

Phosphorus, Sulfur, and Silicon and the Related Elements

ISSN: 1042-6507 (Print) 1563-5325 (Online) Journal homepage: http://www.tandfonline.com/loi/gpss20

Synthesis and Structure of a Kinetically Stabilized 1,3,6-Triphosphafulvene

Hiroki Sugiyama , Shigekazu Ito & Masaaki Yoshifuji

To cite this article: Hiroki Sugiyama, Shigekazu Ito & Masaaki Yoshifuji (2002) Synthesis and Structure of a Kinetically Stabilized 1,3,6-Triphosphafulvene, Phosphorus, Sulfur, and Silicon and the Related Elements, 177:8-9, 2013-2014, DOI: 10.1080/10426500213341

To link to this article: http://dx.doi.org/10.1080/10426500213341

| 0-0- | | | | |
|------|---|--|--|--|
| | | | | |
| | Г | | | |
| | Г | | | |

Published online: 27 Oct 2010.



Submit your article to this journal 🕑

Article views: 6



View related articles 🗹

Full Terms & Conditions of access and use can be found at http://www.tandfonline.com/action/journalInformation?journalCode=gpss20

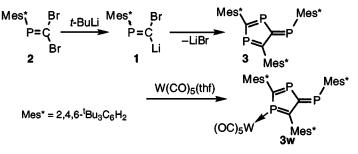


SYNTHESIS AND STRUCTURE OF A KINETICALLY STABILIZED 1,3,6-TRIPHOSPHAFULVENE

Hiroki Sugiyama, Shigekazu Ito, and Masaaki Yoshifuji Tohoku University, Japan

(Received July 29, 2001; accepted December 25, 2001)

Kinetically stabilized phosphanylidene carbenoid **1** is one of the precursors for novel low-coordinated phosphorus compounds. We report the novel trimerization of phosphanylidene carbenoid **1** affording 1,3,6triphosphafulvene **3**. Dibromophosphaethene **2** was allowed to react with 2 molar equivalent of *t*-butyllithium at -78° C to generate the phosphanylidene carbenoid **1**, and the reaction mixture was warmed to 25° C. After purification, 1,3,6-triphosphafulvene **3** was obtained as a deep red solid. The NMR and MS spectrum of **3** supported the 1,3,6-triphosphafulvene structure. Furthermore, triphosphafulvene **3** was treated with an excess amount of W(CO)₅(thf) to afford the pentacarbonyltungsten(0) complex **3w** as a deep red solid. Complex **3w** was recrystallized from toluene at 0°C to afford a suitable crystal for x-ray analysis. The x-ray crystallography of **3w** was confirmed the 1,3,6triphosphafulvene structure with the coordination of tungsten at the P3 position.^{1,*}



SCHEME 1

*In ref. [1], the phosphorus chemical shifts at the P6 position for **3** and **3w** should read $\delta_P = 399.0$ and $\delta_P = 397.1$, respectively.

Address correspondence to Masaaki Yoshifuji, Department of Chemistry, Graduate School of Science, Tohoku University, Aoba, Sendai 980-8578, Japan. E-mail: yoshifj@mail.cc.tohoku.ac.jp

REFERENCE

[1] S. Ito, H. Sugiyama, and M. Yoshifuji, Angew. Chem., Int. Ed., 39, 2781 (2000).