

SUPPLEMENTARY MATERIAL

Polypyridines, Picrates, Lanthanides: A Plethora of Stacks ?

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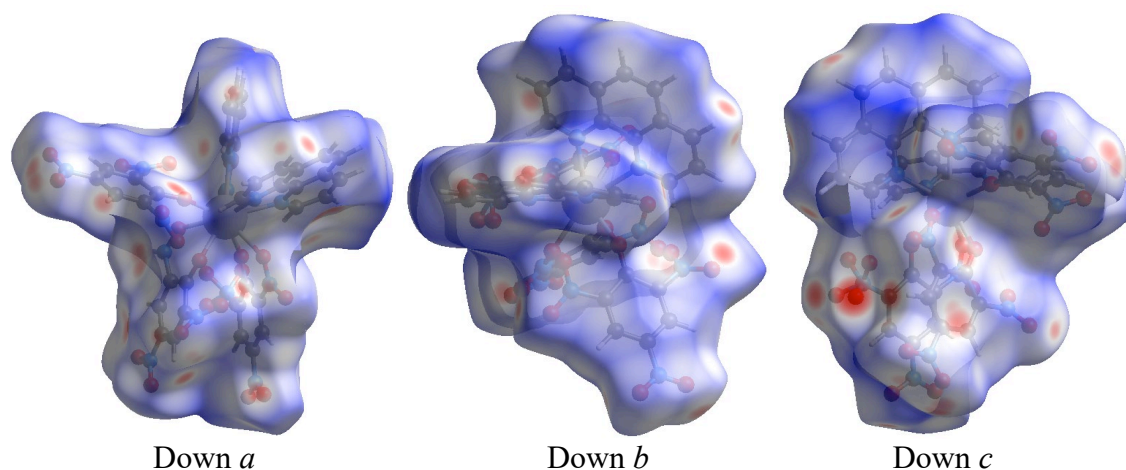


Figure S1 Views of the Hirshfeld surface calculated with CrystalExplorer for the molecule present in the lattice of $[\text{Eu}(\text{phen})_2(\text{pic-O},\text{O}')_2(\text{pic-O})]$. d_{norm} representation, transparent mode; red regions indicate points where interactions with adjacent atoms of a separate molecule exceed dispersion.

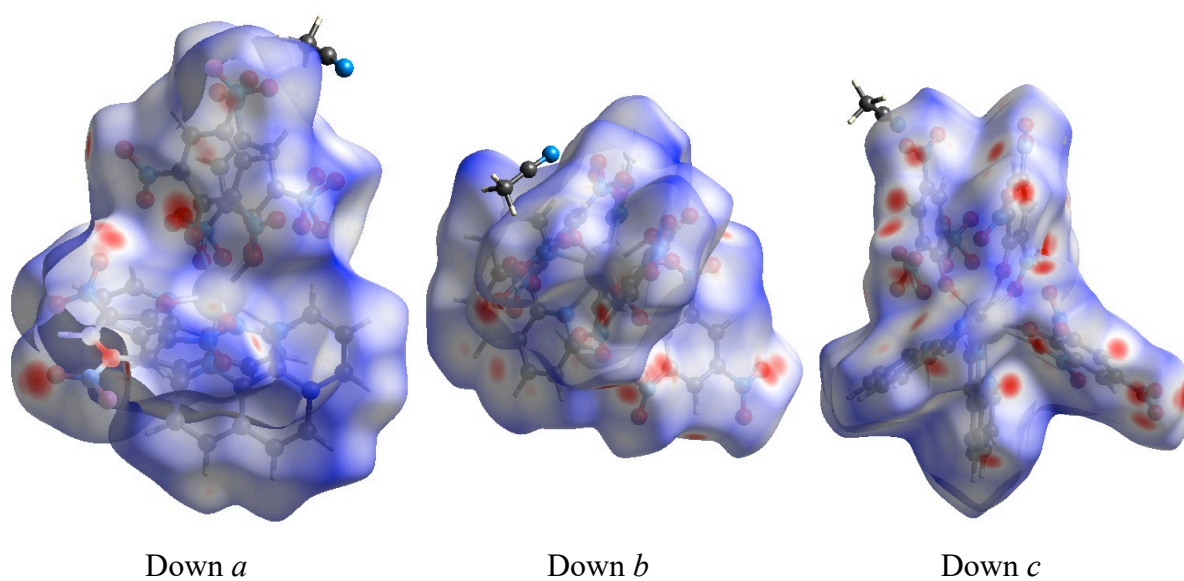


Figure S2(a) Views of the Hirshfeld surface calculated with CrystalExplorer for the molecule present in the lattice of $[\text{Gd}(\text{phen})_2(\text{pic-O},\text{O}')_2(\text{pic-O})]\cdot\text{CH}_3\text{CN}$. d_{norm} representation, transparent mode; red regions indicate points where interactions with adjacent atoms of a separate molecule exceed dispersion.

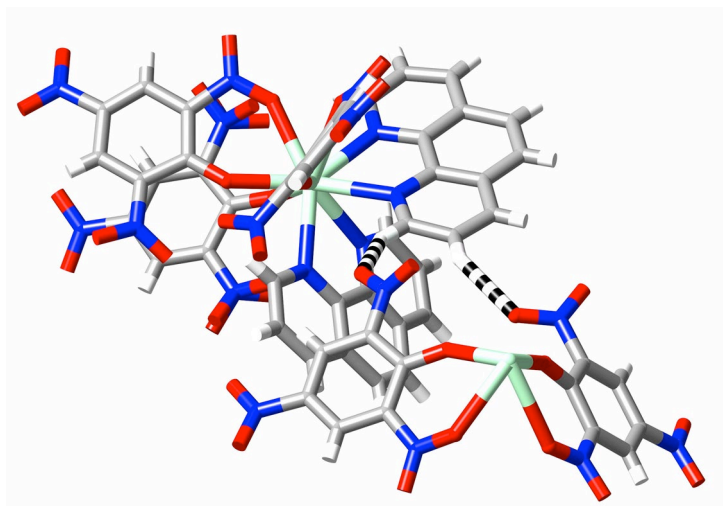


Figure S2(b) *CH...O* interactions (dashed lines) of phen1 in the molecule present in the lattice of $[Gd(phen)_2(pic-O,O')_2(pic-O)].CH_3CN$.

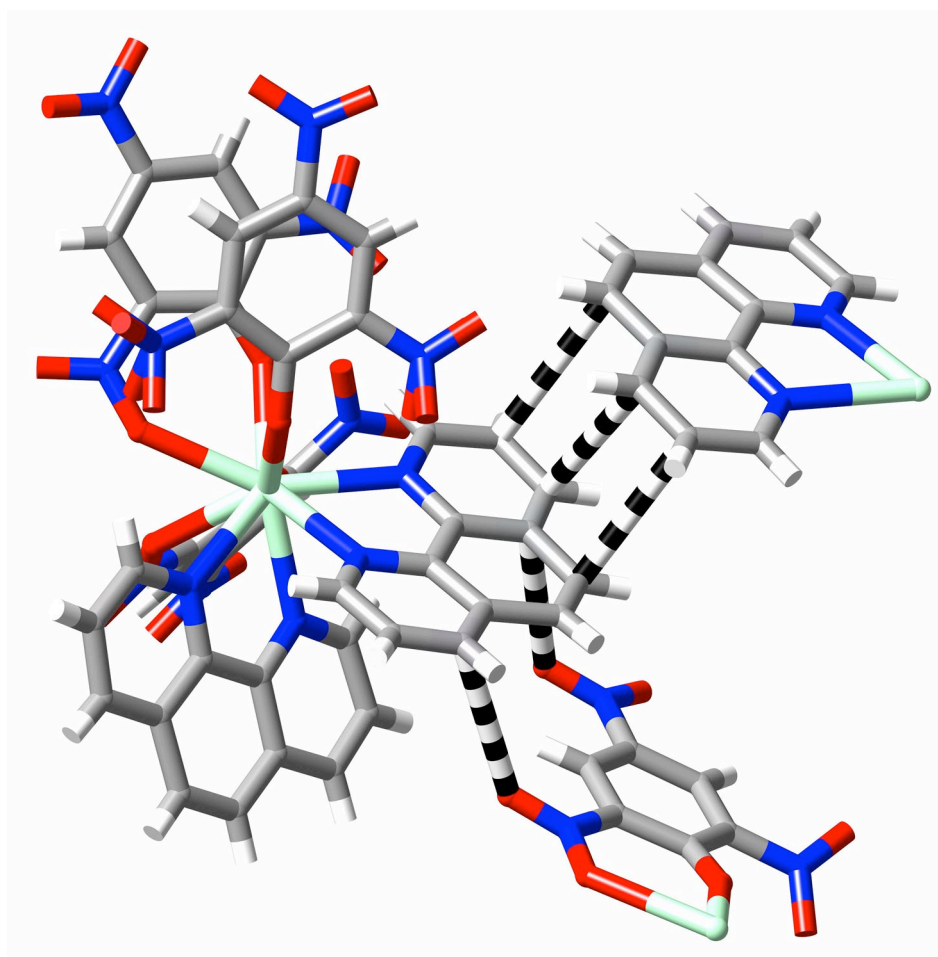


Figure S2(c) *C...C* interactions between phen2 ligands of adjacent molecules in the lattice of $[Gd(phen)_2(pic-O,O')_2(pic-O)].CH_3CN$. Additional *C...O* interactions are shown for one of the stacked pair. All interactions exceeding dispersion are shown as dashed lines.

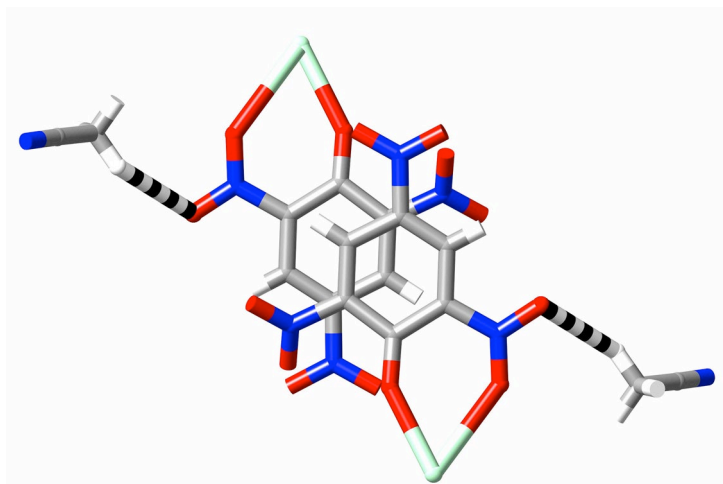


Figure S2(d) The stacked pair of picrate 1 units formed between two molecules in the lattice of $[Gd(phen)_2(pic-O,O')_2(pic-O)].CH_3CN$, projected perpendicular to the mean ring planes.. H-bonds to acetonitrile are shown as dashed lines.

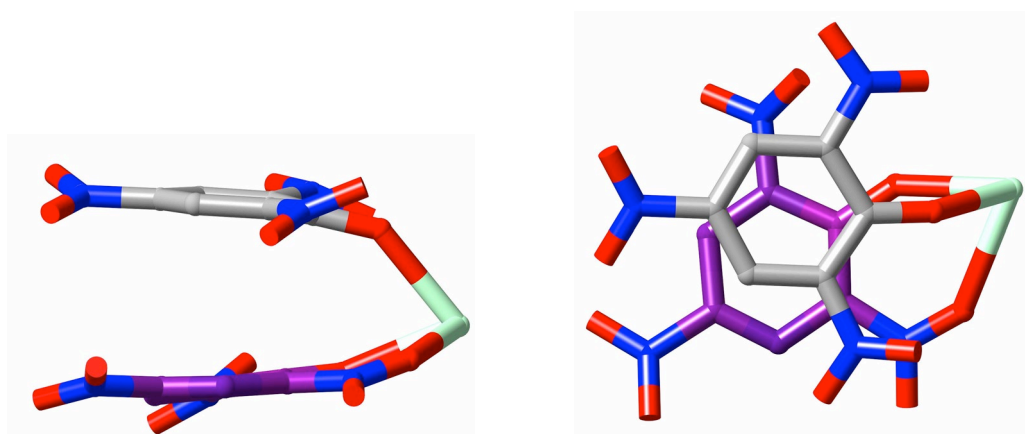


Figure S2(e) Orthogonal views of the intramolecular stacking of picrates 2 and 3 within the coordination sphere of Gd^{III} in $[Gd(phen)_2(pic-O,O')_2(pic-O)].CH_3CN$. C atoms of picrate 3 are shown in violet.

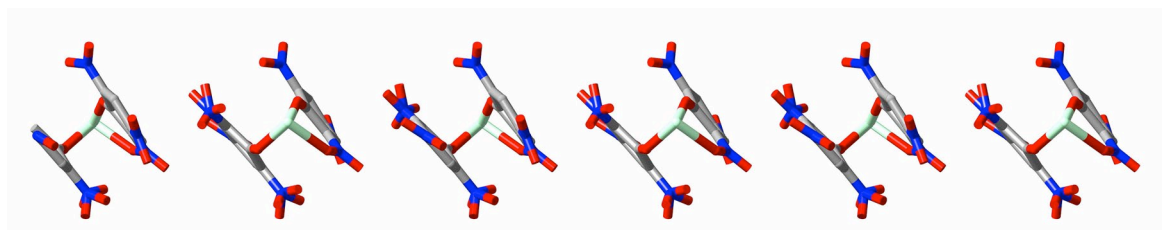


Figure S2(f) The slipped stack single column (running parallel to a) involving picrates 2 and 3 in both intra- and inter-molecular stacking. For clarity, the complete ligand environment of the Gd^{III} ions is not shown.

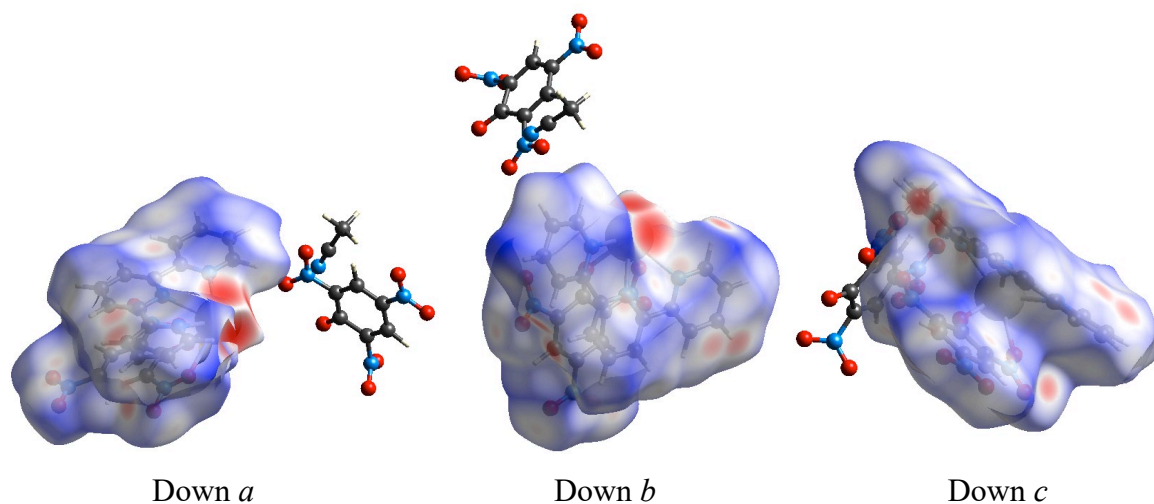


Figure S2(a) Views of the Hirshfeld surface calculated with CrystalExplorer for the asymmetric unit of the cation present in the lattice of [La(terpy)₂(pic)₂](pic).CH₃CN. d_{norm} representation, transparent mode; red regions indicate points where interactions with adjacent atoms of a separate molecule exceed dispersion.