

SUPPLEMENTARY INFORMATIONS

New Ni(II) 1,2-bis(diphenylphosphino)ethane dithiolates: Crystallographic, computational and Hirshfeld surface analyses

Reena Yadav,^a Manoj Trivedi,^b Gabriele Kociok-Köhn,^c Rajendra Prasad^d and Abhinav Kumar^{a*}

a. Department of Chemistry, University of Lucknow, Lucknow 226 007, India.

b. Department of Chemistry, University of Delhi, Delhi 110 007, India.

c. Department of Chemistry, University of Bath, Bath BA2 7AY, UK.

d. Department of Chemistry, S. G. B. University, Amrawati, Maharashtra, India.

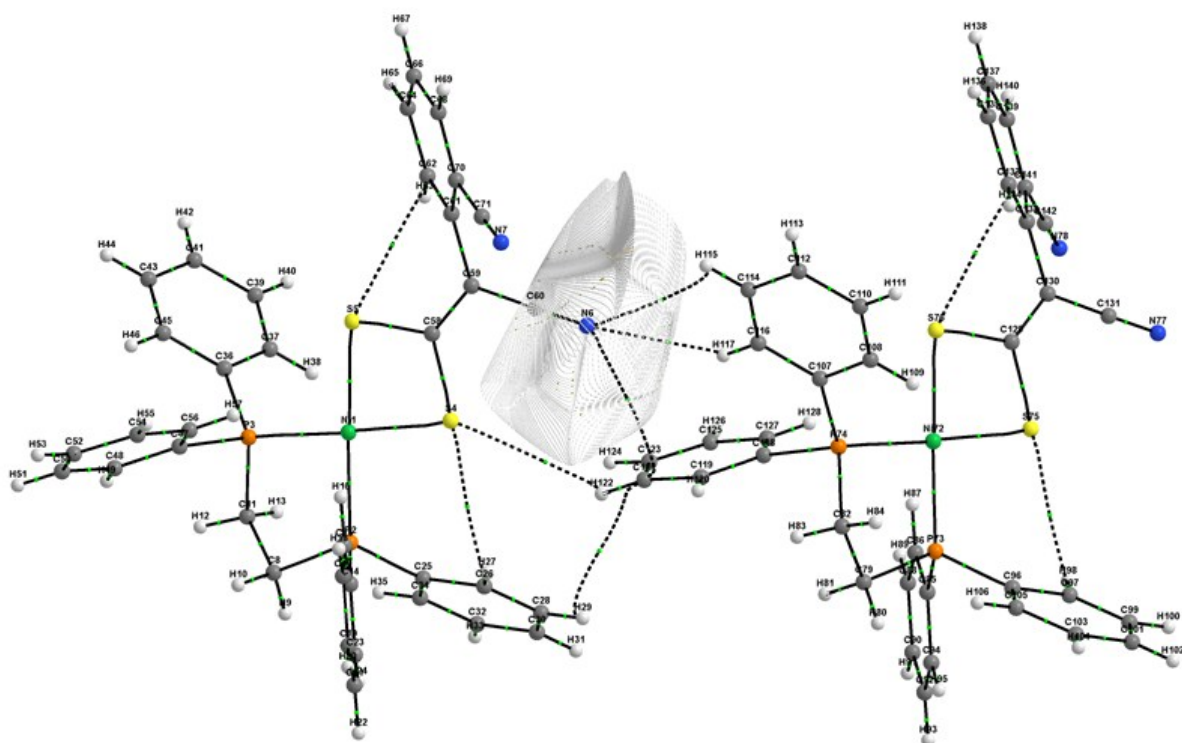


Fig. S1 Molecular graphs for **2** displaying intermolecular (a) C≡N...H interaction along with the interatomic surface of interest.

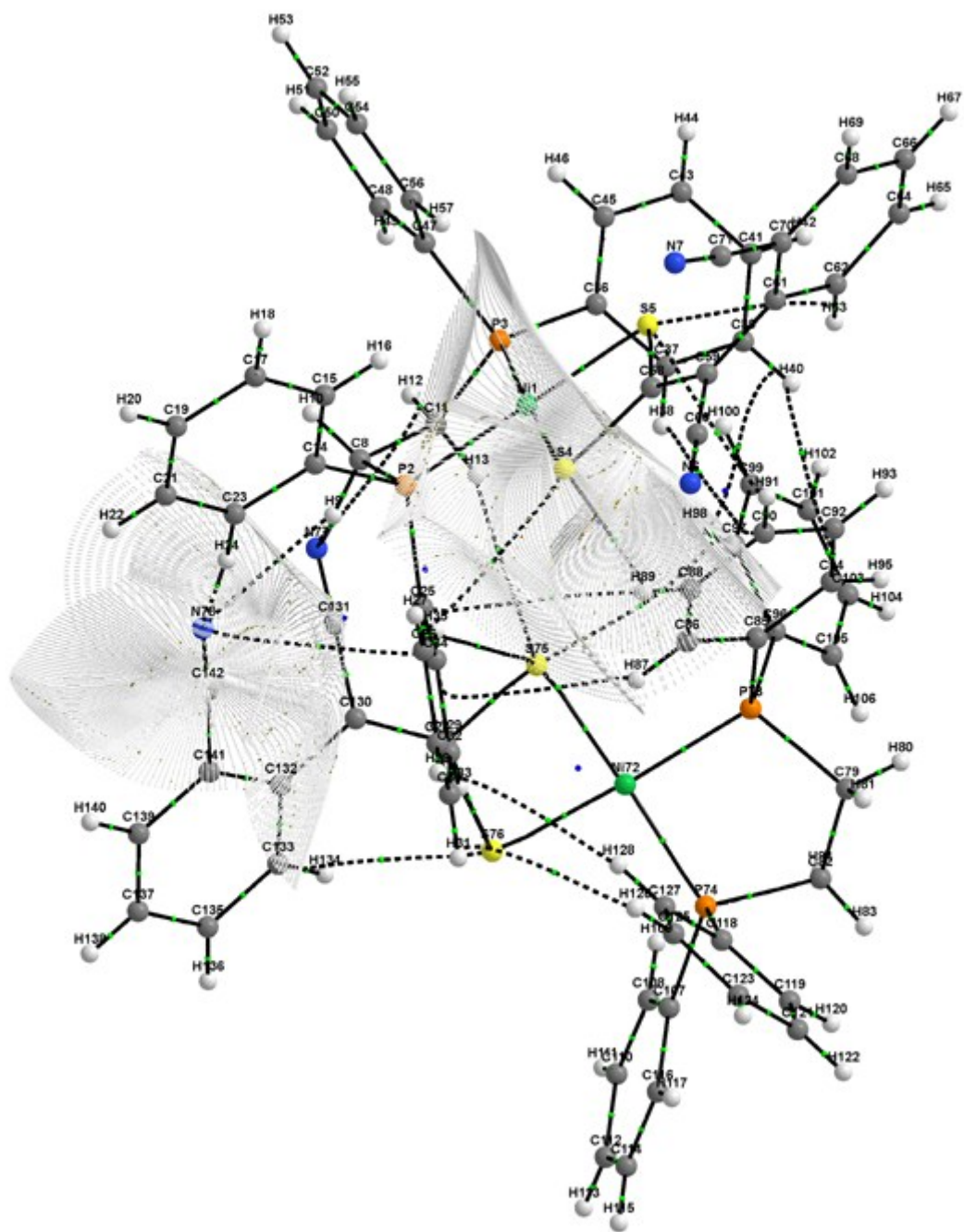


Fig. S2 Molecular graphs for **2** displaying intermolecular C-S...H, C≡N...H interaction alongwith the interatomic surface of interest.

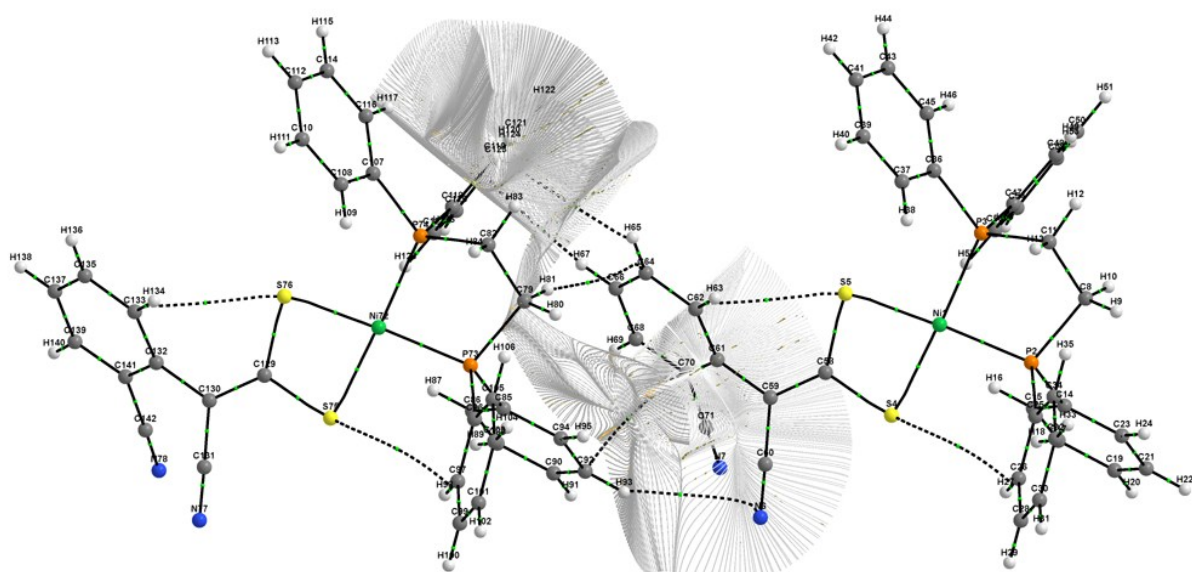


Fig. S3 Molecular graphs for **2** displaying intermolecular $\pi \cdots \pi$ and C-H $\cdots \pi$ interactions along with the interatomic surface of interest.

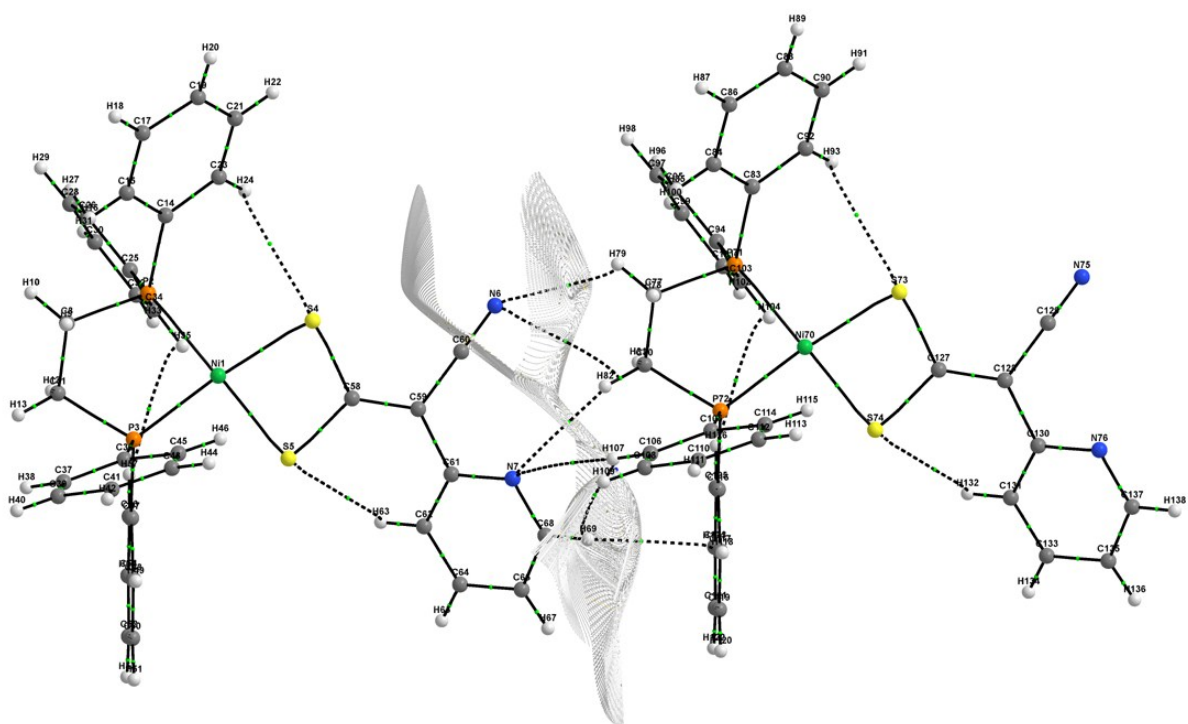


Fig. S4 Molecular graphs for **3** displaying intermolecular C \equiv N \cdots H; (py)N \cdots H; C-H $\cdots \pi$ interactions along with the interatomic surface of interest.

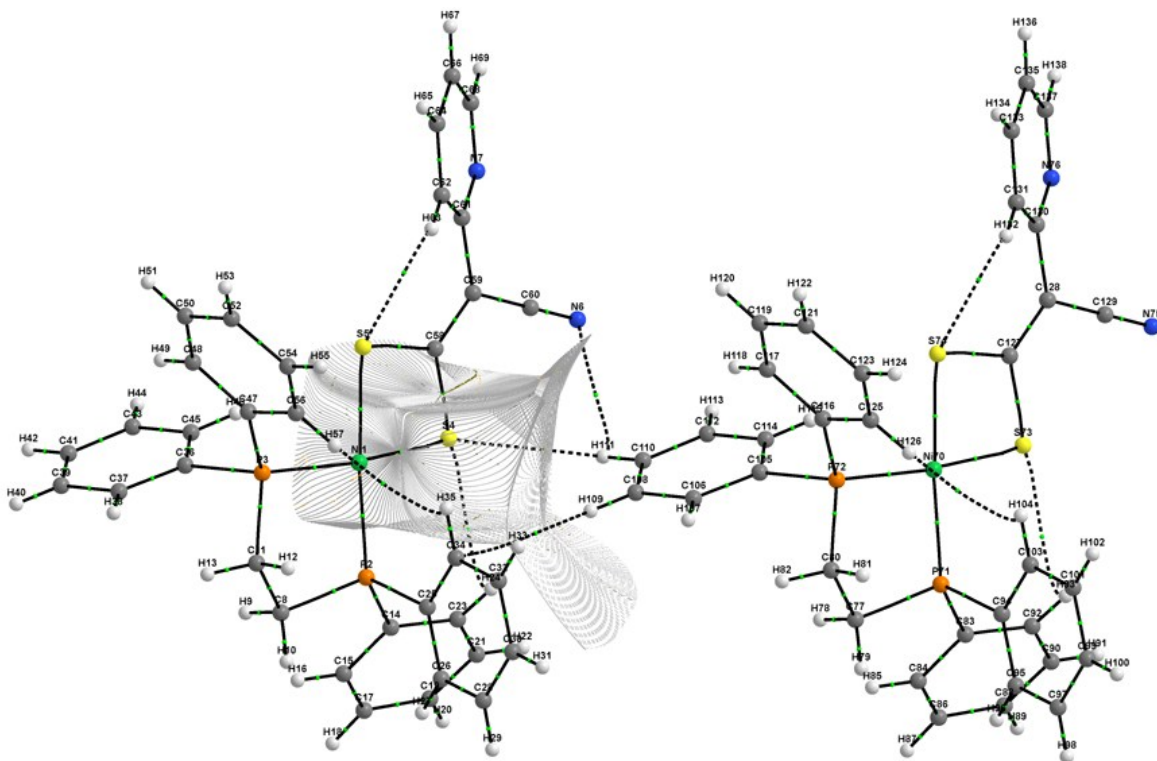


Fig. S5 Molecular graphs for **3** displaying intermolecular C-S...H interaction alongwith the interatomic surface of interest.