HW7

November 14, 2023

WP7 - Exercise 1

Support functions

```
1]: import pandas as pd
   import numpy as np
   import matplotlib.pyplot as plt
   plt.style.use('default') ## reset!
   plt.style.use('paper.mplstyle')
    def shear( sig_n, s1, s2 ) :
        Creates a shear series for given values of principal stresses and sig_n
        center = (s1 + s2)/2
       radius = (s1 - s2)/2
       dx = sig_n - center
       tau_sq = radius * radius - dx * dx
       tau_sq[ tau_sq < 0 ] = None</pre>
       return np.sqrt( tau_sq )
    def mohr2d(s1, s3, ax, title="") :
        Plot Mohr diagram
        # Setup data
       npts = 1000
        [s1,s3] = sorted([s1,s3], reverse=True)
        step = (s1-s3)/npts
        sig_n = np.sort( np.append( np.arange(s3,s1,step), [s1, s3]) )
        s13_tau = shear( sig_n, s1, s3 )
        # Do the plotting stuff
       ax.plot(sig_n, s13_tau, c='gray')
       ax.set_title(title, fontsize=20)
        ax.set_ylabel("Shear stress ($\\tau$)")
       ax.set_xlabel("Effective normal stress ($\sigma_n$)")
       return ax
```

Determine the best fitting parameters of a Coulomb criterion: $\sigma_1 = UCS + q\sigma_3$.

```
import pandas as pd
import matplotlib.pyplot as plt
import numpy as np

df = pd.read_excel("DarleyDaleDataset.xlsx")
df["Sig3_MPa"] = df.S3_MPa - df.Pp_MPa
df["Sig1_MPa"] = df.S1_MPa - df.Pp_MPa
# Filter values beyond the linearity
df = df[df.Sig3_MPa < 100]

df["I1_MPa"] = df.Sig1_MPa + 2*df.Sig3_MPa
s1=df.Sig1_MPa
s3=df.Sig3_MPa
df["J2_MPa2"] = 1/6*( 2 * (s1-s3)**2 )
df["sqrt_J2_MPa"] = np.sqrt(df.J2_MPa2)

fig, ax = plt.subplots()</pre>
```

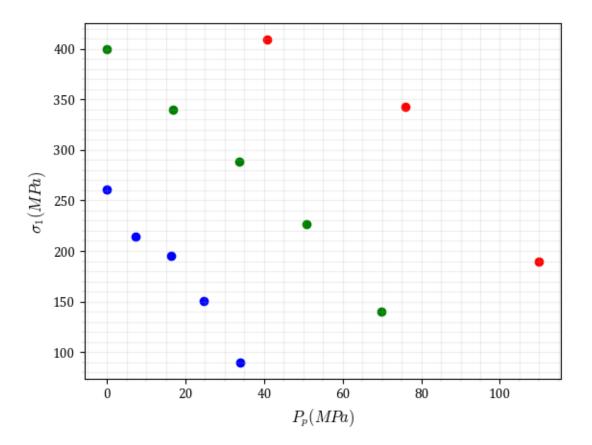
```
for name, g in df.groupby('S3_MPa'):
     ax.scatter(g.Pp_MPa, g.S1_MPa)
 ax.set_xlabel("$P_p (MPa)$")
 ax.set_ylabel("$\sigma_1 (MPa)$")
 from scipy.stats import linregress
 Q, UCS, r_value, p_value, std_err = linregress( df.Sig3_MPa, df.Sig1_MPa )
 Sig3_fit = np.linspace(0,100,10)
 Sig1_fit = Sig3_fit * Q + UCS
 fig, ax = plt.subplots()
 ax.scatter(df.Sig3_MPa, df.Sig1_MPa)
 ax.plot(Sig3_fit, Sig1_fit, ls='--', color='k')
 ax.set_xlabel("$\sigma_3$")
 ax.set_ylabel("$\sigma_1$")
 ax.set_xlim(0,150)
 ax.set_ylim(0,600)
 print(f"UCS={UCS:.2f} MPa q={Q:.4f}")
UCS=80.44 MPa q=4.5881
```

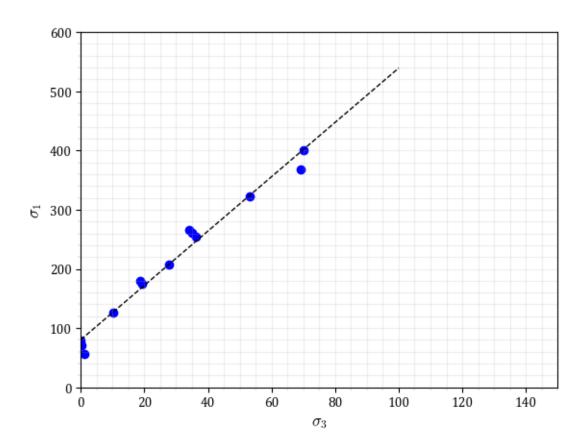
```
2]:
       S3_MPa
                  Pp_MPa
                              S1_MPa Sig3_MPa Sig1_MPa
                                                                  I1_MPa \
   1
         110 40.730779 409.023317 69.269221 368.292538 506.830979
   2
          110
               76.019958 342.219591 33.980042 266.199633 334.159717
   3
          110 110.014132 189.773495 -0.014132 79.759362 79.731098
           70 -0.055399 399.750377 70.055399 399.805776 539.916574
          70 16.932524 339.904257 53.067476 322.971732 429.106683
          70 33.748003 288.242137 36.251997 254.494134 326.998129
   6
           70
   7
               50.737176 226.162799 19.262824 175.425623 213.951271
           70 69.921900 140.320493 0.078100
   8
                                                  70.398593
                                                              70.554794
          35 -0.145370 260.542030 35.145370 260.687400 330.978141
   9
   10
          35
               7.266436 214.585236 27.733564 207.318800 262.785928
   11
          35 16.341879 195.471680 18.658121 179.129801 216.446042

    35
    24.760448
    150.286070
    10.239552
    125.525621
    146.004725

    35
    33.859217
    89.485789
    1.140783
    55.626571
    57.908136

   12
           35
   13
            J2_MPa2 sqrt_J2_MPa
      29804.981366 172.641193
   1
       17975.312781
                     134.072043
        2121.270155
                      46.057249
      36245.103735 190.381469
       24282.769258 155.829295
   6
       15876.543480 126.002157
                      90.160634
        8128.939967
       1648.323920
                      40.599556
   8
   9 16956.402398 130.216752
   10 10750.285693 103.683584
       8583.720043
                      92.648368
   11
   12
        4430.292616
                       66.560443
                     31.457385
        989.567059
   13
```

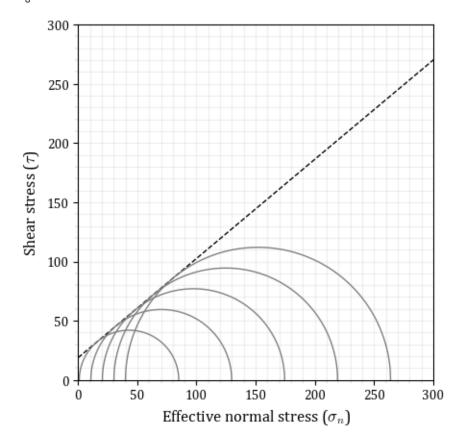




Determine the corresponding values of cohesive strength S_0 and friction angle φ from point (1) and draw the shear failure line together with Mohr circles at failure.

```
SO = UCS / 2 / np.sqrt(Q)
MU = (Q-1) / 2 / np.sqrt(Q)
PHI = np.arctan(MU)
\label{eq:print(f"S0={S0:.2f} MU={MU:.4f} PHI={PHI*180/np.pi:.2f} deg")} \\
sign=np.linspace(0,400,10)
tau = S0 + MU * sign
fig, ax = plt.subplots()
ax.plot(sign, tau, ls='--', color='k')
ax.set_xlim(0,300)
ax.set_ylim(0,300)
sig1_sig3 = Q
for sig3 in np.linspace(1,40,5) :
     sig1 = sig1\_sig3 * (sig3 + S0)
    sig1 = sig1\_sig3 * (sig3 + S0/MU) - S0/MU
    mohr2d(sig1, sig3, ax)
ax.set_aspect('equal', 'box')
```

S0=18.78 MU=0.8376 PHI=39.95 deg



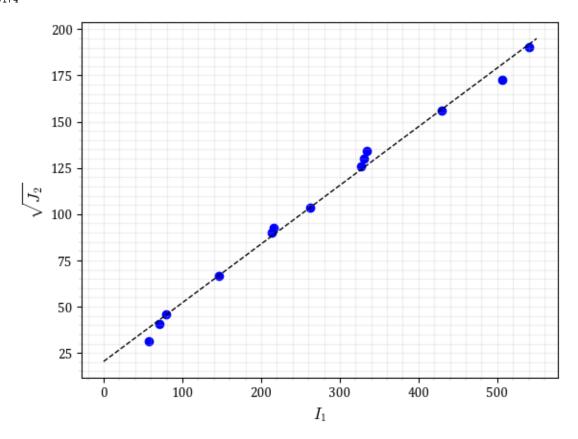
Determine the corresponding fitting parameters for a Drucker-Prager criterion and plot experimental data in a I_1 v.s. $(J_2)^{1/2}$ space.

```
fig, ax = plt.subplots()
ax.scatter( df.I1_MPa, df.sqrt_J2_MPa )

from scipy.stats import linregress
C4, C3, r_value, p_value, std_err = linregress( df.I1_MPa, df.sqrt_J2_MPa )
i1 = np.linspace(0,550,10)
sqrt_j2 = C3 + C4 * i1
ax.plot( i1, sqrt_j2, c='k', ls='--' )
```

```
ax.set_xlabel("$I_1$")
ax.set_ylabel("$\sqrt{J_2}$")
print("Drucker-Prager:")
print(f" C3={C3:.2f} C4={C4:.4f}")
```

Drucker-Prager:
 C3=20.35 C4=0.3174



Compute the corresponding tensile strength assuming a reasonable ratio of unconfined compression strength to tensile strength (browse Jaeger's book).

As of Section 4.5 and 4.6 of Jaeger's book, we can understand that a reasonable ration for $C_0/T_0 = [\sqrt{\mu^2 + 1} + \mu]^2$, or a reference value $C_0/T_0 = 15$.

```
C_0/T_0 = 4.59
UCS = 80.44 T_0 = 17.53 MPa
C_0/T_0 = 15.00
UCS = 80.44 T_0 = 5.36 MPa
```

WP7 - Exercise 2

For this assignment you have to be able to calculate principal stresses on the wall of a wellbore at an arbitrary deviation and azimuth (See Chapter 6 of my notes "Wellbore stability" (https://dnicolasespinoza.github.io/) and Zoback's book Ch. 8).

Develop a script to compute and illustrate graphically (lower hemisphere projection):

- 1. The likelihood of tensile fractures considering wellbore mud pressure is equal to pore pressure in the formation $P_W = P_p$ and a simple tensile strength criterion (independent of intermediate and maximum principal stresses), i.e., required T_s .
- 2. The likelihood of breakouts considering wellbore mud pressure is equal to pore pressure in the formation $P_W = P_p$ and a simple unconfined compression strength criterion, i.e., required UCS.
- 3. The wellbore breakout angle w_{BO} considering a Mohr-Coulomb shear failure criterion and the properties of the Darley Dale Sandstone from Exercise 1.
- 4. The wellbore breakout angle w_{BO} considering a linear I_1 v.s. $(J_2)^{1/2}$ shear failure criterion and the properties of the Darley Dale Sandstone from Exercise 1.
- 5. The wellbore breakout angle w_{BO} considering a Modified Lade shear failure criterion (Zoback's book, p. 100) and the properties of the Darley Dale Sandstone from point 1.

Limit your wellbore breakout angle plots to $0^{\circ} < w_{BO} < 60^{\circ}$ because predictions with linear elasticity over $w_{BO} > 60^{\circ}$ are likely highly inaccurate. Test your script with the three examples given in Zobacks's book in Figures 8.2 and 10.4 (normal faulting, strike-slip faulting, and reverse faulting). When computing breakouts assume $P_W = 32$ MPa (NF), $P_W = 40$ MPa (SSF), and $P_W = 52$ MPa (RF). Use a Poisson ratio of 0.25 for stress calculations.

```
6]:
    def Sp_to_Sg(Sp, alpha, beta, gamma) :
     # ROTATION MATRIX
      CA, SA = np.cos(alpha), np.sin(alpha)
      CB, SB = np.cos(beta), np.sin(beta)
      CG, SG = np.cos(gamma), np.sin(gamma)
      RPG = np.array(
                                     SA*CB
           [ [ CA*CB,
                                                          -SB 1.
                                     SA*SB*SG + CA*CG , CB*SG ],
              [ CA*SB*SG - SA*CG,
              [ CA*SB*CG + SA*SG,
                                     SA*SB*CG - CA*SG , CB*CG ]]
      # MATRIX MULTIPLICATION - STRESS IN GEOGRAPHIC COORDINATES
     Sg = RPG.T @ Sp @ RPG
      return Sg
    def Sg_to_Sw(Sg, delta, phi) :
      # ROTATION MATRIX
      CD = np.cos(delta)
     SD = np.sin(delta)
      CP = np.cos(phi)
     SP = np.sin(phi)
     Rb = np.array(
           [ [ -CD*CP,
                         -SD*CP , SP ],
              ΓSD.
                         -CD , 0
                          SD*SP , CP ]]
              Γ CD*SP.
      Sw = Rb @ Sg @ Rb.T
      return Sw
    def range_levels(min, max, n_steps) :
       n_dec = 0
        _n = n_steps
        while ( True ) :
           step = ( max - min ) / n_steps
            step = round( step, n_dec )
           if ( not step ) :
               n_dec += 1
               continue
            _n = (max - min) / step
           levels = np.arange(np.floor(min*10**n_dec)/10**n_dec, max+step, step)
            if (np.abs(_n - n_steps) < n_steps*.5) : break # Accept 20% deviation
           n_{dec} += 1
        return levels
```

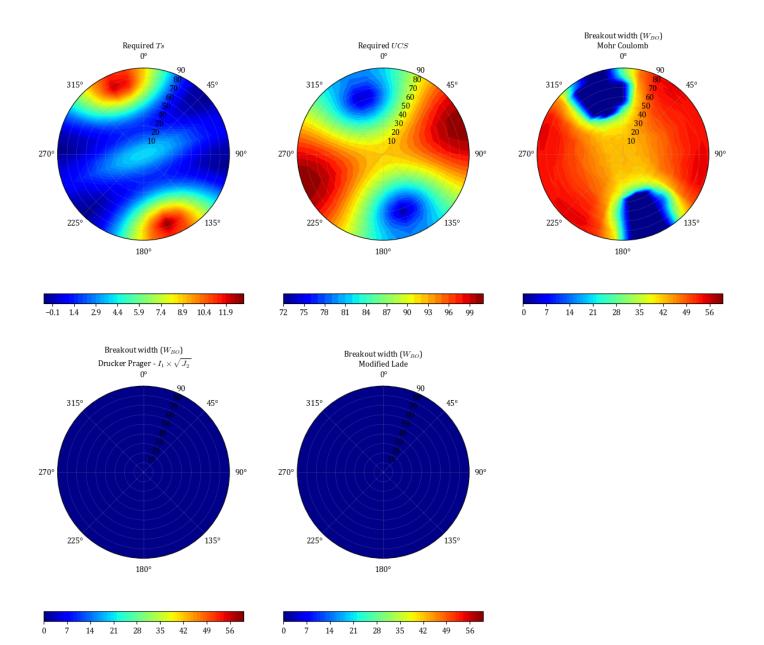
```
7]: import numpy as np
    import pandas as pd
    from numpy import sin, cos, pi
    from numpy.linalg import eigvals
    import matplotlib.pyplot as plt
    def build_polar_plots() :
       global NU
       global S1, S2, S3, AZIMUTH, REGIME
       global PW, PP
       # Mohr coulomb failure criteria: Sig1 = UCS + Q * Sig3
       global MC_UCS, MC_Q
        # I1 x J2 Drucker Parger failure criteria: SQRT_J2 = C3 + V4 * I1
       global DP_C3, DP_C4, DP_I1, DP_SJ2
       DP_I1 = []
       DP_SJ2 = []
       MC_S0 = MC_UCS / 2 / np.sqrt(MC_Q)
       MC_MU = (MC_Q-1) / 2 / np.sqrt(MC_Q)
       MC_PHI = np.arctan(MC_MU)
       # Modified Lade
       ML_S = MC_{SO} / MC_{MU}
       ML_ETA = 4 * MC_MU**2 * (9 - 7 * np.sin(MC_PHI)) / (1 - np.sin(MC_PHI))
       # MPa
       Delta_P = PW - PP
       # Sp
       Sp = [[S1, 0, 0], [0, S2, 0], [0, 0, S3]]
       # Normal
                       : az, pi/2, 0
                      : az, 0, pi/2
        # Strike slip
       # Reverse
                       : az, 0,
       az = AZIMUTH / 180 * np.pi
       if (REGIME == "NORMAL")
                                  : Sg = Sp_to_Sg(Sp, az, pi/2, 0)
       if (REGIME == "STRIKESLIP") : Sg = Sp_to_Sg(Sp, az, 0), pi/2)
       if (REGIME == "REVERSE")
                                  : Sg = Sp_to_Sg( Sp, az, 0
                                                                  , 0 )
       phimesh, deltamesh = np.meshgrid( np.linspace(0,pi/2,30), np.linspace(0, 2*np.pi, 30) )
       Ts = np.zeros_like( deltamesh )
       UCS = np.zeros_like( deltamesh )
       MC_WBO = np.zeros_like( deltamesh )
       DP_WB0 = np.zeros_like( deltamesh )
       ML_WB0 = np.zeros_like( deltamesh )
       print("Calculating ... ")
       for i, j in np.ndindex(deltamesh.shape) :
           delta = deltamesh[i,j] ; phi = phimesh[i,j]
           Sw = Sg_to_Sw(Sg, delta, phi)
            [[sig11, sig12, sig13], [sig21, sig22, sig23], [sig31, sig32, sig33]] = Sw - PP * np.identity(3)
           for theta in np.linspace(0,2*np.pi,180): # Angle around the borehole
               srr = Delta_P
               stt = sig11 + sig22 - 2*(sig11-sig22)*cos(2*theta) - 4*sig12*sin(2*theta) - Delta_P
               szt = 2 * (sig23*cos(theta) - sig13*sin(theta))
               szz = sig33 - 2*NU*(sig11-sig22)*cos(2*theta) - 4*NU*sig12*sin(2*theta)
               ST = [ [ stt, szt ], [ szt, szz] ]
               EV = sorted(eigvals(ST), reverse=True)
               if theta==0 : Ts[i,j] = EV[1]; UCS[i,j] = EV[0] # Initialize
               Ts[i,j] = np.minimum( Ts[i,j], EV[1] )
               UCS[i,j] = np.maximum( UCS[i,j], EV[0] )
                [ sig1, sig2, sig3 ] = sorted( [ EV[0], EV[1], srr ], reverse=True )
                #print(f"{sig1} - {sig2} - {sig3}")
                # Mohr coulomb criteria
               mc\_sig1\_max = MC\_UCS + MC\_Q * sig3
               if sig1 > mc_sig1_max : MC_WBO[i,j] += 1
                # Drucker Prager
```

```
i1 = sig1 + sig3 + sig2
             sqrt_j2 = np.sqrt (1/6*((sig1-sig2)**2 + (sig1-sig3)**2 + (sig2-sig3)**2))
             sqrtj2_max = DP_C3 + DP_C4 * i1
             #print(f"i1:{i1:.2f} sqrt_j2:{sqrt_j2:.2f} max:{sqrtj2_max:.2f}")
            if \ sqrt_j2 > sqrtj2_max : DP_WBO[i,j] += 1
            DP_I1.append(i1)
            DP_SJ2.append(sqrt_j2)
             # Modified Lade
             [ sig1s, sig2s, sig3s ] = np.array([ sig1, sig2, sig3 ]) + ML_S
            i1s = sig1s + sig2s + sig3s
             i3s = sig1s * sig2s * sig3s
            i1s_max_3 = (27 + ML_ETA) * i3s
            i1s_max = np.sign(i1s_max_3) * np.abs(i1s_max_3)**(1/3)
            if i1s > i1s_max : ML_WBO[i,j] += 1
    MC_WBO[MC_WBO > 60] = 60
    DP_WB0[DP_WB0 > 60] = 60
    ML_WB0[ML_WB0 > 60] = 60
    print("Data build ... [ok]")
    fig = plt.figure(figsize=(15, 20))
     def polar_plot( pos, m1, m2, Y, title, range=None, n_levels=50 ) :
        ax = plt.subplot(pos, projection='polar')
        if not range :
            range = [ np.min(Y), np.max(Y)]
        levels = range_levels( range[0], range[1], n_levels )
        CS = ax.contourf( m1, m2, Y, levels, cmap='jet' ) # contourf or pcolormesh
        cbar = fig.colorbar(CS, orientation="horizontal", pad=0.2)
        ax.set_title( title )
        ax.set_theta_zero_location("N")
        ax.set_theta_direction(-1)
    polar_plot(331, deltamesh, phimesh*180/pi, Ts, "Required $Ts$", None, 50)
     polar_plot(332, deltamesh, phimesh*180/pi, UCS, "Required $UCS$", None, 50)
    polar_plot(333, deltamesh, phimesh*180/pi, MC_WBO, "Breakout width ($W_{BO}$)\nMohr Coulomb", [0,60])
    polar_plot(334, deltamesh, phimesh*180/pi, DP_WBO, "Breakout width ($W_{BO}$)\n Drucker Prager - $I_1 \times sqrt\{J_2\}$", [0,60])
    polar_plot (335, deltamesh, phimesh*180/pi, ML_WBO, "Breakout width ($W_{BO}$)\nModified Lade", [0,60])
    print("Done")
# Mohr coulomb failure criteria: Sig1 = UCS + Q * Sig3
MC_UCS=80.44
MC_Q=4.5881
 # I1 x J2 Drucker Parger failure criteria: SQRT_J2 = C3 + V4 * I1
DP_C3=20.35
DP_C4=0.3174
 # Parameters to the Modified Lade
 S1 = 70 \# Sv
 S2 = 67 \# Shmax
S3 = 45 \# Shmin
 AZIMUTH = 160
REGIME = 'NORMAL'
PP = 32
PW = 32
NU = 0.25
print ("ZOBACK - EXAMPLE 1 (NORMAL)")
build_polar_plots()
```

ZOBACK - EXAMPLE 1 (NORMAL)

Calculating ...
Data build ... [ok]

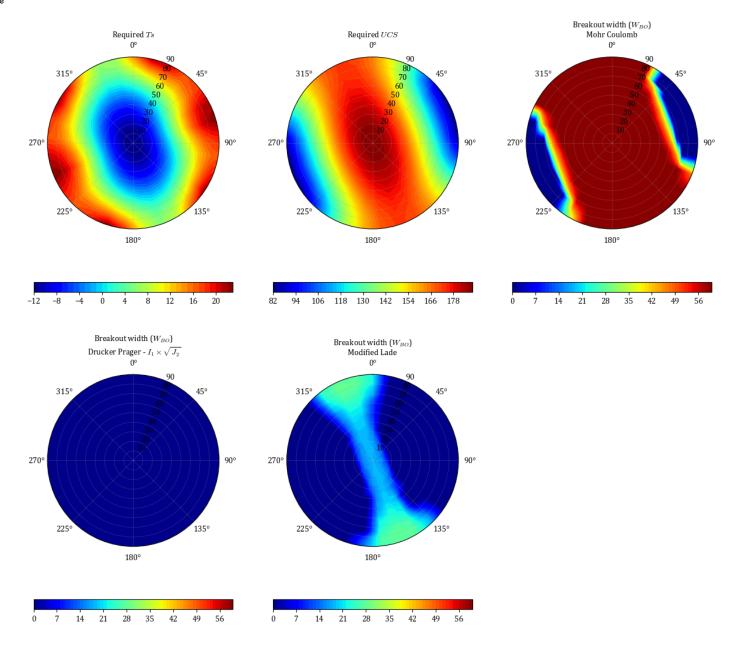
Done



```
# Mohr coulomb failure criteria: Sig1 = UCS + Q * Sig3
MC_UCS=80.44
MC_Q=4.5881
# I1 x J2 Drucker Parger failure criteria: SQRT_J2 = C3 + V4 * I1
DP_C3=20.35
DP_C4=0.3174
# Parameters to the Modified Lade
S1 = 105  # Shmax
S2 = 70
           # Sv
          # Shmin
S3 = 55
AZIMUTH = 70
REGIME = 'STRIKESLIP'
PP = 32
PW = 40
NU = 0.25
print ("ZOBACK - EXAMPLE 2 (STRIKE SLIP)")
build_polar_plots()
```

ZOBACK - EXAMPLE 2 (STRIKE SLIP) Calculating ...

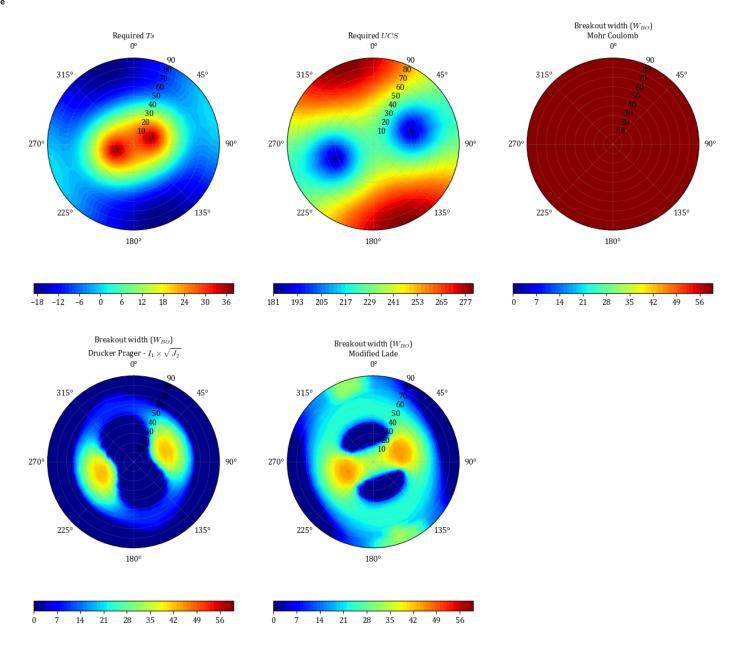
Data build ... [ok] Done



```
# Mohr coulomb failure criteria: Sig1 = UCS + Q * Sig3
MC_UCS=80.44
MC_Q=4.5881
 # I1 x J2 Drucker Parger failure criteria: SQRT_{-}J2 = C3 + V4 * I1
DP_C3=20.35
DP_C4=0.3174
 # Parameters to the Modified Lade
 S1 = 145
          # Shmax
S2 = 125
          # Shmin
S3 = 70
           # Sv
 AZIMUTH = 70
REGIME = 'REVERSE'
PP = 32
PW = 52
NU = 0.25
print ("ZOBACK - EXAMPLE 2 (REVERSE)")
```

build_polar_plots()

ZOBACK - EXAMPLE 2 (REVERSE)
Calculating ...
Data build ... [ok]



```
I]: # Mohr coulomb failure criteria: Sig1 = UCS + Q * Sig3

MC_UCS=80.44

MC_Q=4.5881

# I1 x J2 Drucker Parger failure criteria: SQRT_J2 = C3 + V4 * I1

DP_C3=20.35

DP_C4=0.3174

# Parameters to the Modified Lade
S1 = 70  # Shmax
S2 = 67  # Shmin
S3 = 45  # Sv

AZIMUTH = 70

REGIME = 'NORMAL'

PP = 32
```

```
PW = 60

NU = 0.25

print ("BREAKOUT ANALYSIS")

build_polar_plots()
```

BREAKOUT ANALYSIS
Calculating ...
Data build ... [ok]
Done

