## Novel 'Anti-Prussian Blue' Structure Based on $Zn^{2+}$ Nodes and $[Re_3Mo_3S_8(CN)_6]^{6-}$ Heterometallic Cluster Spacers and its

## Rearrangement to Prussian Blue

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## **Supporting Information**

Fig. S1. IR spectra of compounds 1 (red dotted line) and 2 (black solid line). To prevent decay, spectrum of compound 1 was registered in an ammonia mother solution.

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Fig. S2. Observed (a) and calculated (b) EPR spectra of compound 2.

Bond	d, Å	Bond	d, Å	Bond	d, Å
1		2		*	
M1-M1	2.6248(10)	M1-M1	2.6374(17)	M1-M1	2.6273(8)
M1-M2	2.6230(8)				
M1-S1	2.439(2)	M1-S1	2.426(4)	M1-S1	2.439(2)
M2-S1	2.427(3)				
M1-C1	2.145(14)	M1-C1	2.23(3)	M1-C1	2.173(13)
M2-C2	2.17(3)				
Zn1-N2	2.35(3)	Zn1-N1	2.194(19)		
Zn2-N3	1.994(10)	Zn2-N2	2.16(3)		
Zn2-N4	2.04(2)				

**Table S1.** Bond distances (Å) in 1 and 2 compared to  $CaK_4[Re_3Mo_3S_8(CN)_6] \cdot 8H_2O(*)$ .