## Correlation between gate-dielectric morphology at the nanoscale and charge transport properties in organic field-effect transistors

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## SUPPORTING INFORMATION



**Fig. S1.** Computed mobility as a function of the applied electric field for the four samples of Fig. 2: a) PMMA1; b) PMMA2. Squares and circles correspond to mobilities computed according to the low-field and high-field formula, respectively (see Methods).



**Fig. S2.** Comparison between the electric field dependence of the mobilities of the annealed PMMA1 sample and that obtained by a GDM on a grid with the same (21x21) size. Squares and circles correspond to mobilities computed according to the low-field and high-field formula, respectively (see Methods).



**Fig. S3.** Simulated charge density maps (arbitrary units, logarithmic scale) for the four samples of Fig. 2. The applied electric field  $(1.0 \cdot 10^8 \text{ V m}^{-1})$  is from left to right.



**Fig. S4.** Convergence of mobility values with the number of KMC simulations. Calculations refer to the unannealed samples of **PMMA1** and **PMMA2** (black and red lines, respectively).



**Fig. S5.** Computed mobility as a function of the mean area of interconnected domains for PMMA and PVP samples (exponential fit, R=0.993)

Number of	Bulk density (g/cm <sup>3</sup> )	Rrms	Rp-v	Slab density
monomers		(nm)	(nm)	(g/cm <sup>3</sup> )
32	1.20	0.48	2.8	1.14

0.50

2.8

1.13

1.18

64

**Table S1.** Comparison between structural properties of PMMA as a function of the chain length of oligomers. Calculations refer to the annealed sample of **PMMA1**.