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Title:	Use of Ferrocene Derivatives as Linkers in Forming Metal Organic Frameworks (MOFs): Synthesis, Structural Characterization and Properties
Authors:	Agrawal, Amita (/jspui/browse?type=author&value=Agrawal%2C+Amita)
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Abstract:	Based on ferrocene derivatives as linkers, new metal organic frameworks (MOFs) with a general formula $[M_4(tpxn)_2(FcDC)_2(H_2O)_4]^{4+}$ ($FcDC = 1,1'$ -ferrocenedicarboxylate; $tpxn = tppn, tpbn$, $tpen$ where $tppn = N,N',N'',N'''$ -tetrakis-(2-pyridylmethyl)-1,3-diaminopropane, $tpbn = N,N',N'',N'''$ -tetrakis-(2-pyridylmethyl)-1,4-diaminobutane, $tpen = N,N',N'',N'''$ -tetrakis-(2-pyridylmethyl)-1,5-diaminopentane, $M^{2+} = Cu^{2+}$ (1, 2, 3), Cd^{2+} (4), Co^{2+} (6) and Mn^{2+} (7)) are reported. These are heterometallic MOFs. When Fe^{3+} is chosen as the metal ion a homometallic MOF, $\{[Fe_2(\mu-O)(tpbn)(FcDC)](ClO_4)_2 \cdot 4H_2O\}_n$ (5), is the product. Use of 1,1'-ferrocene disulphonate ($FcDS$) as the linker provides $[Cu_2(tpbn)(FcDS)_2]_n \cdot 4CH_3CN \cdot 2H_2O$ (8a) and $\{[Fe_2(\mu-O)(tpbn)(FcDS)](ClO_4)_2 \cdot 4H_2O\}_n$ (9). All complexes were characterised by elemental analysis, FTIR spectroscopy, UV-visible spectroscopy, mass spectrometry, thermal gravimetric analysis, differential scanning calorimetry, single crystal and powder X-ray diffractometry. 1, 2, 3, 4, 6 and 7 are found to be discrete rectangular MOFs, 9 is a 1D coordination polymer, 5 is a 2D coordination polymer while 8a is a 2D MOF
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