# Data Analysis Exercise

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### **PHY424**

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### 1 Introduction

I analyzed the image "ellipses88.jpg". The sheet of printed paper from which I made the measurements is attached at the end.

# 2 Measurement

## 2.1 Technique

The tools I used are two identical straight rulers, made of metal, with a millimeter scale (the smallest ticks are 1 millimeter apart). The rulers are standard rectangular shape. This is important for the following measurement technique.

To make the measurements, I first made the assumption that the paper is a perfect rectangle, with perfectly straight edges, and the horizontal and vertical axis of the ellipses are perfectly aligned with the edges of the paper when they are printed. A quick examination of the paper attached makes it clear that this is a good assumption, and any error about this assumption is clearly negligible compared to the reading error that follows.

Next, to find the horizontal (or vertical) axis of an ellipse, I needed to align my ruler with two furthest points of the ellipse. I placed the first ruler (I call this the measuring ruler) at the approximate position. Then I take the second ruler (I call this the aligning ruler), align its shorter side with the side of the paper, so that it is perpendicular to the measuring ruler. I then press down on the aligning ruler to fix it on the paper, and slide the measuring ruler along it. This way, I can see the ellipse gradually getting more (or less) covered up by the ruler as I slide. When I see that the amount of ellipse that is covered up by the ruler stops increasing, and just before it starts decreasing, I must be the closest to having my measuring ruler aligned with the horizontal (or vertical) axis of the ellipse. Then I read off the measurement. This method requires that the rulers have perfect 90 degrees angle. Again, this is so accurate that any error is negligible compared to the reading error.

I repeated this measurement for a series of 12 ellipses as well as the calibration square.

### 2.2 Uncertainties

When I read off the reading on the ruler, I estimated the one standard deviation error coming from the human reading error and the blurriness of the edge of the ellipse on the printed paper. This uncertainties are mostly the same for the various ellipse, except for ellipse 1 and 2, whose y- dimensions are especially thin and make the corresponding errors larger.

# 3 Analysis

## 3.1 Initial Data Processing

For each ellipse, I have measured its x- and y- dimensions and estimated their measurement uncertainties. Let's denote these values as  $x_{meas}$ ,  $dx_{meas}$ ,  $dx_{meas}$ ,  $dy_{meas}$ . I also measured the dimensions (with uncertainties) of the calibration square, which converts the measurement unit to the unit used by the given equation. Let's denote these calibration measurements as  $c_x$ ,  $dc_x$ ,  $c_y$ ,  $dc_y$ . Let's denote the calibration square in unit used by the equation as  $c'_x$  and  $c'_y$ , which are exact values known a priori.

The to find (x, y) in the equation unit, I make the calculation

$$x = x_{meas} \cdot s \cdot \frac{c_x'}{c_x} \tag{1}$$

$$y = y_{meas} \cdot s \cdot \frac{c_y'}{c_y} \tag{2}$$

where s is the scale (an exact value) associated with each ellipse. The uncertainty in x and y is a combination of the measurement uncertainty and the calibration uncertainty (which is a systematic uncertainty) via quadrature:

$$dx = s \cdot c_x' \cdot d\left(\frac{x_{meas}}{c_x}\right) \tag{3}$$

$$dx = s \cdot c_x' \cdot \sqrt{\frac{1}{c_x^2} dx^2 + \frac{x_{meas}^2}{c_x^4} dc_x^2}$$
 (4)

and

$$dy = s \cdot c'_y \cdot \sqrt{\frac{1}{c_y^2} dy^2 + \frac{y_{meas}^2}{c_y^4} dc_y^2}$$
 (5)

I made all of the above calculation in the module "process\_data.py" (which also includes the raw measurement data) and copied and pasted all the result of x, dx, y, dy into the text file "ellipse\_data".

#### 3.2 Data Fitting

I first took the sample code "odr\_fit\_to\_data.py" published on the course website. I coded in the three distribution functions of interest: Lorentizian, Log-Normal, and Absolute Sinc. Next, I fitted the data against each of the three fit function, and adjusted the guess parameters based on the program output.

After some trials, I found that if I use all the data, it approximately fits the absolute Sinc function, with the parameters:

$$y_{\mu} = 110 \pm 16, \mu = 24.0 \pm 0.6, \sigma = 4.1 \pm 0.2$$
 (tentative values)

where the absolute Sinc function is given by

$$y(x) = y_{\mu} \left| \frac{\sin(x-\mu)/\sigma}{(x-\mu)/\sigma} \right| \tag{6}$$

A plot of this fit is given below in figure 1. The  $\chi^2$  per degree of freedom is 9.18, which is a reasonable number. However, the corresponding CDF is 0%. In other words, the program is saying that it is impossible to get such a good fit.

I quickly found the cause of the problem. Looking at figure 1, it is clear that most of the data points lie relatively close to the fitted curve, with one exceptional outlier, which is the value with the largest x. The residual is a long horizontal line, which shows how far the point is from the curve. However, the algorithm clearly runs into a problem here. The point should be fitted to the second bump just below it, instead of fitting it to the central bump, which is much further to the left. This program fits the data using the

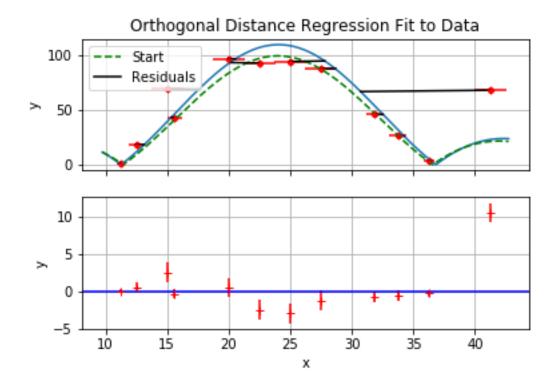


Figure 1: A fit using the absolute Sinc function

orthogonal distance regression, which tries to minimize each point to a "closest" point on the curve. It is know that this algorithm can mistaken which point on the curve is supposed to be the closest.

To remedy this problem, I reran the fitting program but took out the outlier data point. The new fit is shown in figure 2.

This time, my results are

$$y_{\mu} = 102 \pm 5, \mu = 23.7 \pm 0.2, \sigma = 4.0 \pm 0.1$$
 (outlier-adjusted values)

The  $\chi^2$  per degree of freedom is 1.4, which is much smaller than before and indicates a good fit. The CDF is 18.6%, which meas there is a 18.6% chance that the  $\chi^2$  per degree of freedom we found is true. This is a reasonable result. From the plot, the fit is mostly a good fit.

The other two models, Lorentizian and Log-Normal, do not fit the data at all. The fitted parameters have uncertainties of  $10^8$ ; it is unneeded to show the corresponding plots to know that these models are not related to the data.

# 4 Conclusion

Overall, the fit is rather good after I took out the point that not only is an outlier relative to the rest of the data, but also a point that the orthogonal distance regression misses in finding the closest point of the curve. Based on the plots, the uncertainties I estimated were reasonable. If I had more time, I would use a different program to fit the data so the issue with orthogonal distance would not arise.

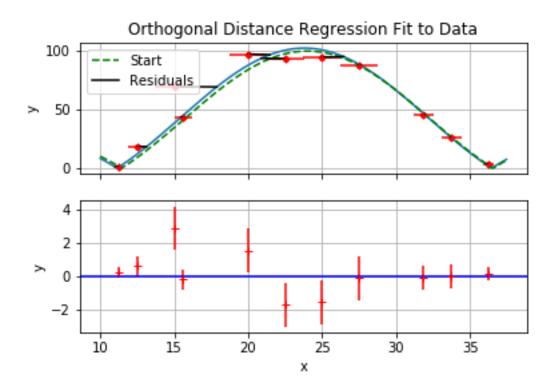


Figure 2: A fit using the absolute Sinc function, but with the outlier removed.

```
# -*- coding: utf-8 -*-
File Name: process_data.py
Purpose: data analysis exercise
Author: Samuel Wong
import numpy as np
scale=np.array([2,2,1,1,2,1,2,2,2,1,1,1]) #exact
#all units below are in cm
#sigma are the uncertainty (1 std deviation)
#y measurements and measurement uncertainty
y meas=np.array([2.65,3.55,2.00,3.50,3.60,3.30,2.60,3.35,3.70,1.40,0.08,0.25])
#x measurements and measurement uncertainty
x_{meas}=np.array([0.60,0.90,2.70,2.55,1.00,1.25,1.65,1.10,0.80,1.00,0.90,2.90])
#calibration and its uncertainty
calibration x=2.00 #25 unit
sigma cal x = 0.02
calibration y=1.90 #25 unit
sigma_cal_y=0.02
#convert to (x,y) unit using the formula:
# unit = measured*scale*(calibration units/calibration)
x = x_meas*scale*(25/calibration_x)
y = y meas*scale*(25/calibration y)
#calculate total uncertainty using quadrature
dx = scale*25*np.sqrt((1/calibration_x**2)*(sigma_x_meas**2) +
                  (x meas**2/calibration x**4)*sigma cal x**2)
dy = scale*25*np.sqrt( (1/calibration_y**2)*(sigma_y_meas**2) +
                  (y_meas**2/calibration_y**4)*sigma_cal_y**2)
for i in range(x.size):
   round(y[i],2), round(dy[i],1)))
```

```
odr_fit_to_data.py
   A simple program to fit x,y data with uncertainties in both x & y
   Uses Orthogonal Distance Regression
   The examples provided are fits of Gaussian, Lorentzian, or linear
         functions to various data files.
   To fit your own data, you need to change:
    (1) func(p,x) to return the function you are trying to fit,
            p is the parameter vector,
            x are the independent variable(s)
            Caution: scipy.odr and scipy.optimize.curve_fit require
                 x & p in opposite orders.
    (2) the name of the data file read in by numpy.loadtxt,
    (3) the initial guesses (p guess) for the fit parameters. The fit
             will not converge if your guess is poor.
   For a more accurate estimate of the Cumulative Distribution
        Function (CDF) by Monte Carlo simulation based on the fitted
        values, set
            Run Monte Carlo CDF Estimator = True
            Number_of_MC_iterations
                                          = 1000
        But note that this can take some time (e.g. > 1s per
            iteration) if the x uncertainties are large.
    For information on scipy.odr, see
   http://docs.scipy.org/doc/scipy/reference/odr.html
   Copyright (c) 2011-2017 University of Toronto
   Modifications:
        7 January 2018 by David Bailey
            Minor bug fixes, e.g. Cauchy example incorrect
        6 August 2017 by David Bailey
            Fixed some Python 2 vs 3 incompatibility, e.g. removed zip
        14 April 2017 by David Bailey
            Fixed Monte Carlo section errors introduced in last two
            modifications. Some reformating and extra comments.
        7 February 2016 by David Bailey
            Made notation and uncertainties compatible with
            curve fit to data, in particular prevented uncertainty
            estimates being reduced if fit is "too good'
        7 January 2016 by David Bailey
            Made printing python 3 compatible
        11 January 2013 by David Bailey
   Original Version
                     : 11 September 2011 by Michael Luzi
   Contact: David Bailey <dbailey@physics.utoronto.ca>
                (www.physics.utoronto.ca/~dbailey)
    License: Released under the MIT License; the full terms are
                appended to the end of this file, and are also found
                at www.opensource.org/licenses/mit-license.php
# Use python 3 consistent printing and unicode
from __future__ import print_function
```

from \_\_future\_\_ import unicode\_literals

```
import inspect
                             # docs.python.org/library/inspect.html
import matplotlib
from matplotlib import pyplot # matplotlib.org/api/pyplot_api.html
import numpy
                             # numpy.scipy.org/
                             # scipy.org/
import scipy
import scipy.odr, scipy.special, scipy.stats
# Decide if Monte Carlo CDF values are desired
Run Monte Carlo CDF Estimator = True
# The default is set to just 11, so all the examples will run quickly
# in a few seconds. For serious work, this should be much larger.
Number_of_MC_iterations
# You can ignore x uncertainties if desired
x_uncertainties_ignored = False
## define some possible functions to be fitted
## You may, of course, define your own.
def Cauchy(p,x) :
    # Cauchy / Lorentzian / Breit-Wigner peak with contant background:
               = p[0] # Constant Background
               = p[1]
                       # Height above/below background
    x1_peak
                       # Central value
    x1
               = p[2]
                       # Half-Width at Half Maximum
    width_x1
               = p[3]
    return ( B + x1 peak*(1/(1+((x-x1)/width x1)**2)) )
def Gauss(p,x):
    # A gaussian or Normal peak with constant background:
              = p[0] # Constant Background
    x1_peak
               = p[1]
                       # Height above/below background
               = p[2]
                         # Central value
    x1
                       # Standard Deviation
    width_x1
              = p[3]
    return ( B + x1_{peak*numpy.exp(-(x-x1)**2/(2*width_x1**2))} )
def Semicircle(p,x) :
    # The upper half of a circle:
               = p[0] # Radius of circle
    R
    x0
               = p[1] # x coordinate of centre of circle
                       # y coordinate of centre of circle
               = p[2]
    from numpy import array, sqrt
    y=[]
    for i in range(len(x)) :
        y_{squared} = R^{**2} - (x[i] - x0)^{**2}
        if y_squared < 0 :</pre>
           y.append(y0)
        else:
           y.append(sqrt(y_squared)+y0)
    return array(y)
def Linear(p,x) :
    # A linear function with:
    # Constant Background
                                    : p[0]
       Slope
                                    : p[1]
    return p[0]+p[1]*x
def Quadratic(p,x) :
    # A quadratic function with:
    # Constant Background
                                    : p[0]
   #
       Slope
                                    : p[1]
    # Curvature
                                    : p[2]
    return p[0]+p[1]*x+p[2]*x**2
```

```
def Cubic(p,x) :
    # A cubic function with:
    # Constant Background
                                     : p[0]
    # SLope
                                     : p[1]
    # Curvature
                                     : p[2]
        3rd Order coefficient
                                     : p[3]
    return p[0]+p[1]*x+p[2]*x**2+p[3]*x**3
def Lorentzian(p,x):
    y_mu = p[0]
    mu = p[1]
    sigma = p[2]
    return y_mu/(1+ (x-mu)**2/sigma**2)
def Log_Normal(p,x):
    y_mu = p[0]
    mu = p[1]
    sigma = p[2]
    denominator = 2*sigma**2
    numerator = -(\text{numpy.log}(x) - \text{numpy.log}(mu))**2
    return y mu*numpy.exp(numerator/denominator)
def Sinc(p,x):
    y_mu = p[0]
    mu = p[1]
    sigma = p[2]
    numerator = numpy.sin((x-mu)/sigma)
    denominator = (x-mu)/sigma
    return y mu*numpy.abs(numerator/denominator)
## Choose function and data to fit:
func = Sinc
if func.__name__ == "Gauss" :
    # Initial guesses for fit parameters
    p_{guess} = (10.,400.,1173.9,0.4,)
    # Gaussian data with large x uncertainties
        This data shows how ODR can fit data that would give great
         trouble to regular chi-squared fits (that ignore x errors).
    data_file = "gauss_xy_errors_ugly.txt"
    # Labels for plot axes
    x label = "energy [KeV]"
    y_label
            = "Counts"
elif func.__name__ == "Semicircle" :
    p_{guess} = (1,1,1)
    # This data is a semicircle, and shows ODR doing the best it can
         with data that is a poor match to the fitting function
    data file = "semicircle.txt"
    x_{abel} = x''
    y_label
             = "y"
elif func.__name__ == "Cauchy" :
    p_{guess} = (0,1,2,1)
    # This data is actually generated from a semicircle distribution,
    # but the fit to a Cauchy is not bad, showing that the same data
       can sometimes be well fit by different distributions.
    data file = "semicircle.txt"
    x label = "x"
    y_label = "y"
elif func.__name__ == "Quadratic" :
    p_{guess} = (0,1,2)
    # This data is parabola
```

```
data file = "parabola xy errors.txt"
    x_label = "x"
            = "y"
    y_label
elif func.__name__ == "Cubic" :
    p_{guess} = (0,1,0,0)
    # This data is a cubic
    data file = "cubic data.txt"
    x_label = "x"
    y_label = "v"
elif func.__name__ == "Lorentzian":
    p_guess = (30,30,10)
    data_file = "ellipse_data.txt"
    x label = "x"
    y_label = "y"
elif func.__name__ == "Log_Normal":
    p_{guess} = (100, 30, 20)
    data file = "ellipse data.txt"
    x_label = "x"
    y_label = "y"
elif func.__name__ == "Sinc":
    p_guess = (100, 24, 4)
    data_file = "ellipse_data.txt"
    x label = "x"
    y_label = "y"
else :
    # default is linear
    func=Linear
    p_guess = (2,2)
    # Linear data with large x uncertainties
    data file = "linear xy errors.txt"
    x_label = "x"
    y_label
            = "y"
## Load data to fit
      Any lines in the data file starting with '#' are ignored
      "data" are values, "sigma" are uncertainties
x_data, x_sigma, y_data, y_sigma = numpy.loadtxt( data_file,
                                                   comments='#',
                                                   unpack = True)
if x uncertainties ignored:
    # x uncertainties cannot be set exactly to zero without crashing the
        fit, but a tiny value seems to do the job.
    x_sigma = len(x_data)*[1e-99]
# Load data for ODR fit
data = scipy.odr.RealData(x=x_data, y=y_data, sx=x_sigma, sy=y_sigma)
# Load model for ODR fit
model = scipy.odr.Model(func)
## Now fit model to data
   job=10 selects central finite differences instead of forward
        differences when doing numerical differentiation of function
   maxit is maximum number of iterations
   The scipy.odr documentation is a bit murky, so to be clear note
         that calling ODR automatically sets full output=1 in the
         low level odr function.
fitted = scipy.odr.ODR(data, model, beta0=p_guess, maxit=5000,job=10)
```

```
output = fitted.run()
                     # 'beta' is an array of the parameter estimates
p = output.beta
cov = output.cov_beta # parameter covariance matrix
#"Quasi-chi-squared" is defined to be the
           [total weighted sum of squares]/dof
#
   i.e. same as numpy.sum((residual/sigma odr)**2)/dof or
#
      numpy.sum(((output.xplus-x)/x sigma)**2
                               +((y_data-output.y)/y_sigma)**2)/dof
   Converges to conventional chi-square for zero x uncertainties.
quasi chisq = output.res var
uncertainty = output.sd_beta # parameter standard uncertainties
  scipy.odr scales the parameter uncertainties by quasi chisq.
    If the fit is poor, i.e. quasi_chisq is large, the uncertainties
    are scaled up, which makes sense. If the fit is too good,
#
    i.e. quasi_chisq << 1, it suggests that the uncertainties have
#
    been overestimated, but it seems risky to scale down the
#
    uncertainties. i.e. The uncertainties shouldn't be too sensistive
    to random fluctuations of the data, and if the uncertainties are
    overestimated, this should be understood and corrected
if quasi chisq < 1.0 :</pre>
   uncertainty = uncertainty/numpy.sqrt(quasi_chisq)
if False: # debugging print statements
   print("sd_beta",output.sd_beta)
   print("cov beta",output.cov beta)
   print("delta",output.delta)
   print("epsilon",output.epsilon)
    print("res_var",output.res_var)
   print("rel_error",output.rel_error)
ORTHOGONAL DISTANCE REGRESSION")
print("
# print out version information for debugging
import sys
print("Python", sys.version)
print("Scipy:",scipy.__version__,", Numpy:",numpy.__version__,
             ", Matplotlib:",matplotlib.__version__, sep="")
print("****************
print("ODR algorithm stop reason: " + output.stopreason[0])
print("\nFit {0} Data points from file: {1}"
                                  .format(len(x data),data file))
print("To Model :")
print(" ",inspect.getsource(func))
if x uncertainties ignored:
   print("** WARNING: x uncertainties set to zero in fit. **\n")
print("Estimated parameters and uncertainties")
for i in range(len(p)) :
   print(("
             p[{0}] = {1:10.5g} +/- {2:10.5g}"+
                     (Starting guess: {3:10.5g})").\
           format(i,p[i],uncertainty[i],p guess[i]))
# number of degrees of freedom for fit
dof = len(x_data) - len(p_guess)
# Note: odr, unlike curve fit, returns the standard covariance matrix.
```

```
print("\nStandard Covariance Matrix : \n", cov, "\n")
print("\nCorrelation Matrix :")
for i,row in enumerate(cov):
   for j in range(len(p)) :
        print("{0:< 8.3g}".</pre>
               format(cov[i,j]/numpy.sqrt(cov[i,i]*cov[j,j])), end="")
            # Newbie Notes: "{0:< 8.3q}" left justifies output with
                space in front of positive numbers, with 3 sig figs;
               end="" suppresses new line.
   print()
# Calculate initial residuals and adjusted error 'sigma odr'
                  for each data point
delta = output.delta # estimated x-component of the residuals
epsilon = output.eps
                       # estimated y-component of the residuals
# (dx star, dy star) are the projections of x sigma and y sigma onto
     the residual line, i.e. he differences in x & y between the data
     point and the point where the orthogonal residual line intersects
     the ellipse' created by x sigma & y sigma.
dx_star = ( x_sigma*numpy.sqrt( ((y_sigma*delta)**2) /
                ((y_sigma*delta)**2 + (x_sigma*epsilon)**2))
dy star = ( y sigma*numpy.sqrt( ((x sigma*epsilon)**2) /
                ((y_sigma*delta)**2 + (x_sigma*epsilon)**2))
sigma odr = numpy.sqrt(dx star**2 + dy star**2)
# residual is positive if the point lies above the fitted curve,
              negative if below
residual = ( numpy.sign(y data-func(p,x data))
              * numpy.sqrt(delta**2 + epsilon**2) )
print(("\nQuasi Chi-Squared/dof
                                = \{0:10.5f\}," +
                            " Chi-Squared CDF = \{1:10.5f\}\%").
       format(quasi chisq,
              100.*float(scipy.special.chdtrc(dof,dof*quasi chisq))))
print("
         WARNING: Above CDF is not valid for large x uncertainties!")
print("\nTo use Monte Carlo simulation to more accurately estimate")
print('
              CDF for large x uncertainties, re-run program with ')
print('
            "Run_Monte_Carlo_CDF_Estimator = True" and')
            "Number of MC iterations >= 1000."', end="")
print(
print(' This may take some time\n')
print("Run Monte Carlo CDF Estimator = {0}"\
                  .format(Run Monte Carlo CDF Estimator))
if Run_Monte_Carlo_CDF_Estimator :
    print("\n**** Running Monte Carlo CDF Estimator ****")
   print("Number of MC iterations = {0}"
                .format(Number_of_MC_iterations), end="")
   # Initialize Monte Carlo output distributions
                     = Number_of_MC_iterations*[None]
   x dist
                     = Number of MC iterations*[None]
   y dist
                    = Number of MC iterations*[None]
   p dist
   quasi_chisq_dist = Number_of_MC_iterations*[None]
   # Initialize timing measurement
   import time
    start time = time.clock()
```

```
for i in range(Number of MC iterations) :
    # Starting with the x and x uncertainty (x sigma) values from
         the data, calculate Monte Carlo values assuming a normal
         gaussian distibution.
    x dist[i] = numpy.random.normal(x data,x sigma)
    # Calculate y using Monte Carlo x, then smear by y uncertainty
    y dist[i] = numpy.random.normal(func(p,x data),y sigma)
    # Fit the Monte Carlo x,y pseudo-data to the original model
    data dist = scipy.odr.RealData( x=x dist[i], y=y dist[i],
                                    sx=x sigma, sy=y sigma)
    model dist = scipy.odr.Model(func)
    fit_dist = scipy.odr.ODR(data_dist, model_dist, p, maxit=5000,
                                    job=10)
    output_dist = fit_dist.run()
    p dist[i] = output dist.beta
    quasi_chisq_dist[i] = output_dist.res_var
end time = time.clock()
print(" simulations in {0} seconds.".\
      format(end_time-start_time))
# Sort the simulated quasi-chi-squared values
quasi_chisq_sort = numpy.sort(quasi_chisq_dist)
# Find the index of the sorted simulated quasi-chi-squared value
# nearest to the data quasi-chi-squared value, to estimate the cdf
print("\nFraction of quasi-chisquared values larger" +
           " than observed value:")
print("
           Monte Carlo CDF = {0:6.1f}%".format(
        100.*(1.0-numpy.abs(quasi chisq sort-quasi chisq).argmin()
                               /float(Number_of_MC_iterations))))
print("
           Minimum, Mean, and Maximum Quasi Chi-Squared values:",
                            end="")
print("{0:6.2g} {1:6.2g} {2:6.2g}".\
      format(numpy.min(quasi_chisq_dist),
                     numpy.average(quasi chisq dist),
                     numpy.max(quasi chisq dist)))
print("\nAverage and Standard Deviation of MC Fit parameters")
ave_p = numpy.average(numpy.array(p_dist),axis=0)
std p = numpy.std(
                      numpy.array(p dist),axis=0)
min_p = numpy.min(
                      numpy.array(p_dist),axis=0)
                      numpy.array(p dist),axis=0)
max p = numpy.max(
for i in range(len(p)) :
              p[\{0\}] = \{1:12.6g\} +/- \{2:12.6g\};"
    print (("
            + " (Min = {3:12.6f}, Max = {4:12.6f}").format(
                  i, ave p[i], std p[i], min p[i], max p[i] ))
print("\nCheck for any Fit Biases in MC Fit parameters'
        +"(Significance and ratio)")
# Check if fit on Monte Carlo data tends to give an average output
    value for the parameters different from the input parameters.
for i in range(len(p)) :
    print("
              p[\{0\}] = \{1: < 6.2f\}  (\{2: <12.7g\}+\/-\{3: <12.7g\}\)"
          .format(i, (ave_p[i]/p[i]-1)/
                  (std_p[i]/p[i]/numpy.sqrt(Number_of_MC_iterations-1)),
                  ave_p[i]/p[i],
                  std_p[i]/p[i]/numpy.sqrt(Number_of_MC_iterations-1)))
```

```
# create figure with light gray background
fig = pyplot.figure(facecolor="0.98")
# 3 rows, 1 column, subplot 1
# 3 rows are declared, but there are only 2 plots; this leaves room
         for text in the empty 3rd row
fit = fig.add subplot(211)
# remove tick labels from upper plot (for clean look)
fit.set xticklabels(())
pyplot.ylabel(y label)
pyplot.title("Orthogonal Distance Regression Fit to Data")
# Plot data as red circles, and fitted function as (default) line.
# For a smooth look, generate many x values for plotting the model
stepsize = (\max(x \text{ data}) - \min(x \text{ data}))/1000.
margin = 50*stepsize
x model = numpy.arange( min(x data)-margin,max(x data)+margin,
                                stepsize)
fit.plot(x_data,y_data,'ro', x_model, func(p,x_model),markersize=4)
# Add error bars on data as red crosses.
fit.errorbar(x_data, y_data, xerr=x_sigma, yerr=y_sigma, fmt='r+')
fit.set_yscale('linear')
    draw starting guess as dashed green line ('r-')
fit.plot(x_model, func(p_guess,x_model), 'g-', label="Start",
                        linestyle="--")
\# output.xplus = x + delta
a = numpy.array([output.xplus,x_data])
\# output.y = f(p, xfit), or y + epsilon
b = numpy.array([output.y,y data])
fit.plot( numpy.array([a[0][0],a[1][0]]),
          numpy.array([b[0][0],b[1][0]]),
          'k-', label = 'Residuals')
for i in range(1,len(y data)):
    fit.plot( numpy.array([a[0][i],a[1][i]]),
              numpy.array([b[0][i],b[1][i]]),
              'k-')
fit.legend(loc='upper left')
fit.grid()
# separate plot to show residuals
residuals = fig.add subplot(212) # 3 rows, 1 column, subplot 2
residuals.errorbar(x=x data,y=residual,yerr=sigma odr,
                            fmt="r+", label = "Residuals")
# make sure residual plot has same x axis as fit plot
residuals.set_xlim(fit.get_xlim())
# Draw a horizontal line at zero on residuals plot
pyplot.axhline(y=0, color='b')
# Label axes
pyplot.xlabel(x label)
# These data look better in 'plain', not scientific, notation, and if
# the tick labels are not offset by a constant (as done by default).
pyplot.ticklabel format(style='plain', useOffset=False, axis='x')
pyplot.ylabel(y_label)
residuals.grid()
pyplot.savefig(func.__name__+"_Fit_Outlier_Adjusted.png")
```

### pyplot.show()

##### END of odr\_fit\_to\_data.py

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