \\ &= M \sum_{q_1=0}^{a-1} \sum_{q_2=0}^{q_1-1} \cdots \sum_{q_k=0}^{q_{k-1}-1} (1) \\ &= M \frac{a^k}{k!}. \end{aligned}$$

Now, extending this to the continuous case is almost trivial.

#### 4.1.2 Continuous Taylor series

First, we write:

$$f(x+a) = f(x) + \int_0^a f'(x+q) dq$$

But now, we note that in turn, we can write

$$f'(x+q) = f'(x) + \int_0^q f''(x+q_2) dq_2$$

Substituting this into ( ) we get:

$$\begin{aligned} f(x+a) &= f(x) + \int_0^a f'(x+q) dq \\ &= f(x) + \int_0^a \left( f'(x) + \int_0^q f''(x+q_2) dq_2 \right) dq \\ &= f(x) + \int_0^a f'(x) dq + \int_0^a \int_0^q f''(x+q_2) dq_2 dq \\ &= f(x) + f'(x)a + \int_0^a \int_0^q \left( f'(x) + \int_0^q f''(x+q_2) dq_2 \right) dq_2 dq \end{aligned}$$

Repeating the process once more:

$$f(x+a) = f(x) + f'(x)a + \int_0^a \int_0^q f''(x+q_2) dq_2 dq$$

$$\begin{aligned}
&= f(x) + f'(x)a + \int_0^a \int_0^q \left( f''(x) + \int_0^{q_2} f'''(x + q_3) dq_3 \right) dq_2 dq \\
&= f(x) + f'(x)a + \int_0^a \int_0^q f''(x) dq_2 dq + \int_0^a \int_0^q \int_0^{q_2} f'''(x + q_3) dq_3 dq_2 dq \\
&= f(x) + f'(x)a + f''(x) \int_0^a \int_0^q dq_2 dq + \int_0^a \int_0^q \int_0^{q_2} f'''(x + q_3) dq_3 dq_2 dq \\
&= \underbrace{f(x) + f'(x)a + f''(x) \frac{a^2}{2}}_{\text{Taylor Polynomial}} + \underbrace{\int_0^a \int_0^q \int_0^{q_2} f'''(x + q_3) dq_3 dq_2 dq}_{\text{Error}}
\end{aligned}$$

We can generalise the Taylor polynomial term by noting that

$$\int_0^a \int_0^{q_1-1} \cdots \int_0^{q_{k-1}-1} dq_k \cdots dq_2 dq_1 = \frac{a^k}{k!}.$$

Thus the  $k^{\text{th}}$  degree Taylor polynomial of  $f(x + a)$  centred at  $x$  is given by:

$$\sum_{q=0}^k \frac{a^q}{q!} f^{(q)}(x).$$

As before, we can give an upper bound for the error term for the  $k - 1^{\text{st}}$  Taylor polynomial. Let

$$M = \max_{q \in [0, a]} \left\{ f^{(k)}(x + q) \right\}.$$

Then:

$$\begin{aligned}
\int_0^a \int_0^{q_1-1} \cdots \int_0^{q_{k-1}-1} f^{(k)}(x + q_k) dq_k \cdots dq_2 dq_1 &\leq \int_0^a \int_0^{q_1-1} \cdots \int_0^{q_{k-1}-1} M dq_k \cdots dq_2 dq_1 \\
&= M \int_0^a \int_0^{q_1-1} \cdots \int_0^{q_{k-1}-1} dq_k \cdots dq_2 dq_1 \\
&= M \frac{a^k}{k!}
\end{aligned}$$

By a similar argument,  $I \frac{a^k}{k!}$  is a lower bound for the error where

$$I = \min_{q \in [0, a]} \left\{ f^{(k)}(x + q) \right\}.$$

Thus, this gives us  $I \frac{a^k}{k!} \leq \text{Error} \leq M \frac{a^k}{k!}$ , and so  $\exists C \in [I, M]$  such that  $\text{Error} = C \frac{a^k}{k!}$ . By the intermediate value theorem,  $\exists q \in [0, a]$  such that  $\text{Error} = f(x + q) \frac{a^k}{k!}$ .

#### 4.1.3 Multivariate Taylor Series

To extend our Taylor series to multivariate continuous functions, it is most convenient to interpret the multivariate function as a univariate function using directional derivatives, rather than using the bottom-up approach used previously.

$$\begin{aligned}
f(\vec{x} + \vec{a}) &= f(\vec{x} + |\vec{a}|\hat{a}) \\
&= f(\vec{x}) + \int_0^{|\vec{a}|} D_{\hat{a}} f(\vec{x} + q\hat{a}) dq \\
&= f(\vec{x}) + \int_0^{|\vec{a}|} D_{\hat{a}} f(\vec{x}) + \int_0^q D_{\hat{a}}^2 f(\vec{x} + q_2\hat{a}) dq_2 dq \quad (5) \\
&= f(\vec{x}) + D_{\hat{a}} f(\vec{x})|\vec{a}| + \int_0^{|\vec{a}|} \int_0^q D_{\hat{a}}^2 f(\vec{x} + q_2\hat{a}) dq_2 dq
\end{aligned}$$

$$\begin{aligned}
& \vdots \\
& = \underbrace{f(\vec{x}) + D_{\hat{a}}f(\vec{x})|\vec{a}| + D_{\hat{a}}^2f(\vec{x})\frac{|\vec{a}|^2}{2}}_{\text{Taylor Polynomial}} + \underbrace{\int_0^{|\vec{a}|} \int_0^q \int_0^{q_2} D_{\hat{a}}^3f(\vec{x} + q_3\hat{a}) dq_3 dq_2 dq}_{\text{Error}}
\end{aligned}$$

Thus, the Taylor polynomial is given by:

$$\sum_{q=0}^k \frac{|\vec{a}|^q}{q!} D_{\hat{a}}^q f(\vec{x}).$$

Using our formula for directional derivatives,

$$\begin{aligned}
D_{\vec{a}}f &= D_f \vec{a} \\
&= \nabla f \cdot \vec{a} \\
&= [\partial_{x_1}f, \dots, \partial_{x_n}f] \begin{bmatrix} a_1 \\ \vdots \\ a_n \end{bmatrix} \\
&= \sum_{i=1}^n a_i \partial_{x_i} f
\end{aligned}$$

and noting that

$$\begin{aligned}
|\vec{a}|D_{\hat{a}} &= |\vec{a}|D_f \hat{a} \\
&= D_f \vec{a} \\
&= D_{\vec{a}}f,
\end{aligned}$$

we may express the Taylor polynomial in terms of partials:

$$\begin{aligned}
\sum_{q=0}^k \frac{|\vec{a}|^q}{q!} D_{\hat{a}}^q f(\vec{x}) &= \sum_{q=0}^k \frac{1}{q!} D_{\vec{a}}^q f(\vec{x}) \\
&= \sum_{q=0}^k \frac{1}{q!} [[D_{\vec{a}}]^q f](\vec{x}) \\
&= \sum_{q=0}^k \frac{1}{q!} \left[ \left[ \sum_{i=1}^n a_i \partial_{x_i} \right]^q f \right](\vec{x})
\end{aligned}$$

As before, we can give an upper bound for the error term for the  $k - 1^{\text{st}}$  Taylor polynomial. Let

$$M = \max_{q \in [0, |\vec{a}|]} \{D_{\hat{a}}^k f(\vec{x} + q\hat{a})\}.$$

Then the upper bound for the error is given by

$$\begin{aligned}
\frac{|\vec{a}|^k}{k!} M &= \frac{|\vec{a}|^k}{k!} D_{\vec{a}}^k f(\vec{x} + q'\hat{a}) \\
&= \frac{1}{k!} D_{\vec{a}}^k f(\vec{x} + q'\hat{a}) \\
&= \frac{1}{k!} \left[ \left[ \sum_{i=1}^n a_i \partial_{x_i} \right]^k f \right](\vec{x} + q'\hat{a})
\end{aligned}$$

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<sup>(5)</sup>Note the potentially confusing notation here.  $D_{\hat{a}}^q f(\vec{x})$  means to first apply the operator  $[D_{\hat{a}}]^q$  to  $f$ , then use the resulting operator on  $\vec{x}$ .



## 4.2 Classification of points in 2 dimensions

If at a point  $\vec{x}$ ,  $D_f(\vec{x})$  is the zero matrix, and the the second directional derivative is:

- positive for all directions, then that point is a local minimum.
- negative for all directions, then that point is a local maximum.
- positive for some directions and negative for others, then that point is a saddle.

So, we want to know when  $[D_{\vec{a}}f]^2(\vec{x})$  is positive or negative. The problem becomes simple when the second directional derivative is re-expressed:

$$\begin{aligned}
 [D_{\vec{a}}]^2 f &= D_{\vec{a}}(D_{\vec{a}}f) \\
 &= D_{\vec{a}}(D_f \vec{a}) \\
 &= D_{\vec{a}}([f_{x_1} \ f_{x_2} \ \cdots \ f_{x_n}] \vec{a}) \\
 &= D_{\vec{a}}([a_1 f_{x_1} + a_2 f_{x_2} + \cdots + a_n f_{x_n}]) \\
 &= D([a_1 f_{x_1} + a_2 f_{x_2} + \cdots + a_n f_{x_n}]) \vec{a} \\
 &= [a_1 f_{x_1 x_1} + \cdots + a_n f_{x_n x_1} \quad a_1 f_{x_1 x_2} + \cdots + a_n f_{x_n x_2} \quad \cdots \quad a_1 f_{x_1 x_n} + \cdots + a_n f_{x_n x_n}] \vec{a} \\
 &= [a_1 \ a_2 \ \cdots \ a_n] \begin{bmatrix} f_{x_1 x_1} & f_{x_1 x_2} & \cdots & f_{x_1 x_n} \\ f_{x_2 x_1} & f_{x_2 x_2} & & f_{x_2 x_n} \\ \vdots & & \ddots & \vdots \\ f_{x_n x_1} & f_{x_n x_2} & \cdots & f_{x_n x_n} \end{bmatrix} \vec{a} \\
 &= \vec{a}^T H_f \vec{a}
 \end{aligned}$$

Thus, the problem reduces to knowing when  $\vec{a}^T H_f \vec{a}$  is positive or negative when applied to a point  $\vec{x}$ . Note however, that  $H_f$  is a symmetric matrix, and so there exists a matrix  $P$  such that  $P^{-1} D P = H_f$ , where  $D$  is diagonal. See 5.2.1 for a proof. (Note that the eigenvalues are functions, as we have left  $H_f$  as an operator.)

**Theorem 4.2.0.1.** For a symmetric matrix  $H_f$  with eigenvalues  $\lambda_i > 0$ ,  $i \in \{1, \dots, n\}$ ,

$$\forall \vec{u} \neq \vec{0}, \quad \vec{u}^T H_f \vec{u} > 0 \iff \forall i \in \{1, \dots, n\}, \quad \lambda_i > 0,$$

and the same statement with  $<$  replaced by  $\leq$ ,  $>$  or  $\geq$  also holds. Also,

$$\exists \vec{u}, \vec{w} \text{ s.t. } \vec{u}^T H_f \vec{u} > 0 \wedge \vec{w}^T H_f \vec{w} < 0 \iff \exists i, j \text{ s.t. } \lambda_i > 0 \wedge \lambda_j < 0.$$

*Proof.* Let  $\vec{v} = P \vec{u}$ , where  $P$  is the appropriate change of basis matrix.  $P^T = P^{-1}$  by orthogonal diagonalisation.

$$\begin{aligned}
 \vec{u}^T H_f \vec{u} &= \vec{u}^T P^T D P \vec{u} \\
 &= \vec{v}^T D \vec{v} \\
 &= [v_1 \ \cdots \ v_n] \begin{bmatrix} \lambda_1 & & \\ & \ddots & \\ & & \lambda_n \end{bmatrix} \begin{bmatrix} v_1 \\ \vdots \\ v_n \end{bmatrix} \\
 &= \sum_{i=1}^n \lambda_i v_i^2
 \end{aligned}$$

So  $\lambda_i > 0 \Rightarrow \forall \vec{u} \neq \vec{0}, \vec{u}^T H_f \vec{u} > 0$ . At least one  $v_i$  is guaranteed to be non-zero as  $P$  is a basis.

Conversely, for an eigenvalue  $\lambda_i$ , consider its unit eigenvector  $\hat{v}$ .  $\hat{v}^T H_f \hat{v} > 0$  by hypothesis but  $\hat{v}^T H_f \hat{v} = \hat{v}^T \lambda_i \hat{v} = \lambda_i$ , so  $\lambda_i > 0$ . The other two statements follow similarly.  $\square$

**Definition 4.2.0.1.** A symmetric matrix is [positive/positive semi-/negative/negative semi-] definite if all eigenvalues are  $[> / \geq / < / \leq] 0$ , respectively. It is indefinite if it is neither positive semi- or negative semi-definite.

From our theorem above, we see that a critical point is a minimum/maximum iff  $H_f$  positive/negative semi-definite on it.

An effective method to check positive/negative definiteness of a Hermitian matrix is through Sylvester's Criterion, to which we give an unintuitive proof below.

**Theorem 4.2.0.2** (Sylvester's Criterion). A matrix  $M$  is positive definite iff all its leading principal minors are positive.

*Proof.* Let  $\delta_k$  denote the  $k^{\text{th}}$  leading principal minor. We prove the theorem by induction. The base case is trivial. Now assuming that the theorem holds for matrices of size  $n \times n$ , let  $A$  be a  $(n+1) \times (n+1)$  matrix with all leading principal minors positive.

To show that  $A$  cannot have 2 or more negative eigenvalues, we consider the unintuitively clever construction  $\vec{w} = v_n \vec{u} - u_n \vec{v}$ , where  $\vec{u}$  and  $\vec{v}$  are eigenvectors corresponding to the two negative eigenvalues. By construction,  $w_n = 0$  so  $\vec{w}^T A \vec{w}$  is equivalent to evaluating the quadratic form of  $A^{(n)}$  (the  $n \times n$  upper-left submatrix) at  $(w_1, \dots, w_n)$ , which is greater than 0 by hypothesis –  $A^{(n)}$  is positive definite. But,

$$\begin{aligned} (v_n \vec{u} - u_n \vec{v})^T A (v_n \vec{u} - u_n \vec{v}) &= (v_n \vec{u}^T - u_n \vec{v}^T) A (v_n \vec{u} - u_n \vec{v}) \\ &= (v_n \vec{u}^T - u_n \vec{v}^T) (A v_n \vec{u} - A u_n \vec{v}) \\ &= v_n \vec{u}^T (A v_n \vec{u} - A u_n \vec{v}) - u_n \vec{v}^T (A v_n \vec{u} - A u_n \vec{v}) \\ &= v_n^2 \vec{u}^T A \vec{u} + u_n^2 \vec{v}^T A \vec{v} \\ &< 0. \end{aligned}$$

Also, there cannot be a single negative eigenvalue by the positivity of  $\delta_{n+1}$ . This also implies that no eigenvalues are 0, and thus all eigenvalues are positive.  $\square$

Note that the determinant of a matrix is the product of its eigenvalues. The proof is as follows. The solutions to the characteristic equation  $\det(A - \lambda I)$  is given by the eigenvalues  $\lambda_i$ . So

$$\det(A - \lambda I) = a \prod_{i=1}^n (\lambda - \lambda_i) \quad (a \in \mathbb{R})$$

From the Laplace algorithm for determinants, it can be inductively seen that the leading coefficient of the characteristic polynomial is  $(-1)^n$ . And so, for the leading coefficients to match on both sides, we require  $a = (-1)^n$ . Now, consider the constant term of the characteristic polynomial, which is given when  $\lambda = 0$ . This gives

$$\det(A) = (-1)^n \prod_{i=1}^n \lambda_i$$

And so for  $2 \times 2$  matrices,  $\det(A) > 0 \wedge a_{1,1} > 0 \Leftrightarrow \lambda_1 > 0$ ,  $\det(A) > 0 \wedge a_{1,1} < 0 \Leftrightarrow \lambda_1 < 0$  and  $\det(A) < 0 \Leftrightarrow \lambda_1 > 0 \wedge \lambda_2 < 0$ .

#### 4.2.1 Lagrange Multipliers

Suppose we wanted to identify the extrema of  $f$ , not on its domain but on  $\vec{x}$  such that  $g(\vec{x}) = 0$ . Then, intuitively it follows that  $\vec{x}_0$  is an extrema iff  $D_{\vec{a}} f(\vec{x}_0) = 0$  for all  $\vec{a}$  such that one is able to travel in the direction of  $\vec{a}$  and still satisfy  $g$  – that is, for some small enough  $\varepsilon$ ,  $g(\vec{x} + t\vec{a}) = 0$ ,  $\forall t \in [0, \varepsilon]$ . An equivalent characterisation is that  $\vec{a}$  satisfies  $D_{\vec{a}} g(\vec{x}_0) = 0$ .

To avoid headaches, I skip any proof of Lagrange multipliers.

### 4.3 Vector Calculus

#### 4.3.1 TNB

**Definition 4.3.1.1.** A scalar/vector field is a function from vectors to scalars/vectors.

**Definition 4.3.1.2.**

$$\nabla := \begin{bmatrix} D_{x_1} \\ D_{x_2} \\ \vdots \\ D_{x_n} \end{bmatrix}.$$

**Definition 4.3.1.3.** The tangent, normal, and binormal vectors  $\vec{T}$ ,  $\vec{N}$  and  $\vec{B}$  of a curve  $\vec{r}(t)$  are defined as:

$$\vec{T} := \frac{dr}{dt} \frac{dt}{ds}$$

Note that  $s(t + \varepsilon) - s(t) = \|\vec{r}(t + \varepsilon) - \vec{r}(t)\|$

Alright, being rigorous here probably makes things worse. We proceed with intuition. Consider a curve  $\vec{r}(t)$ . It is easy to see that the direction of  $\vec{r}'(t)$  signifies the direction of where the curve is heading, while the magnitude is the speed at which the arclength is being covered. Let  $s(t)$  be the arclength as a function of time. Then,

$$s'(t) := \|\vec{r}'(t)\|,$$

so

$$s(t) = \int_{t_0}^t \|\vec{r}'(\tau)\| d\tau.$$

From now on, we also parametrise our curve with arclength. This is convenient when defining concepts such as curvature and torsion, which otherwise depend on the speed at which one travels along the curve. When the curve is parametrised by arclength, we remove this ambiguity by fixing the speed.

To perhaps make the distinction clearer we let  $\vec{\varsigma}$  denote the same curve parametrised by arclength; that is,  $\vec{\varsigma} \circ s = \vec{r}$  (or if the notation is confusing,  $\vec{\varsigma}(s(t)) = \vec{r}(t)$ ). Also,  $\vec{T}_s$  and  $\vec{T}_t$  will be the tangent vector parametrised by arclength and time respectively, with  $\vec{T}_s \circ s = \vec{T}_t$ , with similar conventions for the normal and binormal vectors. Also,  $X'$  denotes the derivative of  $X$  with respect to its parametrisation.

We define the tangent vector  $\vec{T}_t$  to be the unit vector pointing in the direction of  $\vec{r}'$ ; that is,

$$\begin{aligned} \vec{T}_t &:= \frac{\vec{r}'}{\|\vec{r}'\|} \\ &= \frac{(\vec{\varsigma} \circ s)'}{s'} \\ &= \frac{(\vec{\varsigma}' \circ s)s'}{s'} && \text{(Chain Rule)} \\ &= \vec{\varsigma}' \circ s \\ &=: \vec{T}_s \circ s. \end{aligned}$$

Note that we need  $\vec{r}' \neq \vec{0}$ . We quantify how much the direction of the curve is changing by analysing the change in the tangent vector i.e change in the direction of the curve. Thus, the direction of  $\vec{T}_s'$  signifies the direction of where the direction of the curve is heading, while the magnitude quantifies how fast the direction is changing. (Note that by using arclength, we quantify the change in direction per change in unit arclength, instead of change in direction per change in distance travelled in 1 unit of time.) This magnitude is assigned the “curvature”, symbolised by  $\kappa$ . The normal vector  $\vec{N}_s$  is defined to be the unit vector pointing in the direction of  $\vec{T}_s'$ ; that is,

$$\kappa := \left\| \vec{T}_s' \right\|;$$

$$\begin{aligned}\vec{N}_s &:= \frac{\vec{T}_s'}{\|\vec{T}_s'\|} \\ \Rightarrow \vec{T}_s' &= \kappa \vec{N}_s,\end{aligned}$$

and

$$\begin{aligned}(\vec{T}_s \circ s)' &= (\vec{T}_s' \circ s)s' \\ \Rightarrow T_t' &= (\kappa \vec{N}_s \circ s)s' \\ &= s' \kappa \vec{N}_s \circ s \\ &= s' \kappa \vec{N}_t.\end{aligned}$$

We note that a vector valued function is always orthogonal to its derivative, given it has a constant norm. A simple proof is by using the product rule for dot products.

We now want to define the torsion of the curve,  $\tau$ . If one were to keep the tangent vector constant, the curve would travel in a single direction, with 0 curvature; if one were to keep the tangent and normal vectors constant, the curve would travel in a single plane, with 0 torsion.

We define the binormal vector  $\vec{B} := \vec{T} \times \vec{N}$  to characterise the plane on which the curve would live in if the tangent and normal vectors at that instant were to be kept constant. The change in this vector,  $\vec{B}_s'$ , quantifies (...) the change in the normal vector of the plane (per unit change in arclength). It is hard to intuitively reduce this vector into a single word we are familiar with (maybe there is idk). However, it turns out that this direction is equivalent to  $\vec{N}$ .

$$\begin{aligned}\vec{B}_s' &= (\vec{T}_s \times \vec{N}_s)' \\ &= \vec{T}_s' \times \vec{N}_s + \vec{T}_s \times \vec{N}_s' \\ &= \vec{T}_s \times \vec{N}_s'\end{aligned}$$

Noting that  $\vec{B}_s'$  is orthogonal to  $\vec{B}$  (as it is unit (because  $\vec{T}$  and  $\vec{N}$  are unit)) and  $\vec{T}$  (from above), we conclude that it must be pointing in the same direction (up to a factor of -1) as  $\vec{N}$ . (it seems like theres not factor of -1 though?) By convention, the torsion is defined to be the negative of the coefficient of  $\vec{N}_s$ :

$$\begin{aligned}\vec{B}_s' &= -\tau \vec{N}_s; \\ \vec{B}_t' &= -s' \tau \vec{N}_t;\end{aligned}$$

**Theorem 4.3.1.1** (Serret-Frenet Formula). When parametrised by arclength, we have:

$$\begin{bmatrix} \vec{T} \\ \vec{N} \\ \vec{B} \end{bmatrix}' = \begin{bmatrix} 0 & \kappa & 0 \\ -\kappa & 0 & \tau \\ 0 & -\tau & 0 \end{bmatrix} \begin{bmatrix} \vec{T} \\ \vec{N} \\ \vec{B} \end{bmatrix}.$$

*Proof.* We have already shown the top and bottom rows. We show the middle row by proving the matrix must be skew-symmetric. Let

$$Q = \begin{bmatrix} \vec{T} \\ \vec{N} \\ \vec{B} \end{bmatrix}.$$

We note that  $Q$  forms an orthonormal basis, so  $Q^{-1} = Q^T$ . The matrix in question is  $Q'Q^{-1} = Q'Q^T$ . The trick is apply the product rule to  $QQ^T = I$ :

$$\begin{aligned}(QQ^T)' &= Q'Q^T + Q(Q^T)' \\ &= Q'Q^T + Q(Q')^T\end{aligned}$$

$$\begin{aligned}
&= Q'Q^T + (Q'Q^T)^T \\
\Rightarrow Q'Q^T &= -(Q'Q^T)^T
\end{aligned}$$

□

**Corollary 4.3.1.1.** Using the chain rule, we get:

$$\begin{bmatrix} \vec{T}_t \\ \vec{N}_t \\ \vec{B}_t \end{bmatrix}' = s' \begin{bmatrix} 0 & \kappa & 0 \\ -\kappa & 0 & \tau \\ 0 & -\tau & 0 \end{bmatrix} \begin{bmatrix} \vec{T}_t \\ \vec{N}_t \\ \vec{B}_t \end{bmatrix}.$$

The form in the corollary leads to explicit expressions for curvature and torsion in terms of  $\vec{r}$ . A direct approach from trying to calculate  $\kappa$  as  $\|\vec{T}_s'\|$  does not work (for some reason). Instead, we may guess that computing  $\kappa$  and  $\tau$  involves taking the second and third derivatives of  $\vec{r}$ , and compute those expressions first. We try to calculate  $\tau$ .  $\kappa$  will follow easily. We have

$$\begin{aligned}
\vec{r}' &= \|\vec{r}'\| T_t \\
&= s' T_t,
\end{aligned}$$

and so

$$\begin{aligned}
\vec{r}'' &= s'' T_t + s' T_t' \\
&= s'' T_t + (s')^2 \kappa \vec{N}_t,
\end{aligned}$$

and

$$\begin{aligned}
\vec{r}''' &= s''' T_t + s'' T_t' + 2s' s'' \kappa \vec{N}_t + (s')^2 \kappa \vec{N}_t' \\
&= s''' T_t + s'' s' \kappa \vec{N}_t + 2s' s'' \kappa \vec{N}_t + (s')^3 \kappa (-\kappa \vec{T}_t + \tau \vec{B}_t) \\
&= (s''' - (s')^3 \kappa^2) T_t + 3s' s'' \kappa \vec{N}_t + (s')^3 \kappa \tau \vec{B}_t
\end{aligned}$$

We can isolate the  $\tau$  term in the coefficient of  $\vec{B}_t$  by taking a dot product with  $\vec{B}_t$ . Noting that our expressions for  $\vec{r}'$  and  $\vec{r}''$  include  $\vec{T}$  and  $\vec{N}$ , and  $\vec{B} = \vec{T} \times \vec{N}$ , we obtain an expression for the binormal vector:

$$\begin{aligned}
\vec{r}' \times \vec{r}'' &= (s')^3 \kappa \vec{B}_t \\
\Rightarrow \|\vec{r}' \times \vec{r}''\| &= (s')^3 \kappa \\
\Rightarrow (\vec{r}' \times \vec{r}'') \cdot \vec{r}''' &= ((s')^3 \kappa)^2 \tau \\
\Rightarrow \tau &= \frac{(\vec{r}' \times \vec{r}'') \cdot \vec{r}'''}{\|\vec{r}' \times \vec{r}''\|^2}.
\end{aligned}$$

### 4.3.2 Divergence and Curl

**Definition 4.3.2.1.** The *divergence* of a vector field  $\mathbf{F} : \mathbb{R}^n \rightarrow \mathbb{R}^n$  at  $\mathbf{v}$  is

$$\operatorname{div} \mathbf{F}(\mathbf{v}) := \left\langle \frac{1}{|U_{\mathbf{v}}(r)|} \int_{\partial U_{\mathbf{v}}(r)} \mathbf{F} \cdot \hat{\mathbf{n}} \right\rangle_{r \rightarrow 0},$$

The definition is the same if  $U_{\mathbf{v}}(r)$  were to be replaced with any other region containing  $\mathbf{v}$  which approaches 0. Actually I'm not sure but probably.

**Proposition 4.3.2.1.**

$$\operatorname{div} \mathbf{F} = \nabla \cdot \mathbf{F}.$$

*Proof.* The trick is to consider the volumes  $V$  as hypercubes with sidelength, say,  $h$ . Now, consider one of the faces  $S_i$  of the hypercube, say the one with the normal whose only non-zero component is the  $i^{\text{th}}$  one. As we are taking the dot product with the normal, we are only concerned with the  $i^{\text{th}}$  component of  $\mathbf{F}$ . The flux from this face would be:

$$\int_{\mathbf{x} \in S_i} F_i(\mathbf{v} + \mathbf{x}) = \int_{t_n = -h/2}^{h/2} \cdots \int_{t_1 = -h/2}^{h/2} F_i(v_1 + t_1, \dots, v_i + h/2, \dots, v_n + t_n)$$

From this it is clear that

$$\left\langle \int_{\mathbf{x} \in S_i} F_i(\mathbf{v} + \mathbf{x}) \right\rangle_{h \rightarrow 0} = \left\langle h^{n-1} F_i(\mathbf{v} + \frac{h}{2} \mathbf{x}_i) \right\rangle_{h \rightarrow 0}.$$

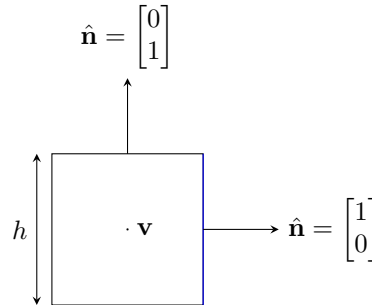
So, we shall just consider the flux of a face with the RHS. Noting that a hypercube has two faces in any given axis, we can write a term in the sequence from the definition of divergence as:

$$\frac{1}{h^n} \sum_{i=1}^n h^{n-1} \left( F_i(\mathbf{v} + \frac{h}{2} \mathbf{x}_i) - F_i(\mathbf{v} - \frac{h}{2} \mathbf{x}_i) \right) = \sum_{i=1}^n \frac{F_i(\mathbf{v} + \frac{h}{2} \mathbf{x}_i) - F_i(\mathbf{v} - \frac{h}{2} \mathbf{x}_i)}{h}$$

But as  $V \rightarrow 0 \iff h \rightarrow 0$ , the summand becomes the partial derivative of  $F_i$  in the  $i^{\text{th}}$  component, and so the expression becomes:

$$\sum_{i=1}^n [\partial_i F_i](\mathbf{v}) = \nabla \cdot \mathbf{F}.$$

To help visualisation, here is part of a 2-dimensional example:



The flux across the blue face is given by:

$$\int_{t=-h/2}^{h/2} F_1(v_1 + \frac{h}{2}, v_2 + t) \approx h F_1(\mathbf{v} + \frac{h}{2} \mathbf{x}_1).$$

□

**Theorem 4.3.2.1** (Gauss' Divergence Theorem). Where  $\mathbf{F} : \mathbb{R}^n \rightarrow \mathbb{R}^n$  is sufficiently differentiable and  $V$  is a compact volume with a piecewise smooth boundary,

$$\int_V \operatorname{div} \mathbf{F} = \oint_{\partial V} \mathbf{F} \cdot \hat{\mathbf{n}}.$$

*Proof.* The key idea of the proof is that the flux of a volume,  $\Phi_F(V)$ , is additive; that is, if  $V = V_1 + V_2$ , we have  $\Phi_F(V) = \Phi_F(V_1) + \Phi_F(V_2)$ . See Wikipedia for a good visualisation. Anyhow, this gives us

$$\begin{aligned} \int_{\partial V} \mathbf{F} \cdot \hat{\mathbf{n}} &= \sum_{V_i \in \text{Slice}(V)} \int_{\partial V_i} \mathbf{F} \cdot \hat{\mathbf{n}} \\ &= \sum_{V_i \in \text{Slice}(V)} \frac{1}{|V_i|} \int_{\partial V_i} \mathbf{F} \cdot \hat{\mathbf{n}} |V_i|. \end{aligned}$$

Now, taking  $V_i \rightarrow 0$ :

$$\begin{aligned} \left\langle \sum_{V_i \in \text{Slice}(V)} \frac{1}{|V_i|} \int_{\partial V_i} \mathbf{F} \cdot \hat{\mathbf{n}} |V_i| \right\rangle_{V_i \rightarrow 0} &= \left\langle \sum_{V_i \in \text{Slice}(V)} \operatorname{div} \mathbf{F} |V_i| \right\rangle_{V_i \rightarrow 0} \\ &= \int_V \operatorname{div} \mathbf{F}. \end{aligned}$$

Note that the continuity of  $\mathbf{F}$  guarantees that the sequence is unchanged in the limit.  $\square$

**Theorem 4.3.2.2** (Stokes' Theorem). Where  $\mathbf{F} : \mathbb{R}^n \rightarrow \mathbb{R}^n$  is sufficiently differentiable and  $V$  is a compact volume with a piecewise smooth boundary,

$$\int_S \operatorname{curl} \mathbf{F} = \oint_{\partial S} \mathbf{F} \cdot \hat{\mathbf{t}}.$$

*Proof.* The proof is extremely similar. The key idea (again) is that the circulation of a surface,  $C_F(S)$ , is additive; that is, if  $S = S_1 + S_2$ , we have  $C_F(S) = C_F(S_1) + C_F(S_2)$ . See Wikipedia for a good visualisation. Anyhow, this gives us

$$\begin{aligned} \oint_{\partial S} \mathbf{F} \cdot \hat{\mathbf{t}} &= \sum_{S_i \in \text{Slice}(S)} \oint_{\partial S_i} \mathbf{F} \cdot \hat{\mathbf{t}} \\ &= \sum_{S_i \in \text{Slice}(S)} \frac{1}{|S_i|} \oint_{\partial S_i} \mathbf{F} \cdot \hat{\mathbf{t}} |S_i|. \end{aligned}$$

Now, taking  $S_i \rightarrow 0$ :

$$\begin{aligned} \left\langle \sum_{S_i \in \text{Slice}(S)} \frac{1}{|S_i|} \oint_{\partial S_i} \mathbf{F} \cdot \hat{\mathbf{t}} |S_i| \right\rangle_{S_i \rightarrow 0} &= \left\langle \sum_{S_i \in \text{Slice}(S)} \operatorname{curl} \mathbf{F} |S_i| \right\rangle_{S_i \rightarrow 0} \\ &= \int_S \operatorname{curl} \mathbf{F}. \end{aligned}$$

Note that the continuity of  $\mathbf{F}$  guarantees that the sequence is unchanged in the limit.  $\square$

### 4.3.3 Change of variables formula

**Theorem 4.3.3.1.** Under appropriate circumstances,

$$\int_{\Phi(V)} f = \int_V f \circ \Phi \det(D\Phi).$$

*Proof.* An alternative characterisation which will help with the change-of-variables theorem is the following:

$$Df(x_0) := \left\langle \frac{|f(L)|}{|L|} \right\rangle_{|L| \rightarrow 0},$$

where each  $L$  is a line/interval containing  $x_0$ . We are guaranteed to have a simple notion of  $L$  as continuous maps preserve intervals, so we need not delve into measure theory. So we can write

$$|f(L)| = \max_{x \in L} f(x) - \min_{x \in L} f(x).$$

The proof of equivalence is as follows. Consider a sequence for

$$\left\langle \hat{\Delta}_h f(x_0) \right\rangle_{h \rightarrow 0}.$$

Now construct a sequence with the effectively the same domain, but expressed in terms of lines, and consider

$$\left\langle \frac{|f(L)|}{|L|} \right\rangle_{|L| \rightarrow 0}.$$

Here is a weird proof that I thought of. I will show that for any  $\delta > 0$  we can find

$|f(L)| = |f(x_0 + h) - f(x_0)|$  with  $0 < |L| < \delta$ , where  $L = [x_0, x_0 + h]$ . Suppose otherwise; that for all  $0 < \delta$ , the maximum value of  $f$  on  $[x_0, x_0 + \delta]$  is not on the endpoints but somewhere in between e.g. on  $x + \xi$ . But then, the maximum of  $f$  on  $[x_0, x_0 + \xi]$  must be on an endpoint, a contradiction.

Anyhow, this shows us that, by choosing a subsequence such that the maximum and minimum values coincide with the endpoints, we can write

$$\begin{aligned} \left\langle \frac{|f(L)|}{|L|} \right\rangle_{|L| \rightarrow 0} &= \left\langle \frac{\max_{x \in L} f(x) - \min_{x \in L} f(x)}{|L|} \right\rangle_{|L| \rightarrow 0} \\ &= \left\langle \frac{|f(x_0 + h) - f(x_0)|}{h} \right\rangle_{h \rightarrow 0} \\ &= Df(x_0). \end{aligned}$$

The existence of the limit at all is guaranteed by differentiability (how?). Actually, I'm not too sure how to generalise this into  $n$  dimensions, but I think the following is true: Where  $f : \mathbb{R}^n \rightarrow \mathbb{R}^n$ ,

$$\left\langle \frac{|f(V)|}{|V|} \right\rangle_{|V| \rightarrow 0} = \det Df.$$

This can be used to prove the change-of-variable formula:

$$\begin{aligned} \int_{\Phi(V)} f &= \left\langle \sum_{\substack{p_i \in P(\Phi(V)) \\ \text{mesh } P(\Phi(V)) \rightarrow 0}} f(p_i^*) |p_i| \right\rangle \\ &= \left\langle \sum_{\substack{p_i \in P(V) \\ \text{mesh } P(V) \rightarrow 0}} f(\Phi(p_i^*)) |\Phi(p_i)| \right\rangle \end{aligned}$$



$$\begin{aligned}
&= \left\langle \sum_{\substack{p_i \in P(V) \\ \text{mesh } P(V) \rightarrow 0}} f(\Phi(p_i^*)) \det(D\Phi)|p_i| \right\rangle \\
&= \left\langle \sum_{\substack{p_i \in P(V) \\ \text{mesh } P(V) \rightarrow 0}} f(\Phi(p_i^*)) \det(D\Phi)|p_i| \right\rangle \\
&= \int_V f \circ \Phi \det(D\Phi).
\end{aligned}$$

The third equality can be justified using the trick used in the definition of the integral – just take the difference of terms in the sequence and one will see that it can be made arbitrarily small. Also, note that the exact sequence of partitions have to be chosen carefully for the equalities to hold; we want  $\Phi(P(V)) = P(\Phi(V))$ .  $\square$

## 5 Linear Algebra

### 5.1 Linear Operators

### 5.2 Eigen

Diagonal matrices are particularly easy to work with, having the effect of scaling axes. In particular, their powers are simply powers of each element. Abusing the change of basis concept from before, we can simplify certain matrices by showing they are equivalent to diagonal matrices under a certain basis.

For example, asdasd. However, not all matrices are diagonal under some basis; consider the rotation matrix asdasdasd.

In general, for a matrix  $A$  to be diagonalisable, there needs to exist a change of basis matrix  $P = P_{S \rightarrow B} = [\vec{p}_1 \quad \vec{p}_2 \quad \dots \quad \vec{p}_n]$  such that  $P^{-1}AP$  is diagonal. This means that

$$\begin{aligned}
AP &= P \begin{bmatrix} \lambda_1 & 0 & \dots & 0 \\ 0 & \lambda_2 & \dots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \dots & \lambda_n \end{bmatrix} \\
&\Rightarrow A\vec{p}_i = \lambda_i \vec{p}_i,
\end{aligned}$$

that is,  $p_i$  are eigenvectors of  $A$ . Notice that the invertibility requirement on  $P$  implies that each vector in  $P$  be independent, that is, each eigenvector of  $A$  be independent.

### 5.2.1 Spectral Theorem

**Definition 5.2.1.1.** A matrix  $A$  is Hermitian iff  $A = A^H$ , where  $A^H$  denotes the conjugate transpose of  $A$ . An equivalent formulation that generalises to operators is that it satisfies  $\langle A\vec{u}, \vec{v} \rangle = \langle \vec{u}, A\vec{v} \rangle$  – that is, it is equivalent to its adjoint.

Note; hermitian operator is an operator from a finite vector space  $V$  to itself.

**Theorem 5.2.1.1** (Hermitian Spectral Theorem). The eigenvectors of a Hermitian matrix can form an orthonormal basis.

*Proof.* The proof uses induction on the size  $n$ . First we note that a Hermitian operator is guaranteed to have at least one eigenvector over complex vector spaces, by the fundamental theorem of algebra. The key step in induction is noting that Hermiticity guarantees preservation of subspace-invariance under orthogonal complementation.

Choose an eigenvector  $\vec{v}$ . Its span is invariant under  $\mathcal{A}$  (the operator corresponding to  $A$ ), so  $\text{span}(\vec{v})^\perp$  is also. But this means that  $\mathcal{A}$  is a Hermitian operator on  $\text{span}(\vec{v})^\perp$ , so we may choose another eigenvector in  $\text{span}(\vec{v})^\perp$ , and continue this process until we exhaust dimensions. Clearly all eigenvectors chosen are orthogonal.

It remains to prove that Hermiticity guarantees preservation of subspace-invariance under orthogonal complementation. Suppose that  $\langle \vec{v}, \vec{u} \rangle = 0$ , where  $\vec{v}$  is an eigenvector of  $A$ . Then,

$$\begin{aligned}\langle A\vec{u}, A\vec{v} \rangle &= \langle A^2\vec{u}, \vec{v} \rangle \\ &= \lambda^2 \langle \vec{u}, \vec{v} \rangle \\ &= 0.\end{aligned}$$

□

**Corollary 5.2.1.1.** The above result holds also for symmetric maps under real inner product spaces.

*Proof.* The proof above need not be modified, as we are guaranteed the existence of eigenvectors as all eigenvalues of symmetric matrices are real. The proof follows from considering two expression equivalent to  $\vec{v}^H A \vec{v}$ , where  $\vec{v}$  is an eigenvector of  $A$ .

$$\begin{aligned}\vec{v}^H A \vec{v} &= \vec{v}^H \lambda \vec{v} \\ &= \lambda \vec{v}^H \vec{v} \\ &= \lambda \|\vec{v}\|^2.\end{aligned}$$

But also,

$$\begin{aligned}\vec{v}^H A \vec{v} &= \vec{v}^H A^H \vec{v} \\ &= (A\vec{v})^H \vec{v} \\ &= (\lambda \vec{v})^H \vec{v} \\ &= \bar{\lambda} \vec{v}^H \vec{v}.\end{aligned}$$

Thus,  $\bar{\lambda} = \lambda$  and we are done.

□

## A Appendix

### A.1 3.1.3

$$\begin{aligned} S \setminus (A \cap S) &= S \cap (A \cap S)' & (1) \\ &= S \cap (A' \cup S') \\ &= S \cap A' \\ &= S \cap (A \cup B) \cap A' \\ &= S \cap B \cap A' \\ &= S \cap B & (B \subseteq A') \end{aligned}$$

## References

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