

## **ELECTRONIC SUPPLEMENTARY INFORMATION**

Donor Disposition and Aliphatic Conformation Effects on Structure in Luminescent Zinc  
Dicarboxylate Coordination Polymers with Dipyridylamide Coligands

Jacob W. Uebler, Julie A. Wilson, Robert L. LaDuca<sup>†\*</sup>

Table S1. Hydrogen bonding parameters for **1–6**.

$D-H\cdots A$	$d(H\cdots A)$	$d(D\cdots A)$	$\angle DHA$	symmetry transformation for $A$
<b>1</b>				
O1W–H1WA $\cdots$ O5	2.43(2)	3.223(3)	155(3)	
O1W–H1WB $\cdots$ O5	2.077(19)	2.879(3)	156(3)	$-x,-y,-z$
N2–H2N $\cdots$ O1W	2.05	2.829(3)	146.7	$-x+1/2,y+1/2,-z+1/2$
<b>2</b>				
N2–H2N $\cdots$ O4	1.95(2)	2.803(4)	156(4)	$x,-y+3/2,z+1/2$
<b>3</b>				
N3–H3 $\cdots$ O2	1.90	2.755(4)	162.4	$-x+1/2,-y+1/2,-z+1$
N3A–H3A $\cdots$ O2	1.90	2.777(10)	174.4	$-x+1/2,-y+1/2,-z+1$
<b>4</b>				
O5–H5A $\cdots$ O1	2.133(17)	2.951(2)	169(3)	$-x+2,-y,-z+1$
N3–H3N $\cdots$ O2	1.939(18)	2.787(3)	166(3)	$-x+1,-y,-z+1$
<b>5</b>				
N2–H2 $\cdots$ O1	2.14	2.885(18)	142.1	$-x,-y+1,-z+1$
N2A–H2A1 $\cdots$ O1	2.07	2.79(2)	138.9	$-x,-y+1,-z+1$
<b>6</b>				
N2–H2N $\cdots$ O3	2.187(17)	3.025(2)	160(2)	

Figure S1. Stacking pattern of  $[\text{Zn}_3(\text{mal})_2(\text{OH})_2(4\text{-pina})_2]_n$  layers in **4**. Hydrogen bonding is shown as dashed lines.

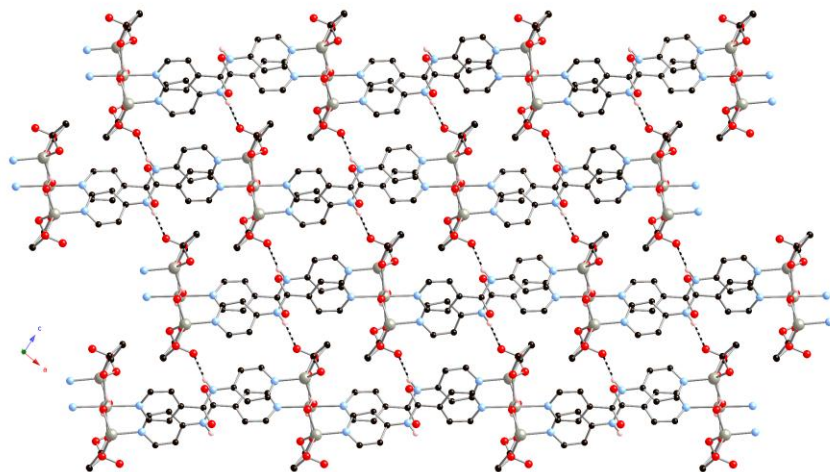


Figure S2. Stacking of  $[\text{Zn}(\text{suc})(4\text{-pna})]_n$  layers in **6**. Hydrogen bonding is shown as dashed lines.

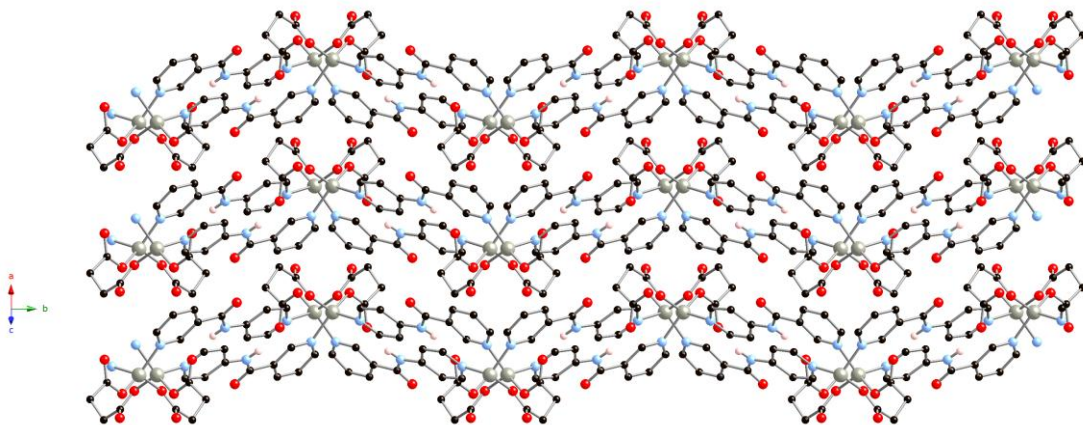


Figure S3. TGA plot for **1**.

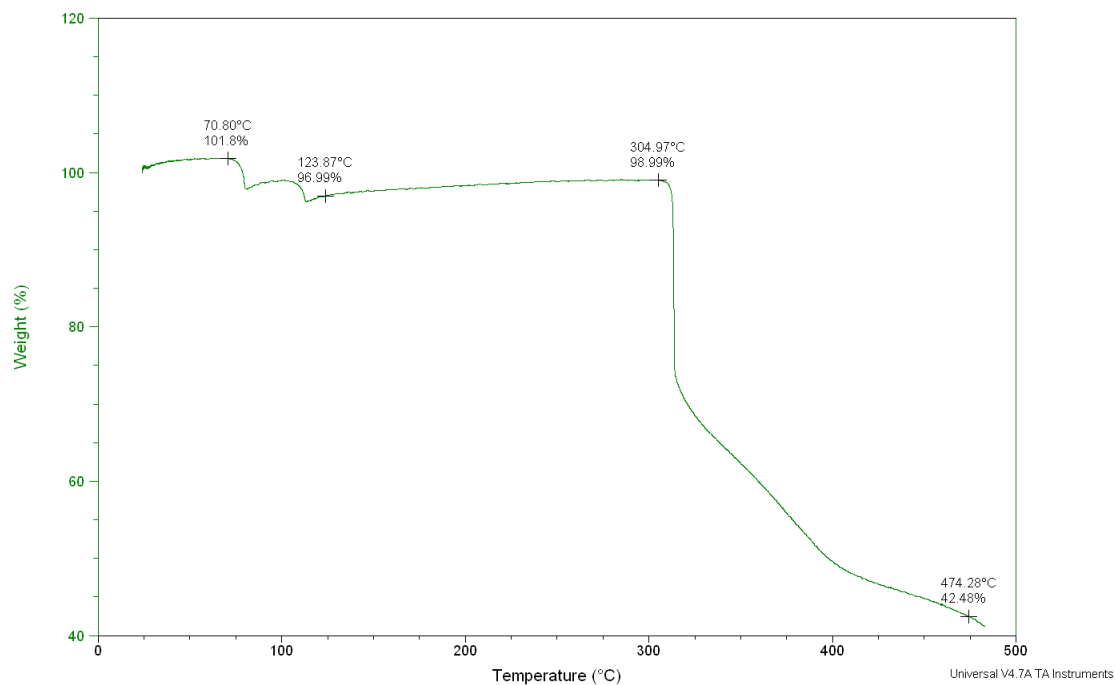


Figure S4. TGA plot for **2**.

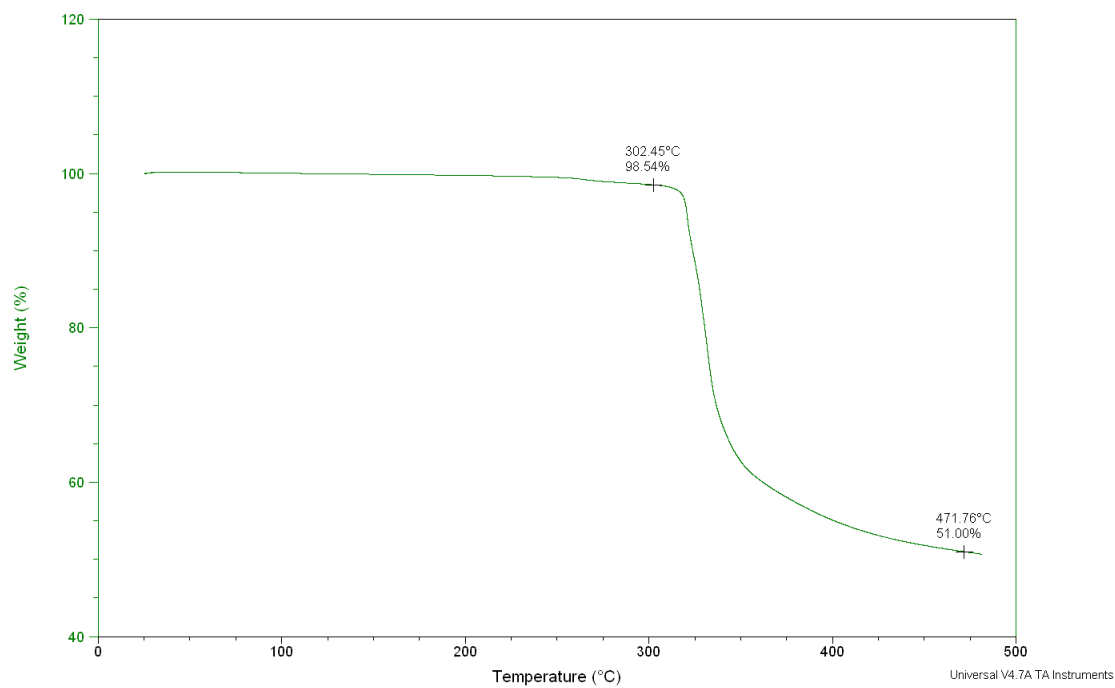


Figure S5. TGA plot for **3**.

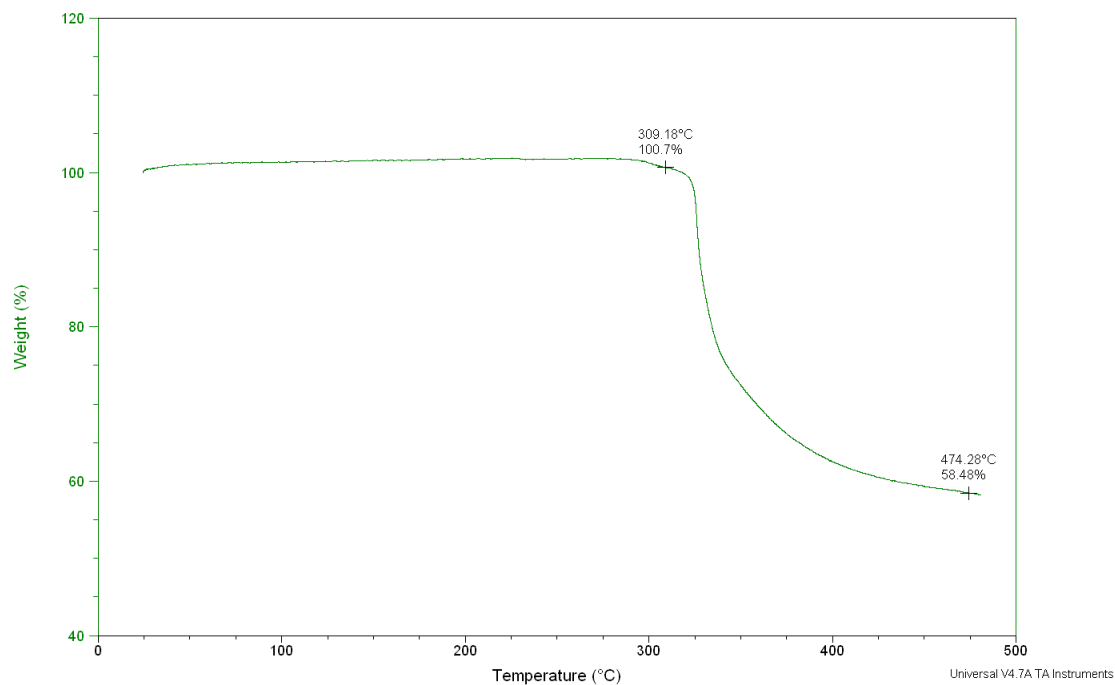


Figure S6. TGA plot for **4**.

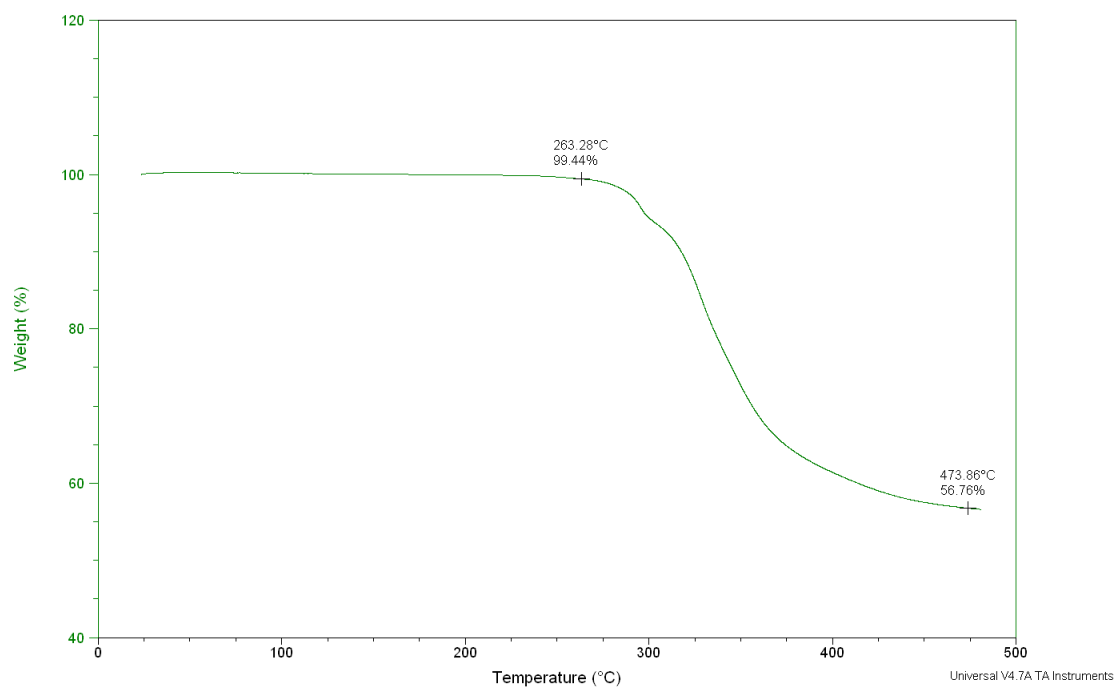


Figure S7. TGA plot for **5**.

