

Gold... π aryl interactions as supramolecular synthons†

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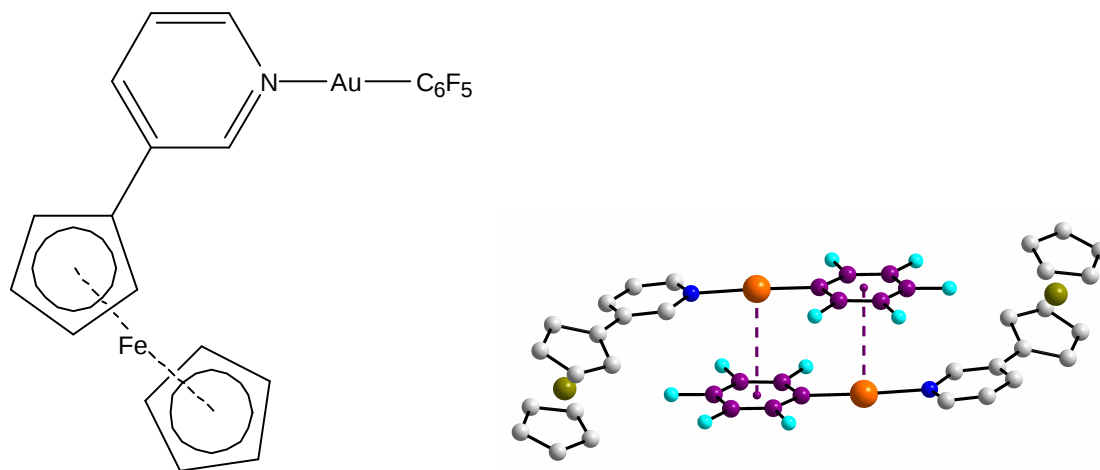
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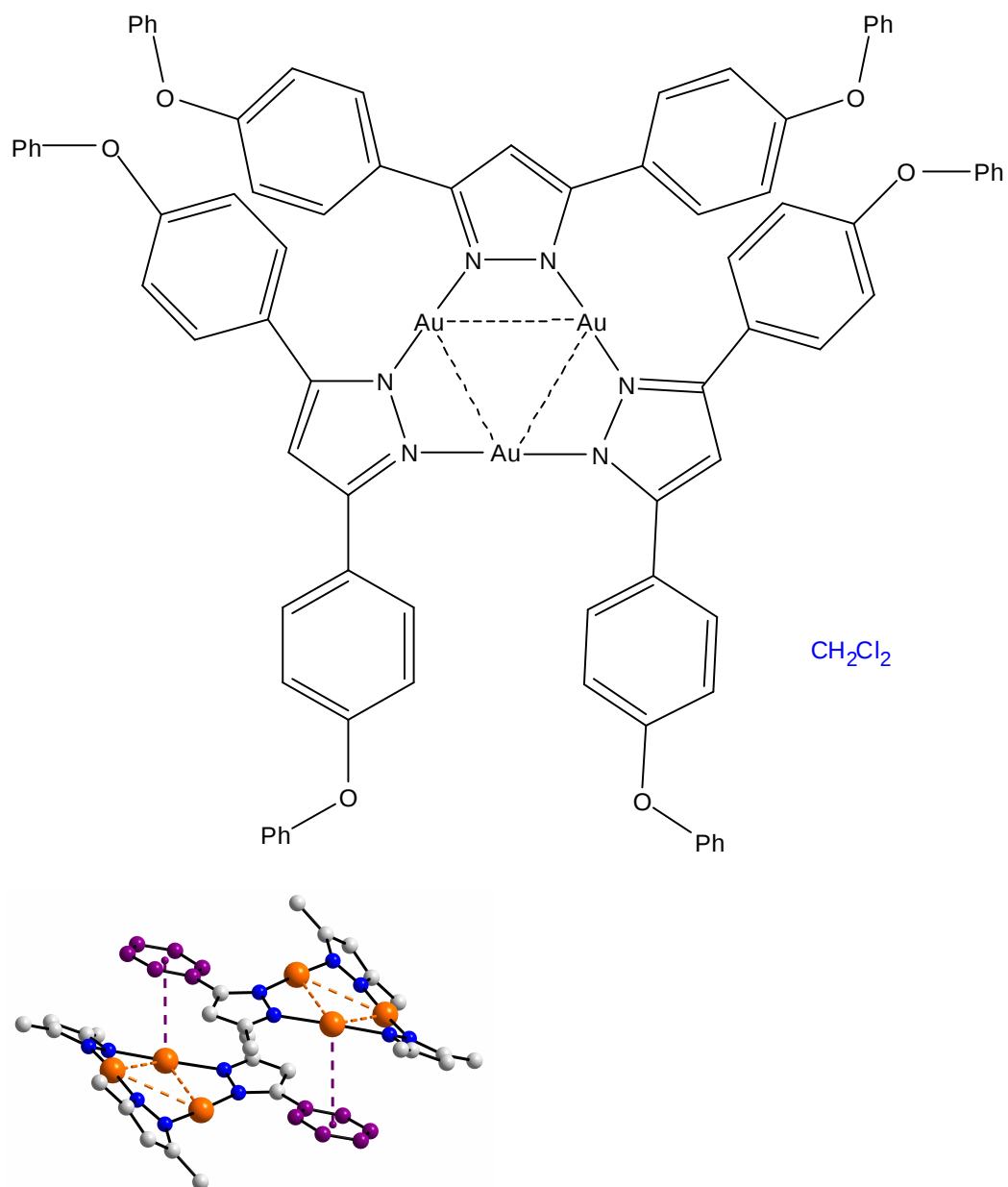
Table S(2) Chemical structures and other descriptors describing intermolecular Au.... π interactions in spherical aggregates, (11) – (33). Species not engaged in Au.... π interactions are shown in blue.

Compound number; Au...C(g), d (Å); Au...C(g)...plane, α^a (°); [ref.] and REFCODE

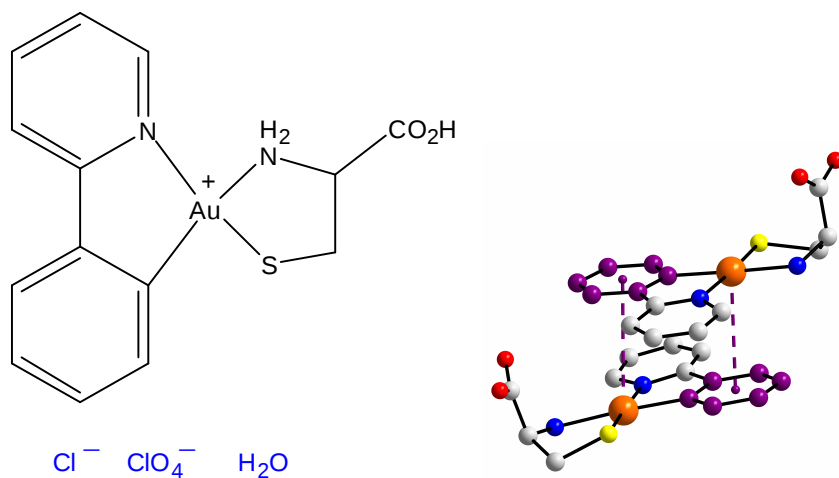
(11), 3.51, 5.1 [E. M. Barranco, O. Crespo, M. C. Gimeno, P. G. Jones, A. Laguna and M. D. Villacampa, *J. Organomet. Chem.*, 1999, **592**, 258] KETLUQ



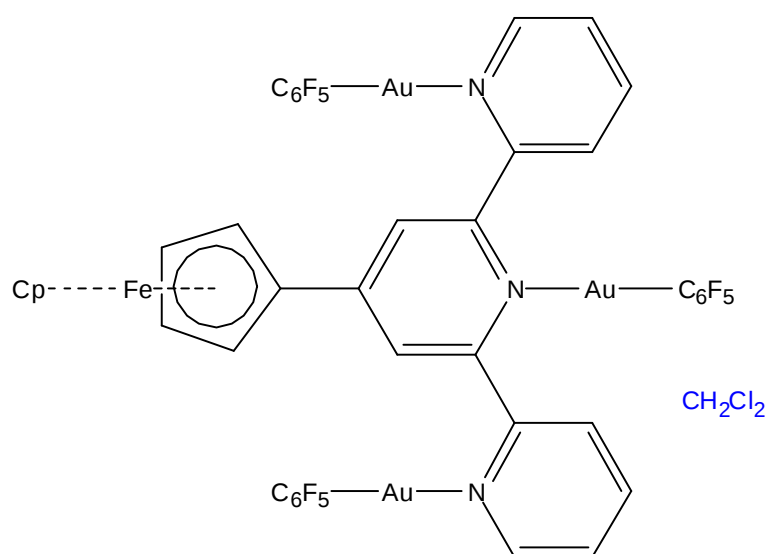
(12), 3.51, 17.1 [M. C. Torralba, P. Ovejero, M. J. Mayoral, M. Cano, J. A. Campo, J. V. Heras, E. Pinilla and M. R. Torres, *Helv. Chim. Acta*, 2004, **87**, 250] ATOWEL



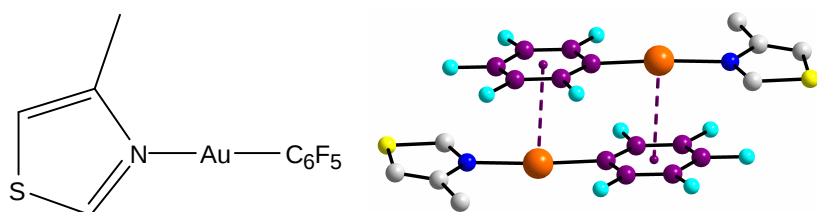
(13), 3.54, 12.7 [D. Fan, C.-T. Yang, J. D. Ranford, J. J. Vittal and P. F. Lee, *Dalton Trans.*, 2003, 3376] ITELEY



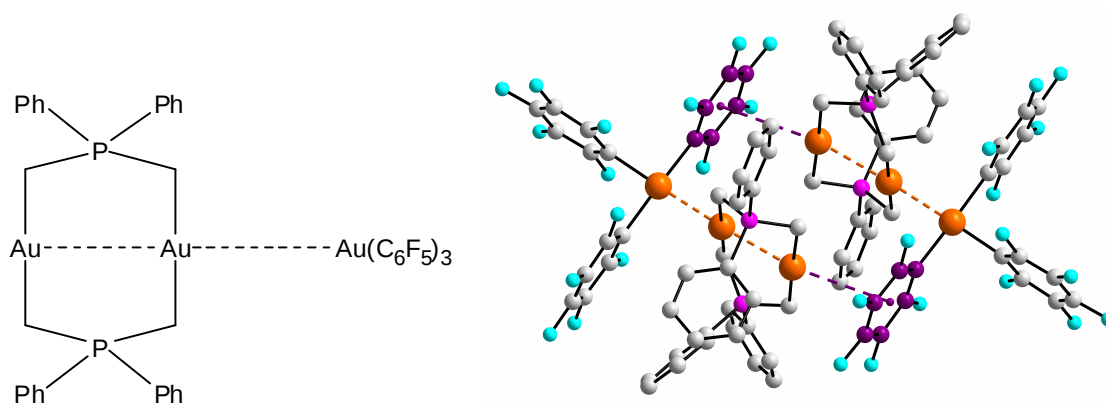
(14), 3.58, 7.5 [E. Aguado, M. J. Calhorda, M. C. Gimeno and A. Laguna, *Chem. Commun.* 2005, 3355] QARFIZ



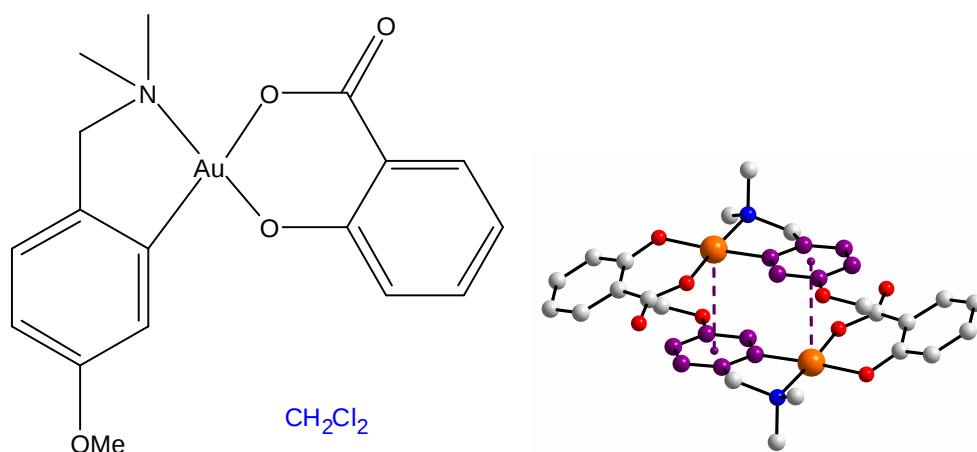
(15), 3.59, 16.6 [S. Cronje, H. G. Raubenheimer, H. S. C. Spies, C. Esterhuysen, H. Schmidbaur, A. Schier and G. J. Kruger, *Dalton Trans.*, 2003, 2859] IKABAX



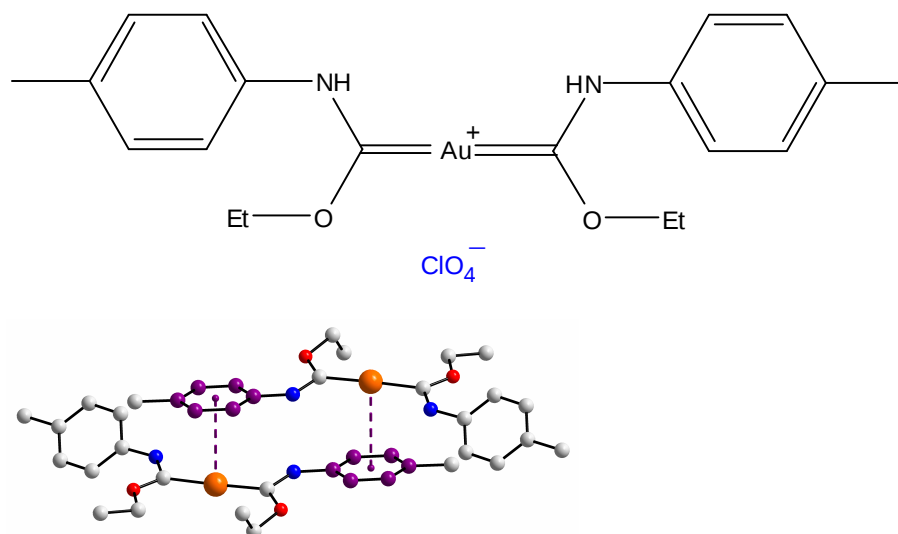
(16), 3.61, 11.3 [R. Uson, A. Laguna, M. Laguna, M. T. Tarton and P. G. Jones, *Chem. Commun.*, 1988, 740] GAVXEG



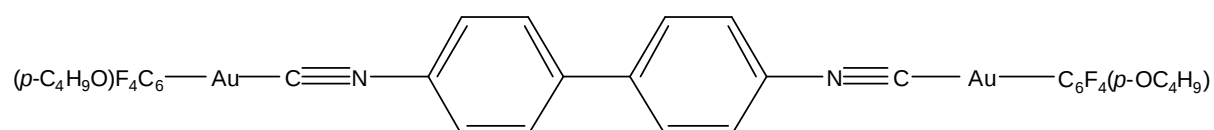
(17), 3.62, 5.2 [M. B. Dinger and W. Henderson, *J. Organomet. Chem.*, 1998, **560**, 233] GOVBUO

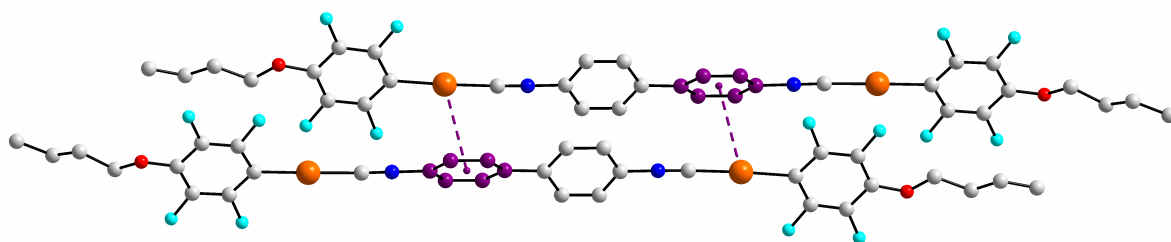


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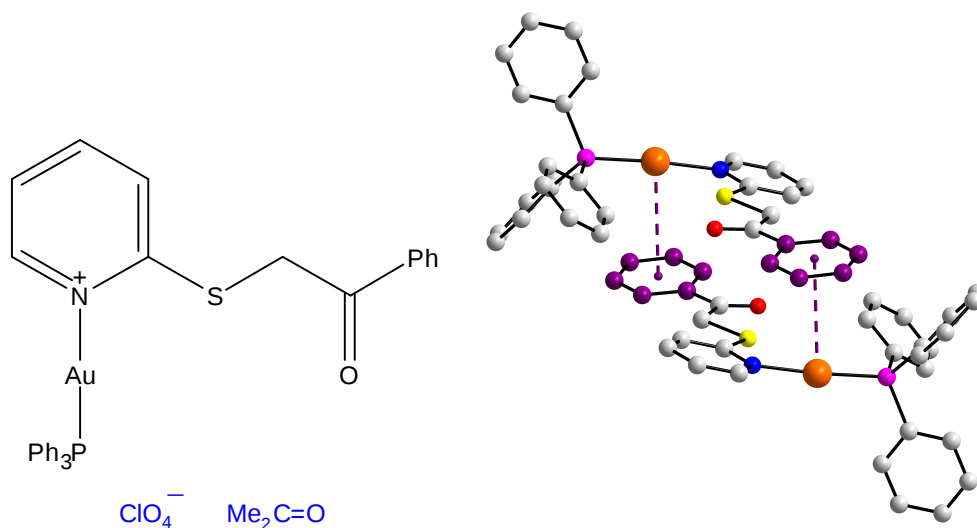


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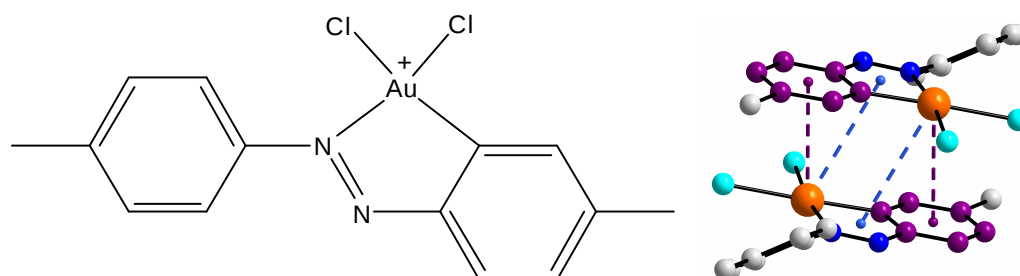




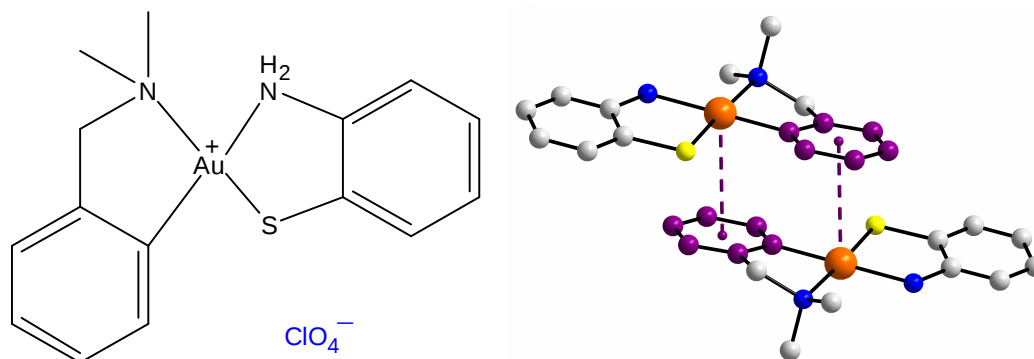
(20), 3.68, 9.1 [J. Vicente, M.-T. Chicote, S. Huertas, M. C. Ramirez de Arellano and P. G. Jones, *Eur. J. Inorg. Chem.*, 1998, 511] NUGGOL



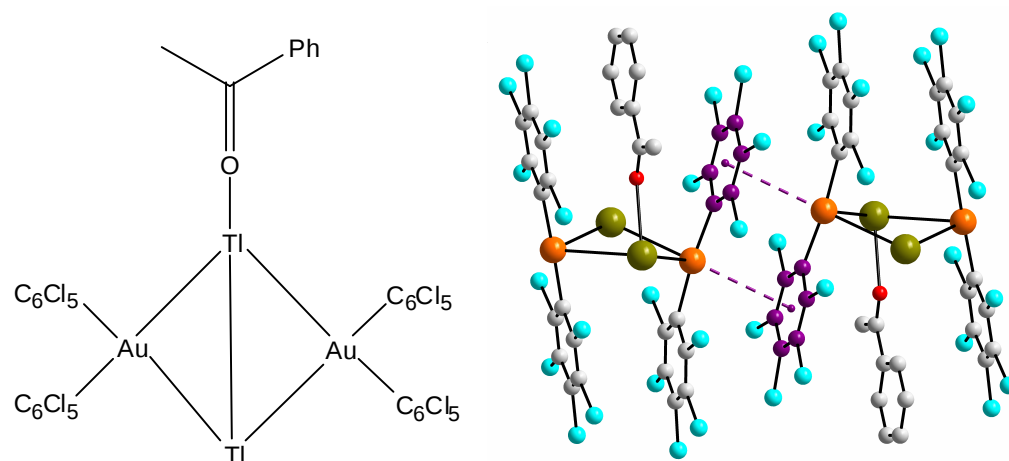
(21), 3.73, 16.8 [J. Vicente, M.-D. Bermudez, M.-P. Carrillo and P. G. Jones, *J. Chem. Soc., Dalton Trans.*, 1992, 1975] KUDSAD



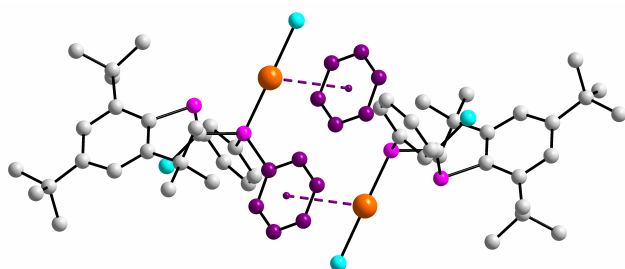
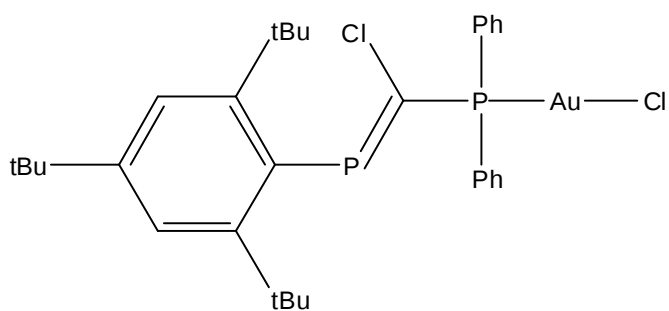
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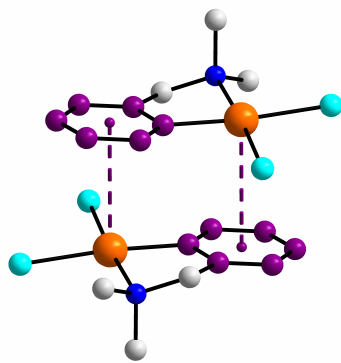
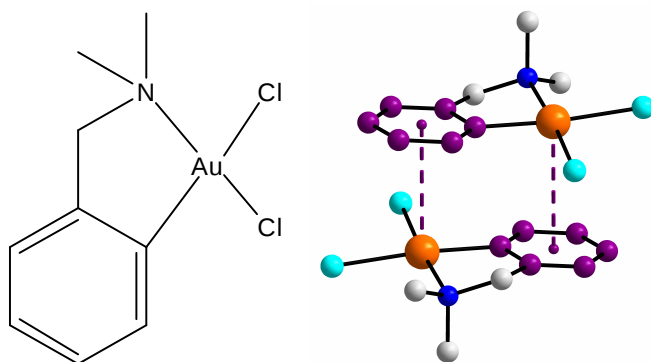
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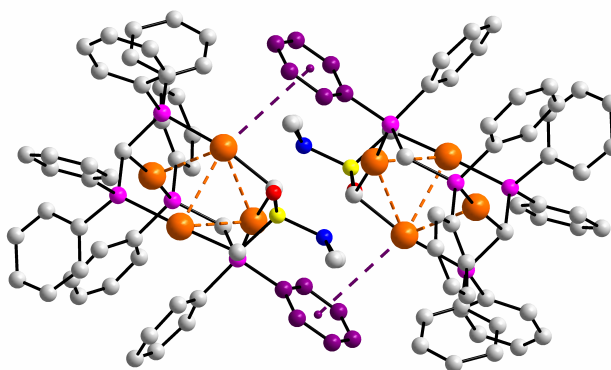
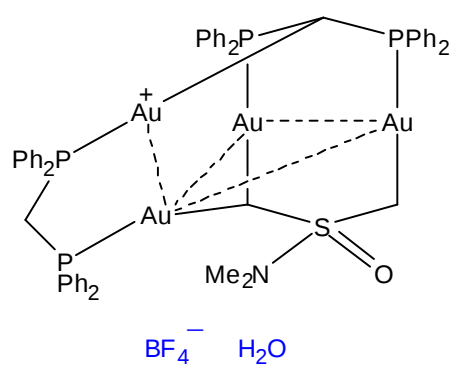
(24), 3.80, 12.8 [M. Freytag, S. Ito and M. Yoshifuji, *Chem. Asian J.*, 2006, **1**, 693] JETCIV



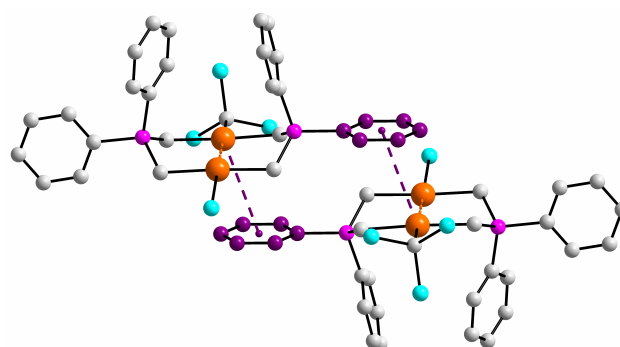
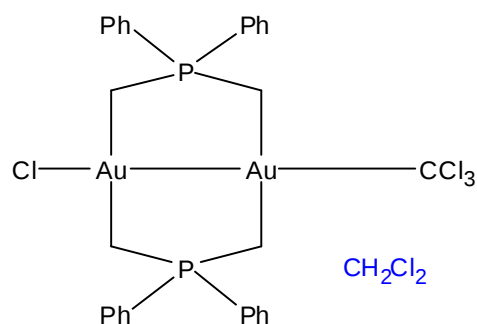
(25), 3.85, 12.2 [J. Mack, K. Ortner, U. Abram and R. V. Parish, *Z. Anorg. Allg. Chem.*, 1997, **623**, 873] RAFZON



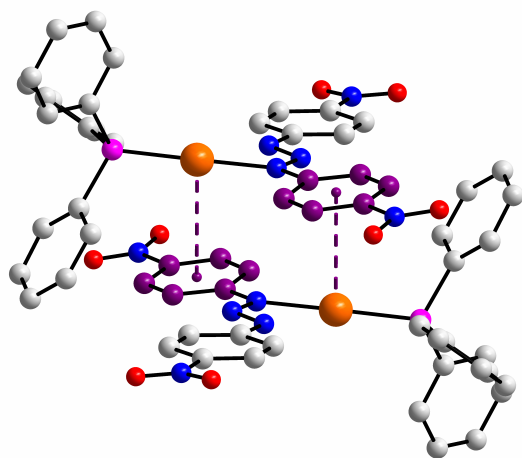
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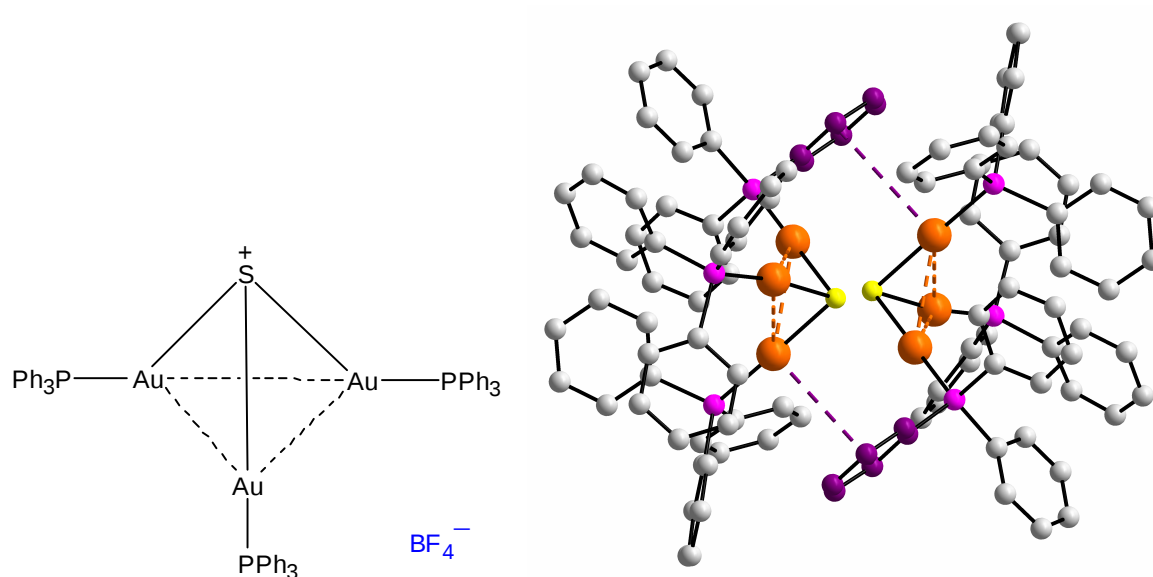


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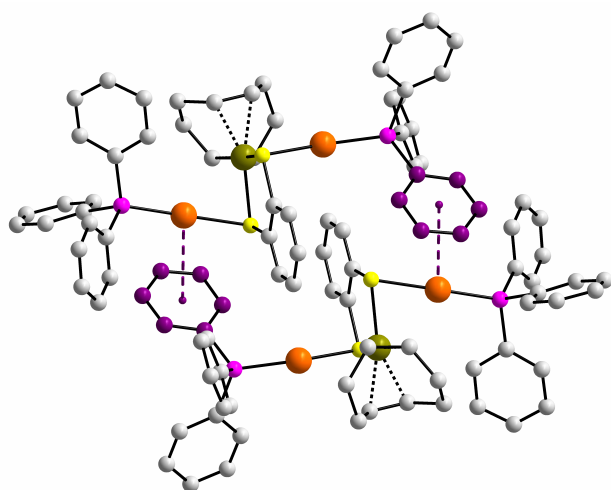
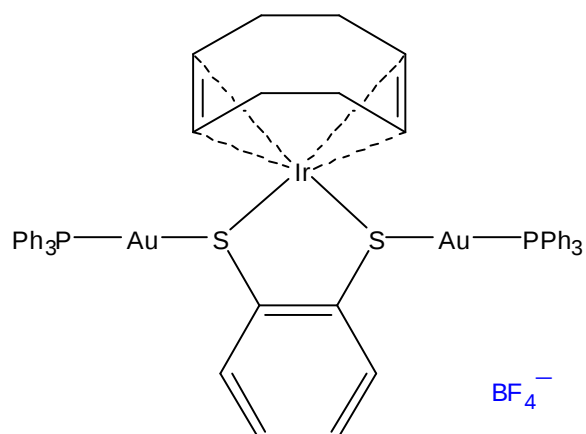


The figure displays the chemical structure and a 3D ball-and-stick model of a gold complex. The chemical structure on the left shows a central gold atom (Au) coordinated to a chlorine atom (Cl), a phenyl group (Ph), and a phosphorus atom (P). The phosphorus atom is further coordinated to two phenyl groups (Ph) and a 2-formylphenyl group (a benzene ring with a formyl group, CHO, at the ortho position). The 3D model on the right illustrates the spatial arrangement of the atoms, with gold (orange), phosphorus (purple), carbon (grey), oxygen (red), and hydrogen (white) atoms. Dashed lines indicate non-covalent interactions, such as hydrogen bonding between the formyl group and the phenyl rings.

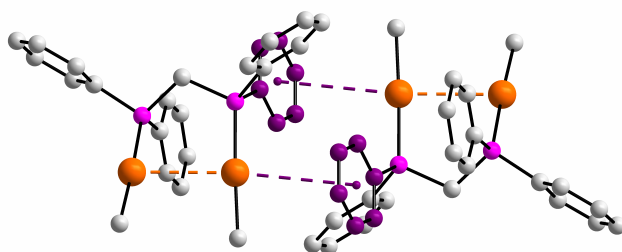
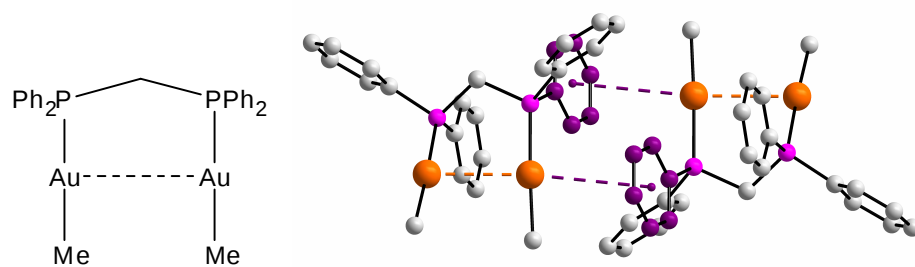
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(32), 3.97, 18.1 [X. Hong, K.-K. Cheung, C.-X. Guo and C.-M. Che, *J. Chem. Soc., Dalton Trans.* 1994, 1867] LEYMAD



(33), 3.99, 19.0 [W. Schneider, A. Bauer and H. Schmidbaur, *J. Chem. Soc., Dalton Trans.*, 1997, 415] REXRUH

