Supporting Information

En Route to Osmium Analogs of KP1019: Synthesis, Structure, Spectroscopic Properties and Antiproliferative Activity of *trans*-$[Os^{IV}Cl_4(Hazole)_2]$

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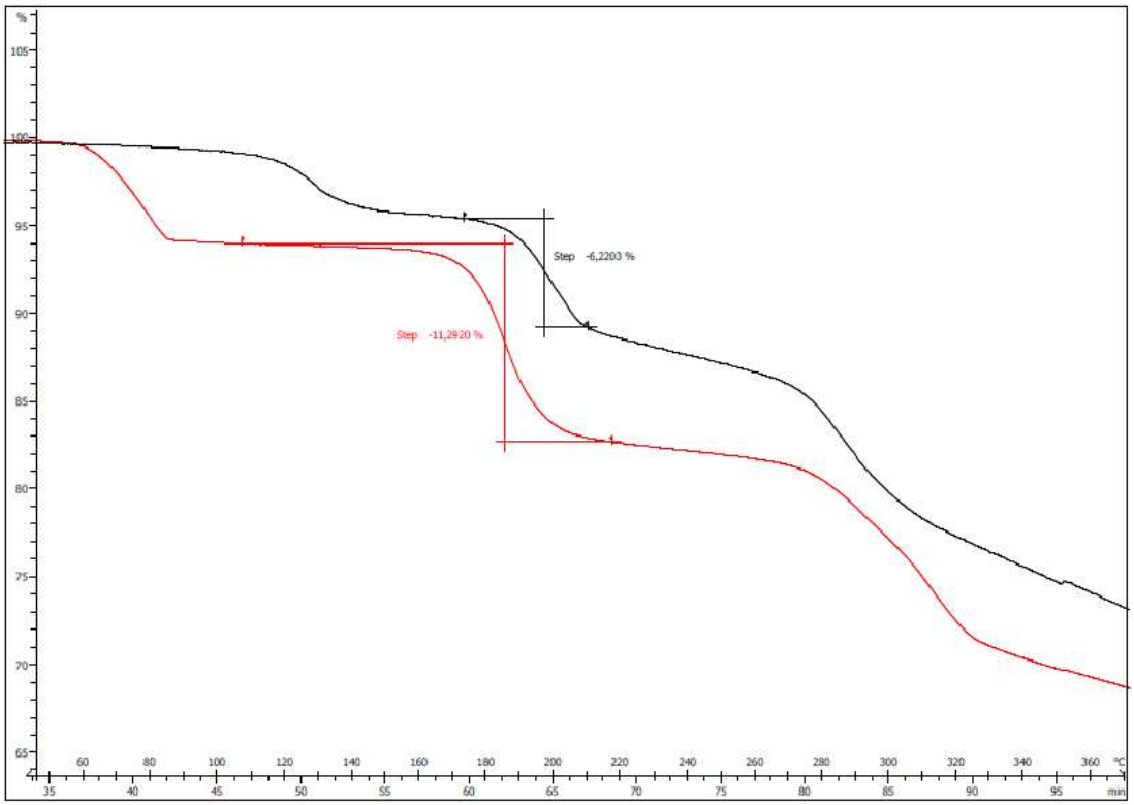

Figure S1: Thermogravimetric data for $(H_2ind)_2[OsCl_6]$ (red) and $(H_2ind)[OsCl_5(\kappa N2\text{-}Hind)]$ (black) showing a weight loss of 11.29% and 6.22%, respectively, attributed to the loss of one (calcd. weight loss 6.03%) and two (calcd. weight loss 11.37%) molecules of HCl.

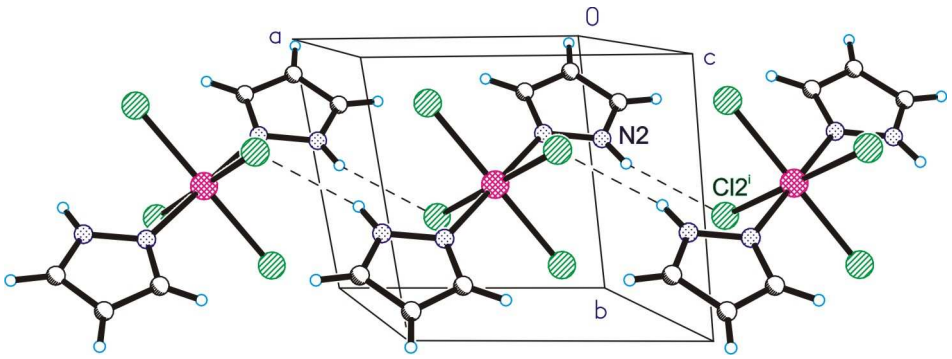

Figure S2. The unit cell of **1** showing the formation of a chain along axis *a* via intermolecular hydrogen bonding interaction N2–H···Cl2^{i} [N2–H 0.88, H···Cl2^{i} 2.511, N2···Cl2^{i} 3.270 Å, N2–H···Cl2^{i} 144.9°] (symmetry code i: *x*–1, *y*, *z*).

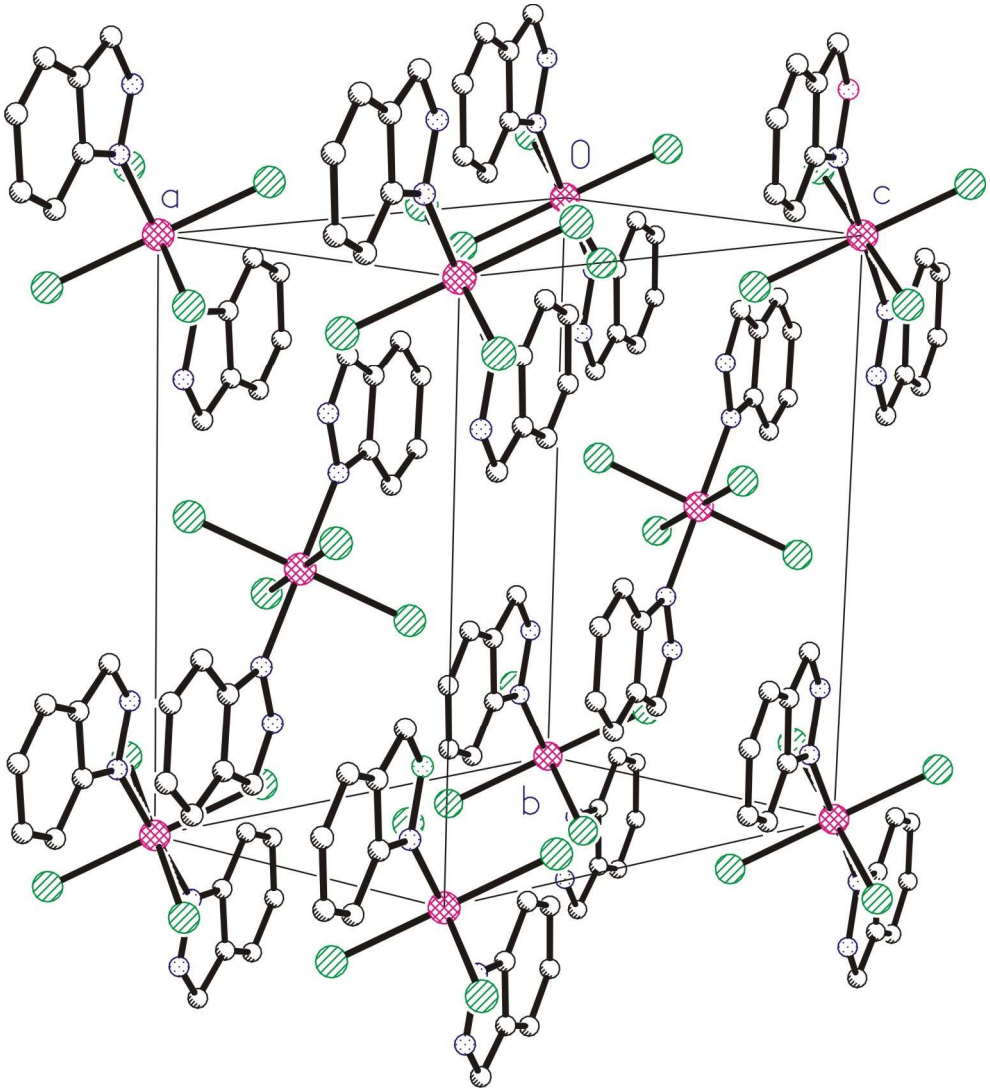

Figure S3. The unit cell of **2**; hydrogen atoms were omitted for clarity.

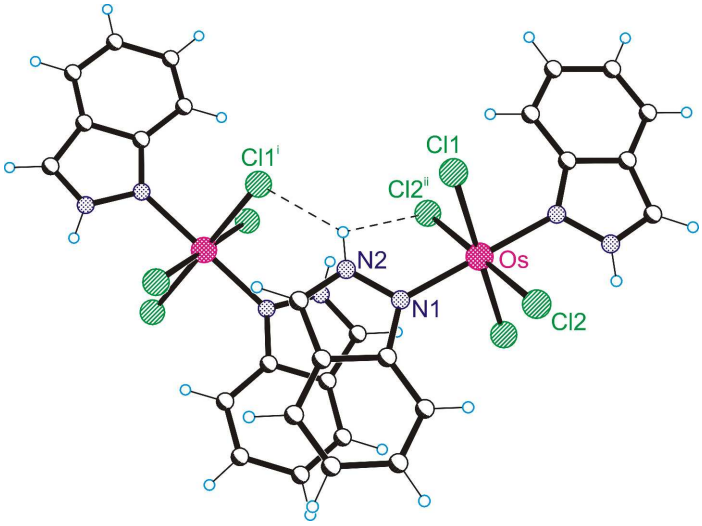

Figure S4. Bifurcated hydrogen bonding interactions N2–H···Cl1^{i} and N2–H···Cl2^{ii} [N2–H 0.88, H···Cl1^{i} 2.798, N2···Cl1^{i} 3.444 Å, N2–H···Cl1^{i} 131.5°; H···Cl2^{ii} 2.713, N2···Cl2^{ii} 3.143 Å, N2–H···Cl2^{ii} 111.4°] (symmetry codes i: *x*, – *y* + 1.5, *z* – 0.5; ii: – *x*, –*y* + 2, –*z*) and π–π stacking interaction in the crystal structure of **2**.

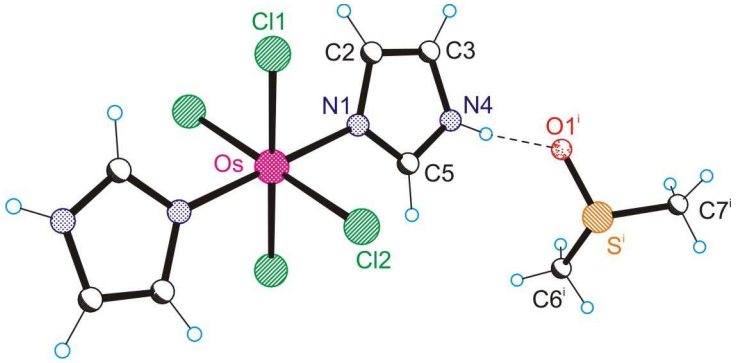

Figure S5. Strong hydrogen bonding N4–H···O1^{i} [N4–H 0.88, H···O1^{i} 1.826, N4···O1^{i} 2.698 Å, N4–H···O1^{i} 170.9°] (symmetry code i: *x* + 1, *y*, *z*) between molecule of **3** and DMSO.

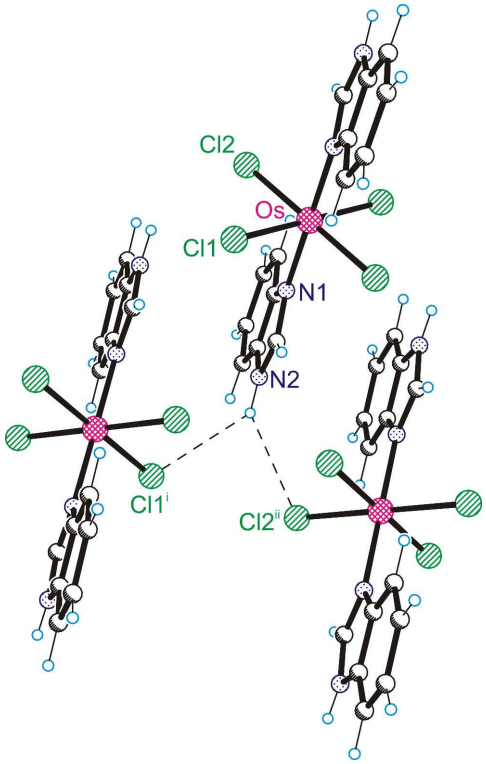

Figure S6. Bifurcated hydrogen bonding interactions N2–H···Cl1^{i} and N2–H···Cl2^{ii} [N2–H 0.88, H···Cl1^{i} 2.867, N2···Cl1^{i} 3.541 Å, N2–H···Cl1^{i} 134.6°; H···Cl2^{ii} 2.770, N2···Cl2^{ii} 3.504 Å, N2–H···Cl2^{ii} 141.9°] (symmetry codes i: $-x + 1$, $y + 0.5$, $-z + 0.5$; ii: x, $-y + 0.5$, $z - 0.5$) and π–π stacking interaction in the crystal structure of **4**.

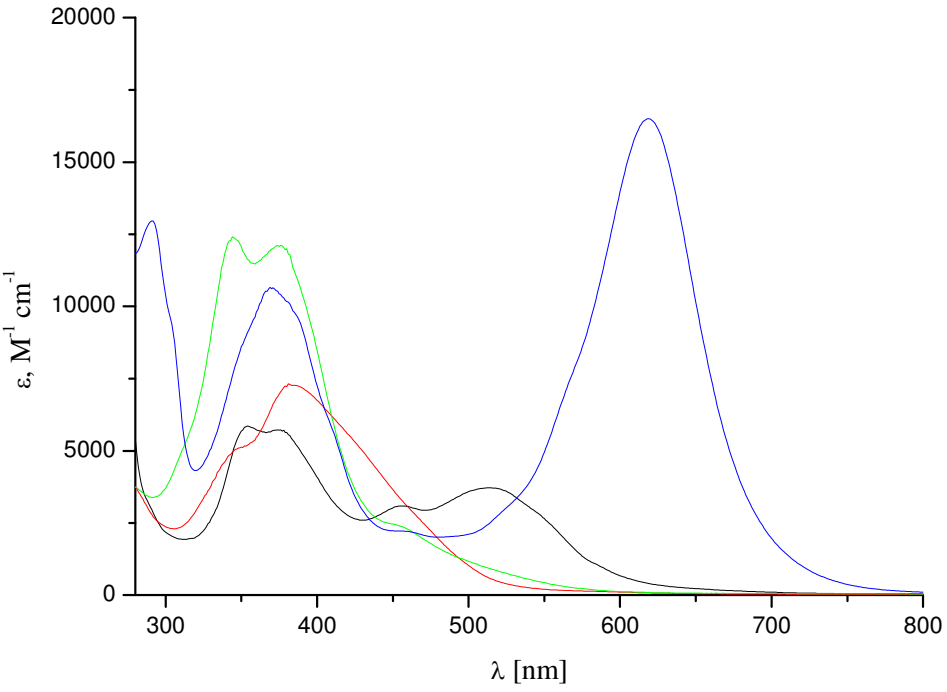

Figure S7: UV−vis spectra of **1** (red), **2** (blue), **3** (green) and **4** (black) in DMSO.