Influence of π – iodide intermolecular interactions on electronic properties of tin(IV) iodide semiconducting complexes

Supporting Information

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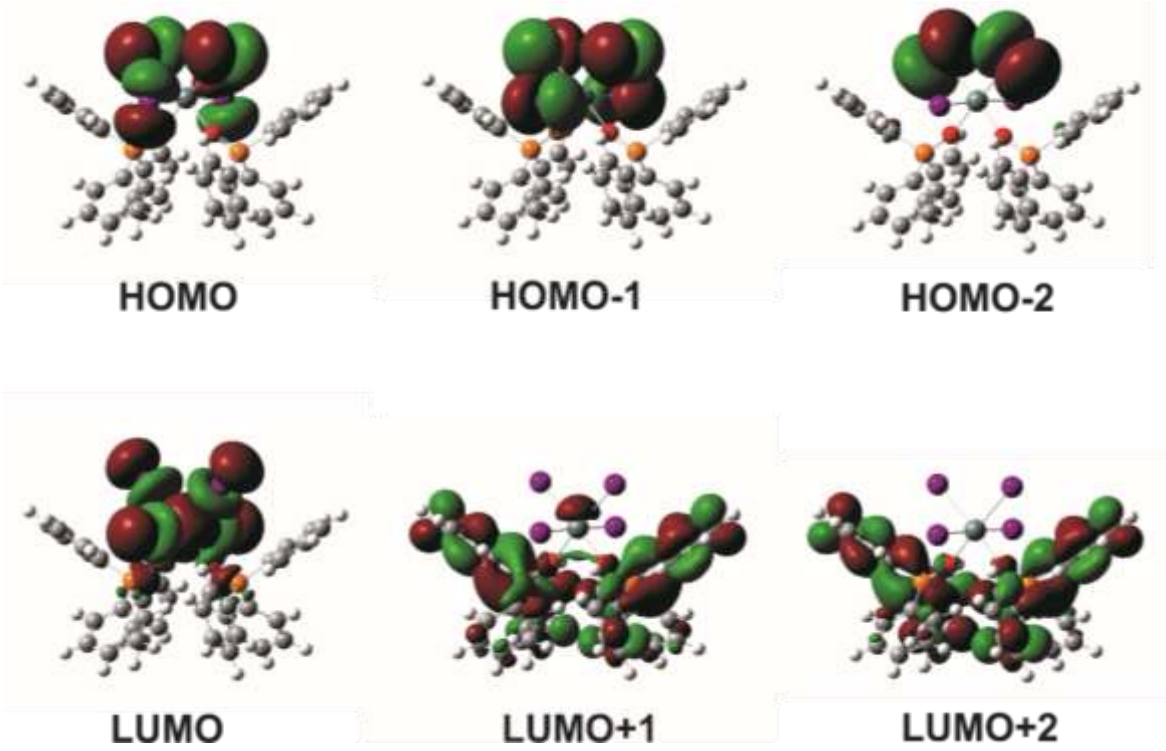

Fig. S1 Visualisation of frontier molecular orbitals of $[(C_6H_5)_3PO]_2SnI_4$

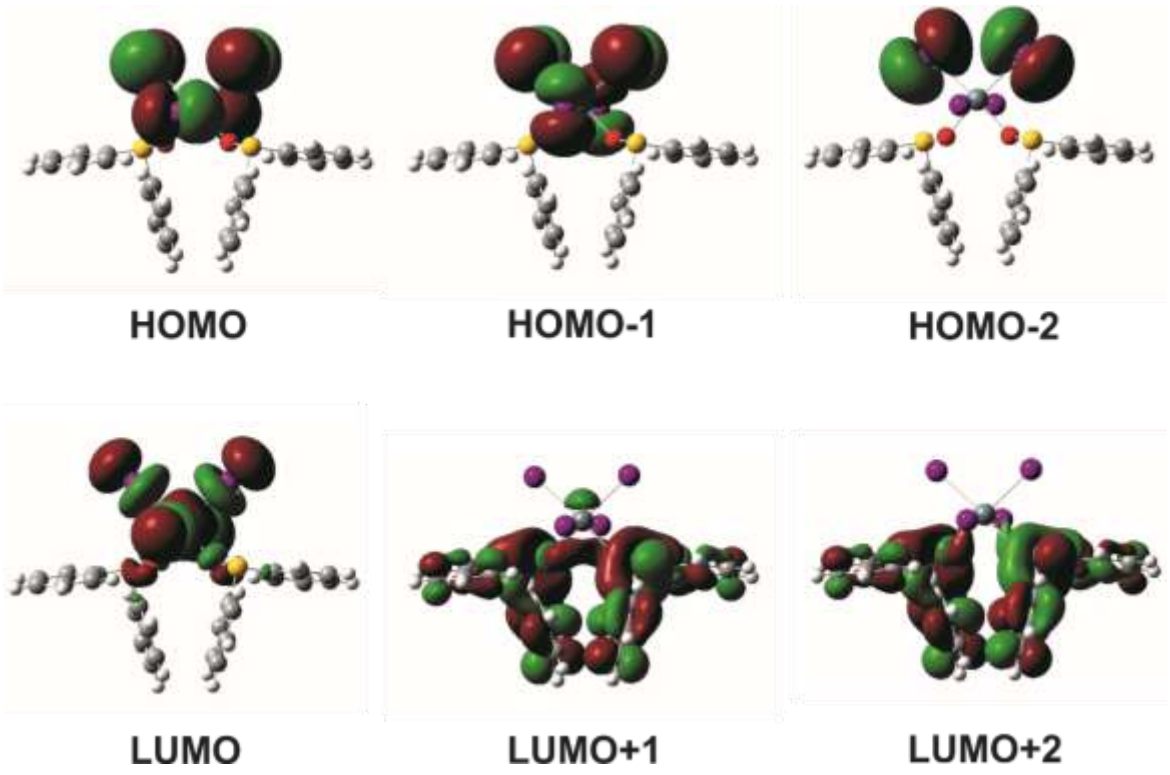

Fig. S2 Visualisation of frontier molecular orbitals of $[(C_6H_5)_2SO]_2SnI_4$

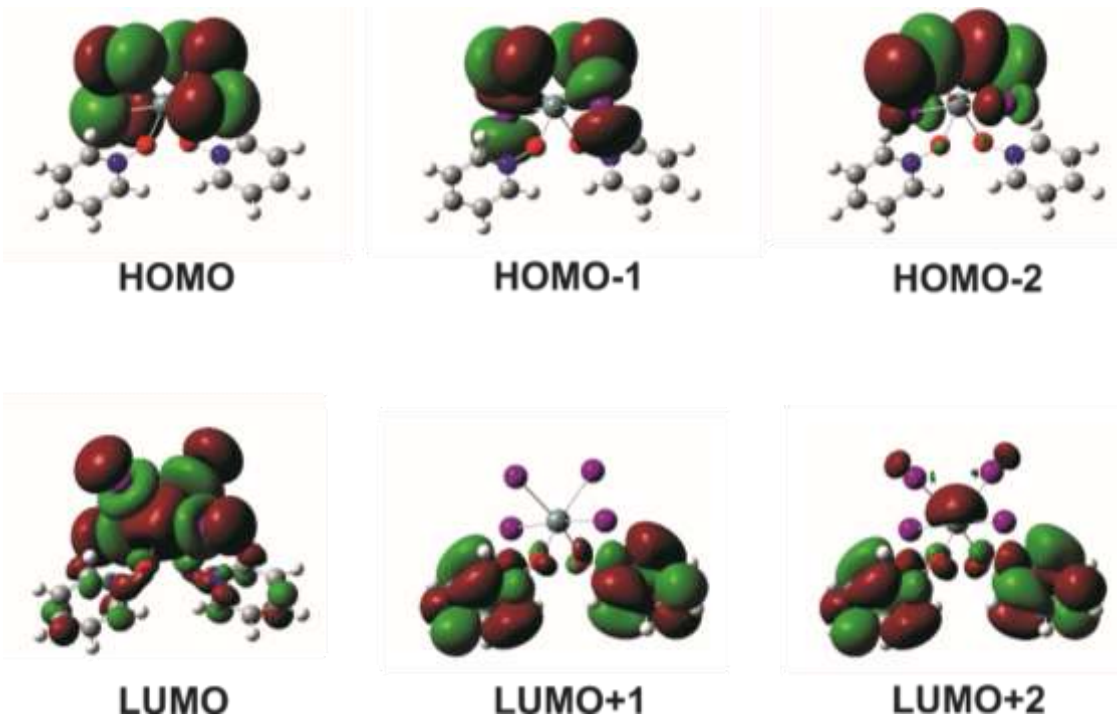

Fig. S3 Visualisation of frontier molecular orbitals of $[C_5H_5NO]_2SnI_4$