

The Fracpole Suite

Ken Kundert

Designer's Guide Consulting, Inc.

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The fracpole suite is a collection of programs used to generate Spectre subcircuits that model components that require a fractional pole impedance. The suite includes a program for generating the fractional pole impedance itself, a program for generating models of inductors that include skin effect, and a program that generates models of capacitors that include dielectric absorption.

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1.0 Background

A *fracpole* is a basic element that can be used to construct models of common components. It exhibits a fractional impedance pole. Such behavior is useful for modeling skin effect in inductors, dielectric absorption in capacitors, and for converting white noise into flicker noise.

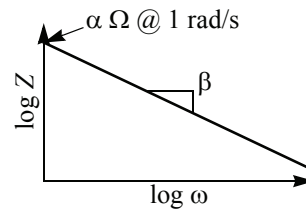
An ideal fracpole exhibits the following impedance,

$$Z(s) = \alpha s^\beta. \quad (1)$$

which is shown in Figure 1.

FIGURE 1.

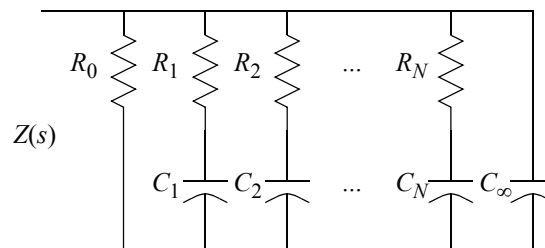
Impedance of an ideal fracpole.



Such a component is distributed if β is not an integer. It has time constants that are spread over an infinite range, from infinitely long to infinitely short. In a transient-based simulator, it must be approximated over a finite range of frequencies, from f_0 to f_1 . The component will be internally constructed as a lumped approximation that consists of a parallel combination of resistors and capacitors, as shown in Figure 2. For this reason, β

FIGURE 2.

A lumped approximation to an ideal fracpole.



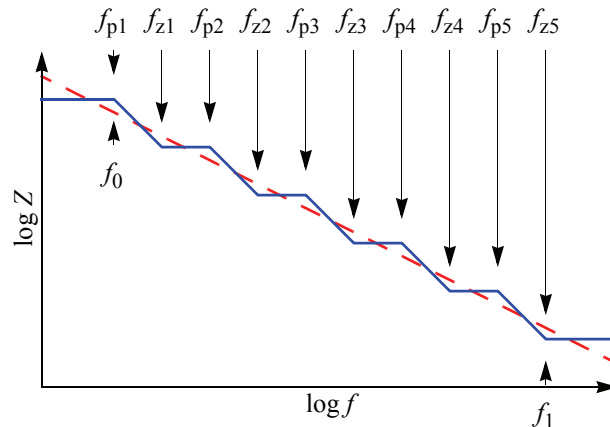
is restricted to fall in the range $-1 < \beta < 0$. In this case, it represents a fractional impedance pole at 0 and is referred to as an impedance fracpole. If instead the approximation is constructed with inductors replacing the capacitors, then β must fall in the range of $0 < \beta < 1$. This represents a fractional admittance pole at 0 and is referred to as an admittance fracpole. However, inductors are somewhat less efficiently modeled than capacitors in most circuit simulators, and the same effect can be accomplished by combining a impedance fracpole with a gyrator.

An ideal fracpole is approximated over a finite range of frequencies by a circuit with real pole and zero pairs such that the poles and zeros alternate and the pairs are evenly

spaced in a logarithmic sense over that range, as shown in Figure 3. The spacing

FIGURE 3.

Approximating an ideal fracpole using a collection of real poles and zeros.

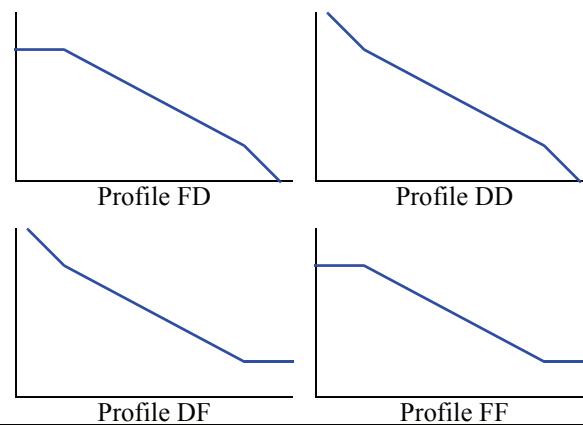


between the pole and zero is the same for each pair, and this spacing relative to the spacing of the pairs themselves determines the slope, β , of the approximation. The number of lumps, N , is chosen to provide a sufficient level of accuracy. The larger N , the closer the impedance of the lumped approximation matches that of the ideal fracpole. Typically, using $1\frac{1}{2}$ lumps per decade of frequency range give a good fit (but 1 lump per decade is often sufficient).

The behavior outside the range of the approximation is determined by whether the first and last critical frequencies are poles or zeros. Four possibilities exist, and they are distinguished by referring to them as different profiles. The profiles are named with a two letter code, where the first letter refers to the low-frequency behavior and the second to the high-frequency behavior. The letters are either d or f , where d signifies ‘down’ or a down-slope and f signifies ‘flat’ or no slope. The profiles are shown in Figure 4.

FIGURE 4.

Log-impedance versus log-frequency for the various profiles.



2.0 Fractional Pole Modeler: FP

The *fp* program takes a series of command line arguments and generates a Spectre sub-circuit that models a fractional pole.

It takes the following arguments:

<i>f0</i>	The low frequency limit for the approximation (Hz).
<i>f1</i>	The high frequency limit for the approximation (Hz).
<i>coef</i>	The unity intercept point for the ideal impedance (the magnitude of the impedance when $\omega=1$ before approximation). The default value is 1.
<i>slope</i>	The slope of the impedance when plotted on a log-log scale (equals the negative of the fraction of a pole desired). Default is -0.5 . Must be less than 0 and greater than -1 .
<i>lumps</i>	The number of lumps used in the approximation (use <i>lumps</i> < 0 to specify number of lumps/decade, in this case it need not be integer).
<i>profile</i>	Specifies whether the extreme critical frequencies of the impedance approximation are poles or zeros.

As an example, entering

```
fp f0=1 f1=1e6 slope=-0.5 lumps=-1.5 profile=dd
```

at the Unix command prompt generates the subcircuit given in Listing 1. It is placed in a file named *fracpole.scs* and has the impedance, resistance, and reactance characteristics shown in Figure 5.

Running the same command, except using *lumps=-1* to get 1 lump (pole/zero pair) per decade, gives the results shown in Figure 6. In this case there is noticeable ripple.

It is recommended that for best accuracy outside the approximation range, the DD profile is used when *slope* < -0.5 and the FF profile is used when *slope* > -0.5 .

3.0 Inductor Modeler: IND

The *ind* program takes a series of command line arguments and generates a Spectre sub-circuit that models an inductor that exhibits skin effect loss. The basic schematic of the inductor is shown in Figure 7 [6]. The component *H* models skin effect and is implemented using a fracpole and a gyrator.

It takes the following arguments:

<i>l</i>	The inductance (H).
<i>rs</i>	The low frequency resistance (the ESR) (Ω).
<i>cp</i>	The shunt parasitic capacitance (F).
<i>rp</i>	The high frequency resistance (Ω).
<i>h</i>	The skin effect parameter.

LISTING 1.

```

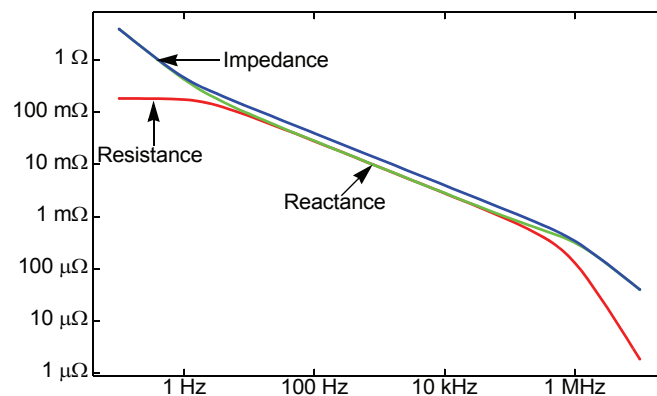
FracPole subcircuit from the file "fracpole.scs".

simulator lang=spectre
// Fractional Impedance Pole
// Impedance has a slope of -0.5 on a log-log scale,
// model is valid from 1 Hz to 1e+06 Hz.
subckt fracpole (p n)
  parameters coef=1

  R1 (p 1) resistor r=0.427978*coef
  C1 (1 n) capacitor c=0.253357/coef
  R2 (p 2) resistor r=0.290855*coef
  C2 (2 n) capacitor c=0.0803174/coef
  R3 (p 3) resistor r=0.143043*coef
  C3 (3 n) capacitor c=0.0351848/coef
  R4 (p 4) resistor r=0.0671815*coef
  C4 (4 n) capacitor c=0.01614/coef
  R5 (p 5) resistor r=0.0312686*coef
  C5 (5 n) capacitor c=0.007471/coef
  R6 (p 6) resistor r=0.0145382*coef
  C6 (6 n) capacitor c=0.00346187/coef
  R7 (p 7) resistor r=0.00678533*coef
  C7 (7 n) capacitor c=0.00159802/coef
  R8 (p 8) resistor r=0.0032317*coef
  C8 (8 n) capacitor c=0.000722864/coef
  R9 (p 9) resistor r=0.00172075*coef
  C9 (9 n) capacitor c=0.000292484/coef
  Cinf (p n) capacitor c=0.000398944/coef
ends fracpole
  
```

FIGURE 5.

Impedance characteristics of fracpole with 1.5 lumps/decade.



- f_0 The low frequency limit for the skin effect approximation (Hz). It is computed automatically to be below the ESR corner frequency if not given.
- f_1 The high frequency limit for the skin effect model (Hz). It is computed automatically to be above the ESL resonant frequency if not given.

FIGURE 6.

Impedance characteristics of fracpole with 1 lump/decade.

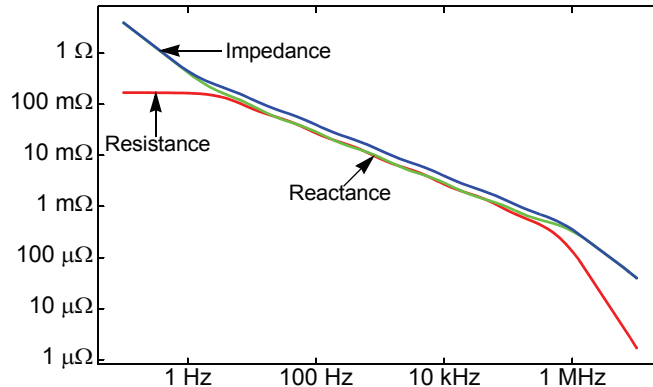
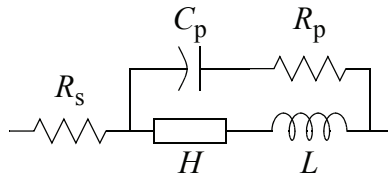


FIGURE 7.

RF inductor model.



lumps The number of lumps used in the skin effect model (use $lumps < 0$ to specify number of lumps/decade, in this case it need not be integer).

profile Specifies whether the extreme critical frequencies of the skin effect approximation are poles or zeros.

name The name that will be used for the subcircuit.

As an example, entering

```
ind name=a01t l=2.6e-9 rp=8 rs=0.001 cp=230e-15 h=704000 lumps=-1.5 \
profile=dd
```

at the Unix command prompt generates the subcircuit given in Listing 2. These are the parameters associated with the Coilcraft *a01t* minispring inductor. The subcircuit name was specified to be *a01t*, so the subcircuit is placed in a file named *a01t.scs*. The subcircuit has the resistance and reactance characteristics shown in Figure 8.

It is recommended that for best accuracy on a full-range model the DD profile be used. It is not necessary to specify $f0$ or $f1$ for a full range model, but the size of the model can be reduced by specifying a restricted range for the skin effect approximation using $f0$ and $f1$ and/or by reducing the number of lumps in the approximation.

It may be necessary to specify the *scaling* parameter on the subcircuit when using it to make the R and C values reasonably sized, though the default value works well in most cases.

LISTING 2.

Inductor subcircuit.

```
simulator lang=spectre
subckt a01t (1 2)
  parameters scaling=1M gmin=1e-12
  L (1 3) inductor l=2.6e-09
  S (3 4) skin_effect s=1/(sqrt(2*M_PI)*704000)
  Cp (1 5) capacitor c=2.3e-13
  Rp (5 4) resistor r=8
  Rs (4 2) resistor r=0.001

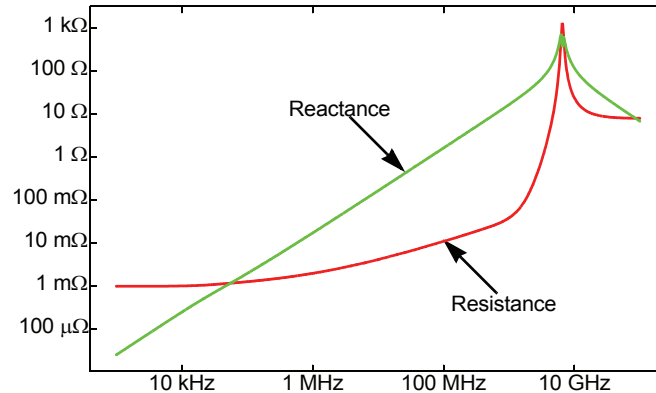
  subckt skin_effect (1 2)
    parameters s=1
    G1 (1 2 3 0) gyrator r=sqrt(scaling)
    A1 (3 0) fracpole coef=scaling/s

    // Gyrator used to convert fractional impedance pole into a fractional admittance
    pole
    subckt gyrator (t1 b1 t2 b2)
      parameters r=1kOhm
      Gm1 (t1 b1 t2 b2) vccs gm=1/r
      Gm2 (b2 t2 t1 b1) vccs gm=1/r
      G1 (t1 b1) resistor r=1/gmin
      G2 (t2 b2) resistor r=1/gmin
    ends gyrator

    // Fractional Impedance Pole
    // Impedance has a slope of -0.5 on a log-log scale (valid from 25kHz to 65GHz.
    subckt fracpole (p n)
      parameters coef=1
      R1 (p 1) resistor r=0.00278394*coef
      C1 (1 n) capacitor c=0.00157399/coef
      R2 (p 2) resistor r=0.00196615*coef
      C2 (2 n) capacitor c=0.000508957/coef
      R3 (p 3) resistor r=0.000999681*coef
      C3 (3 n) capacitor c=0.000228597/coef
      R4 (p 4) resistor r=0.000484109*coef
      C4 (4 n) capacitor c=0.000107801/coef
      R5 (p 5) resistor r=0.000232054*coef
      C5 (5 n) capacitor c=5.13586e-05/coef
      R6 (p 6) resistor r=0.000111005*coef
      C6 (6 n) capacitor c=2.45185e-05/coef
      R7 (p 7) resistor r=5.31306e-05*coef
      C7 (7 n) capacitor c=1.16984e-05/coef
      R8 (p 8) resistor r=2.55506e-05*coef
      C8 (8 n) capacitor c=5.55525e-06/coef
      R9 (p 9) resistor r=1.2561e-05*coef
      C9 (9 n) capacitor c=2.58058e-06/coef
      R10 (p 10) resistor r=6.93017e-06*coef
      C10 (10 n) capacitor c=1.06815e-06/coef
      Cinf (p n) capacitor c=1.56378e-06/coef
    ends fracpole
  ends skin_effect
ends a01t
```

FIGURE 8.

Impedance characteristics of Coilcraft a01t minispring inductor.



The inductor modeler uses the *fracpole* program, which constructs a fracpole as a parallel combination of series RC pairs and so β is restricted to be negative. Skin effect requires that β be positive. To accommodate the limitations of the fracpole program, let $\beta = -1/2$ and then use a gyrator to flip the sign. Choose

$$\alpha = \frac{r^2}{\sqrt{2\pi H}}, \quad (2)$$

where r is the gyrator gain, which is chosen to provide proper scaling for the signal levels. Doing so results in the over all impedance for the combination of the gyrator and the fracpole being

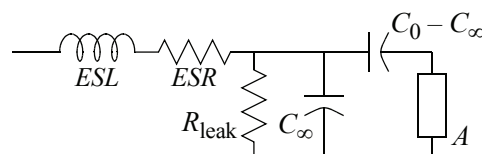
$$Z(f) = \frac{\sqrt{j}f}{H}. \quad (3)$$

4.0 Capacitor Modeler: DA

The *da* program takes a series of command line arguments and generates a Spectre sub-circuit that models a capacitor that exhibits dielectric absorption. The basic schematic of the capacitor is shown in Figure 9 [7]. The component A is a fracpole that models dielectric absorption.

FIGURE 9.

Capacitor model that includes dielectric absorption.



It takes the following arguments:

cinf The high frequency limit for capacitance (F).

<i>c0</i>	The low frequency limit for capacitance (F).
<i>tau0</i>	The average time constant of dielectric dipoles (s).
<i>alpha</i>	Specifies the width of time constant distribution. Must be greater than 0 and less than 1, with larger values implying a larger spread.
<i>esr</i>	The equivalent series resistance (Ω).
<i>esl</i>	The equivalent series inductance (H).
<i>rleak</i>	The leakage resistance (Ω).
<i>f0</i>	The low frequency limit for the dielectric absorption model (Hz).
<i>f1</i>	The high frequency limit for the dielectric absorption model (Hz).
<i>lumps</i>	The number of lumps used in the dielectric absorption approximation (use <i>lumps</i> < 0 to specify number of lumps/decade, in this case it need not be integer).
<i>profile</i>	Specifies whether the extreme critical frequencies of the skin effect approximation are poles or zeros.
<i>name</i>	The name that will be used for the subcircuit.

As an example, entering

```
da name=lossycap cinf=10e-9 c0=22.5e-9 tau0=1 alpha=0.75 f0=0.01 f1=1e6 \  
esr=0.25 esl=8e-9 rleak=1e9 lumps=-1.5 profile=fd
```

at the Unix command prompt generates the subcircuit given in Listing 3. These are the parameters of a 10 nF XR7 multilayer monolithic ceramic capacitor. The subcircuit name was specified to be *lossycap*, so the subcircuit is placed in a file named *lossycap.scs*. The subcircuit has the resistance and reactance characteristics shown in Figure 10.

For a full range model, specify *f0* so that it is below the R_{leak} corner and specify *f1* so that it is above the *ESL* resonance. The the size of the model can be reduced by specifying a restricted frequency range for the dielectric approximation using *f0* and *f1* and/or by reducing the number of lumps in the approximation.

5.0 Conclusion

The *FracPole Suite* is available from www.designers-guide.org/Modeling.

Eventually the fracpole component will be added to Spectre as a built-in primitive. When that occurs, these utility programs will no longer be needed because the inductor and capacitor models can be implemented as simple Spectre parameterized subcircuits. At that point support for the utility programs *fp*, *ind*, and *da* will be retired.

5.1 If You Have Questions

If you have questions about what you have just read, feel free to post them on the *Forum* section of *The Designer's Guide Community* website. Do so by going to www.designers-guide.org/Forum.

LISTING 3.

```

Capacitor subcircuit.

simulator lang=spectre
//
// Capacitor model that include dielectric absorption
//
subckt lossycap (1 2)
  L (1 3) inductor l=8e-09 r=0.25
  C (3 2) capacitor c=1e-08
  RI (3 2) resistor r=1e+09
  Cx (3 4) capacitor c=1.25e-08
  DA (4 2) fracpole

  // Fractional Impedance Pole
  // Impedance has a slope of -0.75 on a log-log scale,
  // model is valid from 0.01 Hz to 1e+06 Hz.
  subckt fracpole (p n)
    parameters coef=8e+07
    R0 (p n) resistor r=7.96829*coef
    R1 (p 1) resistor r=9.42501*coef
    C1 (1 n) capacitor c=0.41576/coef
    R2 (p 2) resistor r=2.40967*coef
    C2 (2 n) capacitor c=0.327729/coef
    R3 (p 3) resistor r=0.703962*coef
    C3 (3 n) capacitor c=0.226085/coef
    R4 (p 4) resistor r=0.210529*coef
    C4 (4 n) capacitor c=0.152355/coef
    R5 (p 5) resistor r=0.0632518*coef
    C5 (5 n) capacitor c=0.102198/coef
    R6 (p 6) resistor r=0.0190214*coef
    C6 (6 n) capacitor c=0.0684891/coef
    R7 (p 7) resistor r=0.00572177*coef
    C7 (7 n) capacitor c=0.0458861/coef
    R8 (p 8) resistor r=0.00172196*coef
    C8 (8 n) capacitor c=0.0307281/coef
    R9 (p 9) resistor r=0.000519356*coef
    C9 (9 n) capacitor c=0.0205325/coef
    R10 (p 10) resistor r=0.000158398*coef
    C10 (10 n) capacitor c=0.0135676/coef
    R11 (p 11) resistor r=5.16126e-05*coef
    C11 (11 n) capacitor c=0.00839165/coef
    Cinf (p n) capacitor c=0.0199735/coef
  ends fracpole
ends lossycap

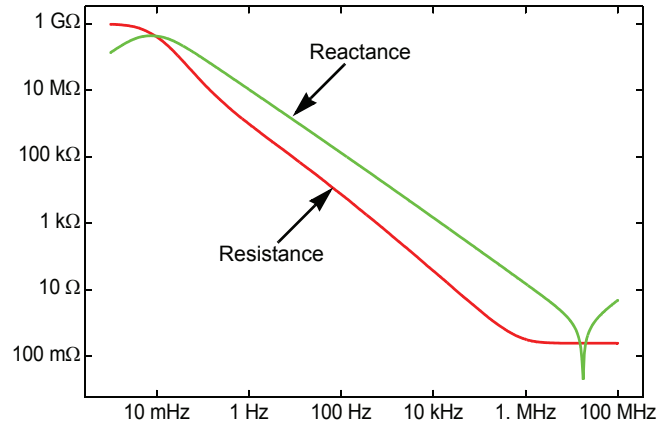
```

Postscript

I would like to thank Cairong (or Charlie) Hu for pointing out a flaw in the *da* program. Prof. J. A. Tenreiro Machado has pointed out that the *fracpole* is an application of fractional calculus (FC) modeling, and that there is a well established community studying

FIGURE 10.

Impedance characteristics of 10 nF XR7 multilayer monolithic ceramic capacitor model.



FC and its applications. He continued by saying that the algorithm for modeling fractional poles with a collection of poles and zeros was first proposed by A. Oustaloup some years ago [9] and that Bohannon also published some interesting work where attempts to exploit the fractional order nature of capacitor, or what he calls ‘fractance’ [3], which is the characteristic exhibited by a ‘fractor’. He references work that presents ideas similar to those presented in this paper, but based on a model of the dielectric developed by Curie that dates back to 1889 [5,10].

Fracpoles also play a role in modeling batteries [2], where a fracpole is referred to as a *Warburg element* and is used to model electrochemical diffusion. There are several diffusion models, but the simplest is the Warburg element [4]. A Warburg impedance element can be used to model semi-infinite linear diffusion, that is, unrestricted diffusion to a large planar electrode. Its impedance is

$$Z_W = \frac{A_W}{\sqrt{j\omega}}. \quad (4)$$

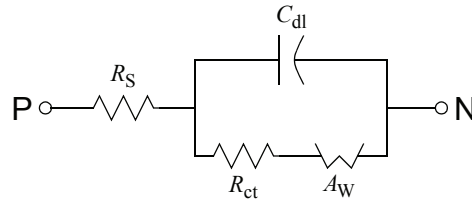
One models a parallel plate electrochemical cell using Randles’ circuit, which employs a Warburg element to represent the electrochemical diffusion. Randles’ circuit is an equivalent electrical circuit that consists of an active electrolyte resistance R_S in series with the parallel combination of the double-layer capacitance C_{dl} and the impedance of a faradaic reaction. The impedance of a faradaic reaction consists of an active charge transfer resistance R_{ct} and a specific electrochemical element of diffusion A_W , which is the Warburg element. [1,8,11]

References

- [1] Bard & Faulkner. *Electrochemical Methods: Fundamentals and Applications*. Second Edition, 2000.
- [2] Baudry, P. et al, Electro-thermal modelling of polymer lithium batteries for starting period and pulse power, *Journal of Power Sources*, vol 54, pp. 393-396, 1995.

FIGURE 11.

Randles circuit model of an electrochemical cell.



- [3] Gary W. Bohannon. *Analog Realization of a Fractional Control Element – Revisited*, October 27, 2002, Wavelength Electronics, Inc. Available on the web from http://mechatronics.ece.usu.edu/foc/cdc02tw/cdrom/additional/FOC_Proposal_Bohannon.pdf.
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- [10] S. Westerlund and L. Ekstam. Capacitor theory. *IEEE Transactions on Dielectrics and Electrical Insulation*, vol. 1, no. 5, pp. 826-839, October 1994.
- [11] Wikipedia. Randles circuit. en.wikipedia.org/wiki/Randles_circuit.