

Table S1. Selected X-ray Data of **3e**

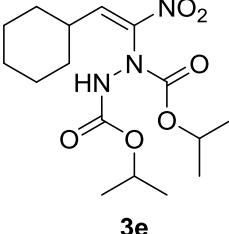
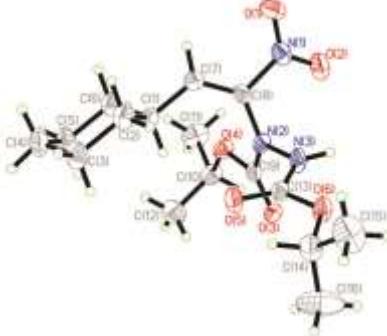
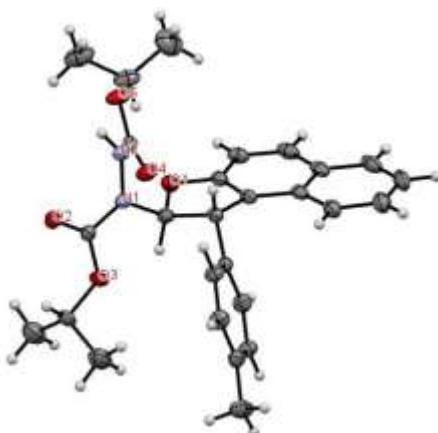
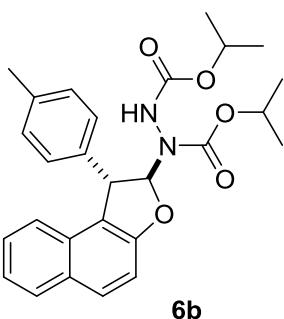
 3e	
Identification code	INN-MVD-CYCLOHEXYL
Empirical formula	C ₁₆ H ₂₇ N ₃ O ₆
Formula weight	357.41
Temperature	293(2) K
Wavelength	0.71070 Å
Crystal system	Triclinic
Space group	P -1
Unit cell dimensions	a = 9.777(5) Å α = 102.265(7) $^{\circ}$ b = 10.159(5) Å β = 100.480(6) $^{\circ}$ c = 10.306(5) Å γ = 90.508(8) $^{\circ}$
Volume	982.4(8) Å ³
Z	2
Density (calculated)	1.208 Mg/m ³
Absorption coefficient	0.093 mm ⁻¹
F(000)	384
Crystal size	0.230 x 0.190 x 0.130 mm ³
Theta range for data collection	3.03 to 29.30 $^{\circ}$
Index ranges	-13 ≤ h ≤ 13, -13 ≤ k ≤ 13, -14 ≤ l ≤ 14
Reflections collected	17915
Independent reflections	5352 (Rint = 0.0759)
Completeness to theta = 29.30 $^{\circ}$	99.4 %
Absorption correction	Numerical
Max. and min. transmission	0.988 and 0.980
Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	5352 / 0 / 226
Goodness-of-fit on F ²	0.872
Final R indices [I>2sigma(I)]	R1 = 0.0497, wR2 = 0.1054
R indices (all data)	R1 = 0.0876, wR2 = 0.1212
Largest diff. peak and hole	0.228 and -0.217 e.Å ⁻³

Table S2. Selected X-ray Data of **6b**

Identification code	INN-MVD-164		
Empirical formula	$C_{27}H_{30}N_2O_5$		
Formula weight	462.53		
Temperature	150 K		
Wavelength	0.71075 Å		
Crystal system	Triclinic		
Space group	P -1		
Unit cell dimensions	$a = 11.116(3)$ Å	$\alpha = 63.523(15)$ °	
	$b = 11.260(3)$ Å	$\beta = 71.584(17)$ °	
	$c = 11.711(3)$ Å	$\gamma = 69.128(17)$ °	
Volume	$1204.2(6)$ Å ³		
Z	2		
Density (calculated)	1.276 Mg/m ³		
Absorption coefficient	0.088 mm ⁻¹		
F(000)	492		
Crystal size	$0.350 \times 0.270 \times 0.130$ mm		
Theta range for data collection	3.064 to 25.000°		
Index ranges	$-13 \leq h \leq 13, -13 \leq k \leq 13, -13 \leq l \leq 13$		
Reflections collected	12275		
Independent reflections	4222 [$R_{(int)} = 0.0404$]		
Completeness to theta = 25.000°	99.6 %		
Absorption correction	Numerical		
Max. and min. transmission	0.972, 0.989		
Refinement method	Full-matrix least-squares on F^2		
Data / restraints / parameters	4222 / 0 / 319		
Goodness-of-fit on F^2	1.040		
Final R indices [I>2sigma(I)]	$R_1 = 0.0421, wR_2 = 0.0969$		
R indices (all data)	$R_1 = 0.0527, wR_2 = 0.1031$		
Largest diff. peak and hole	0.213 and -0.208 e.Å ⁻³		