

## Supporting information

# Nanostructured Fluids from Degradable Nonionic Surfactants for the Cleaning of Works of Art from Polymer Contaminants

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# No kinship exists among these authors.

**Table S1.** Viscosity (Pa·s) of D<sub>2</sub>O/MEK was measured as a function of temperature using a Ubbelhode viscosimeter immersed in a thermostatic bath. D<sub>2</sub>O has a viscosity of 1.11 Pa·s, while MEK has a viscosity of 0.80 Pa·s at 20 °C. The table reports the unusual behavior of the mixture: while the viscosity for a given MEK concentration decreases when temperature is raised, increasing MEK concentration results in a higher viscosity, which is unexpected in view of the viscosities of the two solvents alone.

MEK concentration	25°C	35°C	45°C
4	0.9393	0.744	0.6049
8	1.0413	0.8006	0.6469
13.5	1.3737	1.0404	0.7659
20	1.3134	0.9941	0.7783

**Table S2.** Molecular properties of the chemicals used in SANS experiments. The alkyl portion of the two surfactants was considered as composed of 10 C atoms on average, while no distinction was taken into account between the 5.5 EO groups of NR and the 6 EO groups of BR.

Compound	Formula	Molecular mass (g/mol)	SLD (10 <sup>-6</sup> Å <sup>-2</sup> )
Heavy water	D <sub>2</sub> O	20.04	6.39
BR	C <sub>10</sub> H <sub>21</sub> (CH <sub>2</sub> CH <sub>2</sub> O) <sub>6</sub> O <sub>H</sub>	422.6	0.37
BR tail	C <sub>10</sub> H <sub>21</sub>	141.2	-0.41
BR head	(CH <sub>2</sub> CH <sub>2</sub> O) <sub>6</sub> OH	281.4	0.97
MEK	(C <sub>2</sub> H <sub>5</sub> )CO(CH <sub>3</sub> )	72.11	0.17

**Table S3.** SANS fitting parameters for the BR-based systems.

Fitting parameter	BR – MEK 0%	BR – MEK 4%	BR – MEK 8%	BR – MEK 13.5%	BR – MEK 20%
$\phi$	0.051	0.048	0.045	0.043	0.049
r (Å)	16.6	15.6	15.5	14.4	14.1
poly	0.39	0.38	0.36	0.38	0.40
t (Å)	7.7	8.0	6.1	4.7	4.5
SLD <sub>core</sub> (Å <sup>-2</sup> )	-4.1·10 <sup>-7</sup>	-4.1·10 <sup>-7</sup>	-4.1·10 <sup>-7</sup>	-4.1·10 <sup>-7</sup>	-2.8·10 <sup>-7</sup>
SLD <sub>shell</sub> (Å <sup>-2</sup> )	3.1·10 <sup>-6</sup>	3.5·10 <sup>-6</sup>	2.2·10 <sup>-6</sup>	1.0·10 <sup>-6</sup>	9.1·10 <sup>-7</sup>
SLD <sub>bulk</sub> (Å <sup>-2</sup> )	6.4·10 <sup>-6</sup>	6.0·10 <sup>-6</sup>	5.7·10 <sup>-6</sup>	5.3·10 <sup>-6</sup>	4.9·10 <sup>-6</sup>
K <sub>1</sub>	6.7	5.5	6.5	6.5	10.5
Z <sub>1</sub>	31.0	34.7	34.8	34.8	25.4
K <sub>2</sub>	-1.3	-0.9	-1.7	-1.7	-0.5
Z <sub>2</sub>	11.7	4.0	3.9	3.9	2.3
N <sub>w</sub>	8.5	12.1	3.9	0.1	0.3
N <sub>agg</sub>	62	51	50	40	30

**Table S4.** SANS fitting parameters for the NR-based systems.

Fitting parameter	NR – MEK 0%	NR – MEK 4%	NR – MEK 8%	NR – MEK 13.5%	NR – MEK 20%
$\phi$	0.051	0.048	0.045	0.042	0.038
a (Å)	489.9	-	-	-	317.4
b (Å)	14.4	-	-	-	17.3
r (Å)	-	15.9	16.8	16.2	-
poly	-	0.47	0.44	0.46	-
t (Å)	14.5	8.3	7.6	6.2	15.9
SLD <sub>core</sub> (Å <sup>-2</sup> )	-4.1·10 <sup>-7</sup>	-4.1·10 <sup>-7</sup>	-4.1·10 <sup>-7</sup>	-4.1·10 <sup>-7</sup>	-2.1·10 <sup>-7</sup>
SLD <sub>shell</sub> (Å <sup>-2</sup> )	4.1·10 <sup>-6</sup>	3.6·10 <sup>-6</sup>	2.9·10 <sup>-6</sup>	2.1·10 <sup>-6</sup>	4.0·10 <sup>-6</sup>
SLD <sub>bulk</sub> (Å <sup>-2</sup> )	6.4·10 <sup>-6</sup>	6.0·10 <sup>-6</sup>	5.7·10 <sup>-6</sup>	5.3·10 <sup>-6</sup>	4.9·10 <sup>-6</sup>
K <sub>1</sub>	4.2	21.1	20.2	20.2	10.0
Z <sub>1</sub>	36	25.1	25.1	25.1	35.6
K <sub>2</sub>	-1.3	-2.2	-4.1	-4.0	-3.3
Z <sub>2</sub>	9.0	20.0	20.0	20.0	9.0
N <sub>w</sub>	18.9	12.8	7.7	3.6	25.5
N <sub>agg</sub>	1367	54	64	57	857