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Apêndice A

Arquivo de entrada para o cálculo da substituição de fluidos

Adaptado do código desenvolvido por Kummar (2006)

```
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
% Dados de entrada:
% rho_o = densidade do óleo (API)
% rho_g = densidade específica do gás (API)
% T = Temperatura (oC)
% P = Pressão (psi)
% phi = porosidade (fração)
% VSH = volume de folhelho (fração)
% isw = SWT: saturação inicial de água (fração)
% tsw = saturação final de água
% vp = velocidade de onda P apartir de perfil (ft/s)
% vs = velocidade de onda S apartir de perfil (ft/s)
% rho = densidade total apartir de perfil (gm/cc)
% Dados de saída:
% vp_sat = velocidade de onda P depois da substituição de fluidos(ft/s)
% vs_sat = velocidade de onda S depois da substituição de fluidos(ft/s)
% rho_sat = densidade depois da substituição de fluidos (gm/cc)
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
API_o = 10;
rhoi_g = 1.2;
Ti = xlsread('tempi_3D.xls');
Ti = (5/9)*(Ti-32);
Pi = xlsread('pressi_3D.xls');
T = xlsread('CASE1_DIR1_42_STEP108_T.xls');
T = (5/9)*(T-32);
P = xlsread('CASE1_DIR1_42_STEP108_P.xls');
S = 0.0
phi = 0.32;
vsh = 0.03632;
isg = 0.088;
isw = 0.164;
tsw = xlsread('CASE1_DIR1_42_STEP108_Sw.xls');
tsg = xlsread('CASE1_DIR1_42_STEP108_Sg.xls');
ifluid = 1;
fluid = 2;
vp = 8397.0;
vs = 3759.0;
rho = 2.065 ;
%
% Parâmetros fixos (e.g., Mavko et al., 1998)
%
```

```

k_clay = 20.9;
k_qtz = 36.6;
rho_clay = 2.58;
rho_qtz = 2.65;
%
% Conversões
%
div_mill = 1/1000000;
fs2kms = 0.000305;
kms2fs = 3280.84;
v_clay = vsh*0.70;
v_qtz= 1-v_clay;
ish = 1-isw-isg;
tsh = 1-tsw-tsg;
rhoi_o = 141.5/(API_o+131.5);

Pi = Pi*6.894757*0.001;
P = P*6.894757*0.001;
S = S*div_mill;
vp = vp*fs2kms;
vs = vs*fs2kms;
%
GORi = 2.03*rhoi_g*(0.006894757*Pi.*exp(0.02878*API_o-0.00377*Ti)).^1.205;
GOR = 2.03*rhoi_g*(0.006894757*P.*exp(0.02878*API_o-0.00377*T)).^1.205;
%
% Passo 1: Propriedades da matriz
%
k_voigt = v_clay*k_clay + v_qtz*k_qtz;
k_reuss = 1/(v_clay/k_clay + v_qtz/k_qtz);
k_matrix = 0.5*(k_voigt + k_reuss);
rho_matrix = v_clay*rho_clay+v_qtz*rho_qtz;
%
% Passo 2: Propriedades da salmoura
%
w(1,1) = 1402.85; w(1,3) = 3.437*10^(-3);
w(2,1) = 4.871; w(2,3) = 1.739*10^(-4);
w(3,1) = -0.04783; w(3,3) = -2.135*10^(-6);
w(4,1) = 1.487*10^(-4); w(4,3) = -1.455*10^(-8);
w(5,1) = -2.197*10^(-7); w(5,3) = 5.230*10^(-11);
w(1,2) = 1.524; w(1,4) = -1.197*10^(-5);
w(2,2) = -0.0111; w(2,4) = -1.628*10^(-6);
w(3,2) = 2.747*10^(-4); w(3,4) = 1.237*10^(-8);
w(4,2) = -6.503*10^(-7); w(4,4) = 1.327*10^(-10);
w(5,2) = 7.987*10^(-10); w(5,4) = -4.614*10^(-13);
sumi = 0;
for i=1:5
    for j=1:4
        sumi = sumi+w(i,j).*Ti.^(i-1).*Pi.^(j-1);
    end
end
vi_water = sumi;

sum = 0;
for i=1:5
    for j=1:4

```

```

        sum = sum+w(i,j).*T.^(i-1).*P.^(j-1);
    end
end
v_water = sum;

v1i = 1170-9.6*Ti+0.055*Ti.*Ti-8.5*10^(-5)*Ti.*Ti.*Ti+2.6*Pi-0.0029*Ti....
    *Pi-0.0476*Pi.*Pi;
vi_brine = vi_water+S*v1i+S^1.5*(780-10*Pi+0.16*Pi.*Pi)-1820*S*S;
r1i = 489*Pi-2*Ti.*Pi+0.016*Ti.*Ti.*Pi-1.3*10^(-5)*Ti.*Ti.*Ti....
    *Pi-0.333*Pi.*Pi-0.002*Ti.*Pi.*Pi;
rhoi_water=1+10^(-6)*(-80*Ti-3.3*Ti.*Ti+0.00175*Ti.*Ti.*Ti+r1i);
r2i = 300*Pi-2400*Pi.*S+Ti.*(80+3*Ti-3300*S-13*Pi+47*Pi.*S);
rhoi_brine = rhoi_water+0.668*S+0.44*S*S+10^(-6)*S*r2i;
ki_brine = rhoi_brine.*vi_brine.*vi_brine*div_mill;
%
% Passo 3: Propriedades iniciais do hidrocarboneto
%
if ifluid == 1
    B0i = 0.972+0.00038*(2.495*GORi*sqrt(rhoi_g/rhoi_o)+Ti+17.8).^1.175;
    rhoi_ps = rhoi_o./((1+0.001*GORi).*B0i);
    rhoi_s = (rhoi_o+0.0012*GORi*rhoi_g)./B0i;
    r1i = rhoi_s+(0.00277*Pi-1.71*0.0000001*Pi.*Pi.*Pi)....
        *(rhoi_s-1.15).^2+3.49*0.0001*Pi;
    rhoi_hyc = r1i./(0.972+3.81*0.0001*(Ti+17.8).^1.175);
    vi = 2096*sqrt(rhoi_ps./(2.6-rhoi_ps))-3.7*Ti+4.64*Pi+0.0115*...
        (sqrt(18.33./rhoi_ps-16.97)-1).*Ti.*Pi;
    ki_hyc = rhoi_hyc.*vi.*vi*div_mill;
    R = 8.314;
    Tai = Ti+273.15;
    Ppri = Pi./(4.892-0.4048*rhoi_g);
    Tpri = Tai./(94.72+170.75*rhoi_g);
    E1i = exp(-Ppri.^1.2./Tpri.*(0.45+8*(0.56-1./Tpri).^2));
    Ei = 0.109*(3.85-Tpri).^2.*E1i;
    Z1i = 0.03+0.00527*(3.5-Tpri).^3;
    Zi = Z1i.*Ppri+0.642*Tpri-0.007*Tpri.^4-0.52+Ei;
    rho_hyc_g_i = 28.8*rhoi_g*P./(Zi*R.*Tai);
    dz_dp_i=Z1i+0.109.*(3.85-Tpri).^2.*E1i.*(-1.2*Ppri.^0.2./Tpri.*...
        (0.45+8*(0.56-1./Tpri).^2));
    yoi = 0.85+5.6./(Ppri+2)+27.1./(Ppri+3.5).^2-8.7*exp(-0.65*(Ppri+1));
    k_hyg_i = Pi.*yoi./1000*1.0./(1-Ppri./Zi.*dz_dp_i);
end
%
% Passo 4: Propriedades do fluido
%
ki_fl = 1./(isw./ki_brine+ish./ki_hyc+isg./k_hyg_i);
rhoi_fl = isw.*rhoi_brine+ish.*rhoi_hyc+isg.*rho_hyc_g_i;
%
% Passo 5: Módulo original in-situ (rocha saturada in-situ)
%
dens_poros = 0;
if dens_poros == 1
    rho = phi*rhoi_fl + (1-phi)*rho_matrix;
end
k_sat = rho*(vp*vp-vs*vs^4/3);
g = rho*vs*vs;

```



```

%
% Passo 6: Propriedades do arcaço
%
k1 = k_sat*(phi*k_matrix./ki_fl+1-phi)-k_matrix;
k2 = phi*k_matrix./ki_fl+k_sat/k_matrix-1-phi;
k_frame = k1./k2;
%
% Passo 7: Seleção do tipo de fluido de saída
%
if fluid == 1

elseif fluid == 2

    B0 = 0.972+0.00038*(2.495*GOR*sqrt(rhoi_g/rhoi_o)+T+17.8).^1.175;
    rho_ps = rhoi_o./((1+0.001*GOR).*B0);
    rho_s = (rhoi_o+0.0012*GOR*rhoi_g)/B0;
    r1 = rho_s+(0.00277*P-1.71*0.0000001*P.*P.*P).*(rho_s-1.15).^2+...
        3.49*0.0001*P;
    rho_hyc = r1./(0.972+3.81*0.0001*(T+17.78).^1.175);
    v = 2096*sqrt(rho_ps./(2.6-rho_ps))-3.7*T+4.64*P+0.0115*...
        (sqrt(18.33./rho_ps-16.97)-1).*T.*P;
    k_hyc = rho_hyc.*v.*v*div_mill;
    R = 8.314;
    Ta = T+273.15;
    Ppr = P./(4.892-0.4048*rhoi_g);
    Tpr = Ta./(94.72+170.75*rhoi_g);
    E1 = exp(-Ppr.^1.2./Tpr.*(0.45+8*(0.56-1./Tpr).^2));
    E = 0.109*(3.85-Tpr).^2.*E1;
    Z1 = 0.03+0.00527*(3.5-Tpr).^3;
    Z = Z1.*Ppr+0.642*Tpr-0.007*Tpr.^4-0.52+E;
    rho_hyc_g = 28.8*rhoi_g*P./(Z*R.*Ta);
    dz_dp=Z1+0.109.*(3.85-Tpr).^2.*E1.*(-1.2*Ppr.^0.2./Tpr.*(0.45+8*...
        (0.56-1./Tpr).^2));
    yo = 0.85+5.6./(Ppr+2)+27.1./(Ppr+3.5).^2-8.7*exp(-0.65*(Ppr+1));
    % Salmoura
    v11 = 1170-9.6*T+0.055*T.*T-8.5*10^(-5)*T.*T.*T+2.6*P-0.0029*T.*...
        P-0.0476*P.*P;
    v1_brine = v_water+S*v11+S^1.5*(780-10*P+0.16*P.*P)-1820*S*S;
    r11 = 489*P-2*T.*P+0.016*T.*T.*P-1.3*10^(-5)*T.*T.*T.*P-0.333*P.*...
        P-0.002*T.*P.*P;
    rho1_water=1+10^(-6)*(-80*T-3.3*T.*T+0.00175*T.*T.*T+r11);
    r21 = 300*P-2400*P.*S+T.*(80+3*T-3300*S-13*P+47*P.*S);
    rho1_brine = rho1_water+0.668*S+0.44*S*S+10^(-6)*S*r21;
    k_brine = rho1_brine.*v1_brine.*v1_brine*div_mill;

    k_hyg = P.*yo./1000*1.0./(1-Ppr./Z.*dz_dp);
end
%
% Passo 8: Propriedades do fluido (saturação final) e densidade
% da rocha saturada
%
k_fl = 1./(tsw./k_brine + tsh./k_hyc + tsg./k_hyg);
rho_fl = tsw.*rho1_brine + tsh.*rho_hyc + tsg.*rho_hyc_g;
rho_sat = phi*rho_fl+(1-phi)*rho_matrix;
%

```

```
% Passo 9: Módulo total da rocha saturada
%
k1 = phi./k_fl+(1-phi)/k_matrix-k_frame/(k_matrix*k_matrix);
k_sat_new = k_frame + ((1-k_frame./k_matrix).^2)./k1;
%
% Passo 10: Velocidade sísmica depois da substituição de fluidos
%
vp_sat = sqrt((k_sat_new+g*4/3)./rho_sat)*1000;
vs_sat = sqrt(g./rho_sat)*1000;
AI = vp_sat.*rho_sat;
```