

## Supporting Information

### Diiiodido-bis{N-[2-(diphenylphosphino)benzylidene]benzylamine- $\kappa^2 N,P$ }dicopper(I)

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Schmp. \_\_\_\_\_ °C, Sdp. \_\_\_\_\_ °C, WO<sub>3</sub>- Zugabe  Bemerkungen: \_\_\_\_\_  
Anwesende Elemente: C H N S O Cu, I, P  
Zu bestimmen:      
Erwartete Ergebnisse:  
61,67 % C 4,38 % H 2,77 % N \_\_\_\_\_ %  
Gefundene Ergebnisse:  
Einwaage:  
2,326 mg 55,81 % C 3,97 % H 2,50 % N \_\_\_\_\_ % S  
2,231 mg 55,67 % C 3,96 % H 2,50 % N \_\_\_\_\_ % S  
\_\_\_\_\_ mg \_\_\_\_\_ % C \_\_\_\_\_ % H \_\_\_\_\_ % N \_\_\_\_\_ % S  
\_\_\_\_\_ mg \_\_\_\_\_ % \_\_\_\_\_ % \_\_\_\_\_ % \_\_\_\_\_ %  
\_\_\_\_\_ mg \_\_\_\_\_ % \_\_\_\_\_ % \_\_\_\_\_ % \_\_\_\_\_ %  
Durchgeführt: J.A. Baumann

Figure S1: Elemental analysis of the title compound.

D:\ms\_aug07\vers\fisc46\_070802140658  
tobi Fischer ES-MS (DCM/MeOH + 10 mM NH4Ac) TSQ 7000  
fisc46\_070802140658 #1 RT: 0.00 AV: 1 SM: 3B NL: 3.47E6  
T: + p ESI ms [ 199.96-1799.98]

08/02/2007 02:06:58 PM UM 184

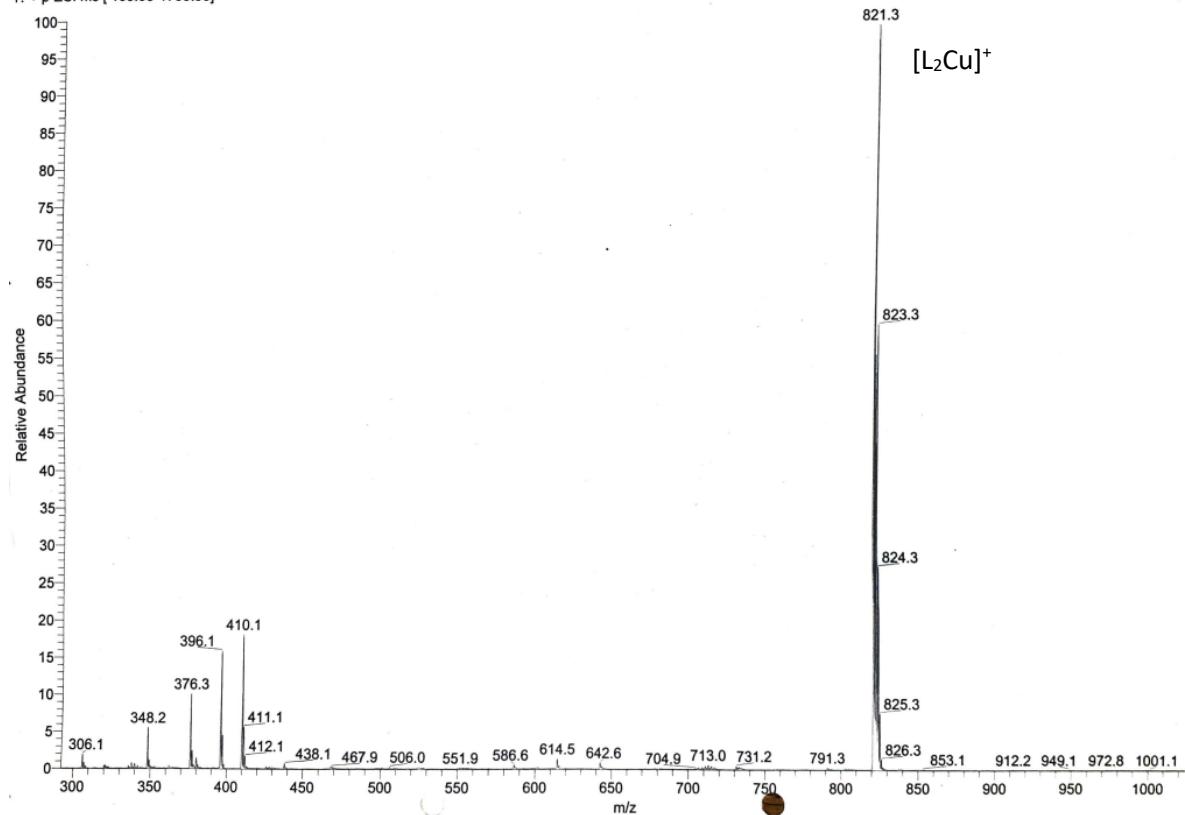


Figure S2: MS(EI) spectrum of the title compound.

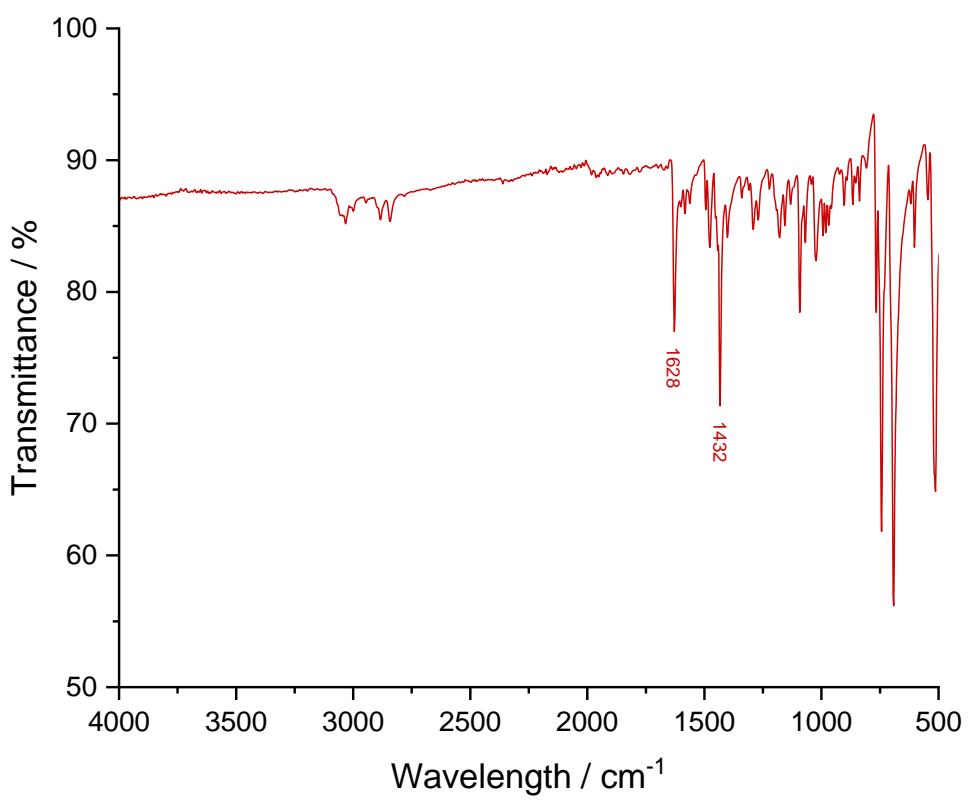


Figure S3: IR-spectrum (ATR) of the title compound.

**Table S1.** Crystal data and data collection and structure refinement details for **1**.

<b>1</b>	
Empirical formula	C <sub>52</sub> H <sub>44</sub> Cu <sub>2</sub> I <sub>2</sub> N <sub>2</sub> P <sub>2</sub>
M <sub>r</sub> , g mol <sup>-1</sup>	1139.73
Size, mm <sup>3</sup>	0.24×0.18×0.08
Crystal system	Monoclinic
Space group	P2 <sub>1</sub> /n
a, Å	15.916(3)
b, Å	9.226(1)
c, Å	15.939(3)
β, deg	95.94(2)
V, Å <sup>3</sup>	2328.0(7)
ρ <sub>calcd.</sub> , g cm <sup>-1</sup>	1.626
Z	2
μ(Mo-K <sub>α</sub> ), mm <sup>-1</sup>	2.346
T, K	297
Θ range, deg	2.80-25.91
Measured reflections	24633
Independent reflections	4505
Reflections with I > 2σ(I)	3751
Absorption correction	Analytical
T <sub>min</sub> /T <sub>max</sub>	0.8257, 0.5272
Restraints/refined	0/274
R1 (I ≥ 2σ(I))	0.0321
wR2	0.0860
ρ <sub>fin</sub> (max/min), e Å <sup>-3</sup>	0.79/-0.29
CCDC no.	2334274