Electronic Supplementary Material (ESI) for Physical Chemistry Chemical Physics.

Gas Phase Complexes of $H_3N\cdots CuF$ and $H_3N\cdots CuI$ Studied by Rotational Spectroscopy and Ab Initio Calculations: The Effect of X (X = F, Cl, Br, I) in $OC\cdots CuX$ and $H_3N\cdots CuX$

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Supplementary Information: High Resolution Fourier Transform Window Function

The program code used to apply the high resolution Fourier transform window function is shown below (as Python script).

```
def f1(x):
        return (1/(numpy.exp(1/x)-1))

def f2(x):
        return (1/x**4.5)

def f(x):
        return f1(x)*f2(x)

def high_res_window(Npoints):
        x = numpy.arange(0,1,1./(Npoints))
        y = f(x)
        y[0] = 0
        maximum = max(y)
        y = y/maximum
        plateau_start = numpy.argwhere(y == max(y))
        y[plateau_start:Npoints] = max(y)
        return y
```