Fundamentals of Artificial Intelligence Programme (2024/25 Q1)

Multiobjective optimization for decision support

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In this assignment, we use two python libraries, one for exploratory modeling analysis <u>The EMA workbench</u> and one for multiobjective optimization <u>Project Platypus</u>, you can install both with pip as follows:

```
1 !pip install ema_workbench platypus-opt
      Requirement already satisfied: ema_workbench in /usr/local/lib/python3.10/dist-packages (2.5.2)
       Requirement already satisfied: platypus-opt in /usr/local/lib/python3.10/dist-packages (1.3.1)
      Requirement already satisfied: numpy in /usr/local/lib/python3.10/dist-packages (from ema_workbench) (1.26.4)
      Requirement already satisfied: pandas in /usr/local/lib/python3.10/dist-packages (from ema_workbench) (2.1.4)
      Requirement already satisfied: scikit-learn in /usr/local/lib/python3.10/dist-packages (from ema workbench) (1.3.2)
      Requirement already satisfied: salib>=1.4.6 in /usr/local/lib/python3.10/dist-packages (from ema_workbench) (1.5.1)
      Requirement already satisfied: matplotlib in /usr/local/lib/python3.10/dist-packages (from ema_workbench) (3.7.1)
       Requirement already satisfied: statsmodels in /usr/local/lib/python3.10/dist-packages (from ema workbench) (0.14.2)
       Requirement already satisfied: seaborn in /usr/local/lib/python3.10/dist-packages (from ema_workbench) (0.13.1)
       Requirement already satisfied: tqdm in /usr/local/lib/python3.10/dist-packages (from ema_workbench) (4.66.5)
      Requirement already satisfied: multiprocess in /usr/local/lib/python3.10/dist-packages (from salib>=1.4.6->ema_workbenc
      Requirement already satisfied: scipy>=1.9.3 in /usr/local/lib/python3.10/dist-packages (from salib>=1.4.6->ema workbenc
      Requirement already satisfied: contourpy>=1.0.1 in /usr/local/lib/python3.10/dist-packages (from matplotlib->ema_workbe
      Requirement already satisfied: cycler>=0.10 in /usr/local/lib/python3.10/dist-packages (from matplotlib->ema_workbench)
       Requirement already satisfied: fonttools>=4.22.0 in /usr/local/lib/python3.10/dist-packages (from matplotlib->ema workb
       Requirement already satisfied: kiwisolver>=1.0.1 in /usr/local/lib/python3.10/dist-packages (from matplotlib->ema workb
      Requirement already satisfied: packaging>=20.0 in /usr/local/lib/python3.10/dist-packages (from matplotlib->ema_workben
      Requirement already satisfied: pillow>=6.2.0 in /usr/local/lib/python3.10/dist-packages (from matplotlib->ema workbench
       Requirement already satisfied: pyparsing>=2.3.1 in /usr/local/lib/python3.10/dist-packages (from matplotlib->ema workbe
      Requirement already satisfied: python-dateutil>=2.7 in /usr/local/lib/python3.10/dist-packages (from matplotlib->ema wo
       Requirement already satisfied: pytz>=2020.1 in /usr/local/lib/python3.10/dist-packages (from pandas->ema_workbench) (20
       Requirement already satisfied: tzdata>=2022.1 in /usr/local/lib/python3.10/dist-packages (from pandas->ema workbench) (
       Requirement already satisfied: joblib>=1.1.1 in /usr/local/lib/python3.10/dist-packages (from scikit-learn->ema_workben
      Requirement\ already\ satisfied:\ threadpoolctl>=2.\ 0.\ 0\ in\ /usr/local/lib/python 3.\ 10/dist-packages\ (from\ scikit-learn->ema\_lib/python 3.\ 10/dist-packages\ (from\ scikit-learn->e
       Requirement already satisfied: patsy>=0.5.6 in /usr/local/lib/python3.10/dist-packages (from statsmodels->ema_workbench
       Requirement already satisfied: six in /usr/local/lib/python3.10/dist-packages (from patsy>=0.5.6->statsmodels->ema_work
      Requirement already satisfied: dill>=0.3.8 in /usr/local/lib/python3.10/dist-packages (from multiprocess->salib>=1.4.6-
```

The Lake model

The goal of this assignment is to demonstrate the use of multiobjective evolutionary optimization, to learn how to visualize and interpret the optimization results, and to use performance metrics to assess the results obtained via multiobjective evolutionary optimization. We will use the lake problem as a test case, this is a classic problem initially developed by Carpenter et al. (1999) where the population of a city has to decide the amount of annual pollution it will release into a lake. In this exercise, we will use the adapted version in Quinn et al. 2017 were the problem is defined as a state-based control problem (where actions are a function of the state of the system). In this case, the 'action' is the Phosphorous (P) emissions which are optimized to balance the economic benefits and the quality of the lake. Since this is a multi-objective problem, we need a flexible function to map the states to actions, so we use radial basis functions to parameterize the emission control policies. In fact, the MOEA will search for the optimal radii, centers and weights that yield good performance for the objectives of the lake model described below. See the paper for more details about the problem formulation.

The model is defined by the following equation:

$$X_{(t+1)} = X_t + a_t + rac{(X_t^q)}{(1+X_t^q)} - bX_t + \epsilon_t$$

where X_t is the pollution at time t, a_t is the rate of anthropogenic pollution at time t, b is the lake's natural removal rate, q is the lake's natural recycling rate, ϵ_t is the rate of natural pollution at time t. The rate of anthropogenic pollution a_t is the decision variable where $a_t \in [0, 0.1]$.

There are four outcomes of interest. The first is the average concentration of phosphor in the lake.

$$f_{phosphorus} = rac{1}{|T|} \sum_{t \in T} X_t$$

where |T| is the cardinality of the set of points in time. The second objective is the economic benefit derived from polluting the lake defined as the discounted benefit of pollution minus the costs of having a polluted lake.

$$f_{economic} = \sum_{t \in T} lpha a_t \delta^t$$

where α is the utility derived from polluting and δ is the discount rate. By default, α is 0.04. The third objective is related to the year over year change in the anthropogenic pollution rate.

$$f_{inertia} = rac{1}{|T|-1}\sum_{t=1}^{|T|}I(|a_t-a_{t-1}|> au)$$

where I is an indicator function that is 0 if the statement is false, and 1 if the statement is true, τ is the threshold that is deemed undesirable, and is for illustrative purposes et to 0.2. Effectively, f_{inertia} is the fraction of years where the absolute value of the change in anthropogenic pollution is larger then τ . The fourth objective is the fraction of years where the pollution in the lake is below the critical threshold.

$$f_{reliability} = rac{1}{|T|} \sum_{t \in T} I(X_t < X_{crit})$$

where I is an indicator function that is 0 if the statement is false, and 1 if the statement is true, X_{crit} is the critical threshold of pollution and is a function of both b and q.

The lake problem is characterized by both stochastic uncertainty and deep uncertainty. The stochastic uncertainty arises from the natural inflow. To reduce this stochastic uncertainty, multiple replications are performed and the average over the replication is taken. Deep uncertainty is presented by uncertainty about the mean μ and standard deviation sigma of the lognormal distribution characterizing the natural inflow, the natural removal rate of the lake β , the natural recycling rate of the lake q, and the discount rate δ . The table below specifies the ranges for the deeply uncertain factors, as well as their best estimate or default values.

Lake model implementation in python

```
1 import math
2 import numpy as np
3 from scipy.optimize import brentq
4
5
6 def
       get_antropogenic_release(xt, c1, c2, r1, r2, w1):
7
8
9
          Parameters
10
11
          xt : float
12
                   polution in lake at time t
13
          c1 : float
14
                   center rbf 1
          c2 : float
15
16
                   center rbf 2
         rl : float
17
18
                   ratius rbf 1
          r2 : float
19
20
                   ratius rbf 2
21
          w1 : float
22
                   weight of rbf 1
23
```

```
24
          Returns
25
26
          float
27
28
          note:: w2 = 1 - w1
29
          , , ,
30
31
32
          rule = w1 * (abs(xt - c1) / r1) ** 3 + (1 - w1) * (abs(xt - c2) / r2) ** 3
33
          at1 = max(rule, 0.01)
34
         at = min(at1, 0.1)
35
36
          return at
37
38
39 \ \# \ \ def \ \ lake\_model(b=0.42, \ \ q=2.0, \ \ mean=0.02,
40 #
                                 stdev=0.001, delta=0.98, alpha=0.4,
41 #
                                 nsamples=100, myears=100, c1=0.25,
42 #
                                 c2=0.25, r1=0.5, r2=0.5,
43 #
                                 w1=0.5, seed=123):
44
45 \; \# adapted from the following request
46 def lake_model(b=0.42, q=2.0, mean=0.02,
47
                               stdev=0.001, delta=0.98, alpha=0.41,
48
                               nsamples=150, myears=100,
49
                              c1=0.25, c2=0.25, r1=0.5, r2=0.5,
50
                               w1=0.5, seed=123):
          "'runs the lake model for nsamples stochastic realisation using
51
          specified random seed.
53
54
          Parameters
55
56
          b : float
57
                 decay rate for P in lake (0.42 = irreversible)
58
          q : float
59
                recycling exponent
60
          mean : float
61
                         mean of natural inflows
62
          stdev : float
63
                         standard deviation of natural inflows
64
          delta : float
65
                         future utility discount rate
66
          alpha : float
67
                         utility from pollution
68
          nsamples : int, optional
69
          myears : int, optional
70
          c1 : float
71
          c2 : float
72
          r1 : float
73
          r2 : float
74
          w1 : float
75
          seed : int, optional
76
                       seed for the random number generator
77
78
          Returns
79
80
          tuple
81
82
83
          np. random. seed (seed)
84
          Pcrit = brentq(lambda x: x ** q / (1 + x ** q) - b * x, 0.01, 1.5)
85
86
          X = np.zeros((myears,))
87
          average_daily_P = np.zeros((myears,))
88
          reliability = 0.0
89
          inertia = 0
90
          utility = 0
91
92
          for _ in range(nsamples):
```

1. Connecting the lake model with the EMA workbench.

max_P = np.max(average_daily_P)

 $\max_{P} = 0.0$

if isinstance(max_P, (bool, np.bool_)):

print(f"max_P: {max_P}, type: {type(max_P)}")

return max_P, utility, inertia, reliability

Given the Python implementation of the lake problem above, adapt the code and connect it to the EMA workbench using the following lever ranges and uncertainty ranges:

Levers	Range	Default val	ue
r1	0.0 - 2.0	(0.5
r2	0.0 - 2.0	0	.5.
c1	-2 - 2	0.	25
c2	-2 - 2	0.	25
w1	0.0-1.0	(0.5
Uncerta	inties	Range	Default value
μ	(0.01 – 0.05	0.02
σ	0.0	01 - 0.005	0.0017
b		0.1 - 0.45	0.42
q		2 - 4.5	2

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127 128 129

You can follow this tutorial for guidance.

The outcomes in the EMA workbench refers to the objectives of the problem, in this case we have four. 1) maximum Phosphorous (to be minimized) 2) utility (to be maximized) 3) intertia (to be maximized), and reliability (to be maximized). Use an alpha value of 0.41, with number of samples= 150, and number of years = 100.

np. arange (my

```
1 # target:
 2 # Maximum phosphorus concentration: should be minimized, because higher phosphorus concentration means deteri
 3 # Economic benefits: should be maximized to achieve long-term economic benefits through pollution control s
 4 # Inertia: should be maximized to maintain consistency and stability of pollution policies between years.
 5 # Reliability: should be maximized to ensure that lake pollution is below the critical value in more yea
 7
 8 from ema_workbench import (Model, RealParameter, ScalarOutcome, MultiprocessingEvaluator)
10 lakemodel = Model('lakemodel', function=lake model)
11
12 # Levers
13 lakemodel.levers = [RealParameter('r1', 0.0, 2.0),
                                        RealParameter ('r2', 0.0, 2.0),
                                        RealParameter ('c1', -2.0, 2.0),
15
                                        RealParameter('c2', -2.0, 2.0),
16
                                        RealParameter('w1', 0.0, 1.0)]
17
18
19 # Uncertainties
20 lakemodel.uncertainties = [RealParameter('mean', 0.01, 0.05),
21
                                                      RealParameter ('stdev', 0.001, 0.005),
22
                                                      RealParameter ('b', 0.1, 0.45),
                                                      RealParameter('q', 2, 4.5),
23
                                                      RealParameter ('delta', 0.93, 0.99)]
24
25
26 # Outcomes
27 lakemodel.outcomes = [ScalarOutcome('max_P', kind=ScalarOutcome.MINIMIZE),
28
                                            ScalarOutcome('utility', kind=ScalarOutcome.MAXIMIZE),
29
                                            ScalarOutcome ('inertia', kind=ScalarOutcome.MAXIMIZE),
30
                                            ScalarOutcome ('reliability', kind=ScalarOutcome.MAXIMIZE)]
```

1 开始借助 AI 编写或生成代码。

2. How would you introduce a constrain within the optimization to reflect a desired performance threshold for a given objective?

You don't actually have to perform this step, simply specify how you would go about establishing a constraint in the optimization formulation, in such a way that it only finds solutions with maximum pollution (max phosphorous) of 0.85.

```
1 # add a constraint to the optimization problem
2 from ema_workbench import Constraint
3 constraints = [Constraint("max pollution", outcome_names="max_P", function=lambda x: max(0.85, x))]
4
```

3. Run the optimization and track the performance metrics

Tip: the EMA Workbench uses <u>Platypus</u> to run the optimization via the evaluator class, you can also collect metrics during runtime specifying the convergence option. Below is a sample snippet on how to run the optimization and collect performance metrics during runtime.

```
9/13/24, 2:48 PM
      8
      9
     11
     12
     14
     15
     16
     17
     18
```

```
10 ema_logging.log_to_stderr(ema_logging.INF0)
13 with MultiprocessingEvaluator(lakemodel) as evaluator:
        results, convergence = evaluator.optimize(
               nfe=10000,
               searchover='levers',
               epsilons=[0.1, 0.1, 0.01, 0.1],
               convergence=convergence_metrics,
19
               constraints=constraints)
20
21
→
    [MainProcess/INFO] pool started with 2 workers
      0%
                                                 | 0/10000 [00:00<?, ?it/s]
      1%
                                         100/10000 [00:09<15:17, 10.79it/s]
      2%
                                         | 200/10000 [00:17<14:17, 11.43it/s]
                                          | 300/10000 [00:25<13:33, 11.93it/s]
      3%
      4%
                                          400/10000 [00:34<13:53, 11.52it/s]
                                          500/10000 [00:43<13:48, 11.46it/s]
      5%
      6%
                                           600/10000 [00:51<13:25, 11.66it/s]
                                           700/10000 [01:00<13:32, 11.44it/s]
      7%
      8%
                                           800/10000 [01:10<13:50, 11.07it/s]
                                            900/10000 [01:17<12:41, 11.95it/s]
      9%
                                            | 900/10000 [01:28<12:41, 11.95it/s]
      9%
     10%
                                           | 1000/10000 [01:28<13:56, 10.76it/s]
                                           | 1100/10000 [01:36<13:04, 11.35it/s]
     11%
     12%
                                            | 1200/10000 [01:45<12:49, 11.44it/s]
     13%
                                            | 1300/10000 [01:52<12:16, 11.82it/s]
     14%
                                            | 1400/10000 [02:01<11:57, 11.99it/s]
                                             | 1500/10000 [02:09<11:50, 11.96it/s]
     15%
     16%
                                             | 1600/10000 [02:16<11:14, 12.45it/s]
     17%
                                             | 1700/10000 [02:26<11:46, 11.75it/s]
     18%
                                             | 1800/10000 [02:33<10:56, 12.49it/s]
     19%
                                             | 1900/10000 [02:42<11:15, 11.99it/s]
     20%
                                              2000/10000 [02:49<10:31, 12.67it/s]
     21%
                                              2100/10000 [02:58<10:51, 12.13it/s]
                                               2200/10000 [03:04<10:04, 12.91it/s]
     22%
                                               2300/10000 [03:14<10:35, 12.12it/s]
     23%
     24%
                                               2400/10000 [03:22<10:37, 11.93it/s]
     25%
                                                2500/10000 [03:31<10:23, 12.03it/s]
```

双击(或按回车键)即可修改

4. Selecting the objectives.

The outputs from the optimization runs will contain the decision variables (i.e. the parameters of the radial basis functions) and objectives combined. Create a data structure that only contains the objective values without the decision variables.

Tip: each row in the output matrix represents a different solution with it's obective values, the first columns are the decision variables and the last columns are the objective values.

1 results

7	r1	r2	c1	c2	w1	max P	utility	inertia	reliability
	· · ·	' -						21101 020	
0	0.475550	0.976891	0.206675	0.183765	0.942480	0.097738	0.267078	0.990000	1.000000
1	0.085196	1.238544	0.240845	-0.261639	0.773910	0.522222	0.729131	0.987867	0.866733
2	0.085196	1.499870	0.240801	-0.405092	0.825415	0.838874	0.910375	0.985933	0.725267
3	0.085196	0.804180	0.242007	-0.012470	0.847954	0.481739	0.707849	0.743333	0.883867
4	0.086045	1.238544	0.240845	-0.261639	0.773910	0.356860	0.637278	0.989067	0.939400
5	0.456780	1.499182	0.327283	-0.298367	0.944601	0.194646	0.534778	0.990000	1.000000
6	0 085106	N 700425	U 34U83E	N 062171	U 83UU34	U 820330	N QN7Q10	U 0888UU	O 805200
4									

1 convergence

```
₹
          hypervolume
                         epsilon_progress
                                                nfe
      0
                    0.0
                                           0
                                                   0
       1
                    0.0
                                           5
                                                100
       2
                                               1100
                    0.0
                                           6
                                           6
                                               2100
       3
                    0.0
       4
                    0.0
                                           6
                                               3100
       5
                    0.0
                                           6
                                               4100
       6
                    0.0
                                           6
                                               5100
      7
                    0.0
                                          10
                                               6100
       8
                    0.0
                                          16
                                               7100
      9
                    0.0
                                          24
                                               8100
      10
                    0.0
                                          29
                                               9100
                                              10021
                    \cap
```

```
1 import pandas as pd
2
3 objective_columns = ['max_P', 'utility', 'inertia', 'reliability']
4
5
6 objectives_df = results[objective_columns]
7
8
```

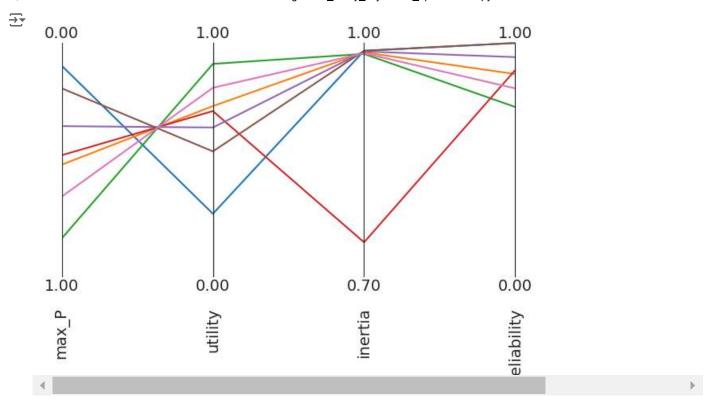
```
9 print(objectives df)
\overline{2}
          max_P
                 utility
                         inertia reliability
    0 0.097738 0.267078 0.990000
                                    1.000000
    1 0.522222 0.729131 0.987867
                                      0.866733
    2 0.838874 0.910375 0.985933
                                    0.725267
    3 0. 481739 0. 707849 0. 743333
                                      0.883867
                                       0.939400
    4 0.356860 0.637278 0.989067
    5 0.194646 0.534778 0.990000
                                      1.000000
    6 0.659329 0.807812 0.986800
                                       0.805200
1 objectives_df.to_csv('objectives_only.csv', index=False)
2 results.to_csv('results.csv', index=False)
3 convergence.to_csv('convergence.csv', index=False)
```

5. Visualizing the results.

Present visually the results of the Pareto optimal solutions (in the objective space), feel free to be creative! Provide a brief discussion the results, are there any tradeoffs observed?

Tip: If you need inspiration check out the EMA workbench parcoords.

```
1 from ema_workbench.analysis import parcoords
2 import matplotlib.pyplot as plt
3
4 data = objectives_df
5
6
7 limits = parcoords.get_limits(data)
8 limits.loc[0, ['max_P', 'utility', 'reliability']] = 0
9 limits.loc[0, ['inertia']] = 0.7
10 limits.loc[1, ['max_P', 'utility', 'inertia', 'reliability']] = 1
11
12
13 paraxes = parcoords.ParallelAxes(limits)
14 paraxes.plot(data)
15 paraxes.invert_axis('max_P')
16
17
18 plt.show()
```

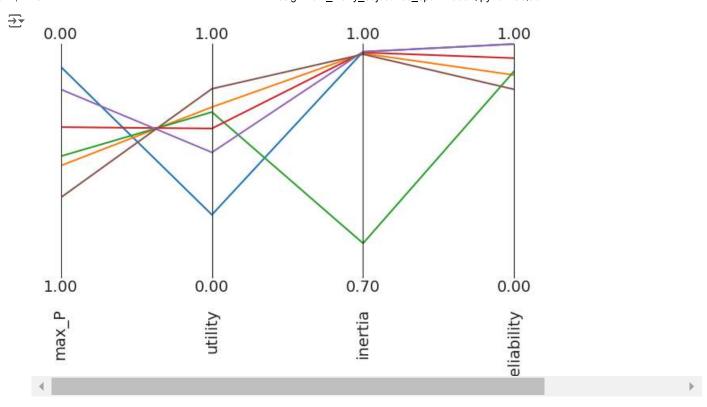


Yes, the parallel coordinates plot reveals several trade-offs among the objectives in the multi-objective optimization problem. Specifically, it shows that reducing maximum phosphorus levels (Max_P) often compromises utility and reliability. Additionally, there's an observable trade-off between inertia and reliability, suggesting that efforts to maintain the status quo (inertia) may adversely affect the reliability of the system.

6. Establishing a performance threshold.

Show visually only the solutions from the Pareto set that yield a reliability above 80%, and briefly discuss the results.

```
1 import plotly.express as px
2
3
4 filtered_data = objectives_df[objectives_df['reliability'] > 0.8]
5
6
7 limits = parcoords.get_limits(data)
8 limits.loc[0, ['max_P', 'utility', 'reliability']] = 0
9 limits.loc[0, ['inertia']] = 0.7
10 limits.loc[1, ['max_P', 'utility', 'inertia', 'reliability']] = 1
11
12
13 paraxes = parcoords.ParallelAxes(limits)
14 paraxes.plot(filtered_data)
15 paraxes.invert_axis('max_P')
16
17
18 plt.show()
19
```



The plot illustrates the balance between Max Phosphorus, Utility, Inertia, and Reliability across different solutions. A tradeoff exists between Max Phosphorus and Reliability, where lower phosphorus doesn't always ensure higher reliability. Similarly, higher utility often corresponds with lower inertia, meaning more flexible, dynamic solutions tend to perform better.

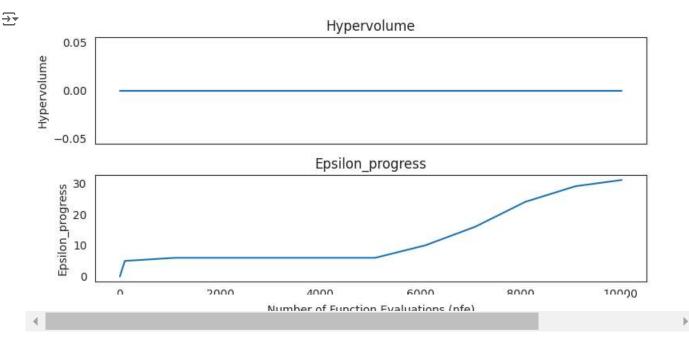
Solutions that achieve high reliability vary in their performance regarding phosphorus and utility, showing multiple paths to reliability. A cluster of solutions also balances high utility and reliability, suggesting an optimal compromise. This visualization helps decision-makers weigh trade-offs based on whether environmental impact or operational performance is prioritized.

7. Performance metrics

Show in a dataframe the results from the metrics (convergence) collected during the optimization. Plot the metrics (e.g hypervolume or epsilon progress) as a function of the number of function evaluations (nfe). Provide a brief discussion of the results.

```
1 import matplotlib.pyplot as plt
2 import seaborn as sns
4 # Set the style
5 sns. set style ("white")
7 # List of possible metrics we want to check if they exist in the DataFrame
8 possible_metrics = ['hypervolume', 'epsilon_progress', 'generational_distance', 'epsilon_indicator', 'inverted_gd',
10 \# Filter out the metrics that exist in the DataFrame
11 available_metrics = [metric for metric in possible_metrics if metric in convergence.columns]
12
13 # Determine the number of plots based on available metrics
14 num_plots = len(available_metrics)
16 # Create subplots dynamically based on the number of available metrics
17 fig, axes = plt.subplots(nrows=num_plots,
                                            figsize=(8,
                                                             num_plots), sharex=True)
18
19 # If only one metric available, wrap the axes object in a list
```

```
20 if num plots == 1:
21
          axes = [axes]
22
23 # Plotting each available metric
24 for ax, metric in zip(axes, available_metrics):
          ax.plot(convergence['nfe'], convergence[metric], label=metric)
26
          ax.set_title(metric.capitalize())
27
          ax. set_ylabel(metric.capitalize())
28
29 # Set common x-axis label
30 axes[-1].set_xlabel('Number of Function Evaluations (nfe)')
32 # Adjust layout
33 plt.tight_layout()
35 # Show the plot
36 plt. show()
37
```



Extra credit: Visualize the phosphorous release from the policy with the highest reliability. If you plot phosphorous release as a function of time, what do you observe?

```
1 def lake_model(b=0.42, q=2.0, mean=0.02, stdev=0.001, delta=0.98, alpha=0.41,
2
                             nsamples=150,
                                          myears=100, c1=0.25, c2=0.25, r1=0.5, r2=0.5, w1=0.5, seed=123):
3
4
         Runs the lake model simulation for a given set of parameters and policies.
5
6
         Parameters:
7
         - b, q, mean, stdev, delta, alpha : float, environmental and economic parameters.
8
         - nsamples : int, number of samples for stochastic realization.
9
         - myears : int, number of years to simulate.
10
          - c1, c2, r1, r2, w1 : float, parameters for the decision-making policy.
11
           seed: int, random seed for reproducibility.
13
         Returns:
14
          - tuple (float, float, float, float, np.array), containing max P, utility, inertia, reliability,
15
             and phosphorous release over time for visualization.
16
17
         np. random. seed (seed)
18
         Pcrit = brentq(lambda x: x ** q / (1 + x **
                                                                           0.01, 1.5)
19
20
         X = np. zeros((myears,))
                                  # Pollution levels
21
          phosphorous_release_over_time = np. zeros((myears,))
                                                              # Store phosphorous release values for each year
          reliability = 0.0
```

```
23
                     inertia = 0
24
                     utility = 0
25
                     for _ in range(nsamples):
26
27
                                    X[0] = 0.0 # Start with no pollution
28
                                    decision = 0.1 # Initial decision
29
30
                                    decisions = np. zeros(myears,)
                                    decisions[0] = decision
31
                                    natural_inflows = np.random.lognormal(math.log(mean ** 2 / math.sqrt(stdev ** 2 + mean ** 2))
32
33
                                                                                                                                                                                         math. sqrt(math. log(1.0 + st
34
                                                                                                                                                                                         size=myears)
35
36
                                     for t in range(1, myears):
37
                                                    decision = get antropogenic release(X[t - 1], c1, c2, r1, r2, w1)
38
                                                    decisions[t] = decision
39
                                                    X[t] = (1 - b) * X[t - 1] + X[t - 1] ** q / (1 + X[t - 1] ** q) + decision
                                                    phosphorous release over time[t] += decision / nsamples # Accumulate average release
41
42
                                    reliability += np. sum(X < Pcrit) / (nsamples * myears)
43
                                     inertia += np.sum(np.absolute(np.diff(decisions) < 0.02)) / (nsamples * myears)</pre>
44
                                    utility += np.sum(alpha * decisions * np.power(delta, np.arange(myears))) / nsamples
45
46
                     max_P = np. max(X)
47
                     return max_P, utility, inertia, reliability, phosphorous_release_over_time
48
49
  1 # Identify the policy with the highest reliability, and it's not only one
  2 max_reliability = results['reliability'].max()
  3 highest_reliability_policies = results[results['reliability'] == max_reliability]
  5 plt.figure(figsize=(12, 6))
  6
  7 for _, policy in highest_reliability_policies.iterrows():
                     max_P, utility, inertia, reliability, phosphorous_release = lake_model(
  9
                                    c1=policy['c1'],
10
                                    c2=policy['c2'],
11
                                    r1=policy['r1'],
                                    r2=policy['r2'],
12
13
                                    w1=policy['w1']
14
15
16
                     plt.plot(np.arange(100), phosphorous_release, label=f'Policy: rl={policy["rl"]}, r2={policy["r2"]}, wl={policy["r2"]}, wl={poli
17
18 plt.xlabel('Year')
19 plt.ylabel ('Phosphorus Release')
20 plt.title('Phosphorus Release Over Time for Policies with Highest Reliability')
21 plt. legend()
22 plt.grid(True)
23 plt. show()
```

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