

AMPSE File Hierarchy

NDA protected

Open Source











- 1- Netlists_Database.py
- 2- APMSE_Graphs.py
- 3- APMSE_CircTest.py

Desanitizer



netlists_desanitized







netlists_sanitized







datasets





reg_files







make_reg





O PyTorch



DARPA AMPSE File Hierarchy

- Python Codes:
 - Netlist_Database.py:
 - Links python to all necessary SPICE testbenches
 - AMPSE_Regsearch.py:
 - Searching the parameters globally in regression models
 - AMPSE_TestCircuit.py:
 - AMPSE Verification on circuit
- Folders:
 - netlists (sanitized vs desantized)
 - datasets
 - reg_files
 - make_reg

DARPA Spectre Testbench (example)

```
0- // Generated by Cadence - spectre
•••
4- // Design view name: schematic
5- simulator lang=spectre
6- global 0
7- parameters rres=2K fnnn=15 fppp=5 VBIAS=0.9 wpppn=670n wnnn=390n lastt=60n
8- include "/technology65nm.scs" section=MC_TT
260- simulator lang=spice // for better compatibility with Xyce
261- .MEAS TRAN_CONT cont_vout1 find v(vbias) when v(oo\<1\>)=0.8 rise=1
262- .MEAS TRAN_CONT cont_tout1 when v(oo\<1\>)=0.8 rise=1
```



DARPA SPICE Output (preprocessed results)

```
$DATA1 SOURCE='AFS'
VERSION='2017_Q1_update4'
Measurement results:1.
cont_tout1, result= 2.2153e-11
cont_tout1, result= 2.7812e-10
cont_tout1, result= 5.3852e-10
cont_tout1, result= 7.9873e-10
cont_tout1, result= 1.0588e-09
cont_tout1, result= 1.3186e-09
cont_tout1, result= 1.5784e-09
cont tout1, result= 1.8382e-09
cont_tout1, result= 2.0982e-09
```



Calculate Metrics from Testbench in Repository

- Four primitive functions:
 - 1. Initialization:
 - Set parameters/metrics name,
 - Parameters range
 - Test-bench location
 - Work-area folder
 - 2. Run SPICE with given parameters
 - Runs Spectre/AFS/Xyce
 - Saves pre-processed data into list of arrays in Python
 - 3. Metrics calculation in Python
 - Calculate metrics from saved preprocessed data
 - 4. Run SPICE + Metrics calculation



Before making the Class import TestSpice and Netlists

from spectreIOlib import TestSpice, Netlists



Primitive function 1. Initialization

```
from spectreIOlib import TestSpice, Netlists
Class YOUR_MODULE_NAME(Netlists):
      def ___init___(self, tech = 65, testfolder = None):
              if tech = 65:
                    self.testbench = YOUR_TESTBENCH
                    self.testfolder = YOUR_TESTFOLDER
                    self.minpar = PARAMETER_MINVALUES # np.array 1D
                    self.maxpar = PARAMETER MAXVALUES # np.array 1D
                    self.stppar = PARAMETER_STEPSIZES # np.array 1D
                    self.parname = PARAMETER_NAMES # list
                    self.par_line_number = LINEINTESTBENCH # Integer
                    self.lst_metrics = ... # List of Dictionary
                    self.runspectre = TestSpice(...) # TestSpice Instance
                    self.finaldataset = YOURDATASET.csv
```



Primitive function 1. Initialization (Example)

```
class VCOSpice(Netlists):
       def init (self, tech = 65):
       if tech ==65:
           self.testbench = home address + '/Netlists/VCO testbenchstatic TT.scs'
           self.testfolder = home address + '/Garbage/TrashVCO 1 1'
           self.minpar = np.array([60e-9, 0.2e-6, 2, 0.2e-6, 2, 2000, 0.9, 1.0])
           self.maxpar = np.array([60e-9, 1.2e-6, 20, 1.2e-6, 20, 2000, 0.9, 1.0])
           self.stppar = np.array([1e-9, 10e-9, 1, 10e-9, 1, 1, 0.1, 0.1])
           self.parname = ['lastt','wnnn','fnnn','wpppn','fppp', 'rres','VBIAS','VDD']
           self.metricname =
['power','vcm','vfs','fnoise','f1','f2','f3','f4','f5','f6','f7','f8']
           self.par_line_number = 7
           self.lst_metrics=[{'read':'c','filename':self.testfolder +
'/test.out/test.measure', 'number':2, 'measurerange':range(9,10)},
{'read':'c','filename':self.testfolder +
'/test_cont_vout1.mt0', 'number':3, 'measurerange':range(4,600)},
{'read':'c','filename':self.testfolder +
'/test_cont_tout1.mt0', 'number':3, 'measurerange':range(4,600)}]
           self.runspectrel=TestSpice(simulator='afs',dict_folder={'testbench':self.testbench,
'trashfolder':self.testfolder},verbose = True)
```



Primitive function 1. Initialization (Example)

```
dict_folder={ 'testbench':self.testbench,
      'trashfolder':self.testfolder},
      verbose = True)
Simulator Type:
1. 'spectre'
2. 'afs'
3. 'apsplus'
4. 'apsplusplus'
5. 'spectre_ascii'
6. 'afs_ascii'
7. 'apsplus_ascii'
8. 'apsplusplus_ascii'
```



Primitive function 2. Run SPICE

```
def normal_run(self, param):
    self.dict_parameters = YOUR_PARAMETER_DICTIONARY(param)
    error_occured = self.runspectrel.runspectre(self.dict_parameters)
        if error_occured:
            out = []
        else:
            out = self.runspectrel.readmetrics(self.lst_metrics)
        return out
```



Primitive function 2. Run SPICE (Example)



Primitive function 3. Metrics calculation

```
def analysis(self, out):
    metrics = YOURFUNC(out)
    return metrics
```



Primitive function 3. Metrics calculation (Example)

```
def analysis(self, lst_out):
        time_edges=np.array(lst_out[2])
       v_edges=np.array(lst_out[1])
       prd_edge=(time_edges[wside:]-time_edges[:-wside])/wside
       v_edges=v_edges[:-wside]
       v_edges=v_edges[lside:-rside]
       prd edge=prd edge[lside:-rside]
       x=v_edges
       y=1/prd edge
        cof = np.polyfit(x, y, polynum)
       vcm=(v edges[0] + v edges[-1])/2
       vfs=(v_edges[-1]-v_edges[0])/2
       vs = np.linspace(vcm-vfs,vcm+vfs,8)
        fout=cof[7]+cof[6]*vs
+cof[5]*vs**2+cof[4]*vs**3+cof[3]*vs**4+cof[2]*vs**5+cof[1]*vs**6+cof[0]*vs**
       return [vcm, vfs]+list(fout)
```



Primitive function 4. Run Spectre + Metrics calculation

```
def wholerun_normal(self,param):
    x = self.normal_run(param)
    w = self.analysis(x)
    return w
```



Primitive function 4. Run Spectre + Metrics calculation

```
def wholerun_normal(self,param):
    x = self.normal_run(param)
    w = self.analysis(x)
    return w
```

4 primitive functions:

```
__init___(tech = YOUR_TECHNOLOGY)

RAW_SPICE_DATA = normal_run(param = TESTBENCH_PARAMETERS)

METRICS = analysis(lst_out = RAW_SPICE_DATA)

METRICS = wholerun_normal(param = TESTBENCH_PARAMETERS)
```



After making the primitive functions you have access to secondary functions:

```
paramset(xmin,xmax,xstp) # set the bounds: max, min, and step
                             # randomly choose parameters within bounds
PARAMETERS = random param()
RAW SPICE DATA = random run() # Runs normal run with random parameter
METRICS = wholerun random()
                             # Runs wholerun normal with random parameters
# BAD PARAMS are parameters that won't match with the given step. Ex.
fingers should be integers and we cannot give finger = 1.5. This is
BAD PARAMS
PARAMETERS = param_std(BAD_PARAMS) # Fixes BAD_NUMBERS
RAW_SPICE_DATA = standard_run(BAD_PARAMS) # runs param_std with normal_run
METRICS = wholerun std(BAD PARAMS) # runs param std with wholerun normal
# Making dataset:
put on csv(self, tedad=NUMBER OF SAMPLES, outcsv=OUTPUTFILE.CSV, do header=True
or False)
# do header : True if you want to have header on dataset, False when outcsv
already exists
```