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a) Prior to the submission of the manuscript, the author(s) should validate and deposit files in Crystallographic Information File (CIF) format.

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Authors should deposit the data corresponding to each structure to be reported in the intended publication with the relevant Date Centre Data for metal organic structures, i.e. those which contain organic carbon in any of the species present in the structure, (including metal carbonyls) should be sent to: The Cambridge Crystallographic Data Centre (CCDC), by e-mail (address:deposit@ccdc.cam.ac.uk), preferably in CIF format. A checklist of data items to be included in the deposition can be obtained from the CCDC home page (http://www.ccdc.cam.ac.uk/), or by e-mail (fileserv@ccdc.cam.ac.uk), with the one line message: sendme checklist.

Data for inorganic compounds should be sent to: Fachinformationszentrum Karlsruhe (FIZ), by e-mail (address: crysdata@FIZ-Karlsruhe.de), by FTP (please contact FIZ if you choose this option), or on disk. No hard copy submissions will be accepted.

Within a day or two of submission the Data Centres will provide deposition codes for each data set. This should be noted in the Supplementary Material section at the end of the text file before the reference list in the manuscript submitted to the journal. The deposited data will be accessed by the journal and checked as part of the refereeing process.

b) When the paper is submitted to the Journal, the following guidelines should be adhered to: The Title should contain reference to the presence in the paper of X-ray crystallographic results. The Abstract should not contain crystal data, but should include a concise statement of the main features of the structural results. The Experimental results must be presented in a concise format.

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(ii) Chemical Formula. This should correspond to the complete chemical unit encompassing crystallographic symmetry (e.g., a centrosymmetric dimer should be included in full form). The use of fractional coefficients (e.g. 0.5H2O), should be used, only, if the unit has partial occupancy of its site. Formulae should be presented in a way that molecules, ionic fragments, solvate molecules etc. are separately identified.
(iii) Unit cell parameters with esds and X-ray wavelength used.
(iv) Crystal system, space group and number of complete chemical units (see ii) per cell.
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Note. Easily derived parameters should not be given unless required in discussion of other data (e.g. number of solvent molecules, etc).

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