Investigation of structure-property-application

relationships of the hydrogel-based solar vapor generator

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Figure S1. The calculated equivalent water evaporation enthalpy in the hydrogel samples at various temperatures by DSC-TGA.



Figure S2. Electrostatic Potential Maps of the polymer repeating unit.



Figure S3. The molecule model and the automatic computed energy of (a) the $-NH_2$ group in PAM connected with two water molecules via hydrogen bonding and (b) the moment of dragging the second water molecule to break the hydrogen bonding between the water molecules.

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Sample name	E_0 (kJ mol ⁻¹)	E_1 (kJ mol ⁻¹)	$E_{\rm hyd}$ (kJ mol ⁻¹)		
PAMPSA-GO	-361.19	-337.25	23.94		
PAA-GO	-190.17	-173.00	17.17		
PHEA-GO	-24.23	-11.72	12.51		
PPEG-GO	150.36	159.99	9.64		
PDMA-GO	-124.32	-115.36	8.96		
PAM-GO	-183.01	-176.33	6.68		
PAETAC-GO	4.32	31.10	26.78		

Table S1. The energy E_0 and E_1 output by the *Avogadro*, and the simulated hydrogen bonding energy E_{hyd} of IW.

	0 min	15 min	30 min	45 min	60 min
PAMPSA-GO	25.60	33.99	36.34	36.66	36.44
	23.40	26.20	31.30	32.10	32.50
PAA-GO	21.78	35.30	37.88	36.94	36.77
	21.00	26.30		29.50	31.30
PHEA-GO	23.63	35.81	37.12	38.35	38.52
	-21.80	26.10	28.20	-30.00	31.00
PPEG-GO	21.77	34.25	æ 37.11	36.09	36.83
	21.10	26.00	29.20	.30.30	31.30
PDMA-GO	27.59	36.92	38.12	38.94	39.22
	*23.10	27.10	28.50	30.00	30.90
PAETAC-GO	26.54	33.93	35.81	36.53	37.65
	22.60	25.80	2 7.8 0	29.00	30.95

Figure S4. Thermal figures of PAMPSA-GO, PAA-GO, PHEA-GO, PPEG-GO, PDMA-GO and PAETAC-GO at the 0 min, 15 mins, 30 mins, 45 mins and 60 mins during the SVG testing.



Figure S5. Evaporation rates of the hydrogels in seawater.

Sample name	Mass of the monomer (mg)	Mass of the cross-linker (mg)
PAMPSA-GO hydrogel	277.11	22.89
PAA-GO hydrogel	242.43	57.57
PHEA-GO hydrogel	261.48	38.49
PPEG-GO hydrogel	300.00	0.00
PDMA-GO hydrogel	258.36	44.22
PAM-GO hydrogel	241.80	58.20
PAETAC-GO hydrogel	344.40	24.49

Table S2 The mass of the monomer and crosslinker used for the synthesis of the corresponding hydrogel.