

## Raman Spectrum of Gaseous Tin(II) Chloride

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*Summary* The first vibrational spectrum of a discrete, monomeric, bent  $\text{MX}_2$  species (where X = halogen) is reported.

THE Raman spectrum of tin(II) chloride was obtained using 5145 Å laser excitation of one atmosphere of compound contained in a silica ampoule at 650°. The spectrum (Figure) is consistent with a bent molecule. Two of the predicted three bands for a  $C_{2v}$  point group were observed, both of them being polarised. The (less intense) asymmetric stretch was probably obscured by the peak assigned

to the symmetric stretch. Linear metal dihalides (*e.g.* zinc chloride) have been observed to give only one polarised band in the Raman spectrum, in agreement with prediction.<sup>1</sup> Vapour density and electron diffraction studies<sup>2,3</sup> on tin chloride have shown that it is predominantly monomeric under the conditions used in these experiments; it was thought that the molecules were bent, but the valence angle was not determined.

Due to extensive polymerisation, the observed Raman spectra of solid and molten tin chloride are quite different to the gas phase spectrum.<sup>4</sup> The frequencies of vibration

of discrete tin chloride molecules have been measured by chemiluminescent emission spectroscopy. However, the method is fraught with difficulty in the assignments of band progressions to particular molecular species, electronic states and vibrational modes, and only in one case is there good agreement with the frequencies obtained by the more direct method of this investigation (Table).

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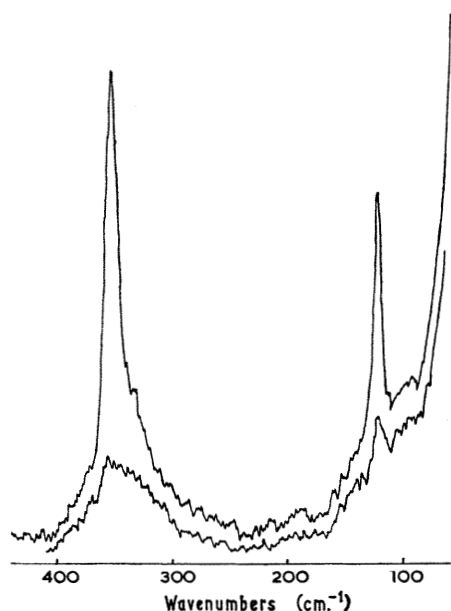
Vibrational mode	Calculated			Observed for a gas		
	Ref. 7	Ref. 5	Ref. 8	Ref. 5	Ref. 6	This work
$\nu_1$	398	435	—	377	$350 \pm 5$	$352 \pm 1s,p$
$\nu_2$	102	205	116	207	$120 \pm 5$	$120 \pm 1m,p$
$\nu_3$	406	—	—	—	—	(356)

Ref. 4	Solid		Melt
		This work	Ref. 4
193s		192sh 162s	228s,p
90s		125m 95sh 83s	112s,p

The vibrations are labelled according to convention, *i.e.*,  $\nu_1$ , symmetric stretch;  $\nu_2$ , deformation;  $\nu_3$ , asymmetric stretch. Frequencies are quoted in wavenumbers ( $\text{cm}^{-1}$ ); p denotes polarised.

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FIGURE

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<sup>6</sup> D. Naegeli and H. B. Palmer, *J. Mol. Spectroscopy*, 1966, **21**, 325.

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