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## Preparation of RSn(I)-Sn(I)R with Two Unsymmetrically Coordinated Sn(I) Atoms and Subsequent Gentle Activation of P-4

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# Preparation of RSn(I)-Sn(I)R with Two Unsymmetrically Coordinated Sn(I) Atoms and Subsequent Gentle $Activation \ of \ P_4$

Shabana Khan, Reent Michel, Johannes M. Dieterich, Ricardo A. Mata, Herbert W. Roesky, Hean-Philippe Demers, Adam Lange, and Dietmar Stalke

<sup>a</sup>Institut für Anorganische Chemie, Georg-August-Universität, Tammannstrasse 4, D-37077, Göttingen, Germany.

<sup>b</sup>Institut für Physikalische Chemie, Georg-August-Universität, Tammannstrasse 6, D-37077, Göttingen, Germany.

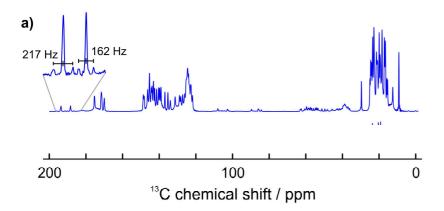
<sup>c</sup>Max-Planck-Institut für biophysikalische Chemie, Am Faßberg11, 37077 Göttingen, Germany

E-mail: hroesky@gwdg.de, dstalke@chemie.uni-goettingen.de, rmata@gwdg.de

#### **Content:**

- (S1). Solid-state NMR of 2
- (S2). Computational details of 2

#### (S1). Solid-state NMR:



**Figure S1.**  $^{13}$ C solid-state cross-polarization NMR spectrum of a bis(stannylene) **2**. The excerpt highlights the one-bond carbon-tin J couplings. The integrals of the satellite peaks (18.9% and 15.9% of total volume) correspond to the combined natural abundance of  $^{117}$ Sn (7.61%) and  $^{119}$ Sn (8.58%). The spectra were recorded at 8.2 KHz MAS on a 9.4 T spectrometer.

#### (S2). Computational Results (Structures and Density Plots)

#### Optimized B3LYP/def2-SVP structure of compound 2

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#### Optimized B3LYP/def2-SVP structure of 3- and 4-coordination model

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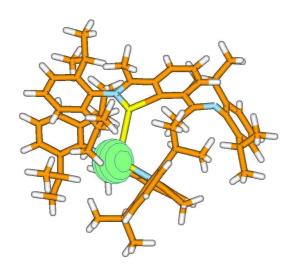
#### Optimized B3LYP/def2-SVP structure of 3- and 4-coordination model

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**Figure S2.** Isosurface plots of the two Sn NLMOs derived from the Sn NBO lone pairs in compound **2**.