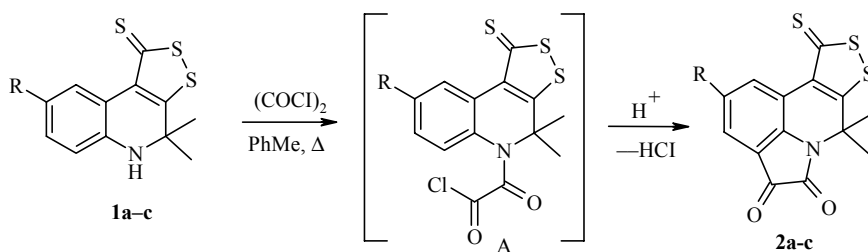


**4,5,7,10-TETRAHYDRO[1,2]DITHIOLO-
[3,4-*c*]PYRROLO[3,2,1-*ij*]QUINOLINE –
A NEW HETEROCYCLIC SYSTEM**

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Keywords: 2,3-dithiolo[3,4-*c*]quinoline-1-thione, 4,5,7,10-tetrahydro[1,2,]dithiolo[3,2,1-*ij*]quinoline, oxalyl chloride.

We have found that the acylation of 8-*R*-4,4-dimethyl-4,5-dihydro[1,2]dithiolo[3,4-*c*]quinoline-1-thiones (**1a-1c**) by oxalyl chloride, as in the case of acyl chlorides [1], proceeds exclusively at the nitrogen atom of the dihydroquinoline ring but is accompanied by spontaneous cyclization as in the Shtolle reaction. Derivatives of a new heterocyclic system, namely, 2-*R*-7,7-dimethyl-10-thioxo-4,5,7,10-tetrahydro[1,2]dithiolo[3,4-*c*]pyrrolo[3,2,1-*ij*]quinoline-4,5-diones (**2a-c**) are obtained in 60-80% yield.



1, 2 a R = H, b R = Me, c R = OMe

This reaction does not require use of a Lewis acid commonly employed as a catalyst in the Shtolle synthesis [2, 3], probably because of the catalytic effect of hydrogen chloride liberated in the first reaction step.

7,7-Dimethyl-10-thioxo-4,5,7,10-tetrahydro[3,4-*c*]pyrrolo[3,2,1-*ij*]quinoline-4,5-dione (2a); mp 300-301°C. ¹H NMR spectrum (DMSO-*d*₆), δ, ppm: 2.12 (6H, s, CMe₂); 6.91 (1H, m, 2-CH arom); 7.11 (1H, m, 1-CH arom); 9.48 (1H, m, 3-CH arom). Mass spectrum, *m/z*: 319 [M⁺]. Found, %: C 52.69; H 2.61; N 4.30; S 30.28. C₁₄H₉NO₂S₃. Calculated, %: C 52.60; H 2.82; N 4.38; S 30.05.

2,7,7-Trimethyl-10-thioxo-4,5,7,10-tetrahydro[3,4-*c*]pyrrolo[3,2,1-*ij*]quinoline-4,5-dione (2b); mp 254-256°C. ¹H NMR spectrum (DMSO-*d*₆), δ, ppm: 2.15 (6H, s, CMe₂); 2.38 (3H, s, Ar-Me); 6.97 (1H, s, 1-CH arom); 9.42 (1H, s, 3-CH arom). Mass spectrum, *m/z*: 333 [M⁺]. Found, %: C 53.69; H 2.65; N 4.37; S 28.98. C₁₅H₁₁NO₂S₃. Calculated, %: C 53.98; H 3.30; N 4.20; S 28.79.

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2-Methoxy-7,7-dimethyl-10-thioxo-4,5,7,10-tetrahydro[3,4-c]pyrrolo[3,2,1-ij]quinoline-4,5-dione (2c); mp 278-279°C. ¹H NMR spectrum (DMSO-d₆), δ, ppm: 2.12 (6H, s, CMe₂); 3.82 (3H, s, OMe); 7.04 (1H, s, 1-CH arom); 9.50 (1H, s, 3-CH arom). Mass spectrum, *m/z*: 349 [M⁺]. Found, %: C 51.67; H 3.41; N 4.23; S 27.78. C₁₅H₁₁NO₃S₃. Calculated, %: C 51.51; H 3.15; N 4.01; S 27.47.

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