

Problem 2 a:

```
C:\Users\lab\Dropbox\Computational Physics\Jinesh_HW3>python -i prob2a.py
>>> Trapezoidal_Rule(2)
0.3252319078064746
>>> Trapezoidal_Rule(4)
0.5122828507233315
>>> Trapezoidal_Rule()
Traceback (most recent call last):
  File "<stdin>", line 1, in <module>
TypeError: Trapezoidal_Rule() takes exactly 1 argument (0 given)
>>> Trapezoidal_Rule(8)
0.40299744847824825
>>> Trapezoidal_Rule(16)
0.43010336929474696
>>> Trapezoidal_Rule(32)
0.4484146657874698
>>> Trapezoidal_Rule(64)
0.45391293121537596
>>>
```

This is the result of inputting the values suggested.

```
>>> calculate_e_2(2)
0.0623503143056
>>> calculate_e_2(4)
-0.036428467415
>>> calculate_e_2(8)
0.00903530693883
>>> calculate_e_2(16)
0.00610376549757
>>> calculate_e_2(32)
0.00183275514264
>>> calculate_e_2(64)
0.000478524385809
>>>
```

These are the errors with the respective number of steps.

Problem 2 b:

The following was the result of the Romberg approximations put in table form.

```

C:\Users\lab\Dropbox\Computational Physics\Jinesh_HW3>python -i Prob2b.py
>>> romberg(16)
[ 0.14797948]
[ 0.32523191 0.38431605]
[ 0.51228285 0.57463317 0.58732097]
[ 0.40299745 0.36656898 0.35269804 0.34897386]
[ 0.43010337 0.43913868 0.44397666 0.44542552 0.44580376]
[ 0.44841467 0.45451843 0.45554375 0.45572735 0.45576775 0.45577749]
[ 0.45391293 0.45574569 0.4558275 0.45583201 0.45583242 0.45583248
0.45583249]
[ 0.4553485 0.45582703 0.45583245 0.45583253 0.45583253 0.45583253
0.45583253 0.45583253]
[ 0.45571127 0.45583219 0.45583253 0.45583253 0.45583253 0.45583253
0.45583253 0.45583253]
[ 0.4558022 0.45583251 0.45583253 0.45583253 0.45583253 0.45583253
0.45583253 0.45583253]
[ 0.45582495 0.45583253 0.45583253 0.45583253 0.45583253 0.45583253
0.45583253 0.45583253]
[ 0.45583064 0.45583253 0.45583253 0.45583253 0.45583253 0.45583253
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[ 0.45583206 0.45583253 0.45583253 0.45583253 0.45583253 0.45583253
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[ 0.45583241 0.45583253 0.45583253 0.45583253 0.45583253 0.45583253
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