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However, empirical assignments of vibrational modes for peaks in the fingerprint region are difficult. The C=C stretching vibrations are generally observed at 1400-1600 cm<sup>-1</sup> in benzene derivatives which are assigned to ring vibrations.

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phine, and of the infrared (3650-200 cm<sup>-1</sup>) and Raman spectra of dichlorophenylphosphine.

Vibrational Spectra of Benzene Derivatives. G. Varsányi. \$69.99; \$69.99; Publisher Description. In organic chemistry a significant amount of knowledge has been obtained since the middle of the last century about the structure of molecules by the use of methods based on chemical properties. The study of the spectra of molecules has led to ...

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In the halogenobenzenes the Raman data support a value of ca. 1320cm<sup>-1</sup>, but in 610 Vibrational spectra of benzene derivatives II benzonitrile the slightly lower value of 1310 cm<sup>-1</sup> seems most likely, the other frequencies in this region being explained as

shown in Table 1.

Genre/Form: Electronic books: Additional Physical Format: Print version: Varsanyi, G. *Vibrational Spectra of Benzene Derivatives*. Oxford : Elsevier Science, ©1969

*Vibrational spectra of benzene derivatives—IX: o ...*

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The value 588  $\text{cm}^{-1}$  has been retained *Vibrational spectra of benzene derivatives IX. o-disubstituted compounds* 1916 [4, 5] for Vig (O2) although the calculated value is 520  $\text{cm}^{-1}$ , and this very weak infrared band could arise from  $2 \times 296$  ( $A^{\wedge}$ ). Small changes have been made in some of the other values.

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Note Continuation of the author's Vibrational spectra of benzene derivatives. Vol. 2 rev. by K. Lempert. "A Halsted Press book." ISBN 0470903309

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