Ethene-*trans*-1,2-bis(4-pyridinium) Dihydrogenphosphite and Dihydrogenphosphate Compounds Exhibiting Cooperative and Directed Hydrogen Bonds between Cations and Anions: H_2 bpe $(H_2PO_3)_2$ and H_2 bpe $(H_2PO_4)_2 \cdot H_2O$

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Z. Naturforsch. 61b, 1067-1071 (2006); received March 27, 2006

 $H_2bpe(H_2PO_3)_2$ (1) and $H_2bpe(H_2PO_4)_2 \cdot H_2O$ (2) ($H_2bpe = ethene-trans-1,2-bis(4-pyridinium)$, $H_2PO_3 = dihydrogenphosphite, and <math>H_2PO_4 = dihydrogenphosphate$) have been prepared and structurally characterized. In compound 1, the dihydrogenphosphite anions form dimers, with a P···P distance of 4.2073(7) Å, by two O–H···H hydrogen bonds, and the dimeric dihydrogenphosphite units interact with the H_2bpe cations by way of N–H···O and O–H···O hydrogen bonds, resulting in a one-dimensional chain. The chains are held together by C–H···O interactions. In compound 2, the phosphate ions are connected by O–H···O hydrogen bonds into an unusual 2D square gridtype framework with P···P separations ranging from 4.7533(7) to 4.9506(8) Å. The H_2bpe cations crosslink the dihydrogen phosphate layers by N–H···O hydrogen bonds, forming a three-dimensional supramolecular network with channels. The water molecules in compound 2 occupy these channels and make O–H···O bonds to adjacent phosphate O atoms and also O–H···O bonds to the next water O atom in the channel.

Key words: Dihydrogenphosphites, Dihydrogenphosphates, Ethene-*trans*-1,2-bis(4-pyridinium) Cations, Hydrogen Bonds, Crystal Structure