

Supplementary Information

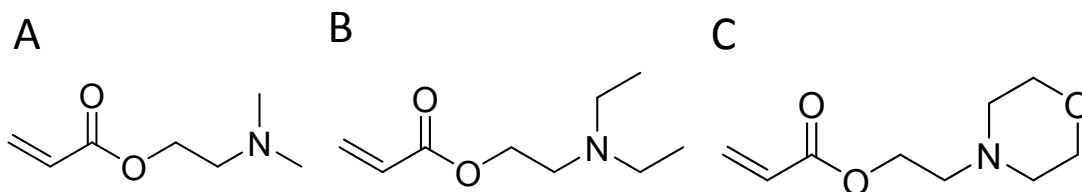


Figure SI 1 Structures of the monomers used: A) Dimethylamino ethyl acrylate; B) Diethylamino ethyl acrylate; C) 2-N-Morpholinoethyl acrylate

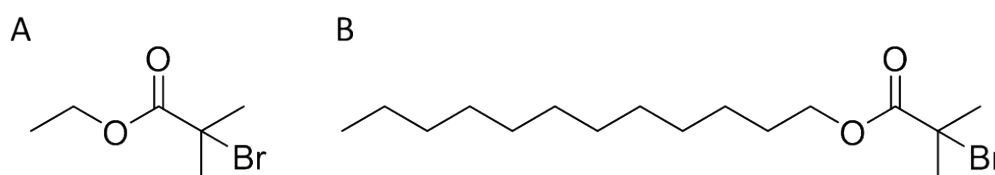


Figure SI 2 Structures of the initiators used: A) EBiB; B) DBiB

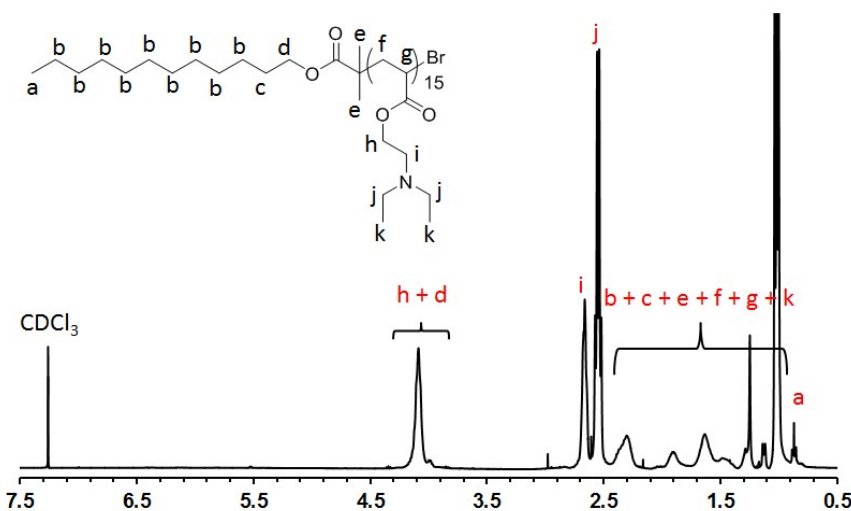


Figure SI 3 ^1H NMR spectra of DEAEA polymer

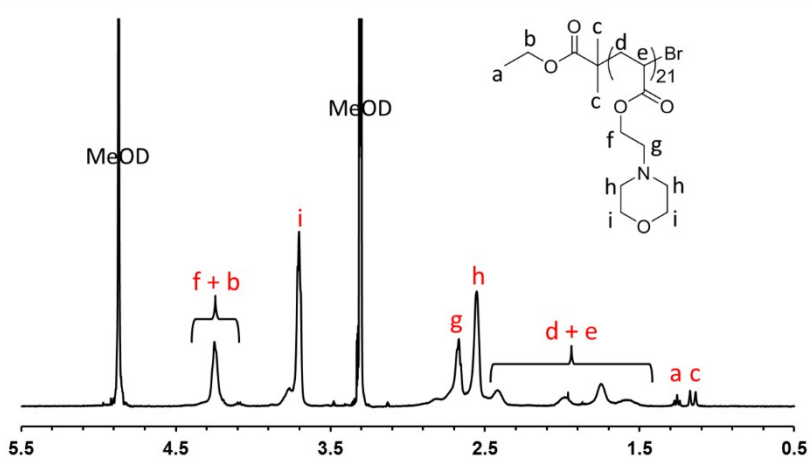


Figure SI 4 ^1H NMR spectra of MEA polymer

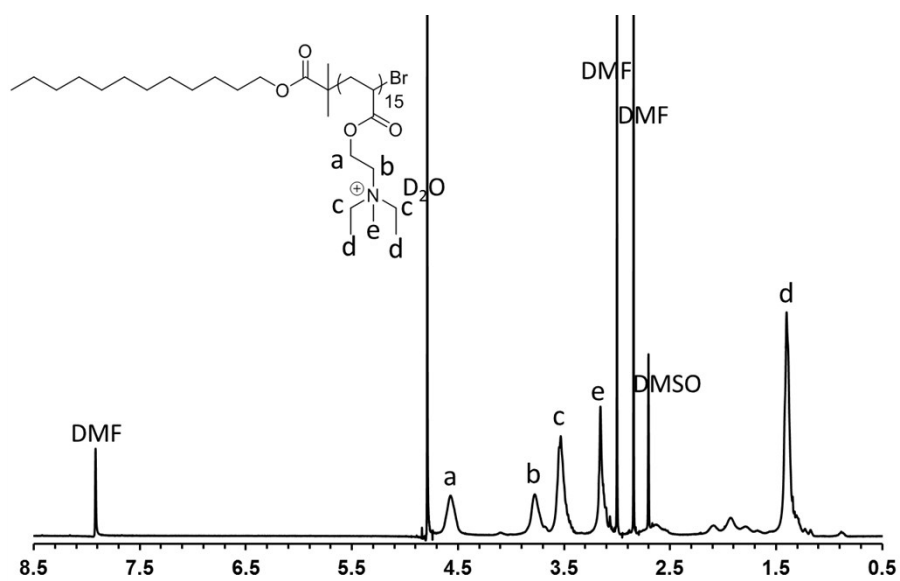


Figure SI 5 ^1H NMR of quaternized DEEA polymer

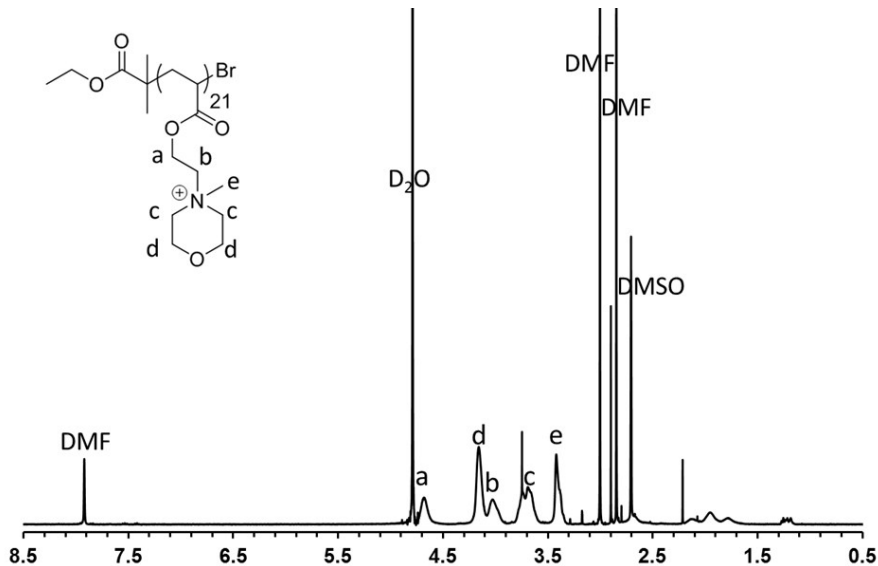


Figure S1 ^1H NMR of quaternized MEA polymer

Table S1 1 Characterization of the polymers using ^1H NMR and GPC

Entry N°	Monomer	Tail length	DP ^a	M _n ^a	M _n ^b	PDI ^b
1	DMAEA	Ethyl	27	4100	2686	1.20
2	DMAEA	Ethyl	5	900	937	1.05
3	DMAEA	Dodecyl	26	4100	2700	1.32
4	DMAEA	Dodecyl	5	1100	900	1.05
5	DEAEA	Ethyl	16	2900	2600	1.33
6	DEAEA	Ethyl	5	1100	1100	1.05
7	DEAEA	Dodecyl	15	2900	2500	1.38
8	DEAEA	Dodecyl	5	1200	1300	1.14
9	MEA	Ethyl	21	4100	4500	1.18
10	MEA	Ethyl	4	900	1300	1.12
11	MEA	Dodecyl	27	5300	4900	1.14
12	MEA	Dodecyl	4	1100	1300	1.13

^aDP (degree of polymerization) and M_n values were determined by ^1H NMR peak integration analysis.

^bM_n and PDI was determined by GPC analysis in DMAC against PSTY standards.

Table S1 2 Table of the compounds that were tested for cell viability in Tables S1 3 and S1 4

	1	2	3	4	5
A	Polymer 1	Polymer 2	Polymer 3	Polymer 4	Polymer 5
B	Polymer 6	Polymer 7	Polymer 8	Polymer 9	Polymer 10
C	Polymer 11	Polymer 12	Polymer 13	Polymer 14	Polymer 15
D	Polymer 16	Polymer 17	Polymer 18	Polymer 19	Polymer 20
E	Polymer 21	Polymer 22	Polymer 23	Polymer 24	

Table SI 3 Cell viability (% of cell survival) in 1% FBS (100 µg/mL of each polymer)

	1	2	3	4	5
A	96.5	97.4	105.9	4.0	82.8
B	95.4	23.8	85.5	82.7	34.1
C	-0.9	-0.1	88.9	91.4	44.9
D	26.3	87.2	105.1	111.6	4.7
E	104.9	97.0	101.3	29.6	

Table SI 4 Cell viability (% of cell survival) in 10% FBS (100 µg/mL of each polymer)

	1	2	3	4	5
A	112.2	127.0	139.3	31.2	84.0
B	90.4	79.1	106.0	98.4	74.0
C	1.7	-3.5	105.6	93.5	80.6
D	104.3	98.0	107.6	113.0	105.6
E	103.3	118.0	108.6	117.0	

Table SI 5 Selectivity index values^a for the lead compounds. MIC values are located in Table 1 and Table 2

Polymer N ^o	HC ₅₀ (µg/mL) ^b	<i>P.</i> <i>aeruginosa</i>	<i>A.</i> <i>baumannii</i>	<i>K.</i> <i>pneumoniae</i>	<i>S.</i> <i>aureus</i>	<i>A. baumannii</i> R
4	187	1	4	1	0	5
5	>1500	>3	>3	>3	>30	>4
7	>1500	>3	>6	>3	>60	>17
12	127	0	1	1	1	8
16	407	1	2	1	8	20

^aThe selectivity index is obtained by dividing the HC₅₀ by the MIC value for a given bacterium; and these values are shown in Table SI 5. The selectivity index is used routinely^{1,2} to describe the selectivity of the compound towards a particular bacterium. ^bHC₅₀ values were calculated by interpolation of results in Figure 4.

2. Y. M. Li, N. Bionda, A. Yongye, P. Geer, M. Stawikowski, P. Cudic, et al., *ChemMedChem.*, 2013, **8**, 1865-1872.