**Supporting Information**

<table>
<thead>
<tr>
<th>Table S1. Summary of X-ray Crystallography Data of 2b</th>
</tr>
</thead>
<tbody>
<tr>
<td>Emp. Formula</td>
</tr>
<tr>
<td>Form. Weight</td>
</tr>
<tr>
<td>Crystal system</td>
</tr>
<tr>
<td>Space group</td>
</tr>
<tr>
<td>a (Å)</td>
</tr>
<tr>
<td>b (Å)</td>
</tr>
<tr>
<td>c (Å)</td>
</tr>
<tr>
<td>α α α</td>
</tr>
<tr>
<td>α α α</td>
</tr>
<tr>
<td>α α α</td>
</tr>
<tr>
<td>Volume (Å$^3$)</td>
</tr>
<tr>
<td>Z</td>
</tr>
<tr>
<td>Density</td>
</tr>
<tr>
<td>F(000)</td>
</tr>
<tr>
<td>Crystal size(mm$^3$)</td>
</tr>
<tr>
<td>Index ranges</td>
</tr>
<tr>
<td>Ref. collected</td>
</tr>
<tr>
<td>Ind. Reflections</td>
</tr>
<tr>
<td>Complete to</td>
</tr>
<tr>
<td>Max. and min transmission</td>
</tr>
<tr>
<td>Data / Restraints/ parameters</td>
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<tr>
<td>GOF</td>
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<td>Final R indices</td>
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<tr>
<td>[I &gt;2 sigma (I)]</td>
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<tr>
<td>R indices</td>
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<tr>
<td>(all data)</td>
</tr>
<tr>
<td>Largest diff. Peak and hole</td>
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<tr>
<td>Temperature</td>
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<tr>
<td>Wavelength</td>
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<tr>
<td>Abs. correction</td>
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<tr>
<td>Refine. Method</td>
</tr>
</tbody>
</table>

$^a$R1=$|(|F_o|-|F_c|)/|F_o||$. 

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br2=\left\{\frac{w(F_o^2-F_c^2)^2}{w(F_o^2)^2}\right\}^{1/2}, w=0.10.

c\text{GoF}=\left[\frac{w(F_o^2-F_c^2)^2}{(N_{\text{refl}}-N_{\text{params}})}\right]^{1/2}.

**Table S2** Selected bond lengths (Å) and angle (°) of 2b

<table>
<thead>
<tr>
<th>Bond</th>
<th>Length (Å)</th>
<th>Angle (°)</th>
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</thead>
<tbody>
<tr>
<td>Zn-O(1)</td>
<td>1.9028(13)</td>
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<tr>
<td>Zn-O(2)</td>
<td>1.9766(13)</td>
<td></td>
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<tr>
<td>Zn-O(2)#1</td>
<td>1.9827(13)</td>
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<tr>
<td>Zn-N</td>
<td>2.1015(15)</td>
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<tr>
<td>O(1)-Zn-O(2)</td>
<td>123.74(6)</td>
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<tr>
<td>O(1)-Zn-O(2)#1</td>
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<tr>
<td>O(2)-Zn-O(2)#1</td>
<td>82.42(5)</td>
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<tr>
<td>O(1)-Zn-N</td>
<td>100.07(6)</td>
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<td>O(2)-Zn-N</td>
<td>124.13(6)</td>
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<tr>
<td>O(2)#1-Zn-N</td>
<td>117.25(6)</td>
<td></td>
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</tbody>
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