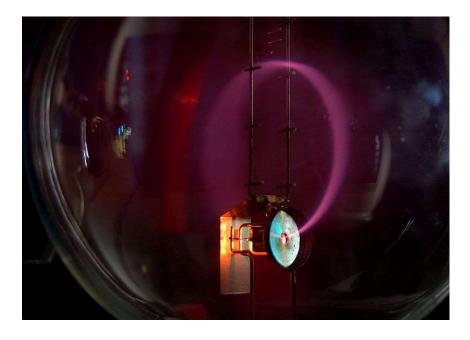
Worksheet 2 — $E/M\,$ Ratio

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1 - Preamble

1A - Modules

Import NumPy, Matplotlib, SymPy, and more

```
In [1]: %pylab inline
    import re
    import sympy
    from IPython.display import display, Markdown, Latex
    import requests
    import scipy.interpolate as interpolate
    import scipy.stats as stats
    from scipy.constants import physical_constants, mu_0
```

Populating the interactive namespace from numpy and matplotlib

1B - Units

Define units to convert measurements to SI

2 - Functions

2A - Model

```
In [3]: def B0(K, I):
    return K*I

In [4]: def B(r, R, B0):
    return

In [5]: def Kr(B, B0, K):
    return B/B0 * K

In [6]: def B_T(Kr, I_s, I_1):
    return Kr * (I_s + I_1) / 2

In [7]: def B_E(Kr, I_s, I_1):
    return Kr * (I_s - I_1) / 2
```

K_r expression:

Interpolate a function from the provided data table

```
# This is the data table provided in the experiment outline.
In [8]:
        radius_vs_field_table = array([
             [0.0, 1],
             [0.1, 0.99996],
             [0.2, 0.99928],
             [0.3, 0.99621],
             [0.4, 0.98728],
             [0.5, 0.96663],
             [0.6, 0.92525],
             [0.7, 0.85121],
             [0.8, 0.73324],
             [0.9, 0.56991],
             [1.0, 0.38007]
        ])
        # Do the interpolation
        radius_vs_field_interp = interpolate.UnivariateSpline(
             # X-values
             radius_vs_field_table[:,0],
             # Y-values
              radius_vs_field_table[:,1]
        def K_r(K, r, R):
             return radius_vs_field_interp(r/R) * K
```

```
In [9]: def e_over_m(V, B, r):
    return 2*V/(B**2*r**2)
```

2B - Data Processing Functions

Subtract 1% from anode voltage readings:

```
In [10]: def correct_anode_voltage(data):
    return data * [0.99, 1]
```

Convert a float into a latex exponential notation: latex_exp(2.3e-3) $ightarrow 2.3 imes 10^{-3}$

```
In [11]: def latex_exp(x, pres=2):
    exp = int(math.log10(abs(x)))
    mant = x / 10**exp * sign(x)
    return ('{:.'+str(pres)+'f} \\times 10^{{{:d}}}').format(mant, exp)
```

Get the strength of the magnetic field in Teslas using the <u>NOAA website (http://www.ngdc.noaa.gov/geomagweb/#igrfwmm)</u>:

```
def get B E(year, month, day, lat, long):
In [12]:
              url = 'http://www.ngdc.noaa.gov/geomag-web/calculators/calculateIgrfwmm'
              response = requests.post(url, data={
                  'browserRequest': 'true',
                  'coordinateSystem': 'M',
                  'dateStepSize': 1.0,
                  'elevation': 1,
                  'elevationUnits': 'K',
                  'endDay': day,
                  'endMonth': month,
                  'endYear': year,
                  'lat1': lat,
                  'lat1Hemisphere': 'N',
                  'lon1': long,
                  'lon1Hemisphere': 'W',
                  'model': 'WMM',
                  'resultFormat': 'csv',
                  'startDay': day,
                  'startMonth': month,
                  'startYear': year
              })
              csv = re.split('[,\n]', response.text)
              return float(csv[-2])*nT, float(csv[-10])*nT
```

2C – Data Analysis Functions

```
In [14]: def analyze_paired_data(diam, V, Il, Is):
    # Calculate Kr
    Kr = K_r(K, diam/2, R)

# Eath's magnetic field
    Be = B_E(Kr, Is, Il)

# Total field
    Bt = B_T(Kr, Is, Il)

# Charge to mass ratio
    em = e_over_m(V, Bt, diam/2)

return Be, Bt, em
```

3 - Data

3A - Readings

```
R = 15.4*cm
R err = 5*mm
K = 7.73e-4 * T/A
K err = 0.04e-4 *T/A
diamter_err = 1.5*mm
voltage_err = 0.5*V
amperage_err = 0.001*A
N = 130
# Store our results by diameter, and by pol
results = {
     8*cm:{},
    10*cm:{}
}
# 10cm radius, pol 1
results[10*cm]['pol1'] = array([
    [150.00, 0.966],
    [170.50, 1.020],
    [190.35, 1.068],
    [210.00, 1.161],
    [230.26, 1.196],
    [249.97, 1.275],
]) * [V, A] # Specify units
# 10cm radius, pol 2
results[10*cm]['pol2'] = array([
    [150.00, 1.109],
    [170.00, 1.157],
    [190.35, 1.221],
    [210.06, 1.312],
    [230.35, 1.335],
    [250.27, 1.402],
]) * [V, A] # Specify units
# 8cm radius, pol 1
results[8*cm]['pol1'] = array([
    [150.55, 1.176],
    [171.63, 1.302],
    [191.29, 1.350],
    [209.58, 1.434],
    [230.36, 1.493],
    [250.66, 1.557]
]) * [V, A] # Specify units
# 8cm radius, pol 2
results[8*cm]['pol2'] = array([
```

```
[150.58, 1.360],
]) * [V, A] # Specify units
```

In tabular form:

10cm, pol1

Out[16]:

Volts	Current
150.0	0.966
170.5	1.02
190.35	1.068
210.0	1.161
230.26	1.196
249.97	1.275

10cm, pol2

```
In [17]: display_tabular_data(
        ['Volts', 'Current'],
        results[10*cm]['pol2']
)
```

Out[17]:

Volts	Current
150.0	1.109
170.0	1.157
190.35	1.221
210.06	1.312
230.35	1.335
250.27	1.402

```
In [18]: display_tabular_data(
        ['Volts', 'Current'],
        results[8*cm]['pol1']
)
```

Out[18]:

Volts	Current
150.55	1.176
171.63	1.302
191.29	1.35
209.58	1.434
230.36	1.493
250.66	1.557

8cm, pol2

```
In [19]: display_tabular_data(
        ['Volts', 'Current'],
        results[8*cm]['pol2']
)
```

Out[19]:

Volts	Current
250.14	1.702
230.57	1.634
210.63	1.578
190.7	1.498
170.12	1.38
150.58	1.36

3A – Accepted Values

```
B_E_accepted,
              B_E_accepted_err
          ) = get_B_E(
              year=2016,
              month=2,
              day=3,
              lat=44.2253523,
              long=76.5009076
         Markdown('NOAA reports that $B E$ is locally {:.4e} $\pm$ {:.2e} T'.format(
              B_E_accepted,
              B_E_accepted_err
          ))
Out[20]: NOAA reports that B_E is locally 5.0522e-05 \pm 1.65e-07 T
In [21]: (
              E_M_ratio_accepted,
              _, # Skip
              E M ratio accepted err
          ) = physical_constants['electron charge to mass quotient']
          E_M_ratio_accepted *= -1 # Make it positive instead of negative
         Markdown('SciPy.constants reports that $e/m$ = {:.9e} $\pm$ {:.1e}'.format(
```

Out[21]: SciPy.constants reports that e/m = 1.758820088e+11 \pm 3.9e+03

E_M_ratio_accepted_err

E M ratio accepted,

4 – Analysis

))

In [20]: (

4A – Correct for anode voltage drop

```
In [22]: for radius, data_by_pol in results.items():
    for pol, data in data by pol.items():
```

4B – Calculate B_E , B_T , and e/m

Pair up data from pol1 and pol2 for each diameter

```
In [23]:
         # View into results: results pol1 & pol2 for each diameter
         results_iterator = [
              (diam, data_by_pol['pol2'], data_by_pol['pol1'])
             for diam, data_by_pol in results.items()
         ]
         B E_results = []
         B T results = []
         E_M_{results} = []
         # Loop through each diameter, with pols paired.
         for diam, data 1, data s in results iterator:
             # Select current from data_l, and data_s
             Il = data_l[:,1]
             Is = data_s[:,1]
             # Voltage is the average between the data sets
             V = (data s[:,0]+data 1[:,0])/2
             Be, Bt, em = analyze_paired_data(diam, V, Il, Is)
             # Append to lists
             B_E_results = np.concatenate((B_E_results, Be))
             B T results = np.concatenate((B T results, Bt))
              E M results = np.concatenate((E M results, em))
```

Now take the average of these results, and the error is the standard error on the mean:

```
In [24]: B_E_averaged = -mean(B_E_results)
B_E_averaged_err = stats.sem(B_E_results, ddof=1)

B_T_averaged = mean(B_T_results)
B_T_averaged_err = stats.sem(B_T_results, ddof=1)

E M averaged = mean(E M results)
```

4C - Charge to Mass Ratio

Using the average of all the readings, we get that the charge-to-mass ratio e/m is:

```
In [25]: Markdown("""
The calculated charge to mass ratio is: $e/m_{{calc}} = {em} \\pm {em_d}$\n
The accepted value is $e/m_{{acc}} = {ema} \\pm {ema_d}$""".format(
    em = latex_exp(E_M_averaged),
    em_d = latex_exp(E_M_averaged_err,0),
    ema = latex_exp(E_M_ratio_accepted),
    ema_d = latex_exp(E_M_ratio_accepted_err,0),
))
```

Out[25]: The calculated charge to mass ratio is: $e/m_{calc}=1.92 imes10^{11}\pm2 imes10^{9}$

The accepted value is $e/m_{acc}=1.76 imes 10^{11} \pm 4 imes 10^3$

Which lies outside the uncertainty bounds.

```
In [31]: E_M_percent_err = abs(E_M_averaged-E_M_ratio_accepted)/E_M_ratio_accepted * 100
Markdown(
    'The percent error relative to the accepted value is ${:.1f} \%$'
    .format(E_M_percent_err))
```

Out[31]: The percent error relative to the accepted value is 9.4%

This is a deviation from the expected result.

4D - Local Magnetic Field

```
In [27]: Markdown("""
The calculated local magnetic field of the Earth is: $B_{{E_,calc}} = {be} \pm
{be_d}$\n
The accepted value is $B_{{E_,acc}} = {bea} \pm {bea_d}$""".format(
    be = latex_exp(B_E_averaged,3),
    be_d = latex_exp(B_E_averaged_err,1),
    bea = latex_exp(B_E_accepted,3),
    bea_d = latex_exp(B_E_accepted_err,0),
))
```

Out[27]: The calculated local magnetic field of the Earth is: $B_{E,calc}=0.544 imes10^{-4}\pm0.2 imes10^{-4}$

The accepted value is $B_E~=0~505 imes10^{-4}\pm0 imes10^{-6}$

In [28]: B_E_percent_err = abs(B_E_accepted-B_E_averaged)/B_E_accepted * 100
Markdown('The percent error in \$B_E\$ is: \${:.1f} \%\$ '.format(B_E_percent_err))

Out[28]: The percent error in B_E is: 7.6%

4E – Derivation of Eq. (4)

The following is adapted from Physics Lab Online (http://dev.physicslab.org/document.aspx? doctype=3&filename=magnetism biotsavartlaw2.xml):

We begin with the Bior-Savart Law:

$$dec{B}=rac{\mu_0 Idec{l} imes \hat{r}}{4\pi r^2}$$

We are going to integrate this around each loop for each coil of wire. Our path around the coils will always be perpendicular to r, so the cross product becomes the magnitude of dl:

$$dec{B} = rac{\mu_0 I dec{l}}{4\pi r^2} = rac{\mu_0 I dec{l}}{4\pi (z^2 + R^2)}$$

Since we are considering the central axis of the coils, it is clear that any horizontal contributions will cancel out due to symetry, do we only need to consider the vertical direction:

$$dB_x = \sin(heta)dB = -rac{R}{r}dB = rac{R}{\sqrt{z^2+R^2}}dB$$

$$dB_x = rac{\mu_0 I dec{l}}{4\pi (z^2 + R^2)^{3/2}}$$

Now we integrate around each coil:

$$egin{align} ec{B}_0 &= \int_0^{2\pi n} rac{\mu_0 I dec{l}}{4\pi (z^2 + R^2)^{3/2}} \hat{z} \ &= rac{2\pi n \mu_0 I}{4\pi (z^2 + R^2)^{3/2}} \hat{z} \ &= rac{n \mu_0 I}{2(z^2 + R^2)^{3/2}} \hat{z} \ \end{aligned}$$

Finally, we have two coils, so we can multiply this final answer by two to get the expected result:

$$ec{B}_0 = rac{n \mu_0 I}{\left(z^2 + R^2
ight)^{3/2}} \hat{z}$$

Which can be simplified to $B_0=rac{8n\mu_0I}{5^{3/2}R}=KI$ and rearranged to find:

$$K=rac{8n\mu_0}{5^{3/2}R}$$

```
In [29]: K_{expr} = 8*N*mu_0/(5**1.5 * R)
          K_expr_err = abs(8*N*mu_0/(5**1.5 * R**2)*R_err)
          Latex('$K_{{expr}} = {} \pm {}$'.format(
               latex_exp(K_expr,3),
               latex_exp(K_expr_err,1)
          ))
Out[29]: K_{expr} = 0.759 	imes 10^{-3} \pm 0.2 	imes 10^{-4}
```

The manufacturer value for K is:

```
In [32]: Latex('$K_{{expr}} = {} \pm {}$'.format(
                latex_exp(K,3),
                latex_exp(K_err,1)
Out[32]: K_{expr} = 0.773 	imes 10^{-3} \pm 0.4 	imes 10^{-5}
```

Which does fall within the uncertainties.

The relative error is:

```
rel_err = abs(K_expr-K)/K*100
In [33]:
         Latex('${:.1f}\%$'.format(rel_err))
```

Out[33]: 1.8%