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Multimodal Transportation Infrastructure Investment and Regional Economic Development: A Structural Equation Modeling Empirical Analysis in China from 1986 to 2011

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ACCEPTED MANUSCRIPT

Structural characterization and physicochemical features of a new arsenate salt

templated by mono and di-protonated 4-aminopyridine cations:

$(C_5H_7N_2)(C_5H_8N_2)[AsO_4]$ ·H₂O

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Abstract

Single crystals of a new organic–inorganic hybrid compound, with the formula (4-APH)(4-APH₂)[AsO₄]·H₂O, was synthesized at room temperature by slow evaporation method and characterized by X-Ray diffraction at 150 K, DSC–TG measurements, FT-IR and Raman spectroscopies. The title salt, $(C_5H_7N_2)(C_5H_8N_2)$ [AsO₄]·H₂O, contains mono and diprotonated 4-aminopyridine cations, an arsenate trianion and one water molecule. The diprotonated 4-ammoniumpyridinium dication $[C_5H_8N_2]^{2+}$ is disordered over two positions with refined site occupancies of 0.73 and 0.27 however the monoprotonated 4-aminopyridinium cation $[C_5H_7N_2]^+$ is ordered. The 4-aminopyridinium rings are essentially planar and occur in stacks along *b* axis. In the crystal, the As^{III} atom is coordinated by four O atoms in a slightly distorted tetrahedral geometry. The arsenate O atoms link the 4-aminopyridinium cations and water molecules into a three-dimensional network via intermolecular O–H…O and N–H…O hydrogen bonds. Additionally, in this structure, the different types and the nature of aromatic–

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