

## **SYNTHESIS AND INFRARED STUDY OF SOME NEW IODATO ADDUCTS AND DERIVATIVES**

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**Abstract:** Three iodato adducts and one derivative have been synthesized and studied by infrared data. The suggested structures are discrete, the iodate behaving as a mono- or bidentate ligand, or an infinite chain with bridging iodate, the environment around the tin centre being trigonal bipyramidal, tetrahedral or octahedral.

**Keywords:** *coordinating iodate, discrete and infinite chain structures, trigonal bipyramidal, tetrahedral, octahedral environments*

## INTRODUCTION

The organotin (IV) molecules are known for their applications as wood preservatives, antifouling paints, drugs etc [1]. That is the reason why several groups including our [2-15] have been focusing their search in the synthesis of new compounds of this family. We have initiated in this work the study of the interactions between  $\text{Et}_4\text{NIO}_3$  and  $\text{SnPh}_3\text{Cl}$ ,  $\text{SnPh}_3\text{Br}$ ,  $\text{SnMe}_2\text{Cl}_2$  or  $\text{SnCl}_4$  which have given four new adducts and derivatives infrared study of which has been carried out then structures suggested on the basis of spectroscopic data.

## MATERIALS AND METHODS

$\text{Et}_4\text{NIO}_3$  has been obtained as a powder on neutralizing  $\text{IO}_3\text{H}$  with a 20% water solution of  $\text{Et}_4\text{NOH}$  and allowing water to evaporate at  $60^\circ\text{C}$ .

**Compound A:** the mixture of 0.7167 g of  $\text{SnPh}_3\text{Cl}$  in acetone and 0.5670 g of  $\text{Et}_4\text{NIO}_3$  in acetone gives after slow evaporation colourless crystals.

**Compound B:** The mixture of 0.8652 g of  $\text{SnMe}_2\text{Cl}_2$  in the ethanol and 1.2010 g of  $\text{Et}_4\text{NIO}_3$  in ethanol gives a white precipitate.

**Compound C:** The mixture of 0.7070 g of  $\text{SnCl}_4$  in benzene and 1.2010 g of  $\text{Et}_4\text{NIO}_3$  in benzene gives a white precipitate.

The analytical data reported below, have allowed to suggest the following formulae (Table 1).

**Table 1.** Suggested formulae of synthesized compounds and the elemental analyses

Compound	Chemical formula	Elemental analysis (%)					
		C		H		N	
		calc.	found	calc.	found	calc.	found
A	$\text{SnPh}_3\text{IO}_3 \cdot \text{SnPh}_3\text{Cl}$	47.49	48.08	3.32	4.19	-	-
B	$\text{SnMe}_2(\text{IO}_3)_2 \cdot 2/5\text{EtOH}$	6.50	6.37	1.64	1.62	-	-
C	$\text{SnCl}_4(\text{Et}_4\text{NIO}_3)_{1.5}$	20.46	20.78	4.34	4.51	2.70	2.59

The elemental analyses were performed by the laboratory of Microanalyses – University of Padova – Italy. The infrared spectra have been recorded at the University of Padova - Italy - by means of a Bruker FTIR spectrometer using CsI windows, the sample being as Nujol mulls. Infrared data are given in  $\text{cm}^{-1}$  (abbreviations: (vs) very strong, (s) strong, (m) medium, (w) weak, (vw) very weak). All the chemicals were purchased from Aldrich – Germany - and used as such.

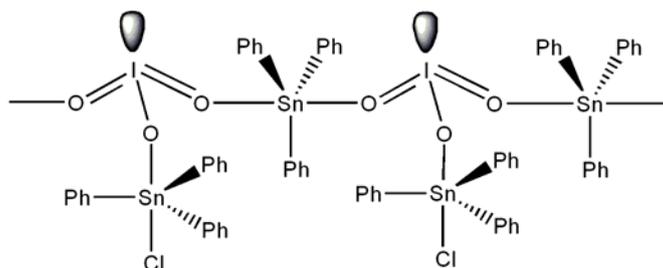
## RESULTS AND DISCUSSION

Let us consider the infrared data (in  $\text{cm}^{-1}$ ) of the studied compounds:

(A):  $\nu_s \text{IO}_3 = 785$  (vs);  $\nu_{\text{as}} \text{IO}_3 = 730$  (vs),  $700$  (vs);  $\delta \text{IO}_3 = 350$  (m);  $\nu_{\text{as}} \text{SnCl}_3 = 270$  (s);  $\nu \text{SnCl} = 235$  (m);  $\nu \text{SnO} = 215$  (m);

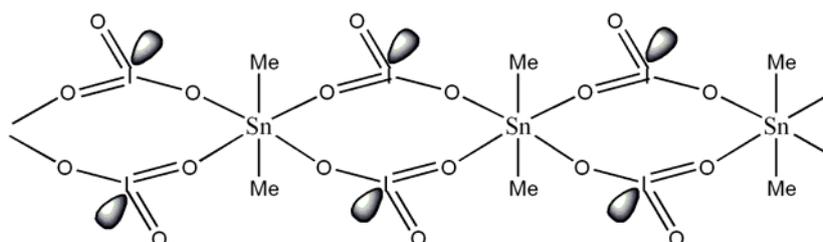
(B):  $\nu_s \text{IO}_3 = 800$  (s);  $\nu_{as} \text{IO}_3 = 745$  (vs),  $700$  (vs);  $\nu_{as} \text{SnC}_2 = 590$  (w);  $\nu \text{SnO} = 230$  (w);  
 (C):  $\nu_s \text{IO}_3 + \nu_{as} \text{IO}_3 = 790$  (br);  $\nu_{as} \text{SnCl}_4 = 294$  (vs).

From all these data we can reasonably suggest for compound **A** an infinite chain of  $\text{SnPh}_3\text{IO}_3$  to which  $\text{SnPh}_3\text{Cl}$  coordinates through the free oxygen atom (Figure 1).



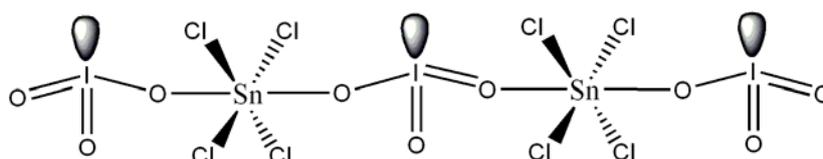
**Figure 1.** Suggested structure for compound **A**

The absence of  $\nu_s \text{SnMe}_2$  in the spectrum of the derivative **B** is an indication of a linear  $\text{SnMe}_2$  residue allowing to deduce an octahedral environment. The suggested structure is an infinite chain with bridging iodates (Figure 2).



**Figure 2.** Suggested structure for compound **B**

$\nu_{as} \text{SnCl}_4$  appears as a sharp and strong band (Eu type according to Group Theory) allowing to conclude to  $D_{4h}$  symmetry for  $\text{SnCl}_4$ . The suggested structure is discrete with one bridging and two monocoordinating iodates (Figure 3).



**Figure 3.** Suggested structure for compound **C**

## CONCLUSIONS

The studied iodato adducts have discrete or chain structures the environment around tin centre being octahedral, trigonal bipyramidal, the iodate behaving as a monocoordinating or a bridging ligand.

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