# Assignment 3 — Math Fundamentals for Robotics 16-811, Fall 2024

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- 1. Consider the function f(x) = sin x 0.5 over the interval  $\left[-\frac{\pi}{2}, \frac{\pi}{2}\right]$ .
  - a) What is the Taylor series expansion for f(x) around x=0 ? Solution:

$$F(x) = -0.5 + \sum_{k=0}^{\infty} \frac{(-1)^k}{(2k+1)!} x^{2k+1}$$
$$\approx -0.5 + x - \frac{x^3}{6} + \frac{x^5}{120} - \frac{x^7}{5040} + \cdots$$

b) Graph f(x) over the interval  $\left[-\frac{\pi}{2}, \frac{\pi}{2}\right]$ . Solution:

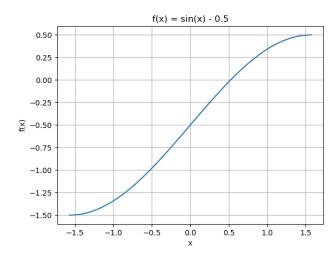


Fig. 1: Graph of f(x) = sin(x) - 0.5

c) Determine the best uniform approximation by a quadratic to the function f(x) on the interval  $[-\frac{\pi}{2}, \frac{\pi}{2}]$ . What are the  $L_{\infty}$  and  $L_2$  errors for this approximation?

# **Solution:**

n=2 so we need n+2=4 points to determine the coefficients of the quadratic. Let them be  $\{x_0,x_1,x_2,x_3\}$ , and let the quadratic function be  $p_2(x)=ax^2+bx+c$ .

 $f^{(n+1)}(x) = f^{(3)}(x) = -\cos(x)$  which doesn't change sign in  $[-\frac{\pi}{2}, \frac{\pi}{2}]$ , so by the theorem taught in class we know

that:

```
\begin{cases} x_0 = -\frac{\pi}{2} \\ x_3 = \frac{\pi}{2} \\ e(x_0) = -e(x_1) = e(x_2) = -e(x_3) = E \end{cases}
+
\begin{cases} e(x_0) = f(x_0) - p_2(x_0) = f(-\frac{\pi}{2}) - p_2(-\frac{pi}{2}) & = -1.5 - \frac{\pi^2}{4}a + \frac{\pi}{2}b - c \\ e(x_1) = f(x_1) - p_2(x_1) & = \sin(x_1) - 0.5 - ax_1^2 - bx_1 - c \\ e(x_2) = f(x_2) - p_2(x_2) & = \sin(x_2) - 0.5 - ax_2^2 - bx_2 - c \\ e(x_3) = f(x_3) - p_2(x_3) = f(\frac{\pi}{2}) - p_2(\frac{\pi}{2}) & = 0.5 - \frac{\pi^2}{4}a - \frac{\pi}{2}b - c \end{cases}
+
\begin{cases} e'(x_1) = f'(x_1) - p'_2(x_1) = \cos(x_1) - 2ax_1 - b = 0 \\ e'(x_2) = f'(x_2) - p'_2(x_2) = \cos(x_2) - 2ax_2 - b = 0 \end{cases}
\downarrow \downarrow
\begin{cases} c = -\frac{\pi^2}{4}a - 1 \\ b = \frac{2E - 2}{\pi} \end{cases}
```

Now we see that the solution is not trivial, so we resort to numerical method Remez Exchange Algorithm to solve for the coefficients.

```
def remez_exchange(
f, n: int, interval: list, max_num_iteration: int = 100, tolerance: float = 1e-5
):
      """Remez Exchange Algorithm
      Args:
            f (_type_): univariate function
            n (int): polynomial degree
            interval (list): interval for uniform approximate
            max_num_iteration (int): max number of iteration
            tolerance (float): tolerance of point change between updates
      11 11 11
      assert len(interval) == 2, "Interval must be a list of 2 elements"
      \# we need n + 2 points to construct a polynomial of degree n, uniform
          initialization
      xi = np.linspace(interval[0], interval[1], n + 2)
      # define x
      x = sp.symbols("x", real=True)
      # numerical function f
      f_numerical = sp.lambdify(x, f, "numpy")
      \# determine if f^{(n + 1)} does not change sign in the interval
      f_n_plus_1 = sp.diff(f, x, n + 1)
      roots = find_all_roots(sp.lambdify(x, f_n_plus_1, "numpy"), interval)
      if len(roots) == 0:
            print("f^(n + 1) does not change sign in the interval")
            xi[0] = interval[0]
            xi[-1] = interval[1]
      # # some turbulence in the initial points
      \# xi[1] = -1.0
      for iter in range(max_num_iteration):
            print(f"\nIteration {iter}\n")
```

```
# construct linear system
      A = np.ones((n + 2, n + 2))
      for degree in range (n + 1):
           A[:, degree] = xi**degree
      A[:, -1] = (-1) ** np.arange(n + 2)
      b = f_numerical(xi).reshape(-1, 1)
      print(f"xi: {xi}")
     print(f"A: {A}")
      print(f"b: {b}")
      # solve linear system
      solution = np.linalg.solve(A, b).flatten()
      print(f"solution: {solution}")
      poly_coeff = solution[:-1]
      error = np.abs(solution[-1])
      # construct polynomial function with sympy
      poly = 0
      for degree in range (n + 1):
            poly += poly_coeff[degree] * x**degree
      print("Polynomial function:")
      sp.pprint(poly)
      # find the maximum error
      error_func = f - poly
      error_func_numerical = sp.lambdify(x, error_func, "numpy")
      max_error_abs, max_point = find_max(error_func, interval)
      print(f"max error: {max_error_abs}")
      print(f"max point: {max_point}")
      # filter out the points with the same sign as the maximum error
      xi_new = xi.copy()
      xi_same_sign = xi_new[
           np.where(
            np.sign(error_func_numerical(xi_new))
            == np.sign(error_func_numerical(max_point))
      # find the closest point to the maximum error point and replace it with the
         maximum error point
      closest_point = xi_same_sign[np.argmin(np.abs(xi_same_sign - max_point))]
      xi_new[np.where(xi_new == closest_point)] = max_point
      print(f"xi_new: {xi_new}")
      # if the error changing is less than tolerance, break the loop
      if np.abs(np.abs(error) - max_error_abs) < tolerance:</pre>
            print(f"Converged after {iter} iterations")
            break
      xi = np.sort(xi_new)
if iter == max_num_iteration - 1:
      print(f"Did not converge after {max_num_iteration} iterations")
return poly, poly_coeff, max_error_abs
```

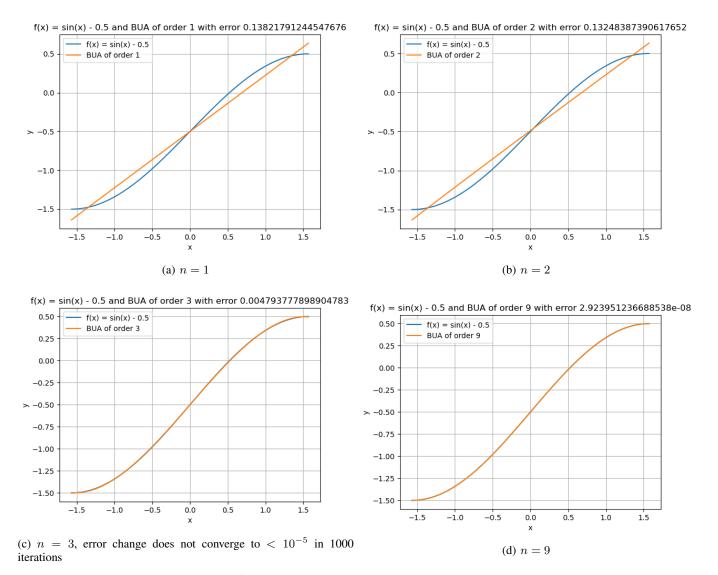


Fig. 2: Best Uniform Approximation using the Remez Exchange Algorithm

Well, I guess odd function approximation can get good results. For n=1:

$$\begin{cases} p_1(x) & \approx 0.72460996019333x - 0.5 \\ L_{\infty} & \approx 0.13821791244547676 \\ L_2 & \approx 0.170396813894351 \end{cases}$$

For n=2:

$$\begin{cases} p_2(x) &\approx -0.500004924940155 + 0.724609736054229x + 1.9960030637173 \times 10^{-6}x^2 \approx 0.724609736054229x - 0.5\\ L_\infty &\approx 0.13822185392370312\\ L_2 &\approx 0.170396982005513 \end{cases}$$

For n = 3:

Its error change does not converge to  $< 10^{-5}$  in 1000 iterations

```
\begin{cases} p_3(x) & \approx -0.5000619896739 + 0.985440080596674x + 0.000139910675817895x^2 + 0.142443200483762x^3 \\ L_\infty & \approx 0.004793777898904783 \\ L_2 & \approx 0.00568482890288639 \end{cases}
```

d) Determine the best least-squares approximation by a quadratic to the function f(x) over the interval  $[-\frac{\pi}{2}, \frac{\pi}{2}]$ . What are the  $L_{\infty}$  and  $L_2$  errors for this approximation?

### **Solution:**

Using Legendre polynomials as basis functions, we can solve for the coefficients of the quadratic. Legendre polynomials modified for interval  $\left[-\frac{\pi}{2},\frac{\pi}{2}\right]$  has the following dot product definition:

$$< f,g> = \int_{-\frac{\pi}{2}}^{\frac{\pi}{2}} f(x)g(x)dx$$

Then with the first 2 Legendre polynomials:

$$\begin{cases} P_0(x) = 1 \\ P_1(x) = x \end{cases}$$

We can derive more Legendre polynomials using the following recurrence relation:

$$P_{n+1}(x) = \left[ x - \frac{\langle xP_n(x), P_n(x) \rangle}{\langle P_n(x), P_n(x) \rangle} \right] P_n(x) - \frac{\langle P_n(x), P_n(x) \rangle}{\langle P_{n-1}(x), P_{n-1}(x) \rangle} P_{n-1}(x), n = 1, 2, \dots$$

More Legendre polynomials:

$$\begin{cases} P_2(x) &= x^2 - \frac{\pi^2}{12} \\ P_3(x) &= x^3 - \frac{3\pi^2}{20}x \\ P_4(x) &= x^4 - \frac{3\pi^2}{14}x^2 + \frac{3\pi^4}{560} \\ P_5(x) &= x^5 - \frac{5\pi^2}{18}x^3 + \frac{5\pi^4}{336}x \\ P_6(x) &= x^6 - \frac{15\pi^2}{44}x^4 + \frac{5\pi^4}{176}x^2 - \frac{5\pi^6}{14784} \end{cases}$$

Now after projecting f(x) onto the first 3 Legendre polynomials for quadratic least square approximation, we can solve for the coefficients of the quadratic.

$$f(x) \approx -0.5P_0(x) + \frac{24}{\pi^3}P_1(x) + 0P_2(x)$$
$$= \frac{24}{\pi^3}x - 0.5$$
$$L_\infty \approx 0.2158542037080533$$
$$L_2 \approx 0.15074041926875908$$

```
def least_square_approximation(f, n: int, interval: list):
    """Least square approximation

Args:
    f (_type_): univariate function
    n (int): polynomial degree
    interval (list): interval for uniform approximate

Returns:
    the polynomial
    """

x = sp.symbols("x", real=True)

# construct the legendre polynomials
legendre_polynomials = [legendre_polynomial(i, interval) for i in range(n + 1)]

print(f"Legendre polynomials: {legendre_polynomials}")

projection_coefficients = [
    legendre_dot(f, P, interval) / legendre_dot(P, P, interval)
    for P in legendre_polynomials
]

print(f"Projection coefficients: {projection_coefficients}")
```

```
poly = 0
for coefficient, P in zip(projection_coefficients, legendre_polynomials):
    poly += coefficient * P

return poly
```

 $f(x) = \sin(x) - 0.5$  and Legendre of order 1 with error 0.1507404192687590{  $f(x) = \sin(x) - 0.5$  and Legendre of order 2 with error 0.1507404192687590{  $f(x) = \sin(x) - 0.5$  and Legendre of order 2 with error 0.1507404192687590{  $f(x) = \sin(x) - 0.5$  and Legendre of order 2 with error 0.1507404192687590{  $f(x) = \sin(x) - 0.5$  and Legendre of order 2 with error 0.1507404192687590{  $f(x) = \sin(x) - 0.5$  and Legendre of order 2 with error 0.1507404192687590{  $f(x) = \sin(x) - 0.5$  and Legendre of order 2 with error 0.1507404192687590{  $f(x) = \sin(x) - 0.5$  $f(x) = \sin(x) - 0.5$  $f(x) = \sin(x) - 0.5$ Legendre of order 1 Legendre of order 2 0.5 0.5 0.0 0.0 -0.5> -0.5 -1.0 -1.0 -1.5 -0.50.5 1.5 -0.5 0.5 1.0 1.5 -1.5-1.00.0 1.0 -1.5-1.00.0 (a) n = 1(b) n = 2 $f(x) = \sin(x) - 0.5$  and Legendre of order 3 with error 0.00491605234430628  $f(x) = \sin(x) - 0.5$  and Legendre of order 9 with error 15.761180424768629 0.50  $f(x) = \sin(x) - 0.5$ 0.50  $f(x) = \sin(x) - 0.5$ Legendre of order 3 Legendre of order 9 0.25 0.25 0.00 0.00 -0.25 -0.25 > -0.50 > -0.50 -0.75-0.75-1.00-1.00-1.25-1.25 -1.50-1.50-1.5 1.5

Fig. 3: Least Square Approximation using Legendre Polynomials

(c) n = 3

Terminology: Suppose an approximation has error function e(x), with x in interval [a,b]. The  $L_{\infty}$  error is  $||e(x)||_{\infty} = \max_{a \leq x \leq b} |e(x)|$  and the  $L_2$  error is  $||e(x)||_2 = \sqrt{\int_a^b |e(x)|^2 dx}$ . For your code submission, submit any code you used.

(d) n = 9, complex error occurs here so they look close

In your pdf, please show all your hand derivations and code results, and replicate any code you would like the TAs to see.

2. Suppose very accurate values of some function f(x) are given at the points  $0 = x_0, x_1, \dots, x_{100} = 1$ , with the  $x_i$  uniformly distributed over the interval [0, 1].

```
(So x_i = \frac{i}{100}, i = 0, \dots, 100.)
```

The values  $\{f(x_i)\}$  are given in the file 'problem2.txt' in sequential order

(so, for example,  $f(0.27) = f(x_{27}) = -0.964603914513021$ ).

Use the method of normal equations discussed in class, to find a description of f(x) as the sum of a very few polynomials and cosines and sines. (You may be able to guess the answer since the function is fairly simple, but please also use the method of normal equations.)

For code, submit any code you used. In your pdf, replicate any code you would like the TAs to see, show your results and explain how you obtained them. If your search for a solution first considered some incorrect solutions, mention those and say how they informed your search for the correct solution.

[Hint: Try to express f(x) as a linear combination of the functions,  $1, x, x^2, \cdots, cos(\pi x), sin(\pi x), cos(2\pi x), sin(2\pi x), \cdots$ . Use the method of normal equations. That method will not be enough by itself, since the underlying functions are redundant (more than a basis). However, that method is useful as a subroutine in a search. Try to find a very simple description of the function f(x) by determining which coefficients in your sum may be set to zero. There may be multiple candidate answers; find one with the fewest nonzero coefficients. Graphing the function may be helpful in your search. You should need at most 3 nonzero coefficients when writing f(x) as a sum of the functions  $1, x, x^2, \cdots, cos(\pi x), sin(\pi x), cos(2\pi x), sin(2\pi x), \cdots$ .]

#### **Solution:**

```
def normal_equation_method(xi: np.array, yi: np.array, n_poly: int, n_trig: int, tol: float
    """Use the method of normal equations to fit a polynomial and trigonometric function to
        the data.
   Aras:
       xi (np.array): xi data points
       yi (np.array): fi data points
       n_poly (int): number of polynomial terms
       n_triq (int): number of trigonometric terms
       tol (float, optional): tolerance for filtering out small coefficients. Defaults to
           1e-5.
    11 11 11
    xi = xi.reshape(-1, 1).copy()
    yi = yi.reshape(-1, 1).copy()
   assert xi.shape == yi.shape, "xi and yi must have the same shape"
   # Construct the design matrix
   A = np.zeros((len(xi), n_poly + 2 * n_trig))
    for i in range(n_poly):
       A[:, i] = (xi**i)[:,
    for i in range(n_trig):
       A[:, n_poly + i] = np.sin((i + 1) * np.pi * xi)[:, 0]
       A[:, n_poly + n_trig + i] = np.cos((i + 1) * np.pi * xi)[:, 0]
    # Use least squares to solve for the coefficients
    coeff, _, _, _ = np.linalg.lstsq(A, yi, rcond=None)
    # Filter out coefficients that are close to zero
    coeff[np.abs(coeff) < tol] = 0
    x = sp.symbols("x", real=True)
   p = 0
    for i in range(n_poly):
       p += coeff[i] * x**i
   for i in range(n_trig):
       p \leftarrow coeff[n\_poly + i] * sp.sin((i + 1) * sp.pi * x)
       p \leftarrow coeff[n_poly + n_trig + i] * sp.cos((i + 1) * sp.pi * x)
    return sp.simplify(p), coeff
```

# My initial guess:

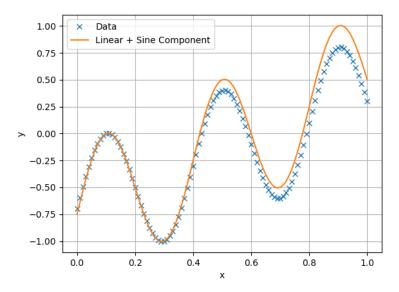


Fig. 4:  $f(x) = 1.25x - 0.75 + 0.625\sin(5\pi x)$ 

# The Method of Normal Equations:

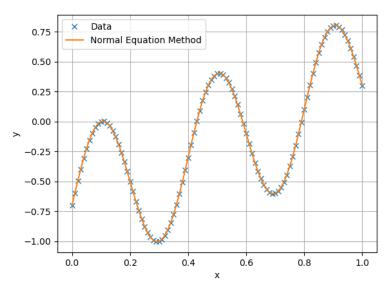


Fig. 5:  $f(x) \approx x - 0.7 + 0.6 \sin(5\pi x)$ 

Accurate result from the method of normal equations:

 $f(x) \approx 0.999999980668295x + 0.600000000000149\sin(5\pi x) - 0.699999999335618$ 

3. The Chebyshev polynomials of the first kind,  $T_n(x)$ , are defined indirectly on [-1, 1] by:

$$T_n(cos\theta) = cos(n\theta)$$
, for  $n \ge 0$ 

Expanding cosine, one finds the recurrence relation  $T_{n+1}(x) = 2xT_n(x) - T_{n-1}(x)$ , for n > 0.

a) Derive  $T_4$  and  $T_5$ . (As always, please show your handwritten work.) **Solution:** 

b) Without actually computing or working out any integrals, prove that  $T_4(x)$  and  $T_5(x)$  are orthogonal polynomials relative to the inner product

$$\langle g, h \rangle = \int_{-1}^{1} (1 - x^2)^{-\frac{1}{2}} g(x)h(x)dx.$$

[Hint: Look carefully at the inner product and use a property of the polynomials.] **Solution:** 

 $x = \cos(\theta)$ , then  $dx = -\sin(\theta)d\theta$ .

$$\langle T_4, T_5 \rangle = \int_{-1}^{1} (1 - x^2)^{-\frac{1}{2}} T_4(x) T_5(x) dx$$

$$= \int_{-\pi}^{0} (1 - \cos^2(\theta))^{-\frac{1}{2}} T_4(\cos(\theta)) T_5(\cos(\theta)) (-\sin(\theta) d\theta)$$

$$= \int_{-\pi}^{0} \frac{\cos(4\theta) \cos(5\theta)}{-\sin(\theta)} (-\sin(\theta) d\theta)$$

$$= \int_{-\pi}^{0} \cos(4\theta) \cos(5\theta) d\theta$$

$$= \int_{-\pi}^{0} \frac{1}{2} \left[ \cos(9\theta) + \cos(\theta) \right] d\theta$$

$$= \frac{1}{2} \left[ \frac{\sin(9\theta)}{9} + \sin(\theta) \right]_{-\pi}^{0}$$

c) Recall that when we have an inner product on a vector space, we may define the *length of vector* v by  $||v|| = \sqrt{\langle v, v \rangle}$ . Here, we may view functions as vectors with an inner product defined by an integral as above. Then the length of  $T_n(x)$  is the number  $\sqrt{\langle T_n, T_n \rangle}$ .

It turns out that all  $T_n(x)$ , with n > 0, have the same length.

Prove this fact by hand-computing the length of  $T_n(x)$ , that is, by working out the relevant integral (leave n symbolic, assume n > 0).

(Hint: You will likely find it useful to make the substitution  $x = cos\theta$  in the integral for  $< T_n, T_n >$ . Don't forget to change the interval of integration as well.)

**Solution:** 

 $x = \cos(\theta)$ , then  $dx = -\sin(\theta)d\theta$ .

$$\langle T_n, T_n \rangle = \int_{-1}^{1} (1 - x^2)^{-\frac{1}{2}} T_n(x) T_n(x) dx$$

$$= \int_{-\pi}^{0} (1 - \cos^2(\theta))^{-\frac{1}{2}} \cos^2(n\theta) (-\sin(\theta) d\theta)$$

$$= \int_{-\pi}^{0} \cos^2(n\theta) d\theta$$

$$= \int_{-\pi}^{0} \frac{1}{2} \left[ \cos(2n\theta) + 1 \right] d\theta$$

$$= \frac{1}{2} \left[ \frac{\sin(2n\theta)}{2n} + \theta \right] \Big|_{-\pi}^{0}$$

$$= \frac{\pi}{2}$$

d) Finally, show that  $\langle T_i, T_j \rangle = 0$  for all i and j such that  $i \geq 0, j \geq 0$ , and  $i \neq j$ . (There are different ways to prove this, e.g., by working out an integral explicitly or by combining known facts from above and lecture.)

### **Solution:**

 $x = \cos(\theta)$ , then  $dx = -\sin(\theta)d\theta$ .

$$\langle T_{i}, T_{j} \rangle = \int_{-1}^{1} (1 - x^{2})^{-\frac{1}{2}} T_{i}(x) T_{j}(x) dx$$

$$= \int_{-\pi}^{0} (1 - \cos^{2}(\theta))^{-\frac{1}{2}} T_{i}(\cos(\theta)) T_{j}(\cos(\theta)) (-\sin(\theta) d\theta)$$

$$= \int_{-\pi}^{0} \frac{\cos(i\theta) \cos(j\theta)}{-\sin(\theta)} (-\sin(\theta) d\theta)$$

$$= \int_{-\pi}^{0} \cos(i\theta) \cos(j\theta) d\theta$$

$$= \int_{-\pi}^{0} \frac{1}{2} \left[ \cos((i+j)\theta) + \cos((i-j)\theta) \right] d\theta$$

$$= \frac{1}{2} \left[ \frac{\sin((i+j)\theta)}{i+j} + \frac{\sin((i-j)\theta)}{i-j} \right]_{-\pi}^{0}$$

$$= 0$$

No code is expected or needed for any part of this problem. In your pdf, please show all your derivations, proofs, and handwritten work.

- 4. After weeks of work you have finally completed construction of a gecko robot. It is a quadruped robot with suctioning feet that allow it to walk on walls. It is also equipped with a Kinect-like sensor, providing a 3D point cloud observation of the world. You want to use these point clouds to reason about the environment and aid in navigation.
  - a) You boot up the robot and place it on a table, taking an initial observation. The observation is saved in the provided clear\_table.txt, and lists (x, y, z) locations in the following format:

$$egin{array}{cccc} x_1 & y_1 & z_1 \ & dots \ x_n & y_n & z_n \end{array}$$

Points are in units of meters and the positive x-direction is right, positive y-direction is down, and positive z-direction is forward. Find the least-squares approximation plane that fits the data. Visualize your fitted plane along with the data. What is the average distance of a point in the data set to the fitted plane?

**Comment:** The phrase "least-squares" is ambiguous. Below are descriptions of two possibilities that might occur to you. **Please use Linear Regression.** That approach is similar to our discussion of the normal equations in lecture.

**Orthogonal-Distance Regression:** In this approach, one computes the SVD decomposition of the  $n \times 3$  matrix whose rows are the data points translated so their centroid is at the origin. It turns out that the third column of V is normal to a plane that minimizes the sum of squared orthogonal distances between the translated points and the plane. (This is a nice result to know.)

**Linear Regression:** This regression is based on the idea that, for perfectly planar data, all the points would satisfy a plane equation of the form ax + by + cz + d = 0. So one has a natural error  $\sum_i (ax_i + by_i + cz_i + d)^2$ , with i indexing the data points. One chooses the coefficients  $\{a, b, c, d\}$  so as to minimize this error. In order to avoid degeneracies, one requires that not all of  $\{a, b, c, d\}$  be 0.

# Please use this approach.

(Additional comments: (i) One convenient approach is to set one of the coefficients  $\{a,b,c,d\}$  to be 1 or -1 while letting the others vary in order to compute the best plane. (ii) Observe that  $|ax_i + by_i + cz_i + d|$  is related to but not necessarily exactly the distance of the ith data point from the plane described by  $\{a,b,c,d\}$ .)

#### **Solution:**

We can set d=1 first, then normalize the plane equation later.

$$ax + by + cz + 1 = 0$$

$$E(a, b, c) = \sum_{i=1}^{n} (ax_i + by_i + cz_i + 1)^2$$
$$= (Ap + 1)^T (Ap + 1)$$

Where:

$$A = \begin{bmatrix} x_1 & y_1 & z_1 \\ \vdots & \vdots & \vdots \\ x_n & y_n & z_n \end{bmatrix}$$

$$p = \begin{bmatrix} a \\ b \\ c \end{bmatrix}$$

$$\mathbf{1} = \begin{bmatrix} 1 \\ \vdots \\ 1 \end{bmatrix}$$

Then:

$$\frac{\partial E}{\partial p} = 2A^{T}(Ap + \mathbf{1}) = 0$$

$$\downarrow \downarrow$$

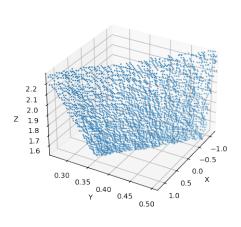
$$A^{T}Ap = -A^{T}\mathbf{1}$$

Solve for p we get  $p^* = [a^*, b^*, c^*]^T$ , then the normalized plane equation is:

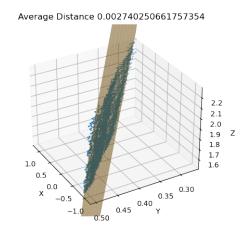
$$\frac{a^*x + b^*y + c^*z + 1}{\|p\|} = 0$$

### Plane Equation:

-0.0951113342629671x - 0.994246831912656y - 0.0492653156527438z + 0.493702932302467 = 0 Average distance = 0.002740250661757354







(b) Fitted Plane, note that Z axis is auto-scaled

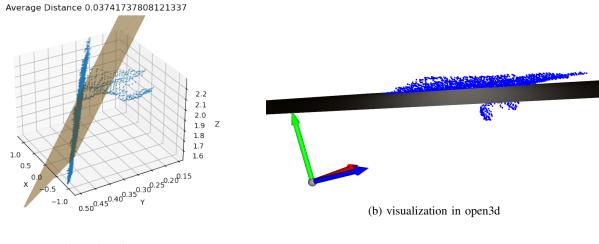
Fig. 6: Fitted Plane on clear\_table.txt

b) Interested in your gecko robot, your cat jumps up on the table. You take a second observation, saved as the provided cluttered table.txt. Using the same method as above, find the least-squares fit to the new data. How does it look? Why? **Solution:** 

Plane equation:

-0.0975611845981769x - 0.973054533289032y - 0.208917903745619z + 0.774119219789838 = 0

Average distance is 0.03741737808121337



(a) cluttered\_table.txt

Fig. 7: Fitted Plane on cluttered\_table.txt

Well, clearly not all the points are close to the ideal plane.

c) Can you suggest a way to still find a fit to the plane of the table regardless of clutter? Verify your idea by writing a program that can successfully find the dominant plane in a list of points regardless of outliers. [Hint: You may assume that the number of points in the plane is much larger than the number of points not in the plane.] Visualize cluttered\_table.txt with your new plane.

#### **Solution:**

#### **RANSAC**

```
def RANSAC_fit_plane(
     points: np.array,
     max_iter: int = 10000,
     sample_ratio: float = 0.01,
     distance_inlier_threshold: float = 0.01,
) -> np.array:
      """Fit a plane to a set of noisy 3D points using RANSAC.
     Aras:
            points (np.array): points in shape (N, 3)
            max_iter (int, optional): maximum iteration to perform random sampling.
               Defaults to 10000.
            sample_ratio (float, optional): the amount of points to sample, sample_ratio
               * N. Defaults to 0.01. Will not be less than 10.
            distance_inlier_threshold (float, optional): inlier points maximum average
               distance to their fitted plane. Defaults to 0.01.
     Raises:
            ValueError: No inlier points found.
     Returns:
           np.array: inlier points
     NUM_POINTS = points.shape[0]
     SAMPLE_SIZE = max(int(points.shape[0] * sample_ratio), 10)
     inlier_index = np.array([], dtype=int)
      for _ in range(max_iter):
            # randomly sample SAMPLE_SIZE points index
            random_index = np.random.permutation(np.arange(NUM_POINTS, dtype=int))[
                  :SAMPLE SIZE
            coeff = fit_plane(points[random_index])
```

```
distance = average_point_to_plane_distance(points[random_index], coeff)
    if distance <= distance_inlier_threshold:
        inlier_index = np.append(inlier_index, random_index)
        inlier_index = np.sort(inlier_index)
        inlier_index = np.unique(inlier_index)

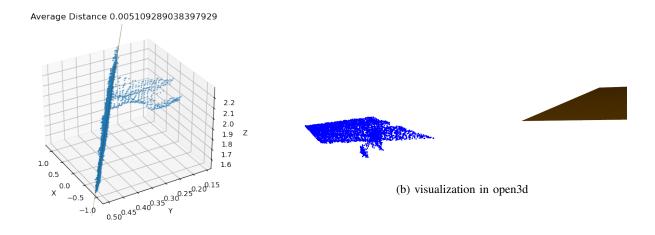
if inlier_index.size == 0:
    raise ValueError("No inlier found.")

print(f"Inliers ratio: {inlier_index.size / NUM_POINTS}")
    return points[inlier_index]</pre>
```

# Plane Equation:

```
-0.10068718707874x - 0.992880691307245y - 0.063639792484089z + 0.518921115704315 = 0\\
```

Average distance to all the points is 0.015754165478591993 Average distance to table points is 0.005143845669654812



(a) cluttered\_table.txt RANSAC

Fig. 8: RANSAC fitted Plane on cluttered\_table.txt

d) Encouraged by your results when testing on a table, you move your geckobot into the hallway and take an observation saved as the provided clean\_hallway.txt. Describe an extension to your solution to part (c) that finds the four dominant planes shown in the scene, then implement it and visualize the data and the four planes.

You may assume that there are roughly the same amount of points in each plane.

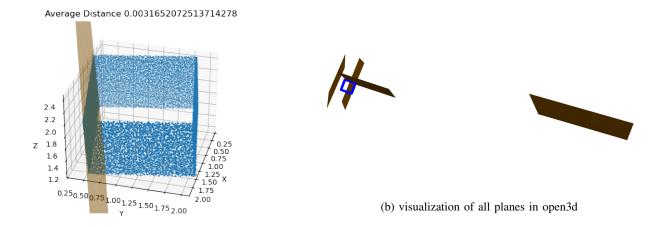
Solution: Extend the inlier criteria of RANSAC:

```
def RANSAC_fit_plane_extended(
     points: np.array,
     num_planes: int,
     max_iter: int = 10000,
     sample_ratio: float = 0.001,
     distance_inlier_threshold: float = 0.01,
      """extended RANSAC to fit multiple planes, inlier points must satisfy:
     1. Itself is a plane with small average distance
     2. Its fitted plane must be the same plane with existing inliers
     3. The fitted plane's close to enough points (1/num_planes * 0.9)
     Args:
           points (np.array): points in shape (N, 3)
           num_planes (int): number of planes to fit, assuming each plane has roughly
               the same number of inliers
           max_iter (int, optional): maximum iteration to perform random sampling.
               Defaults to 10000.
```

```
sample_ratio (float, optional): the amount of points to sample, sample_ratio
         * N. Defaults to 0.0001. Will not be less than 5.
      distance_inlier_threshold (float, optional): inlier points maximum average
         distance to their fitted plane. Defaults to 0.01.
Raises:
     ValueError: No inlier points found.
Returns:
     np.array: inlier points, might not be most points in its plane
NUM_POINTS = points.shape[0]
SAMPLE_SIZE = max(int(points.shape[0] * sample_ratio), 5)
print(f"Number of points: {NUM_POINTS}")
print(f"Sample size: {SAMPLE_SIZE}")
inlier_index = np.array([], dtype=int)
points_homo = np.hstack((points, np.ones((points.shape[0], 1))))
for iter in range(max_iter):
      # randomly sample SAMPLE_SIZE points index
      random_index = np.random.permutation(np.arange(NUM_POINTS, dtype=int))[
            :SAMPLE_SIZE
      extended_index = np.unique(np.append(inlier_index, random_index))
      coeff = fit_plane(points[random_index])
      distance = average_point_to_plane_distance(points[random_index], coeff)
      extended_distance = average_point_to_plane_distance(
            points[extended_index], coeff
      all_distances = np.abs(points_homo @ coeff)
      num_inliers = np.where(all_distances <= distance_inlier_threshold)[0].size</pre>
      inliers_ratio = num_inliers / NUM_POINTS
      # print(f"inliers_ratio = {inliers_ratio}")
      # Must satisfy: itself is a plane, it's the same plane with inliers, and it
         has enough inliers
      if (
            distance <= distance_inlier_threshold</pre>
            and extended_distance <= distance_inlier_threshold</pre>
            and inliers_ratio > 1 / num_planes * 0.9
      ):
            inlier_index = np.append(inlier_index, random_index)
            inlier_index = np.sort(inlier_index)
            inlier_index = np.unique(inlier_index)
if inlier_index.size == 0:
      raise ValueError("No inlier found.")
print(f"Inliers ratio: {inlier_index.size / NUM_POINTS}")
return points[inlier_index]
```

#### Plane Equations:

```
\begin{cases} 0.170391902271525x - 0.984679478280239y - 0.0370529984488184z + 1.80802982578352 & = 0, ||e|| \approx 0.003144184418150 \\ -0.96366746614888x - 0.175264115206761y + 0.201562656776964z + 0.218727644397049 & = 0, ||e|| \approx 0.004516231182443 \\ -0.96403454687357x - 0.170863621367788y + 0.203575576450997z + 1.8162162422856 & = 0, ||e|| \approx 0.007681134237631 \\ 0.16936975663811x - 0.984804871586362y - 0.0383829445993377z + 0.211152350861021 & = 0, ||e|| \approx 0.003350451774268 \end{cases}
```



(a) clean\_hallway.txt extended RANSAC  $1^{st}$  plane

Fig. 9: extended RANSAC fitted plane on clean hallway.txt

More results available in the figs/ folder.

e) You decide it is time to test your gecko robot's suction feet and move it to a different hallway. The feet are strong enough to ignore the force of gravity, allowing the robot to walk on the floor, walls, or ceiling. However, the locomotion of the legs works best on smooth surfaces with few obstacles. Using your solution from part (d), describe how you can mathematically characterize the smoothness of each surface. Load the provided scan cluttered\_hallway.txt, find and plot the four wall planes, describe which surface is safest for your robot to traverse, and provide the smoothness scores from your mathematical characterization.

Note that you may no longer assume that there are roughly the same amount of points in each plane.

### **Solution:**

I characterize the smoothness of each surface by the variance of distance between all the points to the fitted plane, lower variance indicates higher smoothness.

```
def voxel_downsample_point_cloud(points, voxel_size):
    """
    Downsamples the input point cloud using a voxel grid filter.

Args:
    points (np.array): Input points in shape (N, 3).
    voxel_size (float): The size of the voxel grid.

Returns:
    np.array: Downsampled points in shape (M, 3).
    """
    # Create Open3D point cloud
    pcd = o3d.geometry.PointCloud()
    pcd.points = o3d.utility.Vector3dVector(points)

# Downsample
    downsampled_pcd = pcd.voxel_down_sample(voxel_size)

# Extract NumPy array
    downsampled_points = np.asarray(downsampled_pcd.points)

return downsampled_points
```

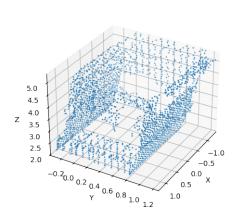
RANSAC is modified to consider inlier points only if their fitted plane is close to enough number of points (threshold is adaptive).

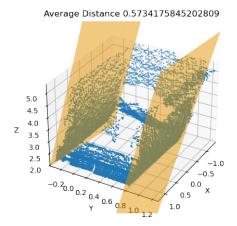
```
def RANSAC_fit_multiple_planes(
    points: np.array,
    num_planes: int,
    max_iter: int = 10000,
```

```
sample_size: int = 5,
      distance_threshold: float = 0.01,
      inlier_ratio: float = 0.9,
):
      """Fit multiple planes to a set of 3D points using RANSAC.
      Args:
            points (np.array): Points in shape (N, 3).
            num_planes (int): Number of planes to fit.
            max iter (int, optional): Maximum iterations for RANSAC. Defaults to 10000.
            sample_size (int, optional): number of points to sample for every RANSAC
                iteration. Defaults to 5.
            distance_threshold (float, optional): Distance threshold for inliers.
                Defaults to 0.01.
            min_inliers (int, optional): Minimum number of inliers to accept a plane.
                Adjust as needed.
      Returns:
      points = points.copy().reshape(-1, 3)
      remaining_points = points.copy().reshape(-1, 3)
      planes = []
      for i in range(num_planes)[::-1]:
            print(f"Fitting plane {i + 1}...")
            min_inliers = int(remaining_points.shape[0] / (i + 1) * inlier_ratio)
            inlier_indices = np.array([], dtype=int)
            # Can no longer assume the same number of inliers for each plane
            for _ in range(max_iter):
                  # Randomly sample points
                  sample_indices = np.random.permutation(
                  np.arange(remaining_points.shape[0], dtype=int)
                  )[:sample_size]
                  extended_index = np.unique(np.append(inlier_indices, sample_indices))
                  sample_points = remaining_points[sample_indices]
                  # Fit a plane to the sample points
                  plane_coefficients = fit_plane(sample_points)
                  fit_distance = average_point_to_plane_distance(
                  sample_points, plane_coefficients
                  inlier_distance = average_point_to_plane_distance(
                  remaining_points[extended_index], plane_coefficients
                  all_distances = np.abs(
                  remaining_points @ plane_coefficients[:3] + plane_coefficients[3]
                  num_inliers = np.where(all_distances <= distance_threshold)[0].size</pre>
                  if (
                  fit_distance <= distance_threshold</pre>
                  and inlier_distance <= distance_threshold</pre>
                  and num_inliers >= min_inliers
                  ):
                  inlier_indices = np.append(inlier_indices, sample_indices)
                  inlier_indices = np.sort(inlier_indices)
                  inlier_indices = np.unique(inlier_indices)
            if inlier indices.size > 0:
                  print(f"Found {inlier_indices.size} inliers.")
                  coeff = fit_plane(remaining_points[inlier_indices])
```

```
# find all points that are close to this plane in points
            distance_all_points = np.abs(points @ coeff[:3] + coeff[3])
            point_close_to_this_plane = points[
            np.where(distance_all_points <= distance_threshold)</pre>
            distance_close_to_this_plane = distance_all_points[
            np.where(distance_all_points <= distance_threshold)</pre>
            print(f"Points close to this plane: {point_close_to_this_plane.shape
                [0] } ")
            planes.append(
                  "coeff": fit_plane(point_close_to_this_plane),
                  "mean_distance": distance_close_to_this_plane.mean(),
                  "var_distance": distance_close_to_this_plane.var(),
                  "var_distance_all": distance_all_points.var(),
                  "points": point_close_to_this_plane,
                  "num_points": point_close_to_this_plane.shape[0],
            \# remove all points close to this plane from remaining points
            distance_remaining_points = np.abs(remaining_points @ coeff[:3] + coeff
                [31)
            remaining_points = remaining_points[
            np.where(distance_remaining_points > distance_threshold)
            print(f"Remaining points: {remaining_points.shape[0]}")
# Sort planes by number of points
planes = sorted(planes, key=lambda x: x["num_points"], reverse=True)
return planes
```

Some results:





(a) cluttered\_hallway.txt downsampled

(b) cluttered\_hallway.txt RANSAC visualized with original point cloud

Fig. 10: RANSAC fitted plane on cluttered\_hallway.txt

The algorithm turns to be very fragile to parameter changes, e.g., sample size, distance threshold. There could be better metric, e.g., project close points to the fitted plane and see if there are holes, to prevent fitting a densely scanned ring instead of a plane.

Anyway, here are the 2 best plane in terms of smoothness and number of points:

Parts (c), (d), and (e) intentionally leave room for some creativity and design. There may be several good approaches. Please submit code for all parts (a), (b), (c), (d), (e) of this problem. In your pdf, explain what you did, how to run your code, and what results you obtained. Describe any design decisions you made. Include as well in your pdf any images you used to visualize data, and explain their meanings. Finally, replicate in your pdf any code you would like the TAs to see.