

Computational Aircraft Prototype Syntheses



Training Session 3 CAPS Analysis ESP v1.19

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- Python Basics
 - Lists, Tuples and Dictionaries
- Accessing/modifying analysis values
 - `analysis.input`
- Analysis execution and outputs
 - `pre/postAnalysis`
 - `analysis.output`
- DIRTY/CLEAN process
 - Tracking changes to inputs that impact outputs
- `capsGroup` attribute
 - Connecting geometry with analysis properties
- Suggested Exercises



Python Basics: Lists

- List: ordered changeable collection. Allows duplicates.
- Created with square brackets []

```
thislist = ["apple", "banana", "banana", "cherry"]
print(thislist)      # Prints ['apple', 'banana', 'banana', 'cherry']
print(thislist[0])  # Prints 'apple'
print(thislist[-1]) # Prints 'cherry'

thislist[1] = "pear" # Change the banana to a pear
thislist.append(42) # Append 42 to the end of the list

for fruit in thislist: # Print each fruit (and 42) in the list via item
    print(fruit)

for i in range(len(thislist)): # Print each fruit (and 42) in the list via index
    print(thislist[i])
```

For more examples: www.w3schools.com/python/python_lists.asp



Python Basics: Tuples

- Tuple: ordered unchangeable collection. Allows duplicates.
- Created with parenthesis ()

```
thistuple = ("apple", "banana", "banana", "cherry")
print(thistuple) # Prints "('apple', 'banana', 'banana', 'cherry')"
print(thistuple[0]) # Prints 'apple'
print(thistuple[-1]) # Prints 'cherry'

for fruit in thistuple: # Print each fruit in the tuple via item
    print(fruit)

for i in range(len(thistuple)): # Print each fruit in the tuple via index
    print(thislist[i])

thistuple[1] = "pear" # Runtime error
```

For more examples: www.w3schools.com/python/python_tuples.asp

- Dictionary: unordered changeable indexed collection. No duplicates.
- Created with curly brackets { }
- key - value pairs separated by colon

```
thisdict = {  
    "status" : "Don't panic",  
    "Dolphin": "So long, and thanks for all the fish.",  
    42       : "The answer"  
    "Years of thought" : 7.5e6  
}  
print(thisdict["status"]) # Prints 'Don't panic'  
print(thisdict[42]) # Prints 'The answer'  
  
# Modify the answer  
thisdict[42] = "The answer to the great question... Of life, the universe and everything..."  
  
for key in thisdict: # Print each key in the dict  
    print(key)  
  
for key in thisdict: # Print each value in the dict  
    print(thisdict[key])
```

More examples: www.w3schools.com/python/python_dictionaries.asp

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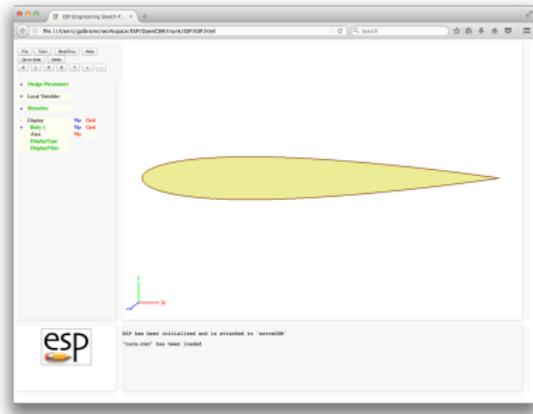


NACA Airfoil Geometry

session03/naca.csm

```
# NACA design parameters
DESPMTR  thick      0.12      #frac of local chord
DESPMTR  camber     0.00      #frac of local chord

# Construct the airfoil
UDPRIM  naca  Thickness thick   Camber  camber
        ATTRIBUTE capsAIM $xfoilAIM;tsfoilAIM
```



- Analysis inputs are set/accessed with analysis.input Sequence of Value Objects

session03/xfoil_1_AnalysisVal.py

```
# Create xfoil aim
print ("\n==> Creating xfoilAIM")
xfoil = myProblem.analysis.create(aim = "xfoilAIM",
                                  name = "xfoil")

# Set Mach number
xfoil.input.Mach = 0.5

# Print the modified mach number
mach = xfoil.input.Mach
print("\n==> Modified Mach =", mach)
```

- Analysis values can be tuples/lists and toggled

session03/xfoil_1_AnalysisVal.py

```
# Print the default value of None
print("\n==> Default Alpha =", xfoil.input["Alpha"].value)

# Set Alpha number
xfoil.input.Alpha = 2.5
print("\n==> Modified Alpha =", xfoil.input["Alpha"].value)

# Set list of Alpha
xfoil.input.Alpha = [0.0, 3.0, 5.0, 7.0, 8.0]
print("\n==> Modified Alpha =", xfoil.input["Alpha"].value)

# Unset Alpha back to None
xfoil.input.Alpha = None
print("\n==> Unset Alpha =", xfoil.input["Alpha"].value)
print()
```

- Available analysis input values in xfoil AIM documentation

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Running analysis with pyCAPS: Inputs

- Create AIM and set analysis values

session03/xfoil_2_Analysis.py

```
# Create xfoil aim
print ("\n==> Creating xfoilAIM")
xfoil = myProblem.analysis.create(aim = "xfoilAIM",
                                 name = "xfoil")

print ("\n==> Setting analysis values")
# Set Mach and Reynolds number
xfoil.input.Mach = 0.5
xfoil.input.Re    = 1.0e6

# Set list of Alpha
xfoil.input.Alpha = [0.0, 3.0, 5.0, 7.0, 8.0]
```

- Run preAnalysis to generate xfoil input files

session03/workDir_2_Analysis/Scratch/xfoil/caps.xfoil

session03/workDir_2_Analysis/Scratch/xfoil/xfoilInput.txt

- “xfoil” in the path is the AIM name

session03/xfoil_2_Analysis.py

```
print ("\n==> Loading geometry from file '"+filename+"...")
myProblem = pyCAPS.Problem(problemName = "workDir_2_Analysis",
                             capsFile = filename,
                             outLevel = 1)
```

```
# Create xfoil aim
print ("\n==> Creating xfoilAIM")
xfoil = myProblem.analysis.create(aim = "xfoilAIM",
                                  name = "xfoil")
```

```
# Run AIM pre-analysis
print ("\n==> Running preAnalysis")
xfoil.preAnalysis()
```

- Execute xfoil in session03/workDir_2_Analysis/Scratch/xfoil

session03/xfoil_2_Analysis.py

```
##### Run xfoil #####
print ("\n\n==> Running xFoil.....")

currentDirectory = os.getcwd() # Get current working directory
os.chdir(xfoil.analysisDir)    # Move into analysis directory

# Run xfoil via system call
os.system("xfoil < xfoilInput.txt > Info.out");

os.chdir(currentDirectory)    # Move back to top directory
#####
```

- CAPS currently does not execute analysis tools
(will execute some tools in next ESP version)
- Driving program responsible for execution



Running analysis with pyCAPS: postAnalysis

- Run postAnalysis to indicate completion and parse output files

session03/xfoil_2_Analysis.py

```
# Run AIM post-analysis
print ("\n==> Running postAnalysis")
xfoil.postAnalysis()
```

- Get outputs with analysis.output Value Object Sequence

```
# Retrieve Alpha, Cl and Cd
print ("\n==> Retrieve analysis results")
Alpha = xfoil.output.Alpha
Cl    = xfoil.output.CL
Cd    = xfoil.output.CD

print()
print("--> Alpha =", Alpha)
print("--> Cl    =", Cl)
print("--> Cd    =", Cd)
print()
```

- Helper function for running the analysis

session03/xfoil_3_Analysis.py

```
def run_xfoil(xfoil):
    # Run AIM pre-analysis
    print ("\n=> Running preAnalysis")
    xfoil.preAnalysis()

##### Run xfoil #####
print ("\n\n=> Running xFoil.....")

currentDirectory = os.getcwd() # Get current working directory
os.chdir(xfoil.analysisDir)    # Move into test directory

# Run xfoil via system call
os.system("xfoil < xfoilInput.txt > Info.out");

os.chdir(currentDirectory)      # Move back to top directory
#####

# Run AIM post-analysis
print ("\n=> Running postAnalysis")
xfoil.postAnalysis()
```

- Compute polar for a range of angles of attack

session03/xfoil_3_Analysis.py

```
print ("\n==> Setting analysis values")
# Set Mach and Reynolds number
xfoil.input.Mach = 0.5
xfoil.input.Re   = 1.0e6

# Set list of Alpha
xfoil.input.Alpha = [0.0, 3.0, 5.0, 7.0, 8.0]

# Run xfoil
run_xfoil(xfoil)

# Retrieve Alpha, Cl and Cd
print ("\n==> Retrieve analysis results")
Alpha = xfoil.output.Alpha
Cl    = xfoil.output.CL
Cd    = xfoil.output.CD

print()
print("--> Alpha =", Alpha)
print("--> Cl    =", Cl)
print("--> Cd    =", Cd)
print()
```

- Switch to compute polar for a range of lift coefficients

session03/xfoil_3_Analysis.py

```
# Unset Alpha, otherwise it will be included in the next analysis
xfoil.input.Alpha = None

# Set specific Cl values instead
xfoil.input.CL = [0.0, 0.1, 0.15, 0.3, 0.4]

# Run xfoil
run_xfoil(xfoil)

# Retrieve Alpha, Cl and Cd
print ("\n==> Retrieve analysis results")
Alpha = xfoil.output.Alpha
Cl    = xfoil.output.CL
Cd    = xfoil.output.CD

print()
print("--> Alpha =", Alpha)
print("--> Cl    =", Cl)
print("--> Cd    =", Cd)
print()
```

- Setup analysis values

session03/xfoil_4_Camber.py

```
# Create xfoil aim
print ("\\n==> Creating xfoilAIM")
xfoil = myProblem.analysis.create(aim = "xfoilAIM",
                                  name = "xfoil")

# Create an alias to the geometry
naca = myProblem.geometry

print ("\\n==> Setting analysis values")
# Set Mach and Reynolds number
xfoil.input.Mach = 0.5
xfoil.input.Re    = 1.0e6

# Set list of Alpha
xfoil.input.Alpha = [0.0, 1.0, 3.0]
```

- Execute sequence of cambers

session03/xfoil_4_Camber.py

```
# List of cambers to analyze
Cambers = [0.00, 0.01, 0.04, 0.07]

Alpha = [] ; Cl = [] ; Cd = []
for camber in Cambers:
    # Modify the camber
    naca.despmtr.camber = camber

    # Run xfoil
    run_xfoil(xfoil)

    # Append Alpha, Cl and Cd
    print ("\n=> Retrieve analysis results")
    Alpha.append(xfoil.output.Alpha)
    Cl.append(xfoil.output.CL)
    Cd.append(xfoil.output.CD)

print()
print("--> Cambers =", Cambers)
print("--> Alpha   =", Alpha)
print("--> Cl     =", Cl)
print("--> Cd     =", Cd)
```

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- Assigning geometry.despmtr marks geometry as DIRTY
 - Geometry always built just-in-time
- Assigning analysis.input or geometry.despmtr marks the AIM as DIRTY
- CAPS does not execute analysis, cannot execute just-in-time (for now)
- Driver is responsible for executing DIRTY AIMs
- Errors reported accessing analysis.output if AIM is DIRTY
- Errors reported running preAnalysis if AIM is CLEAN
 - Avoids inefficiencies with unnecessary calls to preAnalysis

● Execution without errors

session03/xfoil_5_CleanDirty.py

```
print("\n1. No Errors ", "-"*80)

# Set Mach and Reynolds number
xfoil.input.Mach = 0.5
xfoil.input.Re   = 1.0e6

# Set list of Alpha
print("\n==> Setting alpha sequence")
xfoil.input.Alpha = [0.0, 3.0, 5.0, 7.0, 8.0]

# Run xfoil
run_xfoil(xfoil)

# Retrieve Cl
Cl = xfoil.output.CL
print("\n--> Cl      =", Cl)
```



- Trying call analysis.output with DIRTY AIM due to analysis.input change

session03/xfoil_5_CleanDirty.py

```
print("\n2. DIRTY AnalysisVal Error ", "-"*80)

# Set a new alphas
print("\n==> Setting new alpha sequence")
xfoil.input.Alpha = [1.0, 2.0]

# Try to retrieve Cl without executing pre/postAnalysis
print("\n==> Attempting to get Cl")
try:
    Cl = xfoil.output.CL
    print("\n--> Cl      =", Cl)
except pyCAPS.CAPSError as e:
    print("\n==> CAPSError =", e)
```

- Trying get analysis.output with DIRTY AIM due to geometry.despmtr change

session03/xfoil_5_CleanDirty.py

```
print("\n3. DIRTY GeometryVal Error ", "-"*80)

# Modify a geometric parameter
print("\n==> Modifying camber")
myProblem.geometry.despmtr.camber = 0.07

# Try to retrieve Cl without executing pre/postAnalysis
print("\n==> Attempting to get Cl")
try:
    Cl = xfoil.output.CL
    print("\n--> Cl      =", Cl)
except pyCAPS.CAPSError as e:
    print("\n==> CAPSError =", e)
```

- Trying to call analysis.output without calling postAnalysis

session03/xfoil_5_CleanDirty.py

```
print("\n4. DIRTY pre- but no postAnalysis Error ", "-"*80)

# Modify mach number
print("\n==> Modifying Mach")
xfoil.input.Mach = 0.3

# Run AIM pre-analysis
print ("\n==> Running preAnalysis but not running postAnalysis")
xfoil.preAnalysis()

# Retrieve Cl
print("\n==> Attempting to get Cl")
try:
    Cl = xfoil.output.CL
    print("\n--> Cl    =", Cl)
except pyCAPS.CAPSError as e:
    print("\n==> CAPSError =", e)
```

- Calling preAnalysis with a CLEAN AIM

session03/xfoil_5_CleanDirty.py

```
print("\n5. CLEAN Error ", "-"*80)

# Don't modify any analysis or geometry values

try:
    # Run xfoil
    run_xfoil(xfoil)
except pyCAPS.CAPSError as e:
    print("\n==> CAPSError =", e)
```

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capsGroup attribute

- Tags groups of BODY/FACE/EDGE/NODE
 - Entities with same capsGroup value are in the group
- Specific use of capsGroup is AIM dependent

session03/masstran_6_f118_Wing.py

```
filename = "f118-C.csm"
print ("\n==> Loading geometry from file '"+filename+"\\...")
myProblem = pyCAPS.Problem(problemName = "workDir_6_f118_Wing",
                           capsFile = filename,
                           outLevel = 0)

# Create a masstran aim with the wing
masstran = myProblem.analysis.create(aim = "masstranAIM",
                                      name = "masstran",
                                      capsIntent="wing")
```

- Wing FACEs tagged with \$wing:faces
- masstranAIM material and properties for "wing:faces"

session03/f118-C.csm

```
BOX wing:xroot -wing:span/2 wing:zroot   wing:chord  wing:span  wing:chord*wing:thick
SELECT face
ATTRIBUTE capsGroup $wing:faces
```

session03/masstran_6_f118_Wing.py

```
# Define material properties
unobtainium = {"density" : 7850}

# Set the material
masstran.input.Material = {"Unobtainium": unobtainium}

# Define shell property
shell = {"propertyType"      : "Shell",
          "material"        : "Unobtainium",
          "membraneThickness" : 0.2}

# Associate the shell property with capsGroups defined on the geometry
masstran.input.Property = {"wing:faces": shell}
```

- capsGroups on all FACEs

session03/f118-C.csm

```
BOX wing:xroot -wing:span/2 wing:zroot wing:chord wing:span wing:chord*wing:thick
SELECT face
    ATTRIBUTE capsGroup $wing:faces



---



```
BOX htail:xroot -htail:span/2 htail:zroot htail:chord htail:span htail:chord*htail:thick
SELECT face
 ATTRIBUTE capsGroup $htail:faces

```
BOX vtail[4] 0 vtail[5] vtail:chord vtail:chord*vtail[3] vtail:span
SELECT face
    ATTRIBUTE capsGroup $vtail:faces



---



```
BOX 0 -fuse:width/2 -fuse:height/2 fuse:length fuse:width fuse:height
SELECT face 1
 ATTRIBUTE capsGroup $fuse:nose
SELECT face 2
 ATTRIBUTE capsGroup $fuse:tail
SELECT face 3
 ATTRIBUTE capsGroup $fuse:side
SELECT face 4
 ATTRIBUTE capsGroup $fuse:side
SELECT face 5
 ATTRIBUTE capsGroup $fuse:side
SELECT face 6
 ATTRIBUTE capsGroup $fuse:side
```


```


```


```

- Properties assigned to capsGroups

session03/masstran_7_f118.py

```
# Define material properties
unobtainium = {"density" : 7850}
madeupium   = {"density" : 6890}

# Set the materials
masstran.input.Material = {"Unobtainium": unobtainium,
                           "Madeupium": madeupium}

# Define shell properties
shell_1 = {"propertyType"      : "Shell",
           "material"          : "unobtainium",
           "membraneThickness" : 0.2}

shell_2 = {"propertyType"      : "Shell",
           "material"          : "madeupium",
           "membraneThickness" : 0.3}

# Associate the shell property with capsGroups defined on the geometry
masstran.input.Property = {"wing:faces" : shell_1, "htail:faces": shell_1,
                           "fuse:nose"  : shell_1, "fuse:tail"  : shell_1,
                           "vtail:faces": shell_2, "fuse:side"  : shell_2}
```

The capsGroup attribute – AVL Plane Vanilla

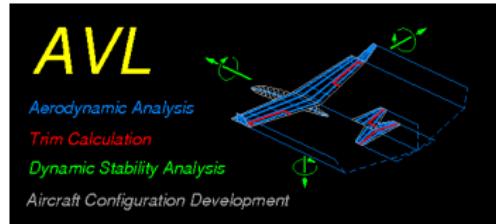
- AVL - Vortex Lattice Method: Geometry defined by airfoils
- capsGroup groups airfoils into surfaces

session03/avlPlaneVanilla.csm

```
UDPRIM    naca Thickness wing:thick Camber wing:camber
SCALE      wing:croot
ROTATEX   90      0      0
TRANSLATE  wing:xroot  0      wing:zroot
SELECT body
ATTRIBUTE capsGroup           $Wing
```

```
UDPRIM    naca Thickness wing:thick Camber wing:camber
SCALE      htail:croot
ROTATEX   90      0      0
TRANSLATE htail:xroot  0      htail:zroot
SELECT body
ATTRIBUTE capsGroup           $Htail
```

```
UDPRIM    naca Thickness vtail:thick
SCALE      vtail:croot
TRANSLATE vtail:xroot  0      vtail:zroot
SELECT body
ATTRIBUTE capsGroup           $Vtail
```



- VLM meshing parameters defined via capsGroups

session03/avl_8_PlaneVanilla.py

```
print ("\n==> Create avlAIM")
avl = myProblem.analysis.create(aim = "avlAIM",
                                name = "avl")

print ("\n==> Setting analysis values")
avl.input.Alpha = 1.0

# Set meshing parameters for each surface
wing = {"numChord" : 4,
         "numSpanTotal" : 24}

htail = {"numChord" : 4,
          "numSpanTotal" : 16}

vtail = {"numChord" : 4,
          "numSpanTotal" : 10}

# Associate the surface parameters with capsGroups defined on the geometry
avl.input.AVL_Surface = {"Wing" : wing ,
                         "Htail": htail,
                         "Vtail": vtail}
```

Thickness

- Plot airfoil polars for a range of airfoil thicknesses
 - Start from a copy of session03/xfoil_4_Camber_Plot.py

New Shells and Material

- Change `capsGroup` value for the top and bottom faces of the fuselage for F-118C (either the same or two different values)
- Starting with session03/masstran_7_f118.py, create a new shell and/or material for the newly created `capsGroup(s)`

F-118C CG Location

- Using session03/masstran_7_f118.py, create an array of F-118C CG x-locations by modifying the wing:xroot location
- Create your own (optionally share it galbramc@mit.edu)