

Synthesis of Fused Cyanopyrroles and Spirocyclopropanes *Via Addition of N-Ylide to Chalconimines*

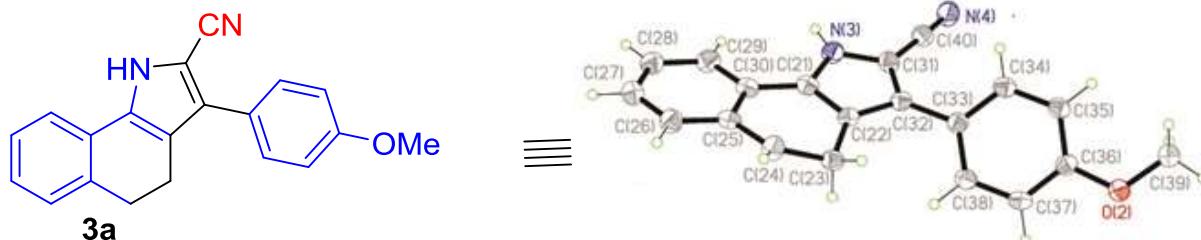
Siba Prasad Midya, Elumalai Gopi, Nishikant Satam and Irishi N. N. Namboothiri*

Department of Chemistry, Indian Institute of Technology, Bombay, Mumbai 400 076

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Selected X-ray data

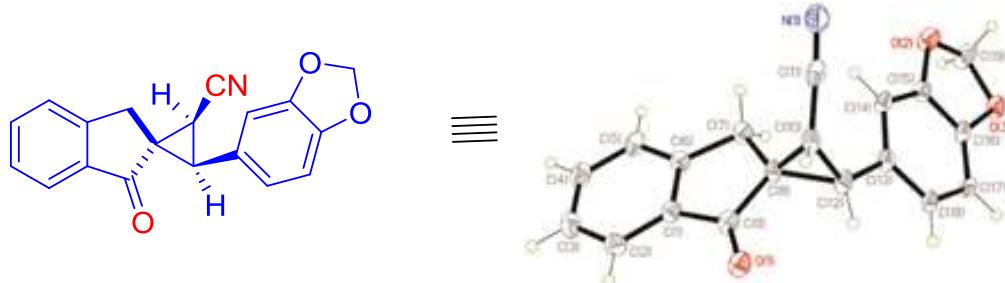
Table S1. Selected X-ray data for **3a**



Crystal data	
Chemical formula	C ₂₀ H ₁₆ N ₂ O
M _r	300.35
Crystal system, space group	Orthorhombic, P2 ₁ 2 ₁ 2 ₁
Temperature (K)	100
a, b, c (Å)	7.199 (3), 13.738 (5), 30.836 (12)
V(Å ³)	3050 (2)
Z	8
Radiation type	Mo Kα
μ (mm ⁻¹)	0.08
Crystal size (mm)	0.23 × 0.18 × 0.15
Data collection	
Diffractometer	CCD diffractometer
Absorption correction	Numerical

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T_{\min}, T_{\max}	0.981, 0.988
No. of measured, independent and observed [$I > 2\sigma(I)$] reflections	23094, 5564, 4866
R_{int}	0.062
$(\sin \theta / \lambda)_{\max} (\text{\AA}^{-1})$	0.602
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Refinement	
$R[F^2 > 2\sigma(F^2)], wR(F^2), S$	0.054, 0.134, 1.02
No. of reflections	5564
No. of parameters	415
H-atom treatment	H-atom parameters constrained
$\Delta\rho_{\max}, \Delta\rho_{\min} (\text{e \AA}^{-3})$	0.17, -0.21

Table S2. Selected X-ray data for 6c



Crystal data	
Chemical formula	C ₁₉ H ₁₃ NO ₃
M_r	303.30
Crystal system, space group	Monoclinic, P2 ₁ /n
Temperature (K)	100
a, b, c (Å)	9.350 (4), 8.514 (3), 18.643 (7)
β (°)	103.789 (18)
V (Å ³)	1441.3 (10)
Z	4
Radiation type	Mo K α
μ (mm ⁻¹)	0.10

Crystal size (mm)	$0.28 \times 0.20 \times 0.16$
Data collection	
Diffractometer	Rigaku Saturn724 diffractometer
Absorption correction	Numerical <i>NUMABS</i> (Rigaku, 1999)
T_{\min}, T_{\max}	0.974, 0.985
No. of measured, independent and observed [$I > 2\sigma(I)$] reflections	12578, 2506, 1318
R_{int}	0.098
$(\sin \theta / \lambda)_{\max}$ (\AA^{-1})	0.595
Refinement	
$R[F^2 > 2\sigma(F^2)], wR(F^2), S$	0.054, 0.150, 0.98
No. of reflections	2506
No. of parameters	208
H-atom treatment	H-atom parameters constrained
$\Delta\rho_{\max}, \Delta\rho_{\min}$ (e \AA^{-3})	0.30, -0.32