

THEORETICAL STUDY ON FERMI RESONANCE OF THE EXCESS PROTON VIBRATION IN BINARY CLUSTERS

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In protonated binary clusters, location and magnitude of delocalization of the excess proton is shown to correlate to the difference of the proton affinities of the two components^a. Theoretical calculations on a series of protonated binary clusters have shown that Fermi resonance can occur in clusters containing amines. For examples, in $(\text{CH}_3)_3\text{N-H}^+-\text{X}$, the N-H^+ stretching frequency is high than the first overtone of N-H^+ bending modes. As the N-H^+ stretching frequency decreases by increasing the proton affinity of X, Fermi resonance inevitably shall occur. At least both N-H^+ bending modes seem to couple with the N-H^+ stretching mode; we performed multi-dimensional mode analyses to understand this coupling.

^aJ. R. Roscioli, L. R. McCunn, M. A. Johnson *Science***316**, 249 (2007).