Supporting Information

Highly Enantioselective Michael Addition of 3-Aryloxindoles with Vinyl Phosphates Catalyzed by Cinchona Alkaloid Derived Thiourea Catalysts

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1. X-Ray Diffraction data of comopound 3u

Table 1. Crystal data and structure refinement for 1683.				
Identification code	1683			
Empirical formula	$C_{30}H_{43}NO_{11}P_2$			
Formula weight	655.59			
Temperature	296(2)			
Wavelength	0.71073 Å			
Crystal system	Orthorhombic			
Space group	P2(1)2(1)2(1)			
Unit cell dimensions	a = 8.8403(5) Å	$\alpha = 90^{\circ}$.		
	b = 11.1099(7) Å	$\beta = 90^{\circ}$.		
	c = 35.459(2) Å	γ = 90°.		
Volume	3482.6(4) Å ³			
Z	4			
Density (calculated)	1.250 Mg/m ³			
Absorption coefficient	0.180 mm ⁻¹			
F(000)	1392			
Crystal size	0.36 x 0.28 x 0.22 mm ³			
Theta range for data collection	1.92 to 25.01°.			
Index ranges	-10<=h<=10, -11<=k<=13, -42<=l<=42			
Reflections collected	40524			
Independent reflections	6107 [R(int) = 0.0337]			
Data / restraints / parameters	6107 / 5 / 397			
Goodness-of-fit on F^2	1.058			
Final R indices [I>2sigma(I)]	R1 = 0.0540, wR2 = 0.1483			
R indices (all data)	R1 = 0.0632, wR2 = 0.1578			
Largest diff. peak and hole	0.312 and -0.337 e.Å ⁻³			
	0			



2. HPLC data for the determination of the ee's of adducts 3 and 5

3a (Table 3, entry 1)



3b (Table 3, entry 2)







3c (Table 3, entry 3)





3d (Table 3, entry 4)







3e (Table 3, entry 5)

PO(OEt)₂ Br PO(OEt)₂ O Boc





3f (Table 3, entry 6)







3g (Table 3, entry 7)

PO(OEt)₂ PO(OEt)₂ O Br





3h (Table 3, entry 8)





























3q (Table 3, entry 17)



3r (Table 3, entry 18)







3s (Table 3, entry 19)







3t (Table 3, entry 20)







3u (Table 3, entry 21)



After Recrystallization of 3u:





3w (Table 3, entry 23)







3x (Table 3, entry 24)



5c (Table 4, entry 4)







5d (Table 5, entry 5)







5e (Table 5, entry 6)







5f (Table 5, entry 7)

ID

1

2

组分名

Σ:

保留时间

9.323

14.138

峰高

120523

238429

358952

峰面积

2814085.4

10304348.7

13118434.1

浓度

21.4514

78.5486

100.0000

拖尾因子



理论塔板

6 (Scheme 2)



8 (Scheme 2)





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1 H NMR of **3b**







¹⁹F NMR of **3b**











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¹H NMR of **3**k







¹⁹F NMR of **3k**







¹³C NMR of **3**l





2



S-50







¹⁹F NMR of **3n**









¹H NMR of **3r**





-100

-150

S-57

PPM

-200

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¹H NMR of **3**t



¹³C NMR of **3**t







¹⁹F NMR of **3t**





















³¹P NMR of **7**

- 26.028 - 25.998 - 25.126 - 25.093









¹H NMR of 8