Exercis One

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1 Applicant Details

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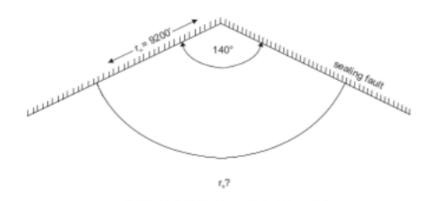
• Assessment Github Repo: Click Here

2 Exercise One

Exercise One (From L.P. Dake Exercise 9.2)

A wedge-shaped reservoir is suspected of having a fair strong natural water drive. The geometry of the reservoir-aquifer system is shown below. The reservoir was at bubble point at the initial condition with no gas cap (m=0). Develop a computer program that will Perform history matching of the reservoir using the production and PVT data given below.

Hint: Use Klins et al paper: "A Polynomial approach to the Van Everdingen - Hurst dimensionless variables for water encroachment"



			Properti	es	and aquifer
		h	=	100 ft	
Reservoir properties			k	=	200 mD
			μ_w	=	0.55 cp
			ø	=	0.25
N =	_	312×10 ⁶ stb	Cw	=	3.0×10 ⁻⁶ /psi
	_		Cf	=	4.0×10 ⁻⁶ / psi
S_{wc}	=	0.05	B _w	=	1.0

2.1 PROJECT ARCHITECTURE

2.1.1 Tools

- Excel
- Python language
- Pycharm IDE
- Flask
- Turtle
- Tkinter
- Jupyter Notebook
- Git and Github

2.1.2 Key points

- Fair strong natural water drive mechanism.
- Wedge-shaped reservoir.
- Initial Conditions are at bubble point(P and T).
- No gas cap(m = 0).

2.1.3 Aim

• To develop a computer program that will perform history matching of the reservoir using the production and PVT data given.

2.1.4 PVT Data:

		Time	Pressure at the	he Plateau Pr	ressure Δρ
Time (years)	N _p (MM/stb)	R _p (scf/stb)	B _o (rb/stb)	R _s (scf/stb)	B _g (rb/scf)
0		650 (R _{si})	1.404(B _{oi})	650 (R _{si})	.00093 (B _{gi})
1	7.88	760	1.374	592	.00098
2	18.42	845	1.349	545	.00107
3	29.15	920	1.329	507	.00117
4	40.69	975	1.316	471	.00128
5	50.14	1025	1.303	442	.00139
6	58.42	1065	1.294	418	.00150
7	65.39	1095	1.287	398	.00160
8	70.74	1120	1.280	383	.00170
9	74.54	1145	1.276	381	.00176
10	77.43	1160	1.273	364	.00182

2.1.5 Production Data

2.1.6 Approach

- Using Klins et al paper: "A Polynomial approach to the Van Everdingen Hurst dimensionless variables for water encroachment"
- 1. The van Everdingen and Hurst dimensionless variables rho/sub D/ and q/sub D/.

3 Importing Necessary Python Libraries

```
[1]: import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
import seaborn as sns
import scipy
```

4 Setting Result Display Margins

```
[2]: pd.set_option("display.max_column", None)
   pd.set_option("display.max_row", None)
   pd.set_option("display.max_colwidth", None)
   #pd.set_option("display.float_format", lambda x: "%.2f" %x)
   plt.style.use("ggplot")
```

```
[3]: # Creating PVT Data table
     Table Features
     1. Time(yrs)
     2. Pressure at Oil water contact(psia)
     3. Plateau Pressure Level(psia)
     4. Change in pressure(psia)
     5. Cummulative Oil Production, Np(MM/stb)
     6. Cummulative Gas Oil Ratio, GOR(scf/stb)
     7. Oil formation volume factor, Bo(rb/stb)
     8. Gas Solubility, Rs(scf/stb)
     9. Gas Formation Volume Factor, Bg(rb/scf)
     The first three(1-3) are independent Variable while 4-9 are dependent variables _{\square}
      ⇒which can be
     empirically determined using there respective equations/correlations.
     Though, the values of the dependent variables are already given in the PVT_{\sqcup}
      \hookrightarrow table shown, we will
     try to automate its determination and append the results into the table.
     11 11 11
```

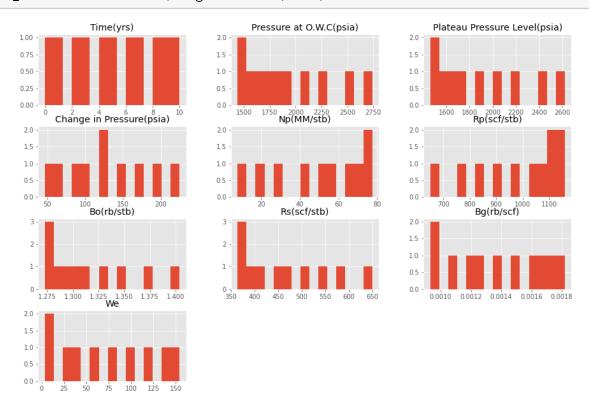
```
pvt_table = {"Time(yrs)": [* range(11)],
              "Pressure at O.W.C(psia)": [2740, 2500, 2290, 2109, 1949, 1818, __
\rightarrow1702, 1608, 1635, 1480, 1440],
              "Plateau Pressure Level(psia)": [np.nan, 2620, 2395, 2199, 2029, __
\rightarrow1883, 1760, 1655, 1571, 1507, 1460],
              "Change in Pressure(psia)": [120, 225, 196, 170, 146, 123, 105, __
\rightarrow84, 64, 47, np.nan],
              "Np(MM/stb)": [np.nan, 7.88, 18.42, 29.15, 40.69, 50.14, 58.42, 65.
\rightarrow39, 70.74, 74.54, 77.43],
              "Rp(scf/stb)": [650, 760, 845, 920, 975, 1025, 1065, 1095, 1120, ___
→1145, 1160],
              "Bo(rb/stb)": [1.404, 1.374, 1.349, 1.329, 1.316, 1.303, 1.294, 1.
\rightarrow287, 1.280, 1.276, 1.273],
              "Rs(scf/stb)": [650, 592, 545, 507, 471, 442, 418, 398, 383, 381, __
→364],
              "Bg(rb/scf)": [0.00093, 0.00098, 0.00107, 0.00117, 0.00128, 0.
\rightarrow00139, 0.00150, 0.00160, 0.00170,
                              0.00176, 0.00182],
              "We": [np.nan, 3.829, 13.460, 26.462, 41.935, 59.207, 77.628, 96.
→805, 116.284, 135.601, 154.401]
pvt_data = pd.DataFrame.from_dict(pvt_table)
pvt_data.head(11)
```

[3]:	Time(vrs)	Pressure at O.	W.C(psia)	Plateau Pressu	re Level(psia) \
0	0	rrobbaro av o.	2740	Tuodaa Tiobba	Na:	
1	1		2500		2620.	
2	2		2290		2395.	
3	3		2109		2199.	
4	4		1949		2029.	
5	5		1818		1883.	
6	6		1702		1760.	
7	7		1608		1655.	
8	8		1635		1571.	0
9	9		1480		1507.	0
10	10		1440		1460.	0
	Changa in	Pressure(psia)	Nn (MM /a+h)	Pn(scf/sth)	Po(rh/g+h)	\
0	Change III	120.0	Np(MM/SCD)	-	1.404	`
1		225.0	7.88	760	1.374	
2		196.0	18.42	845	1.349	
3		170.0	29.15	920	1.329	
4		146.0	40.69	975	1.316	
5		123.0	50.14	1025	1.303	

6	105.0	58.42	1065	1.294
7	84.0	65.39	1095	1.287
8	64.0	70.74	1120	1.280
9	47.0	74.54	1145	1.276
10	NaN	77.43	1160	1.273

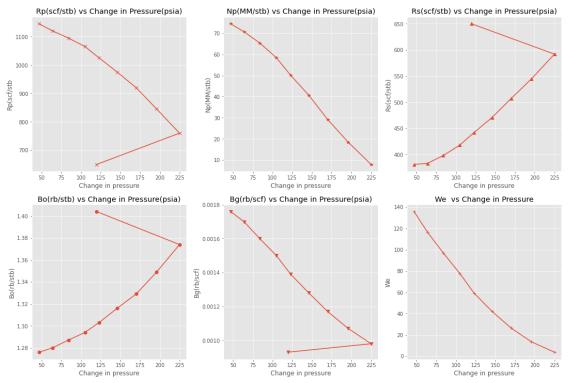
Rs(scf/stb)	Bg(rb/scf)	We
650	0.00093	NaN
592	0.00098	3.829
545	0.00107	13.460
507	0.00117	26.462
471	0.00128	41.935
442	0.00139	59.207
418	0.00150	77.628
398	0.00160	96.805
383	0.00170	116.284
381	0.00176	135.601
364	0.00182	154.401
	650 592 545 507 471 442 418 398 383 381	650 0.00093 592 0.00098 545 0.00107 507 0.00117 471 0.00128 442 0.00139 418 0.00150 398 0.00160 383 0.00170 381 0.00176

[4]: pvt_data.hist(bins = 15, figsize = (15,10));



[5]: f, ax = plt.subplots(nrows=2, ncols=3, figsize=(15,10), constrained_layout=True) #f.tight_layout()

```
ax[0][0].plot(pvt_data["Change in Pressure(psia)"], pvt_data["Rp(scf/stb)"],__
 →marker = "x")
ax[0][0].set_title("Rp(scf/stb) vs Change in Pressure(psia)")
ax[0][0].set(xlabel='Change in pressure', ylabel='Rp(scf/stb')
ax[0][1].plot(pvt_data["Change in Pressure(psia)"], pvt_data["Np(MM/stb)"],
→marker = '*')
ax[0][1].set_title("Np(MM/stb) vs Change in Pressure(psia)")
ax[0][1].set(xlabel='Change in pressure', ylabel='Np(MM/stb)')
ax[0][2].plot(pvt_data["Change in Pressure(psia)"], pvt_data["Rs(scf/stb)"],_u
→marker = '^')
ax[0][2].set_title("Rs(scf/stb) vs Change in Pressure(psia)")
ax[0][2].set(xlabel='Change in pressure', ylabel='Rs(scf/stb)')
ax[1][0].plot(pvt_data["Change in Pressure(psia)"], pvt_data["Bo(rb/stb)"],
→marker = "o")
ax[1][0].set_title("Bo(rb/stb) vs Change in Pressure(psia)")
ax[1][0].set(xlabel='Change in pressure', ylabel='Bo(rb/stb)')
ax[1][1].plot(pvt_data["Change in Pressure(psia)"], pvt_data["Bg(rb/scf)"],
→marker = "v")
ax[1][1].set_title("Bg(rb/scf) vs Change in Pressure(psia)")
ax[1][1].set(xlabel='Change in pressure', ylabel='Bg(rb/scf)')
ax[1][2].plot(pvt_data["Change in Pressure(psia)"],pvt_data["We"], marker = "+")
ax[1][2].set_title("We vs Change in Pressure")
ax[1][2].set(xlabel='Change in pressure', ylabel='We')
plt.show()
```



- The above table represent the production data over time.
- The above table will serve as our observed data or what is usually referred to production history data
- We shall perform history matching by making a comparison between the observed history data and the calculated which is from our Algorithm using Klins et al paper: "Polynomial Approach to the Van Everdigen Hurst Dimensionless variables for water encroachment. The paper is basically using numerical approximation.
- let Calculate the Cumulative water influx using the given production and PVT data
- Since the value of rD wasn't give, we will run a for loop from range 2 25 as stated in Klin et al paper

```
[6]: # Using constant terminal rate, pD
     # Determining tD using rD of 10
    def Klins_params_verification(
    k = int(input("Enter value of permeability in mD")),
    t = int(input("Enter the maximum value of in yrs")),
    phi = float(input("Enter the phase angle value phi")),
    Uw = float(input("Enter the water viscosity value")),
    Cw = float(input("Enter the water compressibility value, Uw")),
    Cf = float(input("Enter the formation compressibility value, Cf")),
    ro = float(input("Enter the reservoir radius value, ro"))):
        Cfw = Cw + Cf
        tD1 = []
        t cross1 = []
        for time in range(t + 1):
            tD = (2.309 * k * time) / (phi * Uw * Cfw * pow(ro, 2))
            print()
            print("========="RESULTS=======")
            print("At time(t):",time, "\t", "The dimensionless time is:", round(tD, __
     →2))
            tD1.append(tD)
        print()
        print("Let Calculate the crossover time to know if the aquifer is finite or_
      →infinite")
```

```
# Calculating crossover time to ascertain if is finite or infinite aquifer
    for rD in range(2, 26):
       bo = 0.0980958
       b1 = 0.100683
       b2 = 2.03863
       t_{cross} = bo * (rD - 1) + b1 * (rD - 1) ** b2
       print()
 →print("==========================")
       print("The dimensionless radius is:",rD, "\t", "The crossover time⊔

→corresponding to tD is:", round(t_cross, 2))
       t cross1.append(t cross)
    print("=====TESTING IF AQUIFER IS FINITE OR INFINITE=====")
    for (val1, val2) in zip(tD1, t_cross1):
       if val1 >= val2:
           print("===Is a finite Aquifer===")
       else:
           print("===Is an infinite Aquifer===")
    pass
    The results shows we have a finite aquifer.
    11 11 11
    print(f"The dimensionless time values are:,\n{tD1}", "\n")
    print("=======END=======")
    print(f"The crossover times are:, \n{t_cross1}", "\n")
    print("=======END========")
    print("==The results shows a finite aquifer===")
params = Klins_params_verification()
params
Enter value of permeability in mD200
Enter the maximum value of in yrs10
Enter the phase angle value phi0.25
Enter the water viscosity value0.55
Enter the water compressibility value, Uw0.000003
Enter the formation compressibility value, Cf0.000004
Enter the reservoir radius value, ro9200
At time(t): 0
              The dimensionless time is: 0.0
```

At time(t): 1 The dimensionless time is: 5.67
======================================
At time(t): 2 The dimensionless time is: 11.34
======================================
At time(t): 3 The dimensionless time is: 17.01
======================================
At time(t): 4 The dimensionless time is: 22.67
======================================
At time(t): 5 The dimensionless time is: 28.34
======================================
At time(t): 6 The dimensionless time is: 34.01
DEGIJI TO
======================================
======================================
At time(t): 8 The dimensionless time is: 45.35
======================================
At time(t): 9 The dimensionless time is: 51.02
At time(t): 10 The dimensionless time is: 56.69
===
Let Calculate the crossover time to know if the aquifer is finite or infinite
===
======================================
The dimensionless radius is: 2 The crossover time corresponding to tD is: 0.2
======================================
The dimensionless radius is: 3 The crossover time corresponding to tD is: 0.61
======================================
The dimensionless radius is: 4 The crossover time corresponding to tD is: 1.24
======================================
The dimensionless radius is: 5 The crossover time corresponding to tD is: 2.09
======================================

The dimensionless radius is: 6 The crossover time corresponding to tD is: 3	3.17
======================================	
The dimensionless radius is: 7 The crossover time corresponding to tD is:	1.47
======================================	
The dimensionless radius is: 8 The crossover time corresponding to tD is: 0	3.01
======================================	
The dimensionless radius is: 9 The crossover time corresponding to tD is:	7.77
======================================	
The dimensionless radius is: 10	tD
======================================	
The dimensionless radius is: 11 The crossover time corresponding to is: 11.99	tD
D Traver Tra	
======================================	
The dimensionless radius is: 12	τD
======================================	
The dimensionless radius is: 13	tD
======================================	
The dimensionless radius is: 14 The crossover time corresponding to is: 20.06	tD
======================================	
The dimensionless radius is: 15 The crossover time corresponding to is: 23.23	tD
The dimensionless radius is: 16 The crossover time corresponding to is: 26.62	tD
======================================	
The dimensionless radius is: 17	tD
======================================	
The dimensionless radius is: 18 The crossover time corresponding to is: 34.13	tD
DEGIT TO	
======================================	

```
is: 38.24
is: 42.59
The dimensionless radius is: 21
                     The crossover time corresponding to tD
is: 47.18
The dimensionless radius is: 22 \, The crossover time corresponding to tD
is: 52.0
is: 57.07
is: 62.38
is: 67.92
====TESTING IF AQUIFER IS FINITE OR INFINITE=====
===Is an infinite Aquifer===
===Is a finite Aquifer===
The dimensionless time values are:,
[0.0, 5.66862249281909, 11.33724498563818, 17.00586747845727, 22.67448997127636,
28.343112464095448, 34.01173495691454, 39.68035744973363, 45.34897994255272,
51.017602435371806, 56.686224928190896]
======END========
The crossover times are:,
[0.19877879999999999, 0.6098529334495365, 1.239718427487128, 2.0919322677795527,
```

3.1690142581459124, 4.472927267793876, 6.0052832825402325, 7.767451993800409, 9.760625355004983, 11.985859190630363, 14.444101628247461, 17.13621340834513,

```
20.06298290822148, 23.225137574375108, 26.623352827591447, 30.258259137021614, 34.13044773467111, 38.240475298846135, 42.58886784130767, 47.17612396953358, 52.00271765163416, 57.06910058044037, 62.37570421090754, 67.92294152855882]
```

```
[7]: # Claculating aquifer constant
     def Aquifer_constant(
     f = float(input('Enter f')),
     phi = float(input('Enter phi')),
     h = float(input("Enter h")),
     Cw = float(input("Enter Cw")),
     Cf = float(input("Enter Cf")),
     ro = float(input("Enter ro"))):
        11 11 11
         f = 0.389
         HHHH
         Cfw = Cw + Cf
         U = 1.119*f*phi*h*Cfw* pow(ro, 2)
         print()
         print(f"The Aquifer constant value is: {round(U, 3)}")
         return U
     U = Aquifer_constant()
```

```
Enter f0.389
Enter phi0.25
Enter h100
Enter Cw0.000003
Enter Cf0.000004
Enter ro9200
```

The Aquifer constant value is: 6447.53

==The results shows a finite aquifer===

[7]: 6447.530292000001

```
[8]: # Calculating Dimensionless water influx, wD using rD = 10

def WrD_10(Aquifer_constant, W_D, t_D):
    We = U * change_in_pressure * W_D*(t_D)
    return We
```

5 Using Klins methods

```
[9]: from math import *
     from cmath import *
     class Klin_et_al:
         \# , rD, k, t, phi, Cw ,Cf, Cfw = Cw+Cf, ro, error
         def __init__(self,
                      rD=float(input("Enter the value of dimensionless radius,rD")),
                      k=float(input("Enter the value of permeability,k")),
                      t=int(input("Enter value of time, t")),
                      phi=float(input("Enter the value of phase angle, phi")),
                      Cw=float(input("Enter the value of water compressibility, ___
      Cf=float(input("Enter the value of pore compressibility, Cf")),
                      ro=float(input("Enter the value of Oil radius, ro")),
                      f=float(input("Enter the value of f")),
                      h=float(input("Enter the value of the pay thickness")),
                      change_in_pressure=float(input("Enter the value of change in_⊔
      →pressure"))):
             11 11 11
             Using rD = 10 as the value of our rD
             k = permeability
             t = range \ of \ time
             phi = phase angle
             Uw = water \ viscosity
             Cfw = Total compressibility
             ro = Reservoir radius
             error = 0.0029\%
             .....
             self.rD = rD
             self.k = k
             self.t = t
             self.phi = phi
             self.Cw = Cw
             self.Cf = Cf
             self.Cfw = self.Cw + self.Cf
             self.ro = ro
             self.f = f
             self.h = h
             self.change_in_pressure = change_in_pressure
                   self.error = 0.000029
         def tD(self):
             # for time in range(self.t):
```

```
tD = 2.309 * self.k * self.t / self.phi * self.Cw * self.Cfw * self.rou
→** 2
       # print("At time(t):", time, "\t", "The dimensionless time is:", "
\rightarrow round(tD, 2))
       return tD
   def t_cross(self):
       bo = -1.767
       b1 = -0.606
       b2 = 0.12368
       b3 = 3.02
       b4 = 2.25
       b5 = 0.50
       t_cross = bo + b1 * self.rD + b2 * self.rD ** b4 + b3 * (log(self.rD, __
→e)) ** b5
       if self.tD > self.t_cross:
           print("=====The Aquifer is Finite Aquifer=====")
           print("======The Aquifer is infinite Aquifer======")
       return t_cross
   # Using the Appendix B approach to determine the qD
   def Apha 1(self):
       bo = -0.00222107
       b1 = -0.627638
       b2 = 6.277915
       b3 = -2.734405
       b4 = 1.2708
       b5 = -1.100417
       alpha_1 = bo + b1 * (1 / sinh(self.rD)) + b2 * self.rD ** b3 + b4 *_{\sqcup}
⇒self.rD ** b5
       return alpha_1
   def Apha_2(self):
       bo = -0.00796608
       b1 = -1.85408
       b2 = 18.71169
       b3 = -2.758326
       b4 = 4.829162
       b5 = -1.009021
       alpha_2 = bo + b1 * (1 / sinh(self.rD)) + b2 * self.rD ** b3 + b4 *_{\sqcup}
⇒self.rD ** b5
       return alpha_2
   # finding Jo amd J1
```

```
HHHH
         Jo = Bessel function of first kind order 0
         J1 = Bessel function of first kind order 1
         Since our aquifer is finite, it will be advisable to find Jo and J1 using
         x in range 0<= x <3. So, our x = 2
         Jo(x) = bo +b1(x/3)**2 +b2(x/3)**4 +b3(x/3)**6 + b4(x/3)**8 +b5(x/3)**10_{\square}
\leftrightarrow +b6(x/3)**12
         (x) - I*J1(x) = bo +b1 (x/3)**2 +b2(x/3)**4 + b3(x/3**6 + b4(x/3)**8 +b5(x/3)**4 + b3(x/3)**6 + b4(x/3)**8 +b5(x/3)**8 +b5(x
\rightarrow 3)**10 + b6(x/3)**12,
         n n n
         def J_o(self):
                     x = 2
                     bo = 1.00
                    b1 = -2.24999997
                     b2 = 1.2656208
                    b3 = -0.3163866
                     b4 = 0.0444479
                    b5 = -0.0039444
                     b6 = 0.0002100
                     Jo = bo + b1 * (x / 3) ** 2 + b2 * (x / 3) ** 4 + b3 * (x / 3) ** 6 + 1
\rightarrow b4 * (x / 3) ** 8 + b5 * (x / 3) ** 10 + b6 * (x / 3)
                     return Jo
         def J_1(self):
                     x = 2
                     bo = 0.5000
                     b1 = -0.56249985
                     b2 = 0.21093573
                    b3 = 0.03954289
                     b4 = 0.00443319
                    b5 = -0.00031761
                    b6 = 0.00001109
                     J1 = (bo + b1 * (x / 3) ** 2 + b2 * (x / 4) ** 4 + b3 * (pow(x / 3, 6))_{\cup}
\rightarrow+ b4 * (x / 3) ** 8 + b5 * (x / 3) ** 10 + b6 * (x / 3) ** 12)*x
                     return J1
         # Finding the dimensionless flow rate
         def dimensionless_rate(self):
                     Alpha_1 = self.Apha_1()
```

```
Alpha_2 = self.Apha_2()
        tD = self.tD()
        J_1 = self.J_1()
        J_o = self.J_o()
        error = 0.000029
        qD_num1 = 2 * e ** (-Alpha_1 ** 2 * tD * J_1 ** 2 * (Alpha_1 * self.rD))
        qD_deno1 = (Alpha_1 ** 2 * (J_o ** 2 * Alpha_1 - J_1 ** 2 * (Alpha_1 *_U))
 ⇒self.rD)))
        qD num2 = 2 * e ** (-Alpha_2 ** 2 * tD * J_1 ** 2 * (Alpha_2 * self.rD))
        qD_{deno2} = Alpha_2 * (J_o ** 2 * Alpha_2 - J_1 ** 2 * (Alpha_2 * self.
 ⊶rD))
        ⊶error
        return qD
    # Calculating aguifer constant
    def Aquifer_constant(self):
        B = 1.119 * self.f * self.phi * self.h * self.Cfw * pow(self.ro, 2)
        return B
    # Calculating Cumulative water influx using klin et al
    def We 10(self):
        B = self.Aquifer_constant()
        qD = self.dimensionless_rate()
        t D = self.tD()
        We = B * self.change_in_pressure * qD * t_D
        return We
method = Klin_et_al()
print(method.We_10())
\# rD = 10, k = 200, t = [* range (11)], phi = 0.25, Cw = 0.000003, Cf = 0.
 \rightarrow000004, ro = 9200, error = 0.000029
Enter the value of dimensionless radius, rD10
Enter the value of permeability, k200
Enter value of time, t10
Enter the value of phase angle, phi0.25
Enter the value of water compressibility, Cw0.000003
Enter the value of pore compressibility, Cf0.000004
Enter the value of Oil radius, ro9200
Enter the value of f0.389
Enter the value of the pay thickness100
Enter the value of change in pressure225
(27231725768.127693+0j)
```

```
[10]: # The time and corresponding We using the Klin et al
      t1 = 2995179062
      t2 = 5158645218
      t3 = 6639422824
      t4 = 7522915487
      t5 = 7839724010
      t6 = 7947626041
      t7 = 7341045140
      t8 = 6326226015
      t9 = 5172722501
      t10 = 0
      We_klins = {"We_klins": [2995179062,5158645218, 6639422824, 7522915487,
      →7839724010, 7947626041,
                              7341045140, 6326226015, 5172722501]}
      df_we = pd.DataFrame.from_dict(We_klins)
      pvt_data["Calculated_We"] = df_we["We_klins"]
      pvt data.head()
[10]:
         Time(yrs)
                    Pressure at O.W.C(psia) Plateau Pressure Level(psia) \
                                        2740
                                                                       NaN
      1
                 1
                                       2500
                                                                    2620.0
      2
                 2
                                       2290
                                                                    2395.0
      3
                 3
                                                                    2199.0
                                        2109
                                        1949
                                                                    2029.0
         Change in Pressure(psia) Np(MM/stb)
                                               Rp(scf/stb) Bo(rb/stb) Rs(scf/stb) \
                                                                  1.404
      0
                            120.0
                                          NaN
                                                        650
                                                                                  650
      1
                            225.0
                                         7.88
                                                        760
                                                                  1.374
                                                                                  592
      2
                            196.0
                                        18.42
                                                        845
                                                                  1.349
                                                                                 545
      3
                                                        920
                                                                  1.329
                                                                                 507
                            170.0
                                        29.15
      4
                            146.0
                                        40.69
                                                        975
                                                                  1.316
                                                                                 471
         Bg(rb/scf)
                         We Calculated We
            0.00093
      0
                        {\tt NaN}
                              2.995179e+09
            0.00098
                     3.829
                              5.158645e+09
      1
      2
            0.00107 13.460
                              6.639423e+09
      3
            0.00117 26.462 7.522915e+09
            0.00128 41.935
                              7.839724e+09
```

6 Performing history matching using graphical representaions

