

## **SUPPORTING MATERIAL**

### **Synthesis and Characterisation of Antimicrobial Metal-Organic Frameworks as Multi-drug Carriers**

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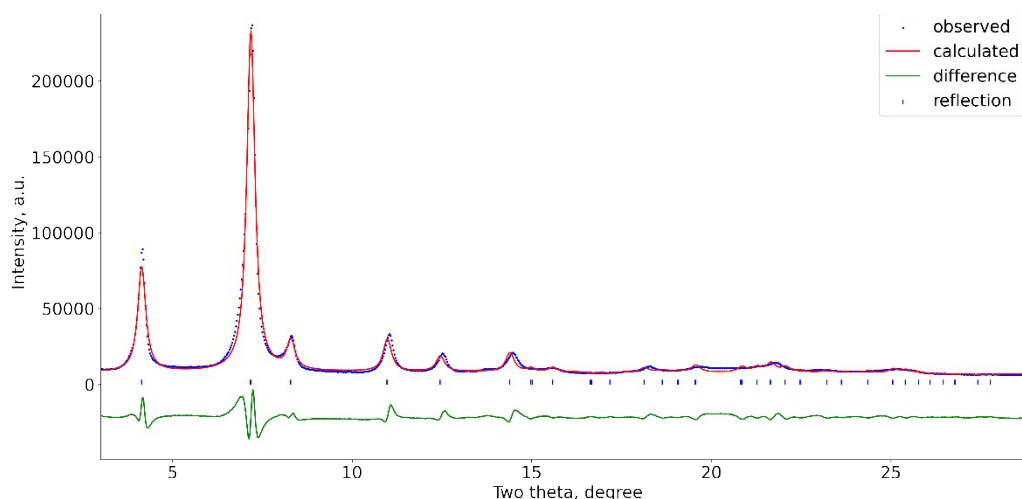
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**Table S1.** Unit cell parameters of the **OnG6** and **Mg<sub>2</sub>(olsalazine)** MOFs. Due to the low number of diffraction peaks the structures of the **OnG6** MOFs were solved in the lowest symmetry space group *P6*, in contrast to the **M<sub>2</sub>(olsalazine)** frameworks, which were solved in the *P3<sub>2</sub>21* space group.

Parameter	<b>OnG6-Zn</b>	<b>OnG6-Mg</b>	<b>OnG6-Cu</b>	<b>OnG6-Co</b>	<b>Mg<sub>2</sub>(olsalazine)</b>
Space Group	<i>P6</i>	<i>P6</i>	<i>P6</i>	<i>P6</i>	<i>P3<sub>2</sub>21</i>
<i>a</i> / Å	24.61(3)	24.32(2)	24.45(3)	24.39(5)	24.97(5)
<i>b</i> / Å	24.61(3)	24.32(2)	24.25(3)	24.39(5)	24.97(5)
<i>c</i> / Å	5.894(7)	5.84(14)	5.81(4)	5.85(3)	6.72(4)
alpha / °	90	90	90	90	90
beta / °	90	90	90	90	90
gamma / °	120	120	120	120	120
volume / Å <sup>3</sup>	3091(5)	2972(8)	2974(18)	3011(14)	3824(3)

**Table S2.** INH loadings for the **OnG6** MOFs.

<b>OnG6</b>	<b>INH Loading</b>	<b>Framework Stoichiometry derived from <sup>1</sup>H-NMR spectrum</b>
<b>OnG6-Zn</b>	4 wt%	<b>Zn<sub>2</sub>(AZD)<sub>1</sub>(INH)<sub>0.12</sub></b>
<b>OnG6-Mg</b>	22 wt%	<b>Mg<sub>2</sub>(AZD)<sub>1</sub>(INH)<sub>0.74</sub></b>
<b>OnG6-Cu</b>	N/A	<b>N/A</b>



**Figure S1.** Pawley profile fit for **OnG7-Mg**.

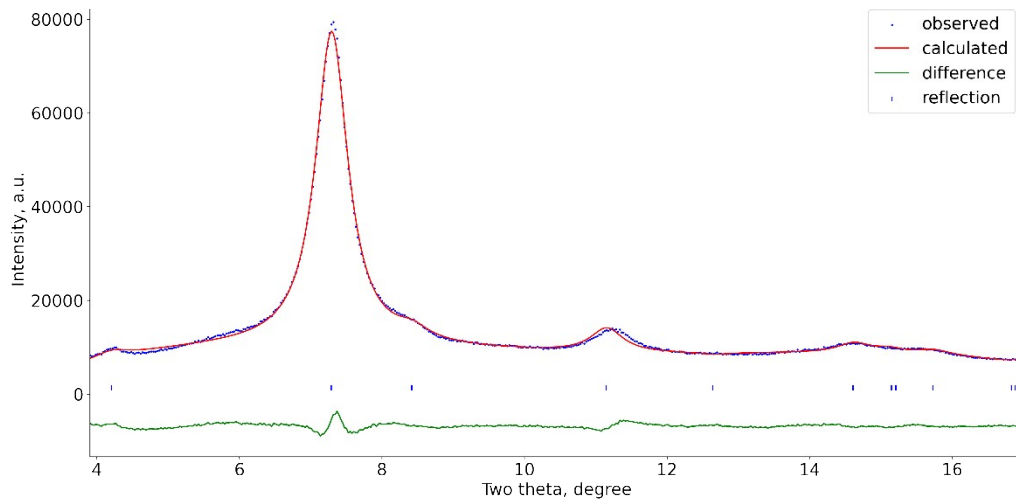


Figure S2. Pawley profile fit for **OnG6-Zn**.

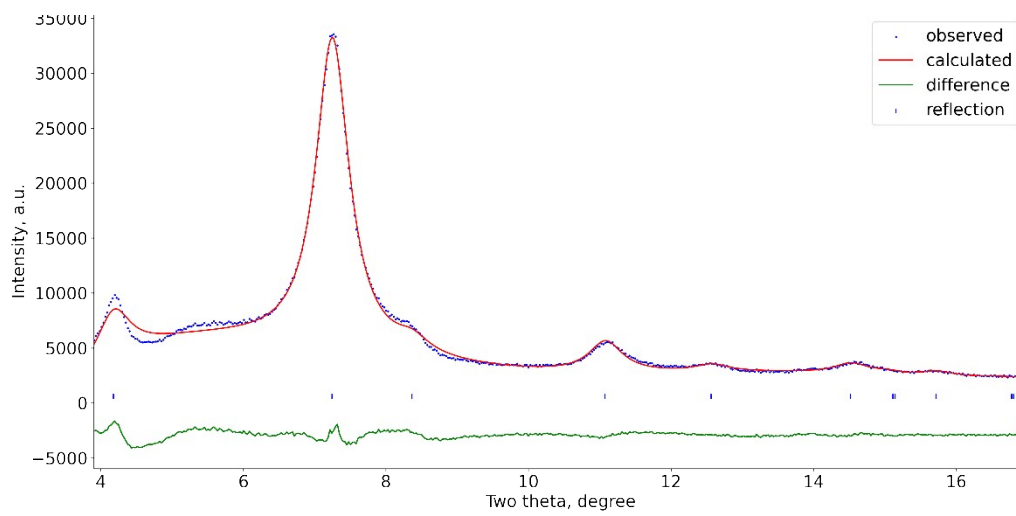


Figure S3. Pawley profile fit for **OnG6-Co**.

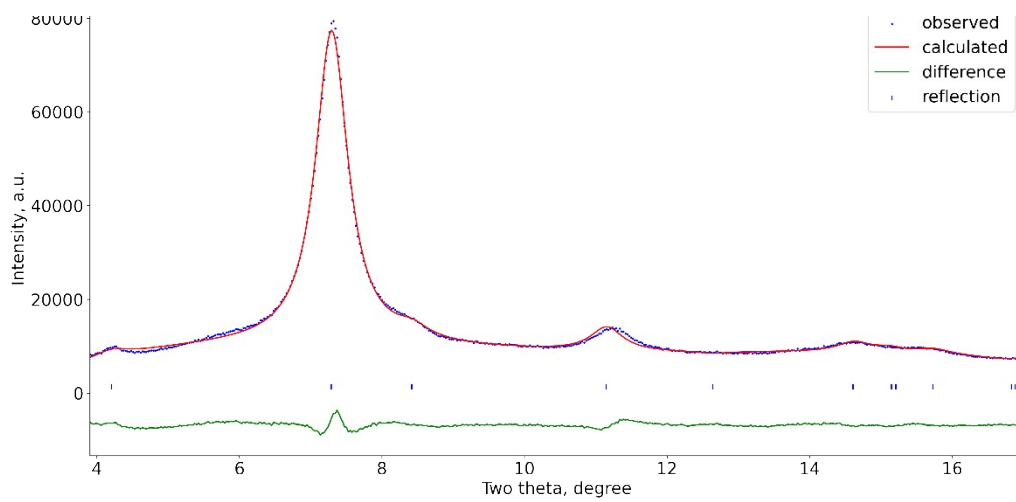


Figure S4. Pawley profile fit for **OnG6-Cu**.

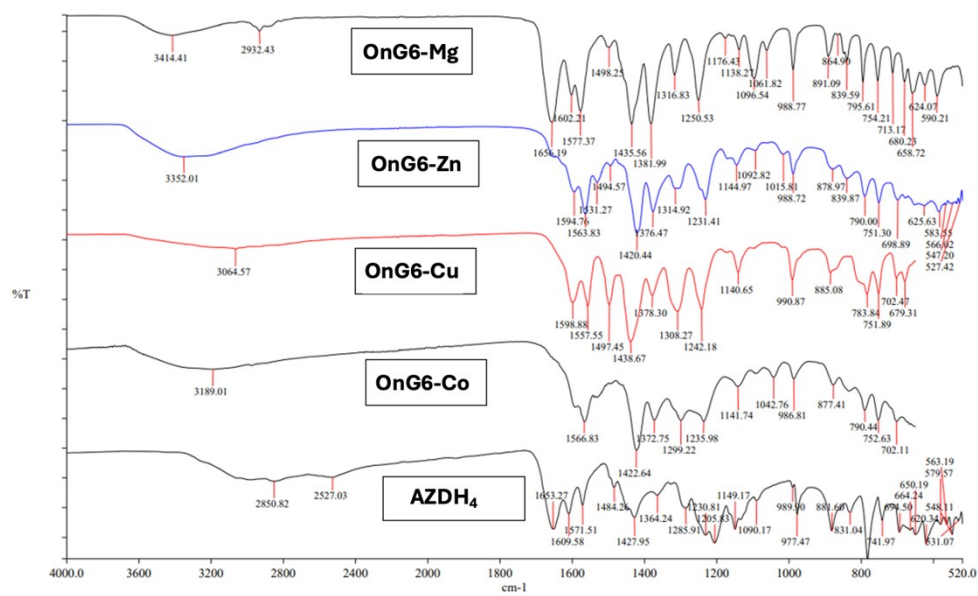


Figure S5. FTIR spectra of the **OnG6** MOFs and the linker **AZDH<sub>4</sub>**.

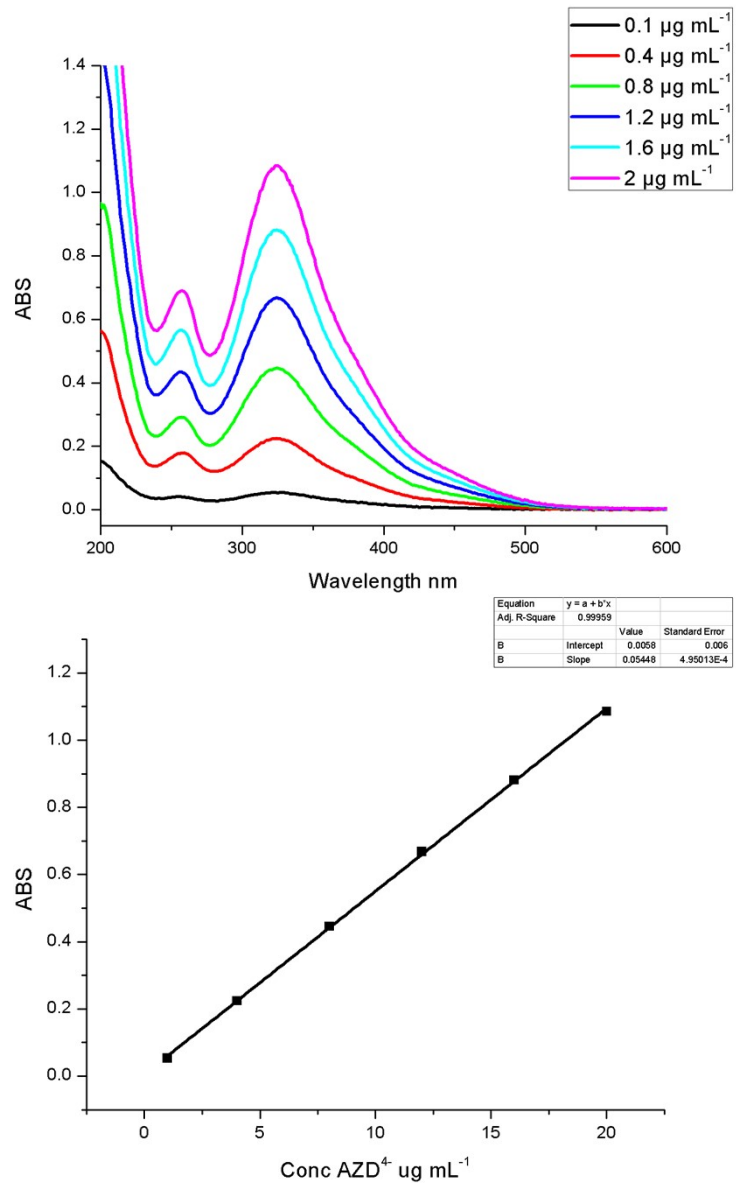


Figure S6. UV-Vis data (top) for Na<sub>4</sub>(AZD) calibration curve (bottom).

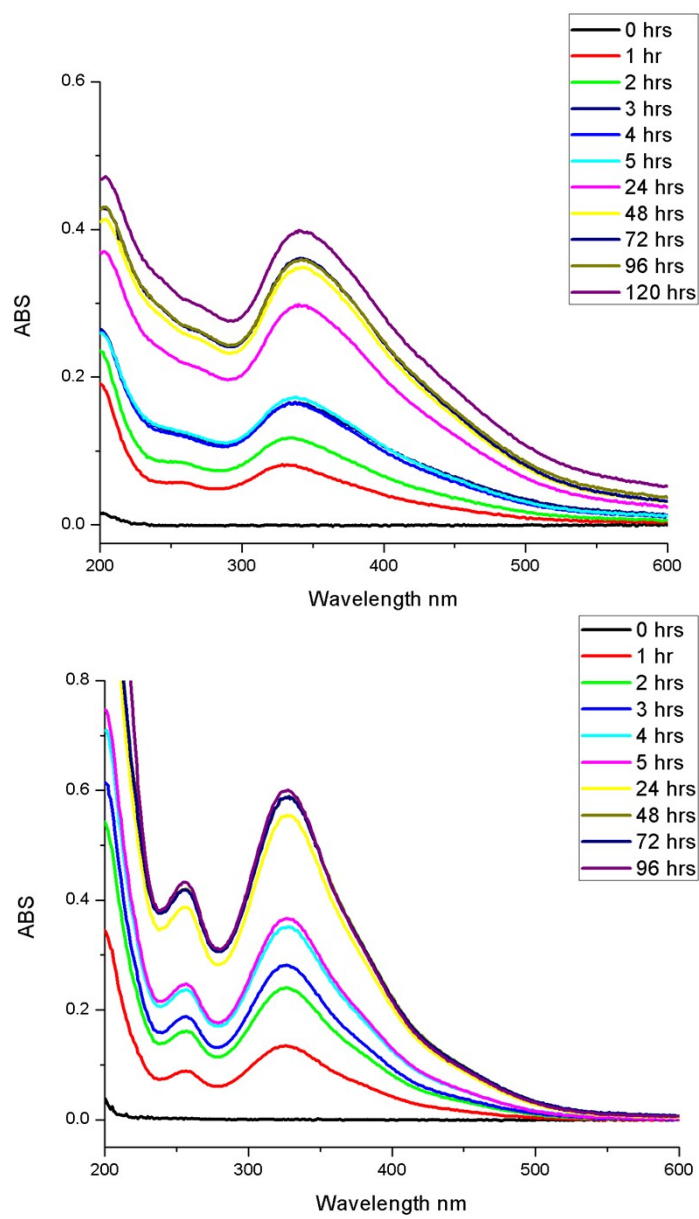


Figure S7. UV-Vis data for **OnG6-Cu** dissolution in H<sub>2</sub>O (top) and PBS solution.

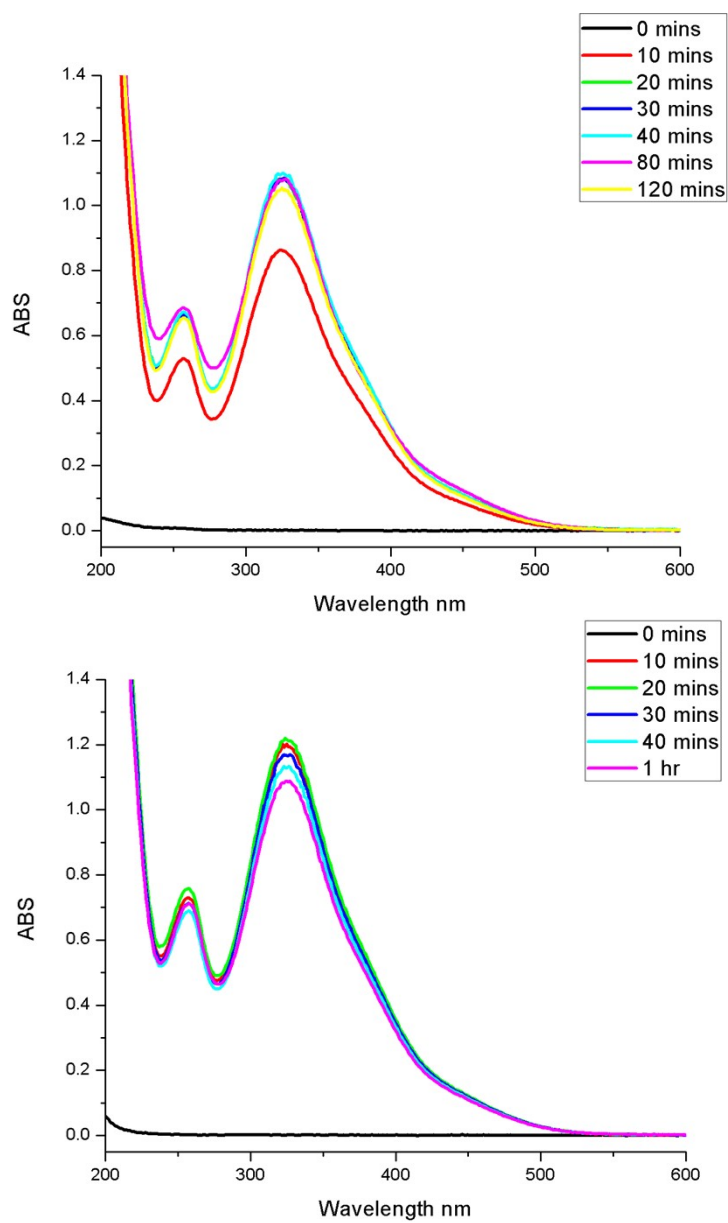


Figure S8. UV-Vis data for **OnG6-Mg** dissolution in H<sub>2</sub>O (top) and PBS solution.

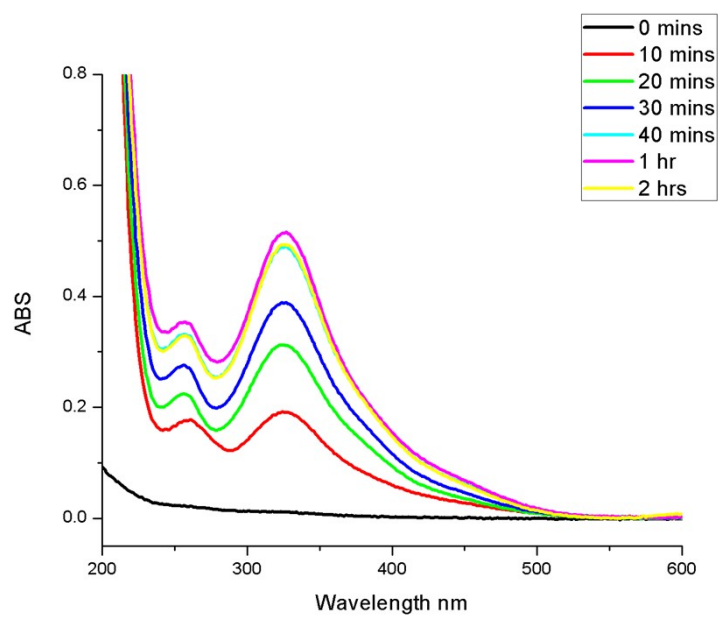


Figure S9. UV-Vis data for **OnG6-Zn** dissolution in water.

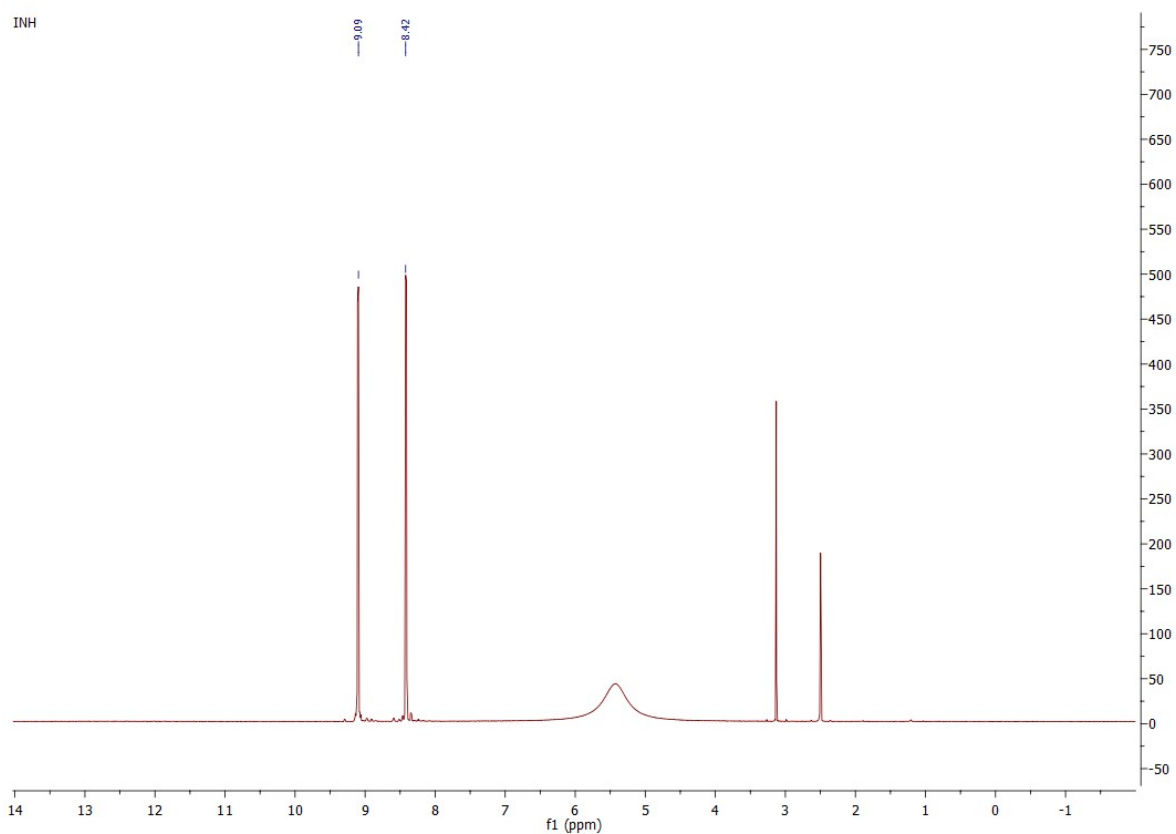


Figure S10. <sup>1</sup>H-NMR of INH in DMSO-d<sub>6</sub>/DCI.



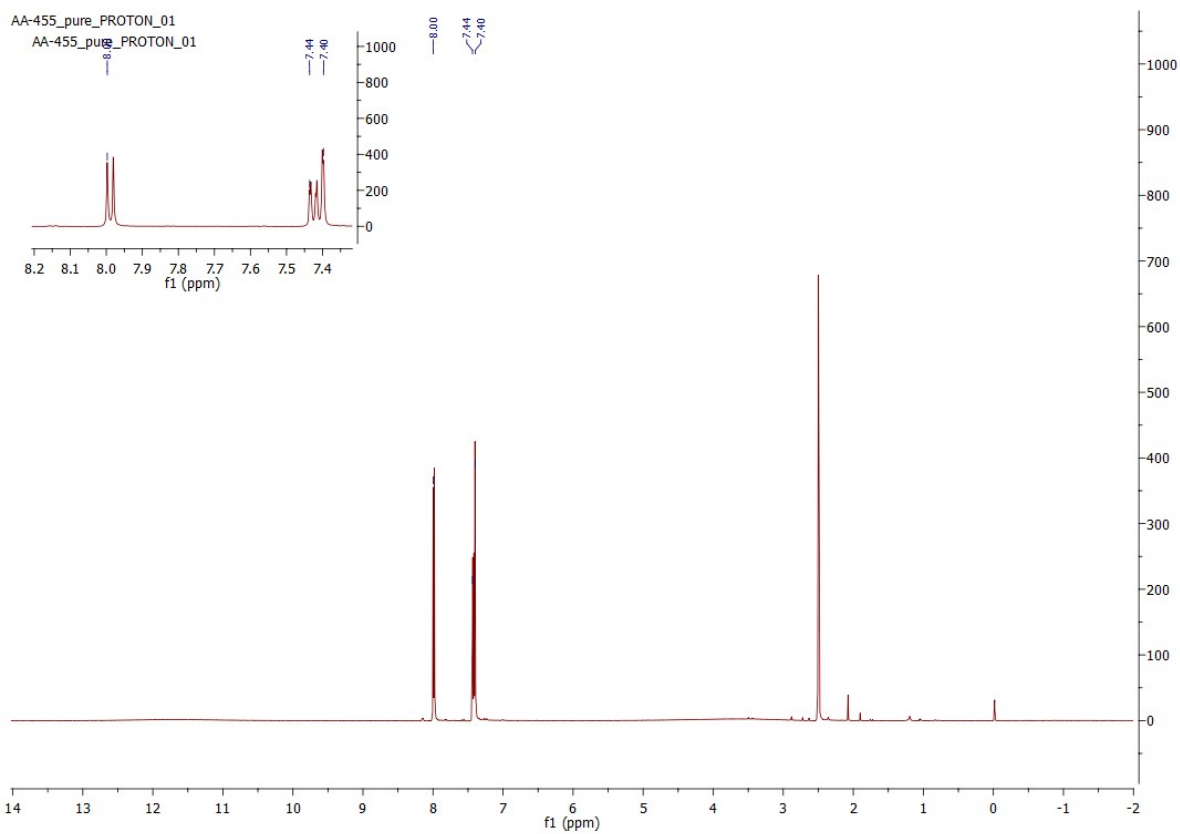


Figure S11.  $^1\text{H-NMR}$  of  $\text{AZDH}_4$  in  $\text{DMSO-d}_6$ .

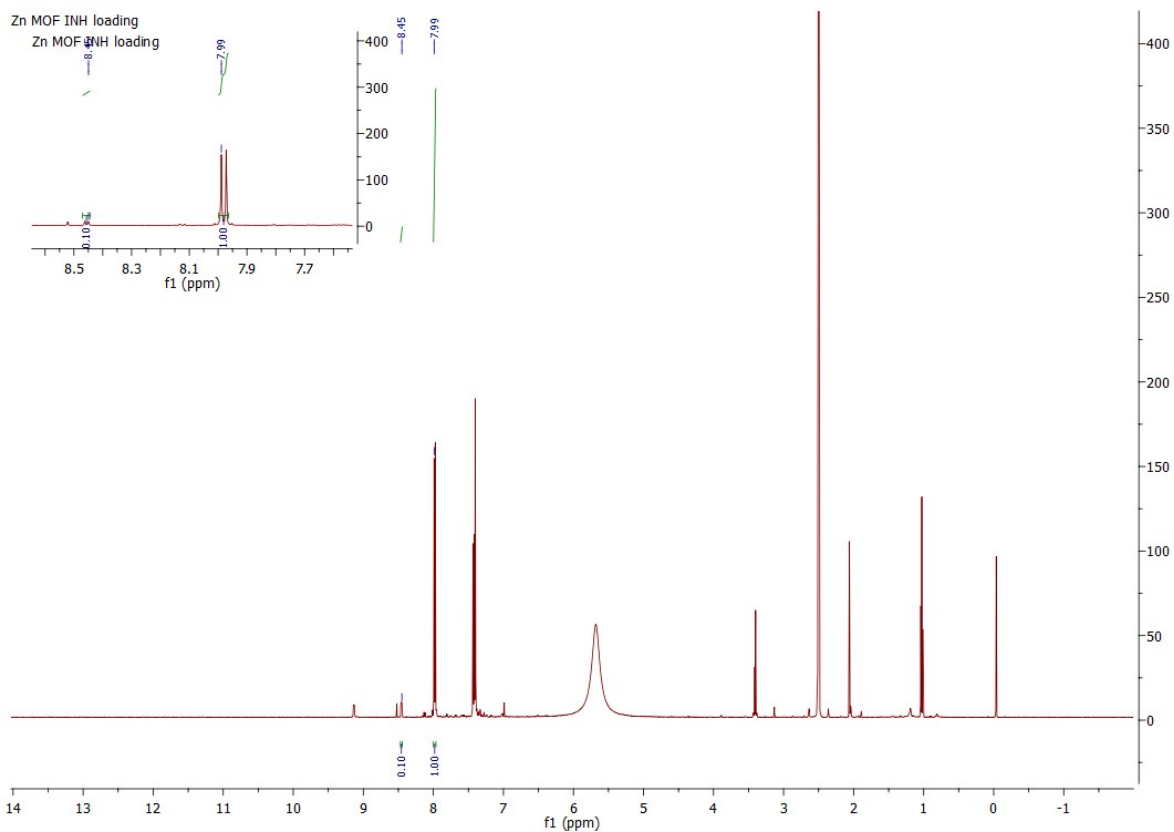


Figure S12.  $^1\text{H-NMR}$  of digested **OnG6-Zn** after INH loading (3 days) in  $\text{DMSO-d}_6/\text{DCI}$ .

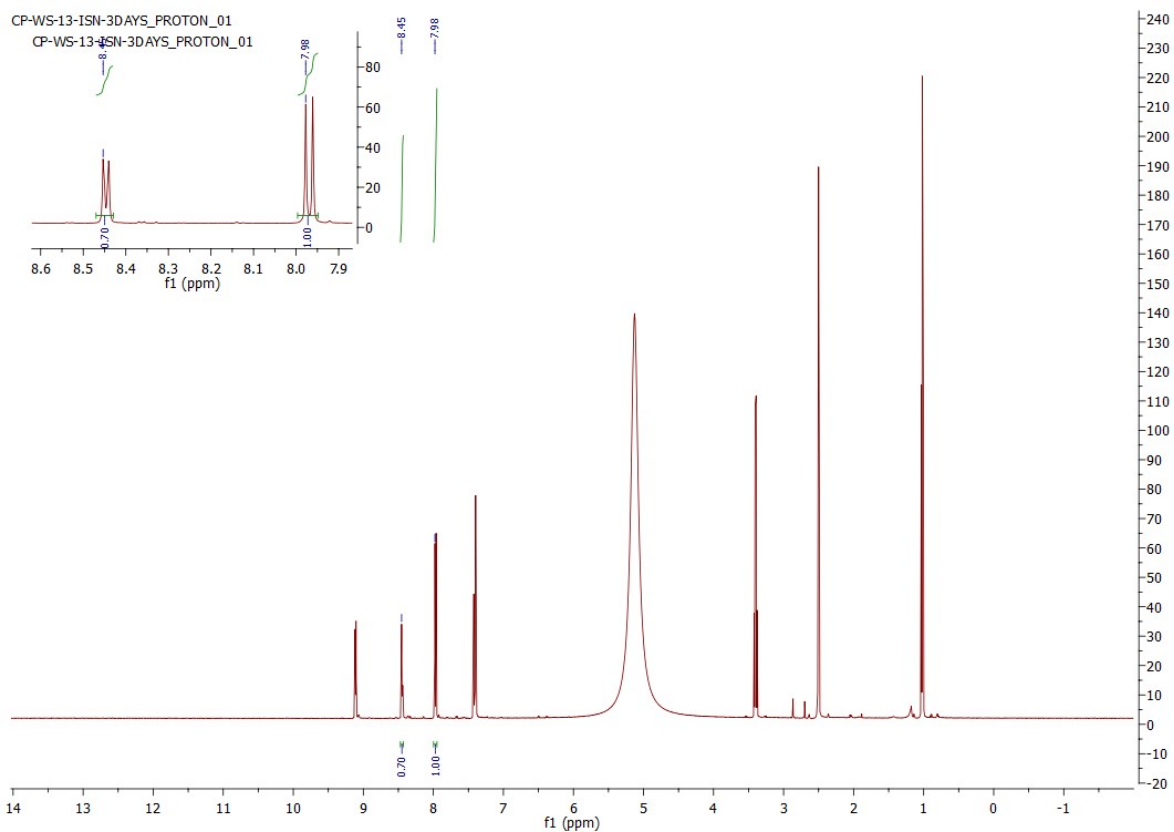


Figure S13.  $^1\text{H-NMR}$  of digested **OnG6-Mg** after INH loading (3 days) in  $\text{DMSO-d}_6/\text{DCI}$ .

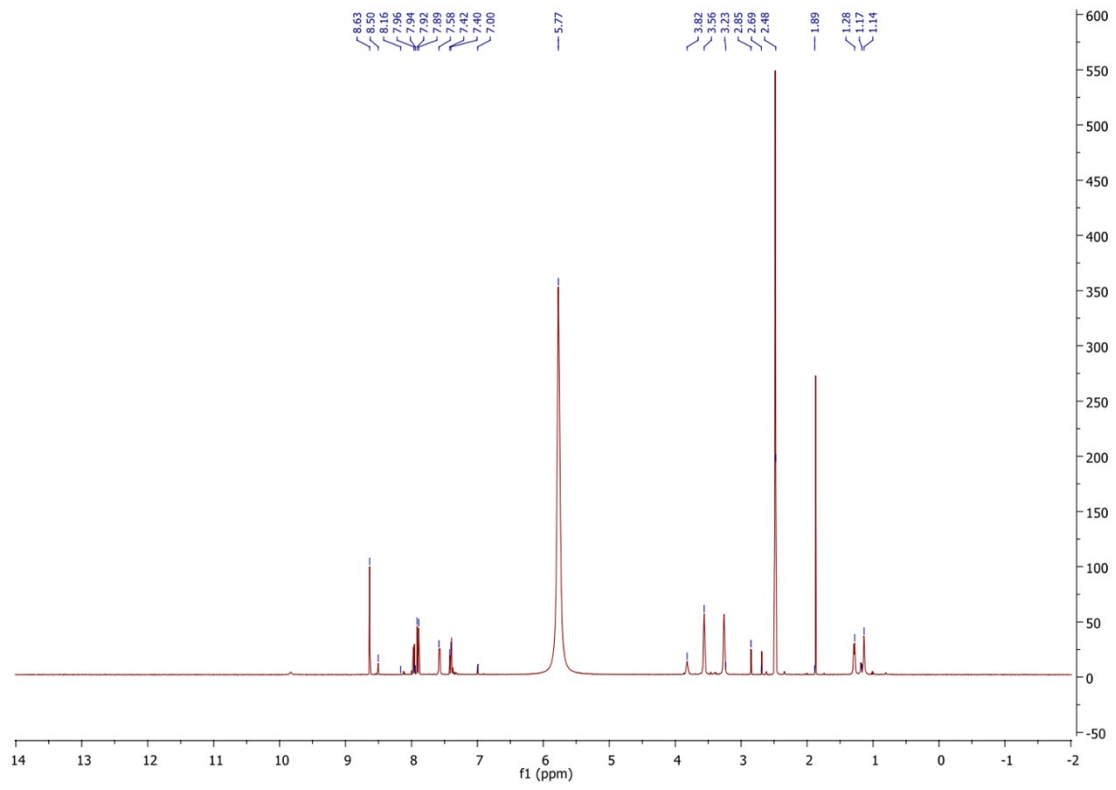


Figure S14.  $^1\text{H-NMR}$  of digested **OnG6-Mg** after CIPRO loading (3 days) in  $\text{DMSO-d}_6/\text{DCl}$ .

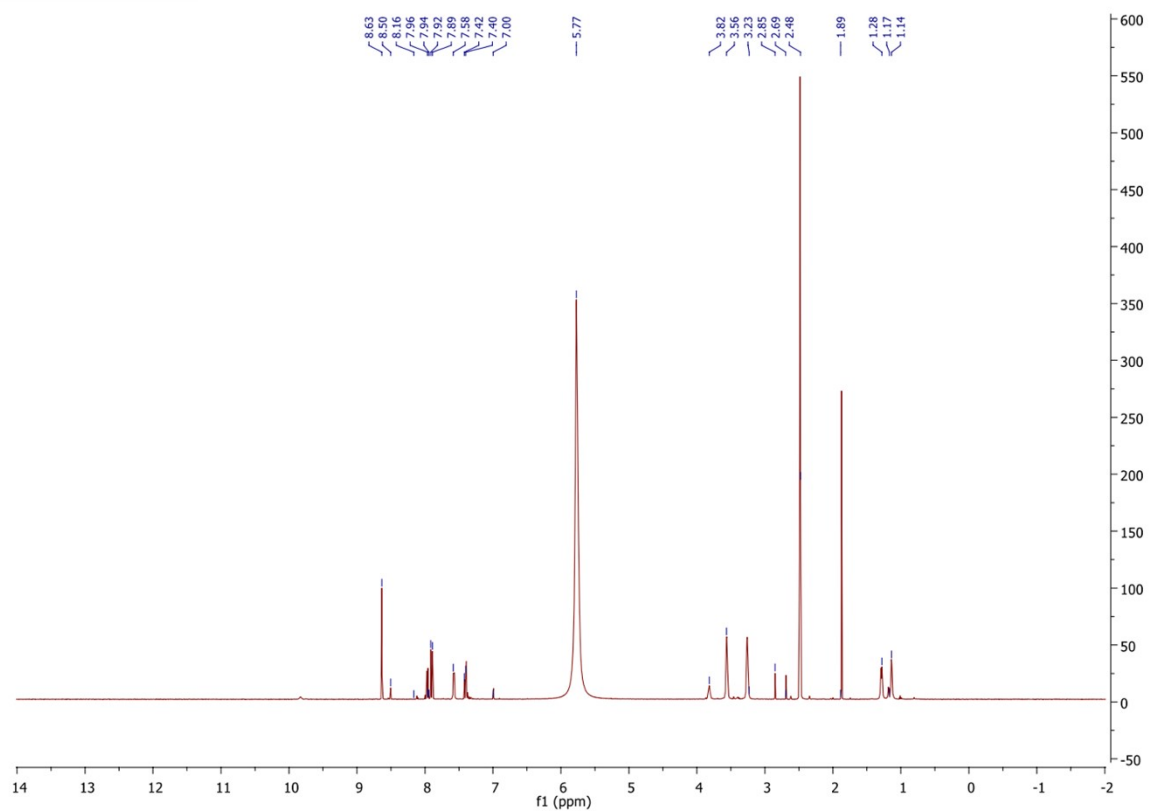


Figure S15. <sup>1</sup>H-NMR of digested **OnG6-Co** after CIPRO loading (3 days) in DMSO-d<sub>6</sub>/DCI.