Electronic Supplementary Information (ESI)

Can multi-biaxial mesogenic mixtures favour biaxial nematics? A computer simulation study

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Figure S1: The homogeneous *side-by-side* and *end-to-end* pair potential energy profiles for some of the various GB ellipsoids (the cutoff radius is $r_c = 5 \sigma_0$).



Figure S2: Representative radial correlation functions for the s = 0.00 species in the isotropic I, nematic N, biaxial nematic N_b, and biaxial smectic S_b phases observed for the $s_{max} = 0.20$ sample.



Figure S3: Snapshots of the N_b phase at $T^* = 2.5$ for the $s_{\text{max}} = 0.20$ sample. The three views have been taken along the l (plate A), m (plate B), and n (plate C), directors. Particles have been colour coded according as s = -0.20 yellow, s = -0.15 cyan, s = 0.00 gray, s = 0.15 pink, s = 0.20 green.



Figure S4: The average potential energy per particle $\langle U_{\rm GB} \rangle$ at various dimensionless (plate A), and reduced (plate B) temperatures for the five non–uniform mixtures with $s_{\rm max} = 0.10$ (blue squares), 0.15 (red circles), 0.20 (green diamonds), 0.25 (orange triangles), and 0.30 (purple diamonds). The data for the reference system with $s_{\rm max} = 0.00$ (black pentagons) are also plotted.



Figure S5: The average number density $\langle \rho \rangle$ at various dimensionless (plate A), and reduced (plate B) temperatures for the five non–uniform mixtures. See caption of Figure S4 for further details.



Figure S6: The phase sequence observed with respect to the average number density values. See caption of Figure 4 for details.



Figure S7: The average number density $\langle \rho \rangle$ plotted with respect to the $\langle R_{00}^2 \rangle$ (plate A), and $\langle R_{22}^2 \rangle$ (plate B) order parameters measured for the reference ellipsoids with s = 0.00 in the five non–uniform mixtures. See caption of Figure S4 for further details.



Figure S8: Pictorial representation of the changes in temperature stability ranges of the N and N_b phases (plate A), and reduction of the T^*_{NbSb} transition temperature (plate B) for the five non–uniform mixtures with respect to the reference system with $s_{\text{max}} = 0.00$.



Figure S9: The average $\langle R_{00}^2 \rangle$ (plate A), and $\langle R_{22}^2 \rangle$ (plate B) order parameters for the reference ellipsoids with s = 0.00, at various reduced temperatures for the five non–uniform mixtures. See caption of Figure S4 for further details.



Figure S10: The average potential energy $\langle U_{\rm GB} \rangle$ plotted with respect to the $\langle R_{22}^2 \rangle$ order parameter measured for the reference ellipsoids with s = 0.00 in the five non–uniform mixtures. See caption of Figure S4 for further details.