Supplementary Information

Figure S1. IR spectra of compound **1**.

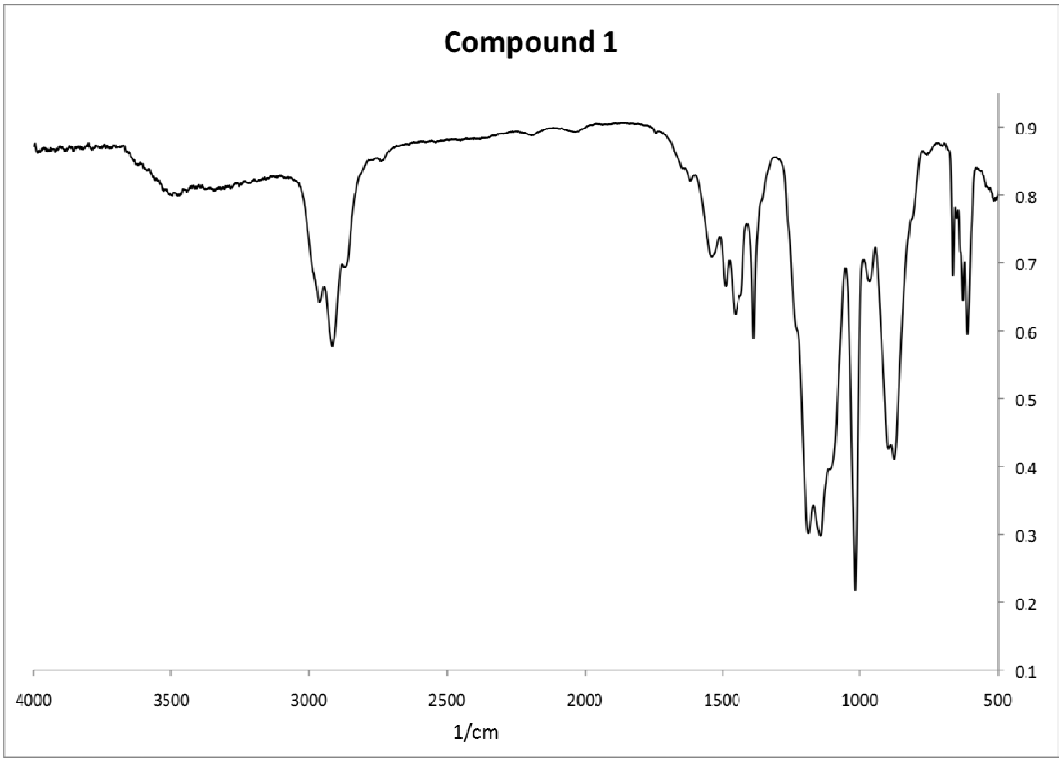

Figure S2. TGA analysis of compound **1**.

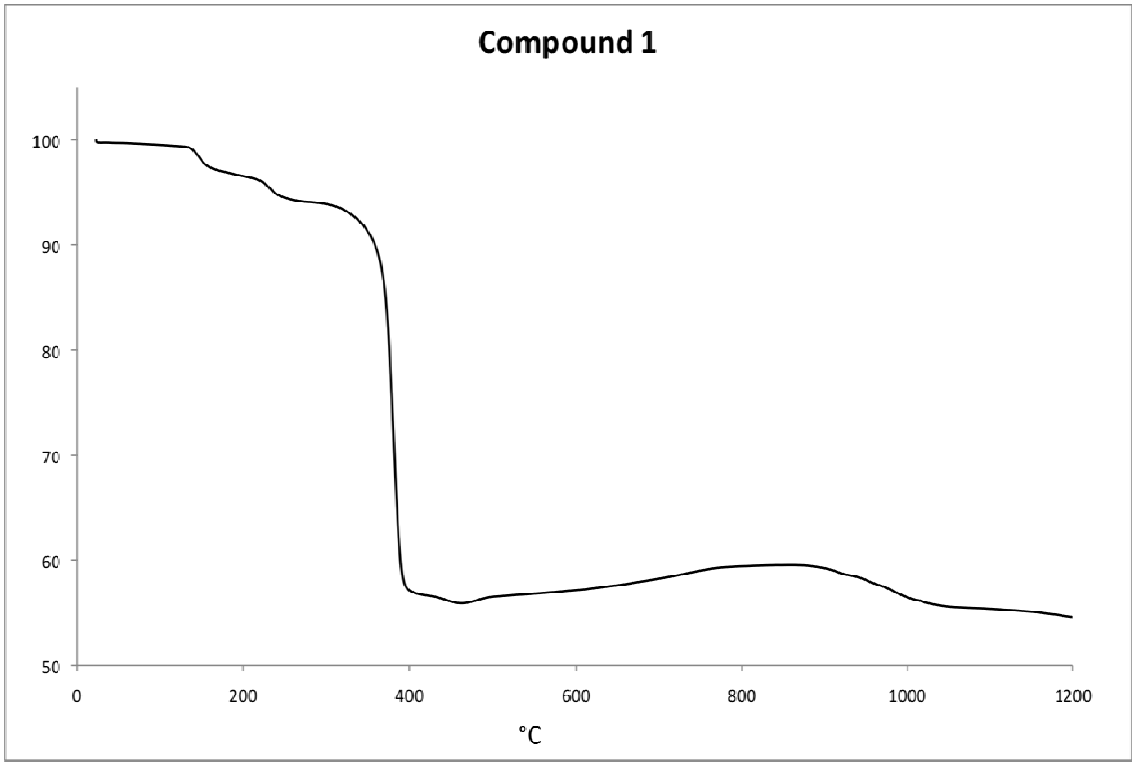

Table S1. Report of the Pawley Refinement.

Pawley method	
Phase name	hkl_Phase
R-Bragg	81.022
Spacegroup	P21/c
Scale	0.814354931
Cell Mass	0
Cell Volume (┼^3)	1143.28167
Wt% - Rietveld	0
Strain	
Strain L	0.0001
Strain G	0.0001
Additional convolutions	
Circles, Constant	0.04765229
Circles, Constant	0.1297621
Hat, Constant	0.0001
Hat, Constant	0.0001
Gaussian, Constant	0.0001
PV_TCHZ peak type	
U	−0.01383141
V	0.00560477
W	−0.00196705
Z	0
X	0.0001
Y	0
Lattice parameters	
a (┼)	11.5957814
b (┼)	10.384685
c (┼)	10.4925736
beta (▒)	115.1966