0.0

-0.5

-1.0

AND. Molecular Meight A Molecular weight Avg. Dipole moment AND Nematic Order A Dipole moment Nematic order Mercept

Avg. Molecular weight

Δ Molecular weight 0.77

Interaction energy

Interaction energy LI

Δ Interaction energy LJ

Avg. Dipole moment 0.84 0.53

Δ Dipole moment **0.27 0.61 0.23**

COF -0.25 -0.039 -0.36 0.15

Δ Nematic order -0.083-0.0075-0.052 0.21 -0.063 0.058 0.14

Δ Interaction energy QQ 0.11 0.029 0.26 0.2 -0.42 0.57 0.07 -0.14 -0.047 0.29

-0.67 -0.29 -0.41 0.1 0.099 0.74 0.37

0.35 0.39 0.34 0.14 0.091 -0.12 -0.52 0.033

Δ Tilt angle 0.37 0.39 0.08 0.065 0.12 -0.76 -0.45 0.073 0.11 -0.69

Diff Tilt 0.65 0.68 0.54 0.3 -0.015 -0.22 -0.2 -0.072 0.53 -0.54 0.2

Avg Nematic order -0.42 -0.38 -0.23 -0.034 -0.03 0.39

Interaction energy A Interaction energy A Interdigitation A Tilk andle Interdigitation Oir Tile

1

0.34 0.19 -0.056 -0.33 -0.17 -0.91 -0.39 -0.042 0.029 -0.72 0.67 0.096 0.013 0.0016 1 -0.68 0.9 -0.47 -0.35 -0.69

0.1 -0.068 0.64 -0.59 -0.18 -0.12 -0.03 -0.69

ALO TILANOLO

Intercept -0.42 -0.21 -0.12 0.31 0.074 1 0.39 0.058 -0.12 0.78 -0.76 -0.22 -0.15 -0.03 -0.91 0.79 -0.92 0.57 0.52 0.74

Avg Tilt angle -0.66 -0.61 -0.34 -0.033 0.081 0.78 0.66 0.13 -0.39 1 -0.69 -0.54 -0.48 0.2 -0.72 0.51 -0.8 0.29 0.58 0.64

Interdigitation 0.43 0.55 0.34 0.37 0.038 -0.15 -0.51 -0.076 0.2 -0.48 0.42 0.38 1 -0.36 0.013 0.014 0.15 0.081 -0.31 -0.12

Δ Interdigitation -0.26 -0.3 -0.21 -0.18 <mark>0.27</mark> -0.03 <mark>0.17</mark> -0.11 <mark>0.087 0.2</mark> -0.076 -0.24 -0.36 **1** 0.0016 -0.16 -0.057 -0.18 <mark>0.13</mark> -0.03

-0.9 -0.51 -0.84 -0.088 0.24 0.52 0.34 0.1 -0.26 0.58 -0.42 -0.49 -0.31 0.13 -0.35 0.36

Δ Interaction energy -0.2 -0.1 0.03 0.2 -0.3 0.79 0.22 -0.067 -0.068 0.51 -0.7 -0.078 0.014 -0.16 -0.68 1 -0.67

Interaction energy QQ 0.66 0.38 0.35 -0.2 -0.24 -0.92 -0.45 -0.079 0.14 -0.8 0.69 0.29 0.15 -0.057 0.9 -0.67

0.53 0.61 -0.039 -0.21 -0.38 -0.0075 0.39 -0.61 0.39 0.68 0.55 -0.3 0.19 -0.1 0.38 0.029 -0.51 -0.29

0.23 -0.36 -0.12 -0.23 -0.052 0.34 -0.34 0.08 0.54 0.34 -0.21 -0.056 0.03 0.35 0.26 -0.84 -0.41

1 0.15 0.31 -0.034 0.21 0.14 -0.033 0.065 0.3 0.37 -0.18 -0.33 0.2 -0.2 0.2 -0.088 0.1

1 0.074 -0.03 -0.063 0.091 0.081 0.12 -0.015 0.038 0.27 -0.17 -0.3 -0.24 -0.42 0.24 0.099

1 0.14 -0.52 0.66 -0.45 -0.2 -0.51 0.17 -0.39 0.22 -0.45 0.07 0.34 0.37

-0.54 0.0011 0.081 -0.18 -0.47 0.9

1 0.033 0.13 0.073 -0.072 -0.076 -0.11 -0.042 -0.067 -0.079 -0.14 0.1 0.1

1 0.38 -0.24 0.096 -0.078 0.29 0.0011 -0.49 -0.18

-0.36

0.6 -0.87

-0.72 0.029

A Interaction energy op Interaction energy op Interaction energy Ly A Interaction energy Ly

0.9 0.36 0.6

-0.36 -0.72 -0.87

0.78

0.029 0.2

0.78

0.2