New Tris(pyrazolyl)triazine and Pyrazolylpyridine Gold(I) and Palladium(II) Derivatives Based on the 3,5-Bis(4-butoxyphenyl)pyrazole Group – Architectures with Different Types of Bonding Interactions

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Keywords: Gold / Hydrogen bonds / Palladium / Supramolecular Chemistry / Triazine

In the original article[1] the footnotes of Tables 1 and 2 on pages 2694 and 2695 are erroneous. The correct version is given below.

Table 1. Selected bond lengths [Å] and angles [°] for **4** with e.s.d.s. in parentheses, including the hydrogen bond geometries

	A	B	C	
Pd−N1	2.10(2)	2.12(1)	2.12(1)	
Pd−N3	2.11(1)	2.12(1)	2.10(1)	
Pd−C9	2.10(1)	2.10(1)	2.10(1)	
Pd−C10	2.13(2)	2.10(1)	2.10(1)	
Pd−C11	2.15(2)	2.12(2)	2.12(2)	
C9−C10	1.382(5)	1.384(5)	1.384(5)	
C10−C11	1.377(5)	1.384(5)	1.386(5)	
N1−Pd−N3	91.1(3)	93.4(4)	93.3(4)	
N1−Pd−C9	98.8(6)	97.9(4)	100.0(4)	
N1−Pd−C10	132.8(6)	130.3(4)	131.1(4)	
N1−Pd−C11	167.1(6)	167.2(4)	169.3(4)	
N3−Pd−C11	100.8(6)	99.3(4)	97.1(4)	
N3−Pd−C9	170.0(6)	168.3(3)	166.1(4)	
N3−Pd−C10	132.9(6)	132.7(4)	131.2(4)	
C9−Pd−C10	38.2(3)	38.3(2)	38.7(2)	
C9−Pd−C11	69.3(5)	69.3(3)	69.8(3)	
C10−Pd−C11	37.6(3)	38.3(2)	38.4(2)	
Hydrogen bond geometries				
D−H···A	D···A	D−H	H···A	D−H···A
N2A−H2A···F9′ [a]	2.81(2)	0.86	1.97	167.9
N4A−H4A···F2	2.89(1)	0.86	2.06	162.9
N2B−H2B···F7	3.02(1)	0.86	2.19	161.7
N4B−H4B···F1	2.81(1)	0.86	1.99	174.0
N2C−H2C···F11	2.81(1)	0.86	1.91	163.8
N4C−H4C···F5	2.81(1)	0.86	2.01	175.1

[a] $x, y, z + 1$.

Table 2. Selected bond lengths [Å] and angles [°] for **5** with e.s.d.s. in parentheses, including the hydrogen bond geometries

N3−O5	1.236(9)			
N3−O3	1.245(8)			
N3−O4	1.213(8)			
Au1−N1	2.077(5)			
Au1−P1	2.239(2)			
P1−C32	1.800(8)			
P1−C26	1.826(8)			
P1−C38	1.822(7)			
Au1···O1′ [a]	3.540(6)			
N1−Au1−P1	178.1(2)			
O5−N3−O3	117.9(8)			
O5−N3−O4	121.0(8)			
O3−N3−O4	121.2(7)			
Hydrogen bond geometries				
D−H···A	D···A	D−H	H···A	D−H···A
N2−H2···O3	2.764(9)	0.95(6)	1.82(7)	171(6)
C4−H4A···O4′′[b]	3.382(9)	0.93	2.63	137.3
C40−H40A···O5′′′[c]	3.31(1)	0.93	2.88	115.1

[b] $-x + 1, -y + 1, -z + 1$. [b] $-x, -y + 1, -z + 1$. [c] $-x, -y + 1, -z$.

[1] R. M. Claramunt, P. Cornago, M. Cano, J. V. Heras, M. L. Gallego, E. Pinilla, M. R. Torres, *Eur. J. Inorg. Chem.*, **2003**, 2693−2704.

Received September 24, 2003
The Editors

 DOI: 10.1002/ejic.200300663