

Supplementary information file S1.

The puckering parameters (Cremer & Pople, 1975) and the lowest displacement asymmetry parameters (Nardelli, 1983) for compounds (I) and (II) – molecules A and B, are as follows:

For the five membered pyrrolidine ring $q_2 = 0.404$ (2)Å, $\varphi = 16.0$ (1)°, $\Delta S(N1)$ is 13.4 (2)° and $\Delta_2(C21)$ is 4.4 (2)° in (I), and $q_2 = 0.400$ (2)Å, $\varphi = 199.7$ (2)°, $\Delta S(N1A)$ is 16.5 (2)° and $\Delta_2(C21A)$ is 1.7 (2)° for molecule A, $q_2 = 0.414$ (2)Å, $\varphi = 23.9$ (2)°, $\Delta S(C12B)$ is 14.0 (2)° and $\Delta_2(C21B)$ is 5.5 (2)° for molecule B in (II).

For the six membered pyran ring $q_2 = 0.448$ (2)Å, $\varphi = 123.4$ (3)°, $\Delta S(C14/C20)$ is 6.2 (2)° and $\Delta_2(C19/C21)$ is 55.0 (2)° in (I), and, $q_2 = 0.435$ (2)Å, $\varphi = 303.0$ (2)°, $\Delta S(C14A/C20A)$ is 6.2 (2)° and $\Delta_2(C19A/C21A)$ is 58.3 (2)° for molecule A, $q_2 = 0.377$ (2)Å, $\varphi = 115.5$ (2)°, $\Delta S(C14B/C20B)$ is 1.0 (2)° and $\Delta_2(O2B/C13B)$ is 60.51 (2)° for molecule B in (II).

For the six membered piperidine ring $q_2 = 0.021$ (2)Å, $\varphi = 55$ (2)°, $\Delta S(N1/C24)$ is 1.9 (2)° and Δ_2 is 115.5 (2)° in (I), and $q_2 = 0.036$ (2)Å, $\varphi = 218$ (2)°, $\Delta S(N1A/C24A)$ is 2.0 (2)° and Δ_2 is 115.3 (2)° for molecule A, $q_2 = 0.019$ (2)Å, $\varphi = 38$ (2)°, $\Delta S(N1B/C24B)$ is 0.7 (2)° and Δ_2 is 116.0 (2)° for molecule B in (II).

Cremer, D. & Pople, J. A. (1975). *J. Am. Chem. Soc.* **97**, 1354--1358.

Nardelli, M. (1983). *Acta Cryst.* **C39**, 1141--1142.