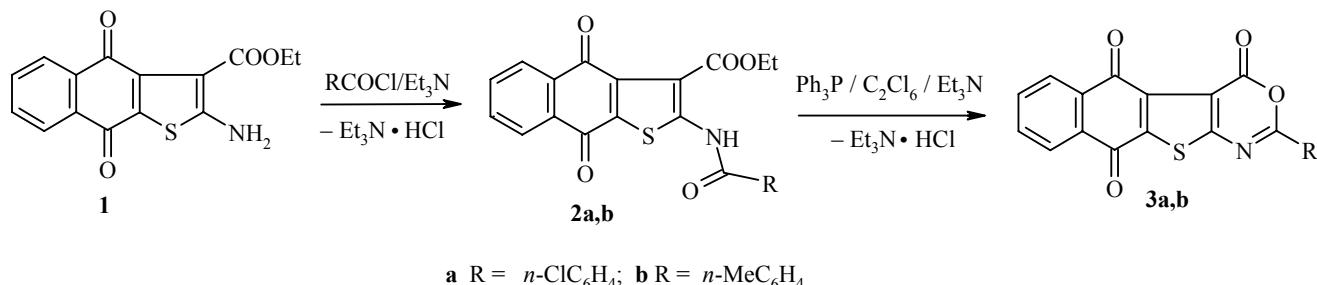


## 2-ARYL-4H-NAPHTHO[2',3':4,5]THIENO+- [2,3-d][1,3]OXAZINE-4,5,10-TRIONES

M. V. Stasevych, S. I. Sabat, M. I. Semenyuk,  
R. Ya. Musyanovych, and V. P. Novikov

**Keywords:** 2-amino-3-ethoxycarbonylnaphtho[2,3-*b*]thiophene-4,9-dione, 2-aryl-4H-naphtho[2',3':4,5]-thieno[2,3-*d*][1,3]oxazine-4,5,10-trione.

The synthetic potential of 2-amino-3-ethoxycarbonylnaphtho[2,3-*b*]thiophene-4,9-dione (**1**) has been little studied until now [1, 2]. We have carried out the reaction of dione **1** with equimolar quantities of aryl chlorides in dioxane in the presence of triethylamine (70–80°C, 5 h) and have obtained 2-arylamino-3-ethoxycarbonylnaphtho[2,3-*b*]thiophene-4,9-diones **2a,b**. Reaction of compounds **2a,b** with dichlorotriphenylphosphorane in the presence of triethylamine in toluene gave the previously undescribed 2-aryl-4H-naphtho[2',3':4,5]thieno[2,3-*d*][1,3]oxazine-4,5,10-triones **3a,b** in yields of 74–78%.



<sup>1</sup>H NMR spectra of DMSO-d<sub>6</sub> solutions with TMS as internal standard were recorded with a Bruker MSL-400 (400 MHz) spectrometer at 25°C.

### 2-(4-Chlorobenzoylamino)-3-ethoxycarbonylnaphtho[2,3-*b*]thiophene-4,9-dione (**2a**).

Yield 61%; mp 184–185°C. <sup>1</sup>H NMR spectrum, δ, ppm (J, Hz): 7.82–7.77 (2H, dd, J = 0.4, J = 7.3, CH<sub>arom</sub>); 8.23–8.09 (2H, m, CH<sub>arom</sub>); 3.30 (1H, s, NH); 7.37–7.82 (4H, dd, J = 2.3, J = 2.1, CH<sub>arom</sub>); 4.24–4.30 (2H, q, J = 7, CH<sub>2</sub>); 1.35 (3H, t, J = 2.9, CH<sub>3</sub>). Found, %: C 59.95; H 2.99; Cl 8.00; N 3.25; S 7.38. C<sub>22</sub>H<sub>14</sub>CINSO<sub>5</sub>. Calculated, %: C 60.07; H 3.21; Cl 8.06; N 3.18; S 7.29.

### 2-(4-Methylbenzoylamino)-3-ethoxycarbonylnaphtho[2,3-*b*]thiophene-4,9-dione (**2b**).

Yield 64%; mp 180–181°C. <sup>1</sup>H NMR spectrum, δ, ppm (J, Hz): 7.81–7.75 (2H, m, CH<sub>arom</sub>); 8.21–8.08 (2H, dd, J = 7.3, J = 0.4, CH<sub>arom</sub>); 3.30 (1H, s, NH); 7.30 (2H, d, J = 2.3, CH<sub>arom</sub>); 7.03 (2H, d, J = 2.2, CH<sub>arom</sub>); 4.24–4.30 (2H, q, J = 6.9, CH<sub>2</sub>); 2.35 (3H, s, CH<sub>3</sub>); 1.35 (3H, t, J = 2.8, CH<sub>3</sub>). Found, %: C 65.97; H 4.15; N 3.20; S 7.38. C<sub>23</sub>H<sub>17</sub>NSO<sub>5</sub>. Calculated, %: C 65.86; H 4.09; N 3.34; S 7.64.

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Lviv Polytechnic National University "Lviv Polytechnic", Lviv 79013, Ukraine; e-mail: vnovikov@polynet.lviv.ua. Translated from Khimiya Geterotsiklicheskikh Soedinenii, No. 7, 1112–1113, July, 2008. Original article submitted January 28, 2008.

**2-(4-Chlorophenyl)-4H-naphtho[2',3':4,5]thieno[2,3-d][1,3]oxazine-4,5,10-trione (3a).** Yield 74%; mp 292-293°C.  $^1\text{H}$  NMR spectrum,  $\delta$ , ppm ( $J$ , Hz): 8.14-7.75 (2H, m, CH<sub>arom</sub>); 8.47-8.41 (2H, t,  $J$  = 7.5, CH<sub>arom</sub>); 7.46 (2H, d,  $J$  = 2.1, CH<sub>arom</sub>); 8.03 (2H, d,  $J$  = 3.3, CH<sub>arom</sub>). Found, %: C 59.94; H 2.10; Cl 8.90; N 3.61; S 8.21. C<sub>21</sub>H<sub>11</sub>NO<sub>4</sub>S. Calculated, %: C 61.00; H 2.05; Cl 9.00; N 3.56; S 8.14.

**2-(4-Methylphenyl)-4H-naphtho[2',3':4,5]thieno[2,3-d][1,3]oxazine-4,5,10-trione (3b).** Yield 75%; mp 292-293°C.  $^1\text{H}$  NMR spectrum,  $\delta$ , ppm ( $J$ , Hz): 8.16-7.74 (2H, t,  $J$  = 7.5, CH<sub>arom</sub>); 8.47-8.41 (3H, m, CH<sub>arom</sub>); 7.21 (2H, d,  $J$  = 3.3, CH<sub>arom</sub>); 8.23 (2H, d,  $J$  = 2.2, CH<sub>arom</sub>); 2.39 (3H, s, CH<sub>3</sub>). Found, %: C 67.71; H 3.02; N 7.02; S 8.40. C<sub>21</sub>H<sub>11</sub>NSO<sub>4</sub>. Calculated, %: C 67.55; H 2.97; N 3.75; S 8.59.

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