Chat history

ChatGPT

New chat

Yesterday

Numerical Integration Methods

Previous 7 Days

Sample Caps for Polished Microscopy

More Time for Research

Error in NumPy division.

Brain Injury Assessment: Questions

Letter of Recommendation Outline

Gaussian Quadrature Function in Python

STL File Composition

Parallelize Python Code

Check Python Version Mac

Square Root in Python

Previous 30 Days

Function Distribution Methods

MPI Rank & Size

Japanese 100 Phrases

Unisex and Gender-Neutral

Top 100 Japanese Nouns

Quantifying Earth's Surface Roughness

February

Generate and Export Matrices

3D Printing Overview

Define Linux Environment File

PhD Research Guidance

Falstaff: Model Explanation

Hazards in Work Environments

Japan 2-Week Itinerary

LinkedIn Message Group

20 Questions Game

Clone Git Repo on HPC

DFW Global Entry Locations

Upgrade planGet GPT-4, DALL·E, and more



James Smith Close sidebar ChatGPT 3.5

You

why am I getting this error in the code that will follow in the next message?

/apps/intel/oneapi/intelpython/latest/lib/python3.7/site-packages/numpy/polynomial/legendre.py:

1581: RuntimeWarning: divide by zero encountered in true_divide

w = 1/(fm * df)

/apps/intel/oneapi/intelpython/latest/lib/python3.7/site-packages/numpy/polynomial/legendre.py:

1455: RuntimeWarning: divide by zero encountered in true divide

scl = 1./np.sqrt(2*np.arange(n) + 1)

Traceback (most recent call last):

File "Homework7.py", line 45, in <module> comm.scatter(N, root=0)

File "mpi4py/MPI/Comm.pyx", line 1267, in mpi4py.MPI.Comm.scatter

File "mpi4py/MPI/msgpickle.pxi", line 730, in mpi4py.MPI.PyMPI scatter

File "mpi4py/MPI/msgpickle.pxi", line 119, in mpi4py.MPI.Pickle.dumpv

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The error message you're encountering seems to be related to numerical calculations involving division by zero. Specifically, it's originating from the numpy.polynomial.legendre module.

Here are a few possibilities as to why you might be encountering this error:

Input Data Issue: Check the input data that you're passing to the function. It's possible that the data is causing division by zero.

Algorithm Issue: There could be a bug in the algorithm or function you're using that's leading to division by zero.

Boundary Issue: Division by zero often occurs when an algorithm or function encounters a boundary condition that isn't handled properly.

Library Version: It's possible that there's a bug in the version of NumPy or related libraries you're using. Updating to the latest version might solve the problem.

To address the issue:

- Double-check your input data to ensure it's valid and doesn't contain any zeros or other problematic values.
- Look into the algorithm or function you're using to see if there are any boundary conditions or special cases that need to be handled differently.
- Consider updating your NumPy or related libraries to the latest version to see if the issue has been addressed in newer releases.

Without seeing the specific code causing this issue, it's challenging to provide a more precise solution. If you could share the relevant part of your code, I might be able to provide more targeted assistance.



```
You
import math
from mpi4py import MPI
import numpy as np
# This code was written by James Smith to perform Gaussian Integration
# Start the MPI Process
comm = MPI.COMM_WORLD
# Determine the Total Number of Tasks
size = comm.Get_size()
# Determine the rank (ID) of this task
rank = comm.Get_rank()
# Function to integrate
def f(x):
  return x*math.exp(x)
N = list(range(1,21))
## --- Master ---
if rank == 0:
  # Generate Quadrature Number
  N = list(range(1,21))
  # Integration Bounds
  x1 = -1
  x2 = 1
  # Weights(w) and Integration Points(x)
  w=[]
  x=[]
  for i in N:
```

```
x_temp,w_temp = np.polynomial.legendre.leggauss(i)
     w.append(w_temp)
     x.append(x_temp)
  ### Compute the Quadrature ###
  # Send Data to Workers
  comm.scatter(N, root=0)
  comm.scatter(x, root=0)
  comm.scatter(w, root=0)
  # Collect the results from each worker
  int results=[]
  for i in range(1,size):
     temp=comm.recv(source=i)
     int_results.append(temp)
  # Compute the table
  print(int results)
## --- Worker ---
else:
  print(f"Process {rank} starts") #Displays worker ID
  # Receive Data
  N_{local} = comm.scatter(N, root=0)
  x local = comm.scatter(x, root=0)
  w_local = comm.scatter(w, root=0)
  # Takes 3 arrays and isolates data list specific to this worker
  data_local=[N_local[rank-1],x_local[rank-1],w_local[rank-1]]
  # Takes Data and separates list into 3 arrays for weights, x, and N
  w_worker = data_local[2]
  x worker = data local[1]
  N_worker = data_local[0]
  # Performs Integration
  sum = 0.0
  for i in range (0,N_worker):
     sum = sum + w_worker[i]*f(x_worker[i])
  result=sum
  # Send Results to Main
  comm.send(result, dest=0)
```

Here is the code[^]

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The error message you're encountering seems to originate from the

numpy.polynomial.legendre.leggauss function. This function computes the Gauss-Legendre quadrature points and weights. The warning indicates that there might be division by zero happening within this function, which suggests that the legendre polynomial computation is encountering some numerical instability.

Given the nature of the legendre polynomial, division by zero can occur if the function is evaluated at certain points where the derivative becomes zero. This can happen when the number of quadrature points is low, or when the interval of integration is not properly specified.

To address this issue, you may consider the following:

Increase Quadrature Points: Increase the number of quadrature points (\mathbb{N}) to see if the error persists. More points might provide better accuracy and avoid division by zero issues.

Check Integration Bounds: Ensure that the integration bounds (x1 and x2) are properly specified. The legendre polynomial is defined on the interval [-1, 1], so if your bounds differ, you may encounter numerical instability.

Handle Numerical Instability: If increasing quadrature points or adjusting bounds doesn't resolve the issue, you may need to handle numerical instability by catching the warning or using a different integration method.

Debug Legendre Polynomial: Debug the legendre polynomial computation to understand where the division by zero is occurring. It might give you insights into what's causing the issue.

Try adjusting the code based on these suggestions and see if the error persists.



You

what is the proper syntax for numpy.polynomial.legendre.leggauss

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The proper syntax for numpy.polynomial.legendre.leggauss is as follows:

python

```
Copy code
```

```
numpy.polynomial.legendre.leggauss(deg)
```

Where:

• deg: int, The degree of the legendre polynomial to use. It determines the number of quadrature points and weights.

This function returns two arrays:

- x: 1-D array, The roots (nodes) of the legendre polynomial of degree deg.
- w: 1-D array, The weights corresponding to the legendre polynomial of degree deg.

Here's an example of how you can use numpy.polynomial.legendre.leggauss:

```
python
Copy code
import numpy as np
# Degree of the legendre polynomial
degree = 5
# Compute Legendre-Gauss quadrature points and weights
x, w = np.polynomial.legendre.leggauss(degree)
print("Quadrature Points (Roots):", x)
print("Weights:", w)
```

This will compute the Legendre-Gauss quadrature points and weights for a Legendre polynomial of degree 5.



You

Assignment:

Write an MPI code with Python that computes the integral of f(x) = xex using quadrature values of n = 1, 2, ..., 20. That is, 20 evaluations of the integral will be obtained, each with a different quadrature. Each integral is to be computed by a different worker process. The Master computes the quadrature (x and y values), passes the quadrature to a worker

process, collects the results from each process, then constructs and prints a table with the results

of the form:

Quadrature no. Integration Result Percent error Run time (s)

1 2 3

... 20

The Python function gauleg.py can be called to obtain the quadrature for any number of points and for arbitrary integration limits. This code can be found in /work/ME5773/sec6/Homework.

The worker processes obtain the quadrature, evaluate the integral, and send the results back to the master process.

Part 1: write the file that is able to execute the MPI program, and run the program with 21 tasks (1 master, and 20 workers). Save the outputs of run in a text file named 'part1.txt' Part 2 (15% extra credit): On a different python file, modify the program for the case where the number of tasks is < 20 but still completes the table for up to 20 integration points. Write the outputs of this run in 'part2.txt'.

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To complete this assignment, you'll need to write two Python files: one for the MPI program itself and another for running the MPI program with different task counts. Below, I'll provide the code for both parts:

Part 1: MPI Program (mpi_integration.py)

```
copy code
import math
from mpi4py import MPI
import numpy as np
import time

# Function to integrate
def f(x):
    return x * math.exp(x)

# Function to compute legendre-gauss quadrature points and weights
def compute_quadrature(n):
    x, w = np.polynomial.legendre.leggauss(n)
```

```
return x, w
# Master process
def master(size):
# Create table header
print("Quadrature no.\tIntegration Result\tPercent error\tRun time (s)")
 # Loop over quadrature points
 for n in range(1, 21):
 # Start timer
 start_time = time.time()
 # Compute quadrature
x, w = compute_quadrature(n)
 # Send quadrature to worker
comm.send((x, w), dest=n, tag=1)
 # Receive result from worker
 result = comm.recv(source=n, tag=2)
 # Stop timer
 end_time = time.time()
 # Compute percent error
 exact_result = math.exp(1) - 1 # Exact integral of f(x) = x * exp(x) from -1 to 1
 percent_error = abs((result - exact_result) / exact_result) * 100
 # Print table row
 print(f"{n}\t\t{result}\t\t{percent_error:.6f}%\t\t{end_time - start_time:.6f}")
# Worker process
def worker(rank):
# Receive quadrature from master
x, w = comm.recv(source=0, tag=1)
 # Compute local integral
 local_result = sum(w[i] * f(x[i]) for i in range(len(x)))
 # Send result to master
 comm.send(local_result, dest=0, tag=2)
if __name__ == "__main__":
 comm = MPI.COMM_WORLD
 size = comm.Get_size()
```

```
rank = comm.Get_rank()

if size != 21:
   if rank == 0:
   print("Error: This program requires exactly 21 processes (1 master and 20 workers).")
   exit()

if rank == 0:
   master(size)
   else:
   worker(rank)
```

Part 2: MPI Program with Fewer Tasks (mpi_integration_part2.py)

```
python
Copy code
import math
from mpi4py import MPI
import numpy as np
import time
# Function to integrate
def f(x):
return x * math.exp(x)
# Function to compute legendre-gauss quadrature points and weights
def compute_quadrature(n):
x, w = np.polynomial.legendre.leggauss(n)
 return x, w
# Master process
def master(size):
# Create table header
print("Quadrature no.\tIntegration Result\tPercent error\tRun time (s)")
# Loop over quadrature points
 for n in range(1, size):
# Start timer
 start_time = time.time()
```

```
# Compute quadrature
 x, w = compute_quadrature(n)
# Send quadrature to worker
 comm.send((x, w), dest=n, tag=1)
 # Receive result from worker
 result = comm.recv(source=n, tag=2)
 # Stop timer
 end_time = time.time()
# Compute percent error
 exact_result = math.exp(1) - 1 # Exact integral of f(x) = x * exp(x) from -1 to 1
 percent_error = abs((result - exact_result) / exact_result) * 100
 # Print table row
 print(f"{n}\t\t{result}\t\t{percent_error:.6f}%\t\t{end_time - start_time:.6f}")
# Worker process
def worker(rank):
# Receive quadrature from master
x, w = comm.recv(source=0, tag=1)
# Compute local integral
 local_result = sum(w[i] * f(x[i]) for i in range(len(x)))
 # Send result to master
 comm.send(local_result, dest=0, tag=2)
if __name__ == "__main__":
comm = MPI.COMM_WORLD
size = comm.Get_size()
 rank = comm.Get_rank()
if rank == 0:
master(size)
else:
 worker(rank)
```

To run these scripts, you need to execute them using MPI. For example:

```
Copy code
```

```
mpirun -np 21 python mpi_integration.py > part1.txt
```

Copy code

```
mpirun -np 10 python mpi_integration_part2.py > part2.txt
```

These commands will execute the MPI programs with 21 and 10 processes respectively and save the outputs to part1.txt and part2.txt.



You

what is the syntax for the comm.scatter command

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The syntax for the comm.scatter command in MPI4py is as follows:

python

Copy code

```
comm.scatter(sendobj, recvobj=None, root=0)
```

Where:

- sendobj: Object to be scattered. This can be a Python list, NumPy array, or any other iterable.
- recvobj: Buffer where the received data will be placed. This argument is required for non-root processes (workers) and can be omitted for the root process (master).
- root: Rank of the sending process (master) that scatters the data. This argument is optional and defaults to 0.

Here's an example of how comm.scatter can be used:

```
python
Copy code
from mpi4py import MPI
comm = MPI.COMM_WORLD
size = comm.Get_size()
```

```
rank = comm.Get_rank()
# Data to be scattered
data = None
if rank == 0:
data = [1, 2, 3, 4, 5]
# Scatter data to all processes
recv_data = comm.scatter(data, root=0)
print(f"Process {rank}: Received data = {recv_data}")
In this example, process 0 scatters the data [1, 2, 3, 4, 5] to all other processes using
comm.scatter. Each process receives one element of the scattered data. The received data is
stored in the recv_data variable.
You
ValueError
                            Traceback (most recent call last)
File /work/ayn868/ME5773SP24/assignment7/Homework7.py:45
        x.append(x_temp)
  42 ### Compute the Quadrature ###
  43
  44 # Send Data to Workers
---> 45 comm.scatter(N, root=0)
  46 comm.scatter(x, root=0)
  47 comm.scatter(w, root=0)
File mpi4py/MPI/Comm.pyx:1587, in mpi4py.MPI.Comm.scatter()
File mpi4py/MPI/msgpickle.pxi:823, in mpi4py.MPI.PyMPI scatter()
File mpi4py/MPI/msgpickle.pxi:161, in mpi4py.MPI.pickle_dumpv()
ValueError: expecting 1 items, got 20
```

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The error message indicates a mismatch between the number of items being scattered and the number of processes in the communicator.

In MPI, when using <code>comm.scatter</code>, the number of items in the send buffer must match the number of processes in the communicator, except for the root process where the send buffer can be larger.

In your case, it seems like you're trying to scatter the list \mathbb{N} , \mathbb{X} , and \mathbb{W} to all processes including the root process. However, you should only scatter to worker processes, not to the root process itself.

Here's how you can modify your code to fix this issue:

```
python
Copy code
## --- Master ---
if rank == 0:
# Generate Quadrature Number
N = list(range(1,21))
# Integration Bounds
 x1 = -1
 x2 = 1
 # Weights(w) and Integration Points(x)
w=[]
 x=[]
 for i in N:
 x_{temp}, w_{temp} = np.polynomial.legendre.leggauss(i)
 w.append(w_temp)
 x.append(x_temp)
 # Send Data to Workers
 comm.scatter(N[1:], root=0) # Send N to all workers except root
 comm.scatter(x[1:], root=0) # Send x to all workers except root
 comm.scatter(w[1:], root=0) # Send w to all workers except root
```

This modification sends N, x, and w to all worker processes (rank > 0), skipping the root process (rank = 0). This should resolve the ValueError you encountered.

You

will the brackets on N,x, and w cause them to only share part of the data?

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Yes, you're correct. Using slicing with brackets ([1:]) will cause only a part of the data to be shared. This is because slicing extracts a portion of the list starting from the index 1 (the second element) to the end of the list.

If you want to send the entire lists $_{\rm N}$, $_{\rm x}$, and $_{\rm w}$ to all worker processes except the root process, you should not use slicing. Instead, you should send the entire lists directly. Here's the corrected code:

```
python
Copy code
## --- Master ---
if rank == 0:
# Generate Quadrature Number
N = list(range(1,21))
# Integration Bounds
 x1 = -1
 x2 = 1
 # Weights(w) and Integration Points(x)
 w=[]
 x=[]
 for i in N:
 x_{temp}, w_{temp} = np.polynomial.legendre.leggauss(i)
w.append(w_temp)
 x.append(x_temp)
 # Send Data to Workers
 comm.scatter(N, root=0) # Send N to all workers including root
 comm.scatter(x, root=0) # Send x to all workers including root
 comm.scatter(w, root=0) # Send w to all workers including root
```

This modification sends the entire lists N, x, and w to all worker processes (including the root process). It ensures that each worker process receives the complete data for N, x, and w.

You

how do I specify to send to all workers except root for comm.scatter?

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I apologize for the confusion. My previous responses were incorrect. The <code>comm.scatter</code> function in MPI doesn't have an option to specify excluding the root process. Instead, you should only scatter to the worker processes and not to the root process.

To achieve this, you can use an if condition to ensure that the root process doesn't scatter data to itself. Here's how you can modify your code:

```
python
Copy code
## --- Master ---
if rank == 0:
# Generate Quadrature Number
N = list(range(1,21))
# Integration Bounds
 x1 = -1
 x2 = 1
 # Weights(w) and Integration Points(x)
 w = []
 x=[]
 for i in N:
x_temp, w_temp = np.polynomial.legendre.leggauss(i)
 w.append(w_temp)
 x.append(x_temp)
# Send Data to Workers
 for i in range(1, size): # Skip root process (rank 0)
 comm.send(N, dest=i, tag=1)
 comm.send(x, dest=i, tag=2)
 comm.send(w, dest=i, tag=3)
```

This modification sends N, x, and w to all worker processes except the root process (rank != 0). The root process iterates over all worker processes and sends the data to each one individually using the comm.send function.



You

how do I receive this tagged data?

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To receive tagged data in MPI4py, you can use the recv method with the appropriate tag. You'll need to ensure that each worker process receives the correct tagged data sent by the master process.

Here's how you can modify your worker process to receive tagged data:

```
python
Copy code
## --- Worker ---
else:
  print(f"Process {rank} starts") #Displays worker ID

# Receive Data
N_local = comm.recv(source=0, tag=1)
x_local = comm.recv(source=0, tag=2)
w_local = comm.recv(source=0, tag=3)

# Perform computations using received data
# ...

# Send Results back to Main
comm.send(result, dest=0, tag=4) # Adjust tag as needed
```

In this modification:

- Each worker process (except the root process) uses <code>comm.recv</code> to receive data from the root process (rank 0) with the specified tag.
- Each piece of data (N, x, and w) is tagged separately so that workers can receive them individually.

• After performing computations using the received data, the worker process can send the results back to the root process with an appropriate tag (e.g., tag=4).

Make sure to adjust the tag numbers according to your application's needs and ensure that they match the tags used in the master process for sending data.



You

can I send without a destination?

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In MPI, when using the <code>comm.send</code> function in Python with MPI4py, you must specify a destination process to which the message will be sent. You cannot send a message without specifying a destination process.

Every communication in MPI involves a sender process and a receiver process. The <code>comm.send</code> function requires you to specify the destination process to which the message will be sent. This ensures that the message is received by the intended process.

Here's an example of how you can use comm. send to send data from one process to another:

```
python
Copy code
from mpi4py import MPI

comm = MPI.COMM_WORLD
rank = comm.Get_rank()

if rank == 0: # If the process is the sender
   data = "Hello, world!"
   destination = 1 # Destination process rank
   comm.send(data, dest=destination) # Sending data to destination process
elif rank == 1: # If the process is the receiver
   source = 0 # Source process rank
   received_data = comm.recv(source=source) # Receiving data from source process
   print("Received:", received_data)
```

In this example, process 0 sends the message "Hello, world!" to process 1 using comm.send, specifying process 1 as the destination (dest=1). Process 1 receives the message using comm.recv, specifying process 0 as the source (source=0).



You

what is comm.bcast

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comm.bcast is a collective communication operation in MPI (Message Passing Interface) that broadcasts data from one process to all other processes in a communicator. It is used to distribute the same data from one process (often referred to as the "root" process) to all other processes in the communicator.

The syntax for comm.bcast in MPI4py is as follows:

```
python
Copy code
comm.bcast(sendobj, root=0)
```

Where:

- sendobj: The data object that will be broadcasted from the root process to all other processes. This can be any picklable Python object (e.g., a list, tuple, NumPy array, etc.).
- root: The rank of the root process that has the data to be broadcasted. This argument is optional and defaults to 0.

Here's an example of how comm.bcast can be used:

```
python
Copy code
from mpi4py import MPI
comm = MPI.COMM_WORLD
rank = comm.Get_rank()
```

```
if rank == 0:
    data = [1, 2, 3, 4, 5] # Data to be broadcasted
else:
    data = None # Placeholder for received data

data = comm.bcast(data, root=0) # Broadcasting data from root process to all other
processes

print(f"Process {rank}: Received data: {data}")
```

In this example, process 0 has the data [1, 2, 3, 4, 5], and it broadcasts this data to all other processes in the communicator using comm.bcast. All processes (including process 0) receive the broadcasted data.



You

how do I receive this broadcast

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To receive a broadcasted message in MPI4py, you can use the <code>comm.bcast</code> function. This function will broadcast the data from the root process to all other processes in the communicator. Each process, including the root process, can receive the broadcasted data.

Here's how you can receive a broadcasted message:

```
python
Copy code
from mpi4py import MPI

comm = MPI.COMM_WORLD
rank = comm.Get_rank()

if rank == 0:
   data = [1, 2, 3, 4, 5] # Data to be broadcasted else:
   data = None # Placeholder for received data
```

data = comm.bcast(data, root=0) # Broadcasting data from root process to all other
processes

```
print(f"Process {rank}: Received data: {data}")
```

In this example, process 0 has the data [1, 2, 3, 4, 5], and it broadcasts this data to all other processes in the communicator using comm.bcast. All processes (including process 0) receive the broadcasted data.

Each process calls <code>comm.bcast</code>, specifying the data to be broadcasted and the root process (rank 0). The return value of <code>comm.bcast</code> is the broadcasted data, so you can assign it to a variable (data in this case) to receive the broadcasted message.



You

what does scattering do vs. broadcasting?

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Scattering and broadcasting are both collective communication operations in MPI, but they serve different purposes and have different behaviors:

Scattering:

- Scattering is a collective communication operation where the root process sends different pieces of data to different processes in the communicator.
- The data to be scattered is divided into segments, and each segment is sent to a different process.
- Each process receives a different segment of the data, and the segments may have different sizes.
- Scattering is useful when you have a large dataset that needs to be divided among different processes for parallel processing.

Broadcasting:

- Broadcasting is a collective communication operation where the root process sends the same piece of data to all other processes in the communicator.
- The root process has the data, and it sends copies of this data to all other processes.
- All processes in the communicator receive the same data.
- Broadcasting is useful when you have a piece of data that needs to be shared among all processes in the communicator.

In summary:

- Scattering divides data into segments and sends different segments to different processes.
- Broadcasting sends the same data from the root process to all other processes.



You

can you please explain this slurm job script #!/bin/bash

#-----

Sample Slurm job script for Arc nodes

#SBATCH -J myjob # Job Name

#SBATCH -o myjob.o%j # Name of the stdout output file

#SBATCH -e myjob.e%j # Name of stderr error file

#SBATCH -p compute1 # Queue (partition) name

#SBATCH -N 2 # Total # of nodes (must be 1 for serial)

#SBATCH -n 40 # Total # of mpi tasks (should be 1 for serial)

#SBATCH -t 00:10:00 # Run time (hh:mm:ss)

#SBATCH --mail-type=ALL

#SBATCH --mail-user=pratik.mitra@my.utsa.edu

Load module

module load openmpi

module swap openmpi intel/oneapi/2021.2.0

#module load libctl/4.5.0

#module load gcc

#module swap intel/oneapi gcc/11.2.0

cd /work/gug303/sources/ME5773SP24/assignment7

#mpirun -np 80 Homework7.py

mpiexec -n 80 python Homework7.py

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Certainly! This is a Slurm job script, which is a shell script used to submit jobs to a

high-performance computing (HPC) cluster managed by the Slurm workload manager. Let's

break down the script and understand each part:

bash

Copy code

#!/bin/bash

This line specifies the interpreter to use, which is Bash in this case. It indicates that the script is written in the Bash scripting language.

```
bash
Copy code
#SBATCH -J myjob
```

This line sets the name of the job to myjob. Jobs submitted to the cluster will be identified by this name.

```
bash
Copy code
#SBATCH -o myjob.o%j
#SBATCH -e myjob.e%j
```

These lines specify the names of the output and error files for the job. %j is a Slurm job ID placeholder, which will be replaced by the actual job ID when the job is submitted.

```
bash
Copy code
#SBATCH -p compute1
```

This line specifies the partition or queue name where the job will be submitted. The job will be scheduled and run on the nodes allocated to this partition.

```
bash
Copy code
#SBATCH -N 2
```

This line specifies the number of nodes requested for the job. In this case, the job requests 2 nodes.

bash

Copy code

#SBATCH -n 40

This line specifies the total number of MPI tasks requested for the job. MPI tasks are typically processes that communicate with each other using the Message Passing Interface (MPI) protocol. In this case, the job requests 40 MPI tasks.

bash

Copy code

#SBATCH -t 00:10:00

This line specifies the maximum run time for the job. It is expressed in hours, minutes, and seconds (hh:mm:ss). In this case, the job is allowed to run for a maximum of 10 minutes.

bash

Copy code

```
#SBATCH --mail-type=ALL
#SBATCH --mail-user=pratik.mitra@my.utsa.edu
```

These lines specify that email notifications should be sent for job events of all types (--mail-type=ALL) to the specified email address (pratik.mitra@my.utsa.edu).

bash

Copy code

```
module load openmpi
module swap openmpi intel/oneapi/2021.2.0
```

These lines load specific software modules. The module load command loads the Open MPI library, and the module swap command swaps the previously loaded Open MPI module with a different version (intel/oneapi/2021.2.0).

bash

Copy code

cd /work/gug303/sources/ME5773SP24/assignment7

This line changes the current working directory to

/work/gug303/sources/ME5773SP24/assignment7. This is where the script expects to find the Python script (Homework7.py) that will be executed.

bash

Copy code

```
#mpirun -np 80 Homework7.py
mpiexec -n 80 python Homework7.py
```

These lines are commented out, but they are options for running the MPI program. The first option (mpirun) is commented out, while the second option (mpiexec) is not. Both commands execute the Python script Homework 7.py with 80 MPI tasks.

Overall, this Slurm job script specifies the job's characteristics, such as its name, requested resources (nodes, MPI tasks, run time), email notification settings, and software environment. It also sets up the execution environment and launches the MPI program with the specified parameters.



You

if I am trying to run the previous code we discussed with 20 workers + the master while using the minimum number of resources, what should I change in this file?

ChatGPT

ChatGPT

To run the previous code with 20 workers (plus the master process) while using the minimum number of resources, you would need to make the following changes to the Slurm job script:

Reduce the number of nodes: Since you only need 21 processes (20 workers + 1 master), you can reduce the number of nodes to 1.

Reduce the number of MPI tasks: Set the total number of MPI tasks to 21.

Adjust the run time: If the job will complete quickly, you can reduce the run time accordingly.

Here's the modified Slurm job script with these changes:

```
Copy code

#!/bin/bash

#SBATCH -J myjob

#SBATCH -o myjob.o%j

#SBATCH -e myjob.e%j

#SBATCH -p compute1

#SBATCH -N 1 # Reduce the number of nodes to 1

#SBATCH -n 21 # Set the total number of MPI tasks to 21 (20 workers + 1 master)

#SBATCH -t 00:05:00 # Adjust the run time as needed

#SBATCH --mail-type=ALL

#SBATCH --mail-user=pratik.mitra@my.utsa.edu
```

cd /work/gug303/sources/ME5773SP24/assignment7

module swap openmpi intel/oneapi/2021.2.0

mpiexec -n 21 python Homework7.py # Run the MPI program with 21 processes

With these modifications, the job script will request one node with a total of 21 MPI tasks (20 workers and 1 master). Adjust the run time as needed based on the expected execution time of your MPI program.



bash

You

how do I show line numbers in vIM

module load openmpi

ChatGPT

ChatGPT

To show line numbers in Vim, you can use the :set number command. This will display line numbers along the left side of the Vim editor window.

Here's how you can enable line numbers in Vim:

Open Vim by typing vim followed by the filename you want to edit.

Copy code

vim filename

Once inside Vim, press the Esc key to ensure you're in normal mode. Type the following command and press Enter:

arduino

Copy code

:set number

This command tells Vim to display line numbers. If you want Vim to remember this setting for future editing sessions, you can add the command to your .vimrc file. Open your .vimrc file (or create it if it doesn't exist) and

add the following line:

arduino

Copy code

set number

Save the .vimrc file and restart Vim for the changes to take effect.

That's it! Now Vim will display line numbers in the editor window, making it easier to navigate through your files.



You

when I run my code, I get an error saying that the x variable is undefined in the $x_{local} = comm.bcast(x,root=0)$ line

ChatGPT

ChatGPT

The error you're encountering suggests that the variable x is undefined at the point where you're attempting to use it as an argument for comm.bcast.

To resolve this issue, you need to ensure that the variable x is defined and initialized before you attempt to broadcast it using comm.bcast. Here's a general approach to ensure the variable x is defined:

Define and initialize the variable x: Ensure that x is defined and assigned a value before attempting to broadcast it. This could involve computing the value of x or loading it from an external source.

Check the scope of x: Make sure that x is defined within the scope of the process where comm.bcast is called. If x is defined in a conditional block or function that is not executed in the process where comm.bcast is called, it will be undefined.

Here's an example of how you can define and initialize x before broadcasting it:

```
Copy code
# Master process
if rank == 0:
    # Generate x
x = generate_x()
```

Broadcast x to all processes

x = comm.bcast(x, root=0)

In this example:

python

else:
 x = None

- In the master process (rank 0), x is generated using a function generate x().
- In all other processes, x is initialized as None.
- Then, comm.bcast is used to broadcast x from the master process to all other processes.

Ensure that $generate_x()$ or the method you're using to define x is properly implemented and that it returns a valid value. Additionally, make sure that the variable x is defined in the scope where $gomm_x$ beast is called.



You

how would I receive this x broadcast in the worker?

ChatGPT

ChatGPT

To receive the broadcasted variable x in the worker process, you can use the comm.bcast function. In MPI4py, all processes in the communicator call comm.bcast, specifying the variable to be broadcasted (x) and the root process (in this case, root=0).

Here's how you can receive the broadcasted variable x in the worker process:

```
python
Copy code
# Worker processes
else:
    # Receive the broadcasted variable x
    x = comm.bcast(None, root=0)

# Now x is available in all worker processes
# Proceed with your computations using x
```

In this code snippet:

- In the worker process (where rank != 0), the comm.bcast function is called to receive the broadcasted variable x.
- The comm.bcast function is called with None as the argument for the variable to be broadcasted. This is because the root process (rank=0) specifies the variable to be broadcasted, and other processes only need to specify the root process.
- After calling comm.bcast, the variable x will contain the broadcasted value in all worker processes, and you can proceed with your computations using x.

Make sure to place this code snippet in the worker process block of your MPI code where you need to use the broadcasted variable \mathbf{x} .



You

if I am broadcasting multiple lists at once, how do I specify which list is received?

ChatGPT

ChatGPT

When broadcasting multiple lists at once in MPI4py, you can use tags to distinguish between the different lists being broadcasted and received. Each list can be assigned a unique tag, allowing processes to identify which list they are receiving.

Here's how you can broadcast multiple lists with different tags and receive them in worker processes:

Master process:

```
python
Copy code
# Master process (rank 0)
if rank == 0:
  # Broadcast list 'x' with tag 1
  comm.bcast(x, root=0, tag=1)
  # Broadcast list 'y' with tag 2
  comm.bcast(y, root=0, tag=2)
  # Broadcast list 'z' with tag 3
  comm.bcast(z, root=0, tag=3)
```

Worker processes:

```
python
Copy code
# Worker processes (rank != 0)
else:
    # Receive list 'x' with tag 1
    x_received = comm.bcast(None, root=0, tag=1)
# Receive list 'y' with tag 2
    y_received = comm.bcast(None, root=0, tag=2)
# Receive list 'z' with tag 3
    z_received = comm.bcast(None, root=0, tag=3)
```

In this example:

• The master process broadcasts three lists (x, y, and z) to all other processes.

- Each list is broadcasted with a unique tag (1, 2, and 3 respectively).
- In the worker processes, the comm.bcast function is called to receive each list with the corresponding tag.
- After calling comm.bcast, the received lists (x_received, y_received, and z_received) contain the broadcasted lists in the worker processes.

By using different tags for each broadcast and receive operation, you can ensure that the worker processes correctly identify and receive each list.



You

what is the significance of the None in the first field of the receiving broadcast

ChatGPT

ChatGPT

In MPI4py, the <code>comm.bcast</code> function is used for broadcasting data from one process to all other processes in the communicator. When receiving a broadcast in a worker process, you might notice that <code>None</code> is passed as the first argument to <code>comm.bcast</code>. Let's discuss the significance of passing <code>None</code> in this context:

```
python
Copy code

# Worker processes (rank != 0)
else:
    # Receive list 'x' with tag 1
    x_received = comm.bcast(None, root=0, tag=1)
```

In this code snippet:

- The comm.bcast function is being called in the worker process to receive the broadcasted data.
- The first argument (None) in the comm.bcast function call is used as a placeholder. It is not used to transmit any actual data.
- The purpose of passing None as the first argument is to indicate to MPI4py that the process is a receiver of the broadcast and to specify additional parameters such as the source root and the tag.
- When comm.bcast is called with None as the first argument, MPI4py expects to receive
 data from the root process specified by the root parameter and identified by the tag
 parameter.

In summary, passing $_{\text{None}}$ as the first argument in the $_{\text{comm.bcast}}$ call when receiving a broadcast in MPI4py indicates to the MPI library that the process is a receiver of the broadcast and expects to receive data from the specified root process. The actual data being broadcasted is transmitted by the root process, and $_{\text{None}}$ serves as a placeholder to trigger the receive operation.



You

when would i put anything other than None in that field?

ChatGPT

ChatGPT

In MPI4py, the first argument of the <code>comm.bcast</code> function when receiving a broadcast is typically set to <code>None</code> because it acts as a placeholder. However, there are situations where you might put something other than <code>None</code> in that field:

Sending Data: If the worker process is also broadcasting data to other processes, you would put the data you want to broadcast in place of None. For example:

python

Copy code

```
# Worker processes (rank != 0)
else:
  # Broadcast data_list to all processes
data_list = [1, 2, 3, 4, 5]
comm.bcast(data_list, root=0, tag=1)
```

In this case, <code>data_list</code> contains the data that the worker process wants to broadcast to all other processes. The <code>comm.bcast</code> function is used to broadcast <code>data_list</code> to all processes in the communicator.

Receiving Non-Root Process Data: If a worker process is receiving data broadcasted by another non-root process, you would put the received data type in place of None. For example:

python

Copy code

```
# Worker processes (rank != 0)
else:
  # Receive data_list broadcasted by process with rank 2
data_list = comm.bcast([], root=2, tag=1)
```

In this case, the worker process expects to receive data broadcasted by the process with rank 2. The empty list [] serves as a placeholder to trigger the receive operation. The actual broadcasted data will be received and stored in the <code>data list</code> variable.

In summary, you would put something other than <code>None</code> in the first argument of <code>comm.bcast</code> when the worker process needs to either send data or receive data broadcasted by another non-root process. Otherwise, when receiving a broadcast from the root process, <code>None</code> is typically used as a placeholder.



You

Traceback (most recent call last):

File "Homework7.py", line 62, in <module>

N local = comm.bcast(None, root=0, tag=1)

File "mpi4py/MPI/Comm.pyx", line 1254, in mpi4py.MPI.Comm.bcast

TypeError: bcast() got an unexpected keyword argument 'tag'

ChatGPT

ChatGPT

The error indicates that the comm.bcast function does not expect the tag keyword argument.

This is because the tag argument is not a valid parameter for the bcast function in MPI4py.

In MPI, communication functions such as beast do not have a tag argument. Tags are used with point-to-point communication functions like send and recv, but not with collective communication functions like beast, scatter, or gather.

Here's how you can modify your code to remove the tag argument:

```
python
Copy code
# Worker processes (rank != 0)
else:
    # Receive broadcasted data from root process
N_local = comm.bcast(None, root=0)
```

In this modified code, the comm.bcast function is called without the tag argument. The worker process simply waits to receive the broadcasted data from the root process specified by the root argument (in this case, root=0). Make sure that the root process broadcasts the data without specifying a tag as well.

If you need to distinguish between different broadcasts in your application, you may consider using different bcast calls for each piece of data, each with its own unique tag, and corresponding recv calls in the worker processes. However, this is only applicable to point-to-point communication, not to collective communication functions like bcast.



You

why the fuck did you tell me to put tag?

ChatGPT

ChatGPT

I apologize for the confusion. It seems there was a misunderstanding. I provided incorrect information in my response.

In MPI4py, the <code>comm.bcast</code> function does not accept a <code>tag</code> argument. Tags are used with point-to-point communication functions like <code>send</code> and <code>recv</code>, not with collective communication functions like <code>bcast</code>.

To correct the issue, simply remove the tag argument from the comm.bcast function call. Here's the corrected code:

```
python
Copy code
# Worker processes (rank != 0)
else:
  # Receive broadcasted data from root process
N_local = comm.bcast(None, root=0)
```

I apologize for any inconvenience caused by the incorrect information. If you have any further questions or need clarification, please feel free to ask.



You

I want to broadcast 3 sets of data at once, how can I do this without mixing them up?

ChatGPT

ChatGPT

To broadcast multiple sets of data at once without mixing them up, you can use a simple approach where you broadcast each set of data separately. This ensures that each set of data is sent and received independently of the others. Here's how you can do it:

```
python
Copy code
# Master process (rank 0)
if rank == 0:
# Broadcast each set of data separately
comm.bcast(data1, root=0)
 comm.bcast(data2, root=0)
comm.bcast(data3, root=0)
python
Copy code
# Worker processes (rank != 0)
else:
 # Receive each set of data separately
data1_received = comm.bcast(None, root=0)
 data2_received = comm.bcast(None, root=0)
data3_received = comm.bcast(None, root=0)
```

In this approach:

- The master process broadcasts each set of data (data1, data2, data3) separately using comm.bcast.
- Each worker process receives each set of data separately using comm.bcast.
- Each set of data is sent and received independently, ensuring that they are not mixed up.

Make sure that data1, data2, and data3 are defined and initialized before calling comm.bcast, and that they are properly received in the worker processes. This approach keeps the broadcast and receive operations separate, avoiding any mixing up of the data sets.



You

how does the worker know which order to receive the data in?

ChatGPT

ChatGPT

In the approach outlined above, the worker processes receive the data sets in the same order in which they are broadcasted by the master process. This order is determined by the sequence of comm.bcast calls made by the master process.

Let's consider the example provided earlier:

```
python
Copy code
# Master process (rank 0)
if rank == 0:
  # Broadcast each set of data separately
  comm.bcast(data1, root=0)
  comm.bcast(data2, root=0)
  comm.bcast(data3, root=0)
```

In this code snippet, data1 is broadcasted first, followed by data2, and then data3.

```
python
Copy code
# Worker processes (rank != 0)
else:
  # Receive each set of data separately
  data1_received = comm.bcast(None, root=0)
  data2_received = comm.bcast(None, root=0)
  data3_received = comm.bcast(None, root=0)
```

In the worker processes, data1 is received first, followed by data2, and then data3. This corresponds to the order in which they were broadcasted by the master process.

The order of broadcast and receive operations should match to ensure that the worker processes receive the data sets correctly. If the order of broadcast and receive operations needs to be different, you would need to adjust the code accordingly. For example, you could modify the master process to broadcast the data sets in a different order, or you could use tags to distinguish between different sets of data.



You

/apps/intel/oneapi/intelpython/latest/lib/python3.7/site-packages/numpy/polynomial/legendre.py: 1455: RuntimeWarning: divide by zero encountered in true_divide scl = 1./np.sqrt(2*np.arange(n) + 1)

/apps/intel/oneapi/intelpython/latest/lib/python3.7/site-packages/numpy/polynomial/legendre.py: 1581: RuntimeWarning: divide by zero encountered in true_divide w = 1/(fm * df)

ChatGPT

ChatGPT

These warnings are coming from the NumPy library, specifically from the numpy.polynomial.legendre module. They indicate that during the computation of Legendre polynomials, a division by zero was encountered.

The warning RuntimeWarning: divide by zero encountered in true_divide suggests that there is division by zero occurring in the computation. This often happens when trying to calculate the Legendre polynomials or associated values with an inappropriate input.

To handle these warnings, you can consider the following steps:

Check your input: Ensure that the input values you are passing to the Legendre polynomial functions are appropriate. Division by zero can occur when using invalid input values, such as when the degree of the polynomial or the range of values is not suitable. Handle special cases: If you expect division by zero to occur in certain cases and you want to handle it gracefully, you can use try-except blocks to catch the warning and handle it appropriately in your code.

Suppress the warning (if appropriate): If you are confident that the division by zero is expected behavior in your specific use case and you do not want to see the warning, you

can suppress it using the warnings module in Python. However, be cautious when suppressing warnings, as they can provide valuable information about potential issues in your code.

Here's an example of how you can suppress the warning:

```
python
Copy code
import numpy as np
import warnings

# Suppress the warning
warnings.filterwarnings("ignore", category=RuntimeWarning)

# Your code that produces the warning
# For example:
# result = np.polynomial.legendre.leggauss(10)
```

Remember to handle warnings appropriately based on your specific use case and whether the warnings indicate potential issues in your code.



You

How do I pair a worker number with the data returned to the main?

ChatGPT

ChatGPT

To pair a worker number with the data returned to the main process, you can use a dictionary or a list to store the data along with the corresponding worker number. Here's a general approach:

Worker Processes:

- Each worker process computes some data.
- After computation, each worker process sends its data along with its worker number (rank) to the main process.

Main Process:

- Collects the data and worker numbers from all worker processes.
- Stores the data and worker numbers in a suitable data structure (such as a dictionary or a list).

Here's an example implementation using a dictionary to pair worker numbers with the data:

```
python
Copy code
from mpi4py import MPI
comm = MPI.COMM_WORLD
rank = comm.Get_rank()
size = comm.Get_size()
# Worker processes
if rank != 0:
# Simulated data computed by each worker process
data = rank * 10 # Example: Worker number multiplied by 10
 # Send data and worker number to main process
 comm.send((rank, data), dest=0)
# Main process
else:
 # Dictionary to store data paired with worker numbers
data_pairs = {}
 # Receive data from each worker process and store it in the dictionary
 for worker_rank in range(1, size):
 worker_data = comm.recv(source=worker_rank)
 worker_number, worker_data = worker_data
 data_pairs[worker_number] = worker_data
 # Print the paired data
 print("Data paired with worker numbers:")
 for worker_number, worker_data in data_pairs.items():
 print(f"Worker {worker_number}: {worker_data}")
```

In this example:

- Each worker process computes some data (in this case, a simulated data value based on its rank).
- After computation, each worker process sends its data along with its rank to the main process using comm.send.
- The main process receives the data from each worker process using comm.recv and stores it in a dictionary, pairing it with the corresponding worker number (rank).
- Finally, the main process prints the paired data.

You can modify this example to suit your specific use case and data types. Depending on your requirements, you might choose a different data structure or communication method.



You

how do I sort a list of tuples (paired worker number with return value) by the worker number ChatGPT

ChatGPT

To sort a list of tuples containing paired worker numbers with their return values by the worker number, you can use the <code>sorted()</code> function in Python. You need to specify the key by which you want to sort the list. In this case, you want to sort by the first element of each tuple, which represents the worker number.

Here's how you can do it:

```
python
Copy code
# List of tuples containing paired worker numbers with return values
paired_data = [(2, 20), (1, 10), (3, 30)]
# Sort the list of tuples by the worker number (first element of each tuple)
sorted_paired_data = sorted(paired_data, key=lambda x: x[0])
# Print the sorted list
print("Sorted paired data by worker number:")
for worker_number, return_value in sorted_paired_data:
    print(f"Worker {worker_number}: {return_value}")
```

Output:

yaml

Copy code

Sorted paired data by worker number:

Worker 1: 10 Worker 2: 20 Worker 3: 30

In this example:

- sorted() function sorts the list of tuples (paired_data) based on the worker number (first element of each tuple).
- The key parameter specifies the function to be called on each element for sorting. Here, we use a lambda function to access the first element (x[0]) of each tuple.
- The sorted list (sorted_paired_data) is printed, displaying the worker numbers and their corresponding return values in sorted order.

```
You
## --- Worker ---
else:
  print(f"Process {rank} starts") #Displays worker ID
  # Receive Data
  N local = comm.bcast(None, root=0)
  x local = comm.bcast(None, root=0)
  w_local = comm.bcast(None, root=0)
  # Takes 3 arrays and isolates data list specific to this worker
  data_local=[N_local[rank-1],x_local[rank-1],w_local[rank-1]]
  # Takes Data and separates list into 3 arrays for weights, x, and N
  w_worker = data_local[2]
  x worker = data local[1]
  N_worker = data_local[0]
  # Performs Integration
  sum = 0.0
  for i in range (0,N worker):
     sum = sum + w worker[i]*f(x worker[i])
  result=sum
  # Send Results to Main
```

ChatGPT

ChatGPT

The code you provided seems to be part of an MPI program where worker processes are performing some computation and then sending the results back to the main process. Let's address the issue you mentioned, where worker 1 returns 0:

comm.send((rank, result), dest=0 Why does this return 0 for worker 1?

The likely reason why worker 1 returns 0 could be due to an indexing issue in your code when accessing the local data ($N_{local}, x_{local}, w_{local}$). In Python, indexing starts from 0, so accessing $N_{local}[rank-1], x_{local}[rank-1]$, and $w_{local}[rank-1]$ might be incorrect, especially if rank is 1-based.

Assuming N_local, x_local, and w_local are lists of data broadcasted from the main process, and rank is the worker process's rank, you should access the data using rank directly, without subtracting 1.

Here's how you should modify the code:

```
python
Copy code
else:
print(f"Process {rank} starts") # Display worker ID
# Receive Data
 N_local = comm.bcast(None, root=0)
 x_{local} = comm.bcast(None, root=0)
 w_local = comm.bcast(None, root=0)
# Access local data directly using rank
N_{worker} = N_{local}[rank]
x_{worker} = x_{local}[rank]
 w_worker = w_local[rank]
# Performs Integration
 result = 0.0
 for i in range(N_worker):
 result += w_worker[i] * f(x_worker[i])
 # Send Results to Main
 comm.send((rank, result), dest=0)
```

In this modified version:

• We access N local, x local, and w local using the worker process's rank directly.

- The loop for integration starts from 0 up to N_worker 1, where N_worker is the number of integration points for the specific worker process.
- Each worker process sends its rank along with the computed result to the main process.



You

Here is the master process Last login: Thu Mar 28 12:37:49 on ttys000

(base) james@Jamess-MacBook-Pro-2 ~ % ssh ayn868@arc.utsa.edu

0

@ WARNING: REMOTE HOST IDENTIFICATION HAS CHANGED! @

0

IT IS POSSIBLE THAT SOMEONE IS DOING SOMETHING NASTY!

Someone could be eavesdropping on you right now (man-in-the-middle attack)!

It is also possible that a host key has just been changed.

The fingerprint for the ED25519 key sent by the remote host is

SHA256:fDPYc5IZ38NmQ2lcyk2gmytzbxsnAOKBmt/H9anJtjo.

Please contact your system administrator.

Add correct host key in /Users/james/.ssh/known_hosts to get rid of this message.

Offending ECDSA key in /Users/james/.ssh/known_hosts:3

Host key for arc.utsa.edu has changed and you have requested strict checking.

Host key verification failed.

(base) james@Jamess-MacBook-Pro-2 ~ % ssh ayn868@arc.utsa.edu

(ayn868@arc.utsa.edu) Password:

(ayn868@arc.utsa.edu) Duo two-factor login for ayn868

Enter a passcode or select one of the following options:

- 1. Duo Push to XXX-XXX-8036
- 2. Phone call to XXX-XXX-8036
- 3. SMS passcodes to XXX-XXX-8036

Passcode or option (1-3): 1

Success. Logging you in...

Last login: Thu Mar 28 12:38:26 2024 from 129.222.77.154

Welcome to Arc!

Do *not* execute code directly on the login nodes. SBATCH should be used to submit jobs to the Slurm job scheduling engine.

If you need to run an interactive job on a compute node, use the following syntax to connect to one:

srun -p compute1 -n 1 -t 00:05:00 --cpus-per-task=1 --pty bash

Starting in work directory /work/ayn868

[login001: ayn868]\$ Is

HDFView ME5773SP24 sec3 source code sources

[login001: ayn868]\$ cd ME5773SP24/

[login001: ME5773SP24]\$ cd assignment7

[login001: assignment7]\$ ls

gauleg.py job_script.slurm myjob.e364679 myjob.o364679 part2.py Homework7.py myjob.e364675 myjob.o364675 part1.py part2.txt

int-trap.py myjob.e364678 myjob.o364678 part1.txt

[login001: assignment7]\$ ml vim

The vim module version 8.1 is loaded.

[login001: assignment7]\$ vim job_script.slurm

[login001: assignment7]\$ vim part1.py

[login001: assignment7]\$ vim

[login001: assignment7]\$ Is

gauleg.py int-trap.py myjob.e364675 myjob.e364679 myjob.o364678 part1.py part2.py Homework7.py job_script.slurm myjob.e364678 myjob.o364675 myjob.o364679 part1.txt part2.txt

[login001: assignment7]\$ vim Homework7.py

[login001: assignment7]\$ Is

gauleg.py int-trap.py myjob.e364675 myjob.e364679 myjob.o364678 part1.py part2.py Homework7.py job_script.slurm myjob.e364678 myjob.o364675 myjob.o364679 part1.txt part2.txt

[login001: assignment7]\$ client loop: send disconnect: Broken pipe

(base) james@Jamess-MacBook-Pro-2 ~ % ssh ayn868@arc.utsa.edu

@ WARNING: REMOTE HOST IDENTIFICATION HAS CHANGED! @

IT IS POSSIBLE THAT SOMEONE IS DOING SOMETHING NASTY!

Someone could be eavesdropping on you right now (man-in-the-middle attack)!

It is also possible that a host key has just been changed.

The fingerprint for the ED25519 key sent by the remote host is

SHA256:fDPYc5IZ38NmQ2Icyk2qmytzbxsnAOKBmt/H9anJtjo.

Please contact your system administrator.

Add correct host key in /Users/james/.ssh/known_hosts to get rid of this message.

Offending ECDSA key in /Users/james/.ssh/known hosts:3

Host key for arc.utsa.edu has changed and you have requested strict checking.

Host key verification failed.

(base) james@Jamess-MacBook-Pro-2 ~ % ssh ayn868@arc.utsa.edu

@ WARNING: REMOTE HOST IDENTIFICATION HAS CHANGED! @

IT IS POSSIBLE THAT SOMEONE IS DOING SOMETHING NASTY!

Someone could be eavesdropping on you right now (man-in-the-middle attack)!

It is also possible that a host key has just been changed.

The fingerprint for the ED25519 key sent by the remote host is

SHA256:fDPYc5IZ38NmQ2Icyk2qmytzbxsnAOKBmt/H9anJtjo.

Please contact your system administrator.

Add correct host key in /Users/james/.ssh/known hosts to get rid of this message.

Offending ECDSA key in /Users/james/.ssh/known hosts:3

Host key for arc.utsa.edu has changed and you have requested strict checking.

Host key verification failed.

(base) james@Jamess-MacBook-Pro-2 ~ % ssh ayn868@arc.utsa.edu

@ WARNING: REMOTE HOST IDENTIFICATION HAS CHANGED! @

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Add correct host key in /Users/james/.ssh/known hosts to get rid of this message.

Offending ECDSA key in /Users/james/.ssh/known hosts:3

Host key for arc.utsa.edu has changed and you have requested strict checking.

Host key verification failed.

(base) james@Jamess-MacBook-Pro-2 ~ % ssh ayn868@arc.utsa.edu

0

@ WARNING: REMOTE HOST IDENTIFICATION HAS CHANGED! @

0

IT IS POSSIBLE THAT SOMEONE IS DOING SOMETHING NASTY!

Someone could be eavesdropping on you right now (man-in-the-middle attack)!

It is also possible that a host key has just been changed.

The fingerprint for the ED25519 key sent by the remote host is

SHA256:fDPYc5IZ38NmQ2Icyk2qmytzbxsnAOKBmt/H9anJtjo.

Please contact your system administrator.

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Offending ECDSA key in /Users/james/.ssh/known_hosts:3

Host key for arc.utsa.edu has changed and you have requested strict checking.

Host key verification failed.

(base) james@Jamess-MacBook-Pro-2 ~ % ssh ayn868@arc.utsa.edu

@ WARNING: REMOTE HOST IDENTIFICATION HAS CHANGED! @

0

IT IS POSSIBLE THAT SOMEONE IS DOING SOMETHING NASTY!

Someone could be eavesdropping on you right now (man-in-the-middle attack)!

It is also possible that a host key has just been changed.

The fingerprint for the ED25519 key sent by the remote host is

SHA256:fDPYc5IZ38NmQ2Icyk2qmytzbxsnAOKBmt/H9anJtjo.

Please contact your system administrator.

Add correct host key in /Users/james/.ssh/known hosts to get rid of this message.

Offending ECDSA key in /Users/james/.ssh/known hosts:3

Host key for arc.utsa.edu has changed and you have requested strict checking.

Host key verification failed.

(base) james@Jamess-MacBook-Pro-2 ~ % ssh ayn868@arc.utsa.edu

(avn868@arc.utsa.edu) Password:

(ayn868@arc.utsa.edu) Duo two-factor login for ayn868

Enter a passcode or select one of the following options:

- 1. Duo Push to XXX-XXX-8036
- 2. Phone call to XXX-XXX-8036
- 3. SMS passcodes to XXX-XXX-8036

Passcode or option (1-3): 1 Success. Logging you in...

Last login: Fri Mar 29 13:42:18 2024 from 129.222.77.154

Welcome to Arc!

Do *not* execute code directly on the login nodes. SBATCH should be used to submit jobs to the Slurm job scheduling engine.

If you need to run an interactive job on a compute node, use the following syntax to connect to one:

srun -p compute1 -n 1 -t 00:05:00 --cpus-per-task=1 --pty bash

Starting in work directory /work/ayn868

[login001: ayn868]\$ ls

HDFView ME5773SP24 sec3 source_code sources

[login001: ayn868]\$ cd ME5773SP24/

[login001: ME5773SP24]\$ Is

assignment3 assignment4 assignment5 assignment6 assignment7 README.md

[login001: ME5773SP24]\$ cd assignment7

[login001: assignment7]\$ Is

gauleg.py int-trap.py myjob.e364675 myjob.e364679 myjob.o364678 part1.py part2.py Homework7.py job_script.slurm myjob.e364678 myjob.o364675 myjob.o364679 part1.txt part2.txt

[login001: assignment7]\$ vim Homework7.py

-bash: vim: command not found [login001: assignment7]\$ ml vim

The vim module version 8.1 is loaded.

[login001: assignment7]\$ vim Homework7.py

[login001: assignment7]\$ Is

gauleg.py int-trap.py myjob.e364675 myjob.e364679 myjob.o364678 part1.py part2.py Homework7.py job_script.slurm myjob.e364678 myjob.o364675 myjob.o364679 part1.txt part2.txt

[login001: assignment7]\$ vim job_script.slurm

[login001: assignment7]\$ cp job_script.slurm James_script.slurm

[login001: assignment7]\$ vim James_script.slurm

[login001: assignment7]\$ Is

gauleg.py int-trap.py job_script.slurm myjob.e364678 myjob.o364675 myjob.o364679 part1.txt part2.txt

Homework7.py James script.slurm myjob.e364675 myjob.e364679 myjob.o364678

part1.py part2.py

[login001: assignment7]\$ sbatch James script.slurm

Submitted batch job 365235

[login001: assignment7]\$ squeue -u ayn868

JOBID PARTITION NAME USER ST TIME NODES NODELIST(REASON)

[login001: assignment7]\$ Is

gauleg.py Jamesjob.e365235 job_script.slurm myjob.e364679 myjob.o364679 part2.py Homework7.py Jamesjob.o365235 myjob.e364675 myjob.o364675 part1.py part2.txt int-trap.py James_script.slurm myjob.e364678 myjob.o364678 part1.txt

[login001: assignment7]\$ vim Jamesjob.e365235

[login001: assignment7]\$ rm Jamesjob. Jamesjob.e365235 Jamesjob.o365235 [login001: assignment7]\$ rm Jamesjob.*

Do you really want to run this delete command?

rm Jamesjob.e365235 Jamesjob.o365235

(y/n)? y

[login001: assignment7]\$ Is

gauleg.py job script.slurm myjob.e364678 myjob.o364675 myjob.o364679 int-trap.py part1.txt part2.txt Homework7.py James script.slurm myjob.e364675 myjob.e364679 myjob.o364678 part1.pv part2.py [login001: assignment7]\$ vim James script.slurm [login001: assignment7]\$ pwd /work/ayn868/ME5773SP24/assignment7 [login001: assignment7]\$ vim James script.slurm [login001: assignment7]\$ Is job_script.slurm myjob.e364678 myjob.o364675 myjob.o364679 gauleg.py int-trap.py part1.txt part2.txt Homework7.py James_script.slurm myjob.e364675 myjob.e364679 myjob.o364678 part1.py part2.py [login001: assignment7]\$ vim Homework7.py [login001: assignment7]\$ sbatch James script.slurm Submitted batch job 365236 [login001: assignment7]\$ squeue -u ayn868 **JOBID PARTITION** NAME USER ST TIME NODES NODELIST(REASON) 365236 compute1 Jamesjob ayn868 R 0:03 1 c022 [login001: assignment7]\$ squeue -u ayn868 JOBID PARTITION NAME **USER ST** TIME NODES NODELIST(REASON) 365236 compute1 Jamesjob ayn868 CG 0:05 1 [login001: assignment7]\$ squeue -u ayn868 JOBID PARTITION NAME **USER ST** TIME NODES NODELIST(REASON) [login001: assignment7]\$ ls gauleg.py Jamesjob.e365236 job script.slurm myjob.e364679 myjob.o364679 part2.py Homework7.py Jamesjob.o365236 myjob.e364675 myjob.o364675 part1.py part2.txt int-trap.py James script.slurm myjob.e364678 myjob.o364678 part1.txt [login001: assignment7]\$ vim Jamesjob.e365236 [login001: assignment7]\$ vim Jamesjob.o365236 [login001: assignment7]\$ Is Jamesjob.e365236 job script.slurm myjob.e364679 myjob.o364679 part2.py gauleg.py Homework7.py Jamesjob.o365236 myjob.e364675 myjob.o364675 part1.py part2.txt int-trap.py James script.slurm myjob.e364678 myjob.o364678 part1.txt [login001: assignment7]\$ rm Jamesjob.* Do you really want to run this delete command? rm Jamesjob.e365236 Jamesjob.o365236 (v/n)? v [login001: assignment7]\$ ls job script.slurm myjob.e364678 myjob.o364675 myjob.o364679 int-trap.py gauleg.py part1.txt part2.txt Homework7.py James_script.slurm myjob.e364675 myjob.e364679 myjob.o364678

part1.pv

part2.pv

[login001: assignment7]\$ vim Homework7.py

[login001: assignment7]\$ vi Homework7.py [login001: assignment7]\$ vim Homework7.py

[login001: assignment7]\$ sbatch James_script.slurm

Submitted batch job 365237 [login001: assignment7]\$ Is

gauleg.py Jamesjob.e365237 job_script.slurm myjob.e364679 myjob.o364679 part2.py Homework7.py Jamesjob.o365237 myjob.e364675 myjob.o364675 part1.py part2.txt

int-trap.py James_script.slurm myjob.e364678 myjob.o364678 part1.txt

[login001: assignment7]\$ vim Jamesjob.e365237

Press ENTER or type command to continue

[login001: assignment7]\$ vim Jamesjob.o365237

[login001: assignment7]\$ rm Jamesjob.*

Do you really want to run this delete command?

rm Jamesjob.e365237 Jamesjob.o365237

(y/n)? y

[login001: assignment7]\$ ls

gauleg.py int-trap.py job_script.slurm myjob.e364678 myjob.o364675 myjob.o364679 part1.txt part2.txt

Homework7.py James_script.slurm myjob.e364675 myjob.e364679 myjob.o364678

part1.py part2.py

[login001: assignment7]\$ rm myjob.*

Do you really want to run this delete command?

rm myjob.e364675 myjob.e364678 myjob.e364679 myjob.o364675 myjob.o364678

myjob.o364679

(y/n)? y

[login001: assignment7]\$ Is

gauleg.py Homework7.py int-trap.py James_script.slurm job_script.slurm part1.py part1.txt part2.py part2.txt

[login001: assignment7]\$ vim Homework7.py

[login001: assignment7]\$ sbatch James script.slurm

Submitted batch job 365239

[login001: assignment7]\$ squeue -u ayn868

JOBID PARTITION NAME USER ST TIME NODES NODELIST(REASON)

[login001: assignment7]\$ [login001: assignment7]\$ Is

gauleg.py int-trap.py Jamesjob.o365239 job_script.slurm part1.txt part2.txt

Homework7.py Jamesjob.e365239 James_script.slurm part1.py part2.py

[login001: assignment7]\$ vim Jamesjob.e365239

[login001: assignment7]\$ vim Jamesjob.o365239

[login001: assignment7]\$ Is

gauleg.py int-trap.py Jamesjob.o365239 job_script.slurm part1.txt part2.txt Homework7.py Jamesjob.e365239 James_script.slurm part1.py part2.py

```
[login001: assignment7]$ vim Homework7.py
[login001: assignment7]$ vi Homework7.py
  x2 = 1
  # Weights(w) and Integration Points(x)
  w=[]
  x=[]
  for i in N:
    x temp,w temp = np.polynomial.legendre.leggauss(i)
    w.append(w temp)
    x.append(x_temp)
  ### Compute the Quadrature ###
  # Send Data to Workers
  comm.bcast(N, root=0)
  comm.bcast(x, root=0)
  comm.bcast(w, root=0)
  # Collect the results from each worker
  int_results=[]
  for i in range(1,size):
     temp=comm.recv(source=i)
     int_results.append(temp)
  # Compute the table
    # Sort Reults
  SortedResults = sorted(int_results, key=lambda x: x[0])
  print(SortedResults)
```

ChatGPT

ChatGPT

It seems like you're trying to submit jobs using Slurm on a remote machine, specifically the Arc cluster at UTSA. However, you seem to encounter issues with SSH host key verification. This typically happens when the remote host's key changes, which can occur due to various reasons such as system updates or changes in the server configuration.

To resolve this issue, you can do the following:

Update known_hosts file: Remove the offending key from your known_hosts file. You can do this manually or use the ssh-keygen command with the -R option followed by the hostname to remove the entry.

bash

Copy code

```
ssh-keygen -R arc.utsa.edu
```

Replace arc.utsa.edu with the hostname of the remote server.

Retry SSH connection: After removing the offending key, try to SSH into the remote machine again. You may be prompted to confirm the authenticity of the host, and upon confirmation, the new host key will be added to your known hosts file.

Ensure correct hostname: Double-check that you're connecting to the correct hostname or IP address of the remote server.

Contact system administrator: If the issue persists or if you suspect a security concern, it's a good idea to contact the system administrator of the remote server for assistance.

Once you've resolved the SSH host key verification issue, you should be able to submit your Slurm jobs without encountering further SSH errors.

```
You
## --- Master ---
if rank == 0:
  # Generate Quadrature Number
  N = list(range(1,21))
  # Integration Bounds
  x1 = -1
  x2 = 1
  # Weights(w) and Integration Points(x)
  w=[]
  x=[]
  for i in N:
    x_temp,w_temp = np.polynomial.legendre.leggauss(i)
    w.append(w temp)
    x.append(x_temp)
  ### Compute the Quadrature ###
```

```
# Send Data to Workers
comm.bcast(N, root=0)
comm.bcast(x, root=0)

# Collect the results from each worker
int_results=[]
for i in range(1,size):
    temp=comm.recv(source=i)
    int_results.append(temp)

# Compute the table

# Sort Reults
SortedResults = sorted(int_results, key=lambda x: x[0])
print(SortedResults)
```

Here is the master process

ChatGPT

ChatGPT

The code snippet you provided represents the master process in your MPI program. Here's a breakdown of what it does:

Generate Quadrature Number: Creates a list ${\tt N}$ containing quadrature numbers from 1 to 20.

Integration Bounds: Sets the integration bounds x1 and x2.

Compute Quadrature Points and Weights: Computes the Legendre quadrature points and weights for each quadrature number in $_{\rm N}$.

Send Data to Workers: Broadcasts the quadrature numbers (\mathbb{N}), quadrature points (\mathbb{N}), and quadrature weights (\mathbb{N}) to all worker processes.

Collect Results from Workers: Receives the integration results from each worker process and appends them to the <code>int_results</code> list.

Compute the Table:

- Sorts the integration results based on the quadrature number.
- Prints the sorted results.

Overall, this code orchestrates the distribution of data to worker processes, collects the results, and prints them in sorted order based on the quadrature number. Make sure the worker processes are correctly handling the received data and sending back the integration results to

the master process. Additionally, ensure that the worker processes are properly synchronized with the master process to avoid deadlocks or race conditions.



You

Given the master and worker processes that I have provided, why is worker returning 0?

ChatGPT

ChatGPT

The issue with the worker process returning 0 might be due to incorrect indexing when accessing the broadcasted data. Let's analyze the worker process code snippet you provided:

```
python
Copy code
else:
 print(f"Process {rank} starts") #Displays worker ID
# Receive Data
 N_local = comm.bcast(None, root=0)
 x_{local} = comm.bcast(None, root=0)
w_local = comm.bcast(None, root=0)
 # Takes 3 arrays and isolates data list specific to this worker
 data_local=[N_local[rank-1], x_local[rank-1], w_local[rank-1]]
 # Takes Data and separates list into 3 arrays for weights, x, and N
 w_worker = data_local[2]
 x_worker = data_local[1]
 N_{worker} = data_{local}[0]
# Performs Integration
 sum = 0.0
 for i in range (0, N_worker):
 sum = sum + w_worker[i]*f(x_worker[i])
 result=sum
 # Send Results to Main
comm.send((rank, result), dest=0)
```

The issue lies in how you are isolating the data specific to each worker. You are using the worker's rank to index the broadcasted arrays ($N_local, x_local, w_local$). However, in

Python, the rank of the processes in MPI starts from 0, not 1. So, when you are trying to access N_local[rank-1], it's accessing the 0th element for the worker with rank 1, which corresponds to the first element of the broadcasted arrays.

To fix this issue, you should directly use the received broadcasted arrays without subtracting 1 from the rank. Here's the corrected version of the code:

```
python
Copy code
else:
print(f"Process {rank} starts") #Displays worker ID
 # Receive Data
 N_local = comm.bcast(None, root=0)
 x_{local} = comm.bcast(None, root=0)
w_local = comm.bcast(None, root=0)
\# Takes Data and separates list into 3 arrays for weights, x, and N
 w_worker = w_local
x_worker = x_local
 N_{worker} = N_{local}
# Performs Integration
sum = 0.0
 for i in range (0, N_worker):
 sum = sum + w_worker[i] * f(x_worker[i])
 result = sum
 # Send Results to Main
comm.send((rank, result), dest=0)
```

This way, each worker will correctly use the received broadcasted data without any indexing issues, and it should return the correct results to the master process.



You

This will not work

ChatGPT

ChatGPT

I apologize for the oversight. Let's correct the code.

Since you're broadcasting the entire lists N, x, and w to all workers, each worker should receive the entire list, not just a part of it. The indexing rank - 1 is not appropriate here because it's intended for distributing data among workers, not for accessing the received broadcasted data.

Here's the corrected worker process code:

```
python
Copy code
else:
    print(f"Process {rank} starts") # Displays worker ID

# Receive Data
    N_local = comm.bcast(None, root=0)
    x_local = comm.bcast(None, root=0)
    w_local = comm.bcast(None, root=0)

# Performs Integration
    sum = 0.0
    for i in range(0, len(N_local)):
    sum += w_local[i] * f(x_local[i])
    result = sum

# Send Results to Main
    comm.send((rank, result), dest=0)
```

In this corrected version, each worker receives the entire broadcasted lists \mathbb{N} , \times , and \mathbb{W} , and then performs the integration using those lists. There's no need for indexing or isolating specific data for each worker since all workers receive the same data broadcasted from the master process.



You

how do I print a table?

ChatGPT

ChatGPT

To print a table in Python, you can use various libraries such as tabulate, prettytable, or simply format the output using string formatting. Here's an example using the tabulate library:

```
python
```

```
Copy code
```

```
from tabulate import tabulate
```

```
# Assuming SortedResults contains the results in the format (quadrature_no,
integration_result)
headers = ["Quadrature no.", "Integration Result", "Percent error", "Run time
(s)"]

# Compute percent error and run time (assuming they are available)
table_data = []
for quadrature, result, _, _ in SortedResults:
   percent_error = compute_percent_error(result) # You need to implement
compute_percent_error function
   run_time = compute_run_time(quadrature) # You need to implement compute_run_time
function
   table_data.append([quadrature, result, percent_error, run_time])

# Print the table
print(tabulate(table_data, headers=headers, floatfmt=".6f")) # Adjust floatfmt as
needed
```

In this example, <code>sortedResults</code> is assumed to be a list of tuples containing the quadrature number, integration result, percent error, and run time. You would replace the placeholders <code>compute_percent_error</code> and <code>compute_run_time</code> with your actual functions for computing these values. Adjust the headers and data accordingly based on your specific requirements.



You

How do you time something in python

ChatGPT

ChatGPT

In Python, you can time the execution of code using the time module or the timeit module.

Here's how you can do it with each:

Using the time module:

python

Copy code

```
import time
start_time = time.time()
# Code to be timed
# For example:
# result = some_function()
end_time = time.time()
execution_time = end_time - start_time
print("Execution time:", execution_time, "seconds")
      Using the timeit module:
python
Copy code
import timeit
# Code to be timed, wrapped in a function
# For example:
# def timed_function():
# return some_function()
execution_time = timeit.timeit("timed_function()", setup="from __main__ import
timed_function", number=1)
print("Execution time:", execution_time, "seconds")
```

Both methods measure the time taken to execute the code between the <code>start_time</code> and <code>end_time</code> markers. The <code>timeit</code> module is more suitable for measuring shorter execution times or for benchmarking purposes, while the <code>time</code> module is more general-purpose.



You

Traceback (most recent call last):

File "Homework7.py", line 4, in <module>
from tabulate import tabulate

ModuleNotFoundError: No module named 'tabulate'

ChatGPT

ChatGPT

It seems that you are trying to use the tabulate module, but it is not installed in your Python environment. You need to install the tabulate module before using it. You can install it using pip, which is the package installer for Python.

You can install tabulate by running the following command in your terminal or command prompt:

Copy code

pip install tabulate

Once installed, you can import and use the tabulate module in your Python script without any errors.



You

can I pip install into a slurm file?

ChatGPT

ChatGPT

No, you cannot directly run pip install commands within a Slurm script (.slurm file). Slurm scripts are meant for submitting and managing job execution on a cluster's compute nodes, and they are typically used to specify job parameters such as resources, runtime, and the command to be executed.

To install packages using pip, you typically need access to the command line interface of the system where you want to install the package. This means you would need to run pip install commands outside of the Slurm script, either directly on the command line or in a separate script.

If you need certain Python packages installed on the compute nodes where your Slurm jobs will run, you should communicate with your system administrator or cluster manager to have those packages installed system-wide or within a specific Python environment accessible to your jobs. Once the packages are available in the environment, you can use them in your Slurm scripts without needing to install them again within the script.

You

what are my other options for making a table?

ChatGPT

ChatGPT

python

If you're unable to use the tabulate module, you have a few other options for creating tables in Python:

Using format Method: You can use the str.format() method to format strings and create a simple table-like structure. While this method requires more manual formatting, it can be effective for basic tabular output.

```
python
    Copy code

data = [("Column1", "Column2", "Column3"),
    (value1, value2, value3),
    (value4, value5, value6)]

for row in data:
    print("{:<15} {:<15} {:<15}".format(*row))</pre>
```

Using String Manipulation: You can manually manipulate strings to create a table-like structure. This approach gives you full control over the formatting but can be more cumbersome for larger tables.

```
Copy code

data = [("Column1", "Column2", "Column3"),
  (value1, value2, value3),
  (value4, value5, value6)]

for row in data:
  print(" | ".join(str(item).ljust(15) for item in row))
```

Using Pandas Library: If you have Pandas installed, you can use it to create tables with more flexibility and customization options.

python

Copy code

```
import pandas as pd

data = [("Column1", "Column2", "Column3"),
  (value1, value2, value3),
  (value4, value5, value6)]

df = pd.DataFrame(data[1:], columns=data[0])
print(df)
```

Each of these methods has its own advantages and limitations, so you can choose the one that best fits your requirements and constraints.



You

how an I add a tuple to a list of tuples?

ChatGPT

ChatGPT

You can add a tuple to a list of tuples using the append() method or by concatenating lists.

Here's how you can do it:

Using append() method:

```
python
Copy code
# Existing list of tuples
list_of_tuples = [(1, 'a'), (2, 'b'), (3, 'c')]
# Tuple to add
new_tuple = (4, 'd')
# Adding the tuple to the list
list_of_tuples.append(new_tuple)
```

```
print(list_of_tuples)
```

Using concatenation:

```
python
Copy code
# Existing list of tuples
list_of_tuples = [(1, 'a'), (2, 'b'), (3, 'c')]
# Tuple to add
new_tuple = (4, 'd')
# Adding the tuple to the list
list_of_tuples += [new_tuple]
print(list_of_tuples)
```

Both methods will produce the same result, adding the tuple (4, 'd') to the list of tuples.



You

can I append to the beginning?

ChatGPT

ChatGPT

Yes, you can append to the beginning of a list in Python using the <code>insert()</code> method. Here's how you can do it with a list of tuples:

```
python
Copy code
# Existing list of tuples
list_of_tuples = [(1, 'a'), (2, 'b'), (3, 'c')]
# Tuple to add
new_tuple = (0, 'z')
# Inserting the tuple at the beginning of the list
list_of_tuples.insert(0, new_tuple)
```

```
print(list_of_tuples)
```

This will insert the tuple (0, |z|) at the beginning of the list, shifting the existing elements to the right.



You

show line numbersi n ivm

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In Vim, you can display line numbers by typing :set number and pressing Enter. To disable line numbers, you can type :set nonumber and press Enter. Alternatively, you can use the shortcut :set nu to enable line numbers and :set nonu to disable them.



You

what is the syntax error in these lines? Traceback (most recent call last):

File "Homework7.py", line 4, in <module>

from tabulate import tabulate

ModuleNotFoundError: No module named 'tabulate'

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The error message indicates that the Python interpreter couldn't find the tabulate module, which suggests that it's not installed in your Python environment. To resolve this error, you need to install the tabulate module using pip.

You can install tabulate using the following command in your terminal or command prompt:

Copy code

pip install tabulate

After installing tabulate, you should be able to import it in your Python script without encountering the ModuleNotFoundError.



You

```
what is the syntax error in these lines? ## --- Worker ---
else:
  # print(f"Process {rank} starts") #Displays worker ID
  StartTime = time.time()
```

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The syntax error in the provided code snippet is the absence of the if statement before the else statement. In Python, the else statement must be preceded by an if statement to form an if-else block. Here's the corrected version of the code:

```
python
Copy code
## --- Worker ---
if rank != 0:
# print(f"Process {rank} starts") #Displays worker ID
StartTime = time.time()
```

With this change, the else block is correctly associated with the preceding if statement, and the code should execute without syntax errors.

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You
```

```
What about now? ## --- Master ---
if rank == 0:
  # Generate Quadrature Number
  N = list(range(1,21))
  # Integration Bounds
  x1 = -1
  x2 = 1
  # Weights(w) and Integration Points(x)
  w=[]
  x=[]
  for i in N:
    x temp,w temp = np.polynomial.legendre.leggauss(i)
    w.append(w temp)
    x.append(x_temp)
```

```
### Compute the Quadrature ###
  # Send Data to Workers
  comm.bcast(N, root=0)
  comm.bcast(x, root=0)
  comm.bcast(w, root=0)
  # Collect the results from each worker
  int_results=[]
  for i in range(1,size):
     temp=comm.recv(source=i)
     int_results.append(temp)
  # Compute the table
     # Sort Reults
  SortedResults = sorted(int_results, key=lambda x: x[0])
  # print(SortedResults)
    # Print Table
  TableHeaders = ["Quadrature No.", "Integration Result", "Percent Error", "Run Time (sec)"]
  SortedResults.insert(0,TableHeaders)
  print(pd.DataFrame(SortedResults[1:], columns=data[0])
## --- Worker ---
else:
  # print(f"Process {rank} starts") #Displays worker ID
  StartTime = time.time()
```

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There are a couple of issues in the provided code snippet:

The import pandas as pd statement is missing. Without it, the code will raise a NameError when trying to use pd. DataFrame.

There's a syntax error in the print statement. The closing parenthesis is missing.

Here's the corrected version of the code:

python

Copy code

```
import pandas as pd # Add this line to import pandas module
## --- Master ---
if rank == 0:
# Generate Ouadrature Number
N = list(range(1,21))
# Integration Bounds
x1 = -1
x2 = 1
# Weights(w) and Integration Points(x)
w=[]
x=[]
 for i in N:
 x_temp, w_temp = np.polynomial.legendre.leggauss(i)
 w.append(w_temp)
x.append(x_temp)
### Compute the Quadrature ###
 # Send Data to Workers
 comm.bcast(N, root=0)
 comm.bcast(x, root=0)
 comm.bcast(w, root=0)
 # Collect the results from each worker
 int_results=[]
 for i in range(1, size):
temp=comm.recv(source=i)
int_results.append(temp)
 # Compute the table
# Sort Results
SortedResults = sorted(int_results, key=lambda x: x[0])
# Print Table
 TableHeaders = ["Quadrature No." , "Integration Result" , "Percent Error" , "Run
Time (sec)"]
SortedResults.insert(0, TableHeaders)
 df = pd.DataFrame(SortedResults[1:], columns=SortedResults[0]) # Corrected this
line
print(df) # Corrected this line
```

```
## --- Worker ---
else:
 # print(f"Process {rank} starts") #Displays worker ID
 StartTime = time.time()
```

With these changes, the code should work correctly, printing the table with the sorted results. Make sure you have the pandas library installed (pip install pandas) to use pd. DataFrame.



You

is there anything wrong with the else: line?

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No, there is nothing inherently wrong with the else: line. However, the else: statement should be followed by an indented block of code that will be executed if the condition in the if statement above it evaluates to False. In your provided code snippet, the else: statement is followed by a comment, which means there is no actual code executed in the else block.

If you intend to execute some code in the else block, you should remove the comment # print(f"Process {rank} starts") or replace it with the actual code you want to execute. Otherwise, if there is no code to execute in the else block, it's perfectly fine to leave it as is.



```
You
```

SyntaxError: invalid syntax File "Homework7.py", line 75 else: ٨ SyntaxError: invalid syntax File "Homework7.py", line 75 else: SyntaxError: invalid syntax File "Homework7.py", line 75 else: SyntaxError: invalid syntax File "Homework7.py", line 75

```
else:
    ٨
SyntaxError: invalid syntax
 File "Homework7.py", line 75
  else:
    ٨
SyntaxError: invalid syntax
 File "Homework7.py", line 75
  else:
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SyntaxError: invalid syntax
 File "Homework7.py", line 75
  else:
    ٨
SyntaxError: invalid syntax
 File "Homework7.py", line 75
  else:
    ٨
SyntaxError: invalid syntax
                                import math
from mpi4py import MPI
import numpy as np
import pandas
import warnings
import time
# This code was written by James Smith to perform Gaussian Integration
# Suppress Leggauss warning about dividing by Zero
warnings.filterwarnings("ignore", category=RuntimeWarning)
# Start the MPI Process
comm = MPI.COMM WORLD
# Determine the Total Number of Tasks
size = comm.Get_size()
# Determine the rank (ID) of this task
rank = comm.Get rank()
# Function to integrate
def f(x):
  return x*math.exp(x)
```

```
N = list(range(1,21))
## --- Master ---
if rank == 0:
  # Generate Quadrature Number
  N = list(range(1,21))
  # Integration Bounds
  x1 = -1
  x2 = 1
  # Weights(w) and Integration Points(x)
  w=[]
  x=[]
  for i in N:
    x_temp,w_temp = np.polynomial.legendre.leggauss(i)
    w.append(w temp)
    x.append(x_temp)
  ### Compute the Quadrature ###
  # Send Data to Workers
  comm.bcast(N, root=0)
  comm.bcast(x, root=0)
  comm.bcast(w, root=0)
  # Collect the results from each worker
  int_results=[]
  for i in range(1,size):
     temp=comm.recv(source=i)
     int_results.append(temp)
  # Compute the table
    # Sort Reults
  SortedResults = sorted(int_results, key=lambda x: x[0])
  # print(SortedResults)
    # Print Table
  TableHeaders = ["Quadrature No.", "Integration Result", "Percent Error", "Run Time (sec)"]
  SortedResults.insert(0,TableHeaders)
  print(pd.DataFrame(SortedResults[1:], columns=data[0])
```

```
## --- Worker ---
else:
  # print(f"Process {rank} starts") #Displays worker ID
  StartTime = time.time()
  # Receive Data
  N local = comm.bcast(None, root=0)
  x local = comm.bcast(None, root=0)
  w_local = comm.bcast(None, root=0)
  # Takes 3 arrays and isolates data list specific to this worker
  data_local=[N_local[rank-1],x_local[rank-1],w_local[rank-1]]
  # Takes Data and separates list into 3 arrays for weights, x, and N
  w worker = data local[2]
  x_worker = data_local[1]
  N_worker = data_local[0]
  # Performs Integration
  sum = 0.0
  for i in range (0,N worker):
    sum = sum + w_worker[i]*f(x_worker[i])
  result=sum
  # Calculate Work Time
  Work_Time = time.time() - StartTime
  # Calculate % Error
  PercentError =100*(result - 0.735758882342885)/0.735758882342885
  # Send Results to Main
  comm.send((rank, result, PercentError, Work Time), dest=0)
  # Debugging, Delete if code is working I guess
  #if rank==1:
  # print(N_worker,x_worker,w_worker)
  # print(sum)
```

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The issue seems to be with the missing closing parenthesis in the print statement just before the else: block. You should add a closing parenthesis) at the end of that line to resolve the syntax error.

Here's the corrected version:

python

Copy code

print(pd.DataFrame(SortedResults[1:], columns=data[0]))

After fixing this, the if-else block structure should be syntactically correct.



You

fuck this

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I'm sorry to hear that you're frustrated. If you have any specific questions or need further assistance, please feel free to ask. I'm here to help.



You

is there anything else wrong with the code?

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There are a few syntax errors and missing imports in the code you provided. Here are the issues I noticed:

In the master process section, you are missing a closing parenthesis ")" for the print (pd.DataFrame (SortedResults[1:], columns=data[0]) line.

In the worker process section, there is a missing closing parenthesis ")" for the comm.send((rank, result, PercentError, Work Time), dest=0) line.

You have a syntax error in the worker process section where the <code>else</code>: statement is not properly aligned. It should be aligned with the <code>if</code> statement.

You are missing an import statement for the pandas library.

The comment "Sort Reults" contains a typo. It should be "Sort Results".

After fixing these issues, your code should run without syntax errors. If you encounter any further issues or need assistance with specific parts of the code, please let me know!



You

thanks

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You're welcome! If you have any more questions or need further assistance, feel free to ask.

Good luck with your code!



Υοι

how do I save this conversation into a pdf?

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To save this conversation as a PDF, you can follow these steps:

Select All Text: Scroll through the conversation and select all the text. You can do this by clicking and dragging your mouse cursor over the entire conversation or by using keyboard shortcuts like Ctrl+A (Cmd+A on Mac) to select all.

Copy the Text: Once all the text is selected, right-click and choose the "Copy" option from the context menu. Alternatively, you can use the keyboard shortcut Ctrl+C (Cmd+C on Mac) to copy the text.

Open a Text Editor or Word Processor: Open a text editor like Notepad (Windows), TextEdit (Mac), or any word processor like Microsoft Word or Google Docs.

Paste the Text: Click inside the text editor or word processor and paste the copied text into a new document. You can right-click and choose the "Paste" option from the context menu or use the keyboard shortcut Ctrl+V (Cmd+V on Mac) to paste the text.

Format the Text (Optional): You can format the text as needed, adjust font sizes, styles, or add headings to organize the conversation.

Save the Document: Once you have pasted and formatted the text, you can save the document. In most text editors or word processors, you can go to the "File" menu and choose "Save As" to save the document. Choose the PDF format when saving.

Name and Choose Save Location: Give your PDF document a name and choose where you want to save it on your computer.

Save the Document: Click the "Save" button to save the document as a PDF.

Following these steps will allow you to save the conversation as a PDF document on your computer. If you encounter any issues or need further assistance, feel free to ask!

ChatGPT can make mistakes. Consider checking important information.

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