

Fig. S1: (a) Packing of the molecules of **6b** in a projection along [010]; (b) arrangement of the molecules in the layer, parallel to (100) at $x = 0.250$, formed due to weak intermolecular interactions: C24-H24...Cg2 ($x, y+1, z$), marked by dashed lines, with the geometrical parameters: H...A $y+ = 2.92\text{Å}$, D...A = 3.797Å , and $\angle\text{DHA} = 157^\circ$.

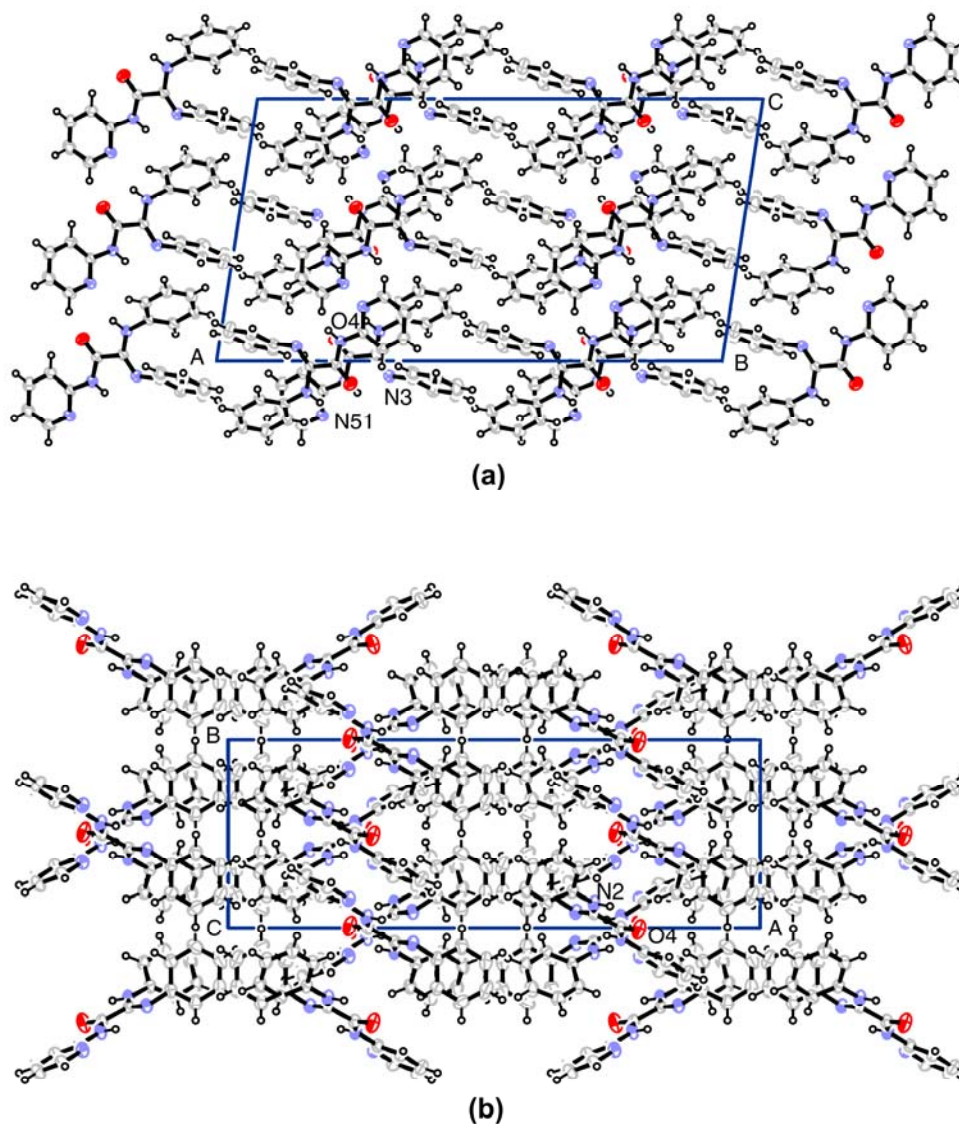


Fig. S2: The packing of the molecules **7a**: (a) in a projection along [010]; (b) in a projection along [001].

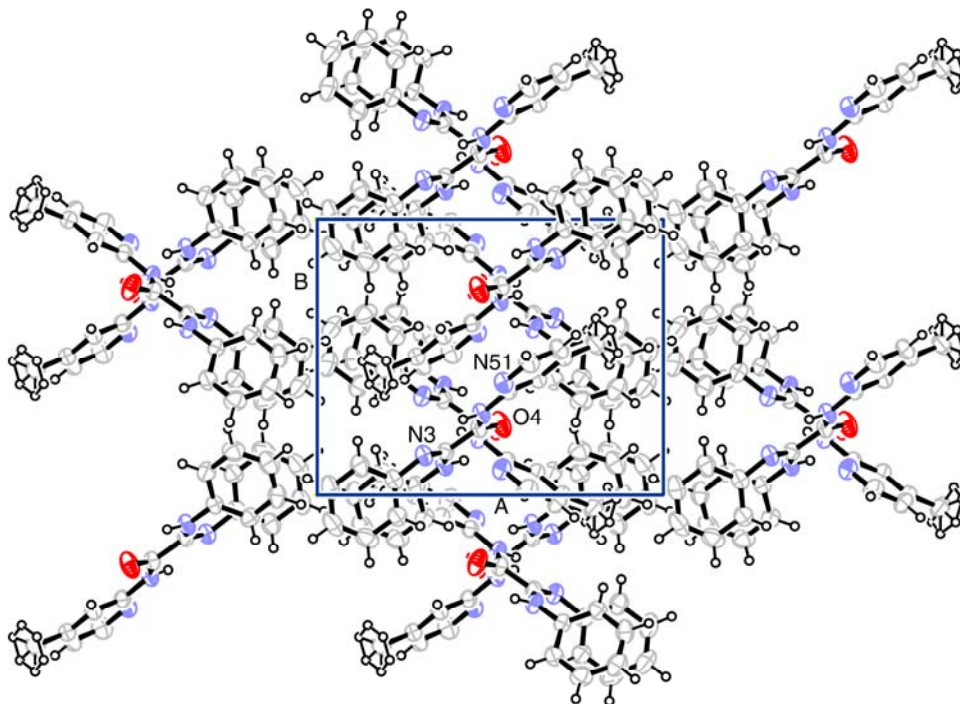


Fig. S3: The packing of the molecules of **7b** in a projection along [001].

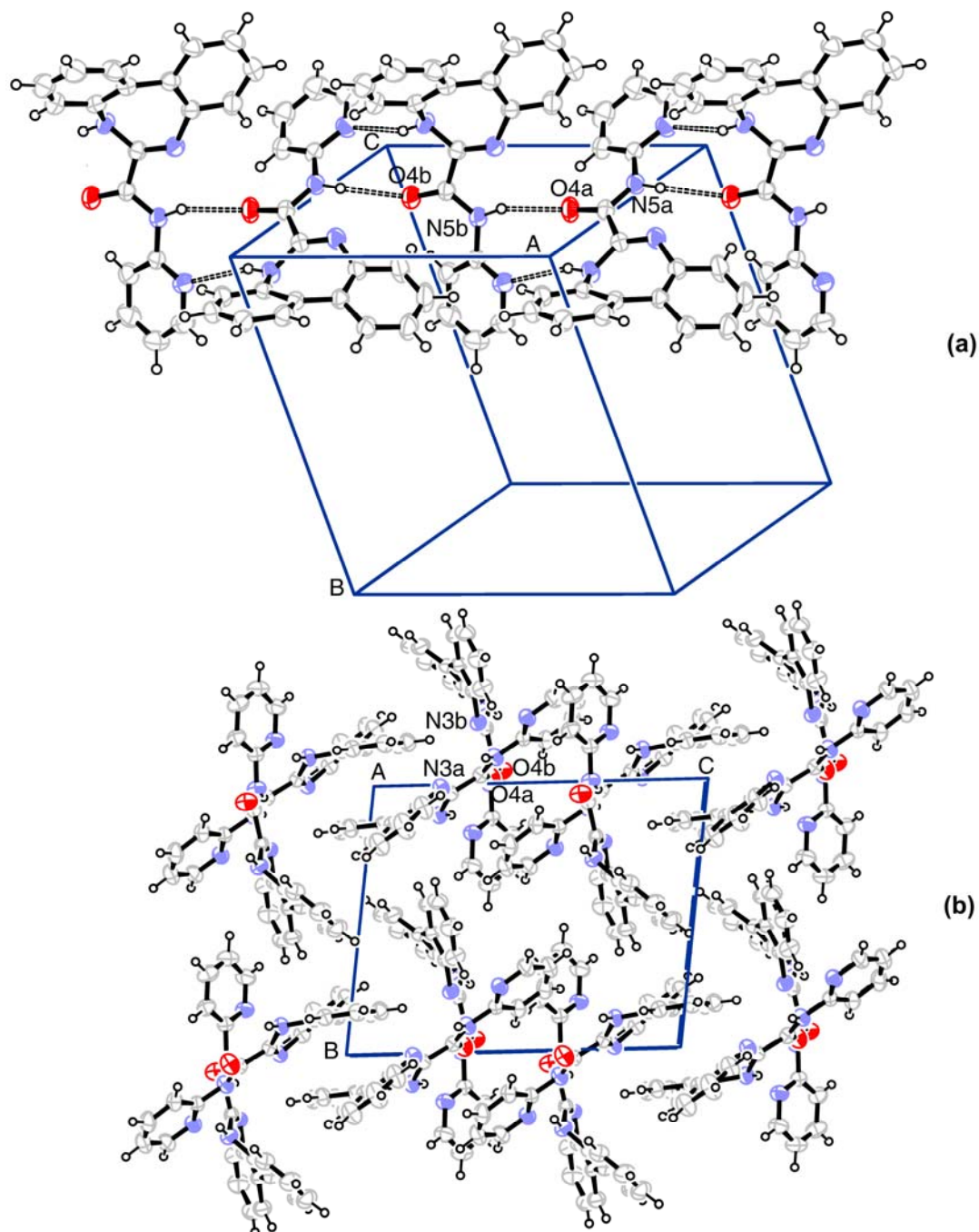


Fig. S4: The packing of the molecules of **11a**: (a) the view of a tape (consisting of atoms at $x, y-1, z$) running along [100] and built of the alternate diastereomeric molecules of **11a_a** and **11a_b**, joined by the H-bonds of N-H...O and N-H...N types, defined in Table 3; (b) in a projection along [100] i.e. along the axis of the tape.

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Table S1 Geometric parameters of weak interactions. Cg1 and Cg2 denote the gravity centres of the C21^C26 and C31^C36 rings, respectively

D-H...A	D-H/Å	H...A/Å	D...A/Å	∠DHA/°
6a				
C54-H54...O4 ⁱⁱ	0.93	2.58	3.420	151
O4...N5 ⁱ			3.261	
O4...Cg2			3.360	
6b				
C24-H24...O4 ⁱ	0.93	2.61	3.261	127
	0.93	2.92	3.797	157
			3.238	
7a				
O4...N5 ⁱ	C24-		3.233	
	H24...Cg2 ⁱ			
Cg1...Cg2	C3...N5 ⁱⁱ		3.955	
7b				
O4...N5 ⁱ			3.193	
Cg1...Cg2			3.658	
11a				
C34b-H34b...Cg2a ⁱⁱ	0.93	2.77	3.589	148
C55a-H55a...Cg2b ⁱⁱⁱ	0.93	3.07	3.913	152
C56b-H56b...Cg1b ^{iv}	0.93	3.08	3.858	143
O4a...N51b			3.221	
O4b...N51a ⁱ			3.155	

Symmetry codes: **6a**: (i) -x+3, -y, -z+1; (ii) -x+1, -y+1, -z+1; **6b**: (i) x, y+1, z; **(ii)** -x, -y-1, -z-1; **7a**: (i) x, -y, z+1/2; **7b**: (i) x, -y+1/2, z-1/2; **11a**: (i) x-1, y, z; (ii) x, y-1, z; (iii) -x+2, -y+1, -z+1; (iv) x, y+1, z.