

H NMR data

isobutyl POSS chloride [$(C_4H_9)_7Si_8O_{12}$]–Cl

1H NMR: 1.87 (m, 7H), 0.95 (m, 42H), 0.62 (m, 14H).
Analysis Calculated for $C_{28}H_{63}ClO_{12}Si_8$: C 39.47, H 7.45.
Found: C 39.02, H 7.58.

cyclopentyl POSS chloride [$(c-C_5H_9)_7Si_8O_{12}$]–Cl

1H NMR: 1.77 (m, 14H), 1.54 (m, 42H), 1.04 (m, 7H).
Analysis Calculated for $C_{35}H_{63}ClO_{12}Si_8$: C 44.91, H 6.79.
Found: C 44.37, H 6.88.

isobutyl POSS mono-ol [$(C_4H_9)_7Si_8O_{12}$]–OH

1H NMR: 1.85 (m, 7H), 0.96 (m, 42H), 0.56 (m, 14H).
Analysis Calculated for $C_{28}H_{64}O_{13}Si_8$: C 40.35, H 7.74.
Found: C 39.96, H 7.68.

Cyclopentyl POSS mono-ol [$(c-C_5H_9)_7Si_8O_{12}$]–OH

1H NMR: 1.74 (m, 14H), 1.56 (m, 42H), 0.98 (m, 7H).
Analysis Calculated for $C_{35}H_{64}O_{13}Si_8$: C 45.84, H 7.03.
Found: C 44.16, H 7.10.

4-methyl phenyl (trioxyisobutyl POSS) silane [$(C_4H_9)_7Si_8O_{12}-O]_3-Si-ArCH_3$

1H NMR: 7.60 (dd, 2H), 7.15 (dd, 2H), 2.38 (s, 3H),
1.81 (m, 21H), 0.94 (m, 126H), 0.59 (m, 42H). Analysis
Calculated for $C_{91}H_{196}O_{39}Si_{25}$: C 41.94, H 7.54. Found: C
41.17, H 7.68.

4-methyl phenyl (trioxycyclopentyl POSS) silane [$(C_5H_9)_7Si_8O_{12}-O]_3-Si-ArCH_3$

1H NMR: 7.63 (dd, 2H), 7.12 (dd, 2H), 2.35 (s, 3H),
1.73 (m, 42H), 1.54 (m, 126H), 0.99 (m, 21H). Analysis
Calculated for $C_{112}H_{196}O_{39}Si_{25}$: C 47.06, H 7.54. Found: C
41.17, H 6.87.

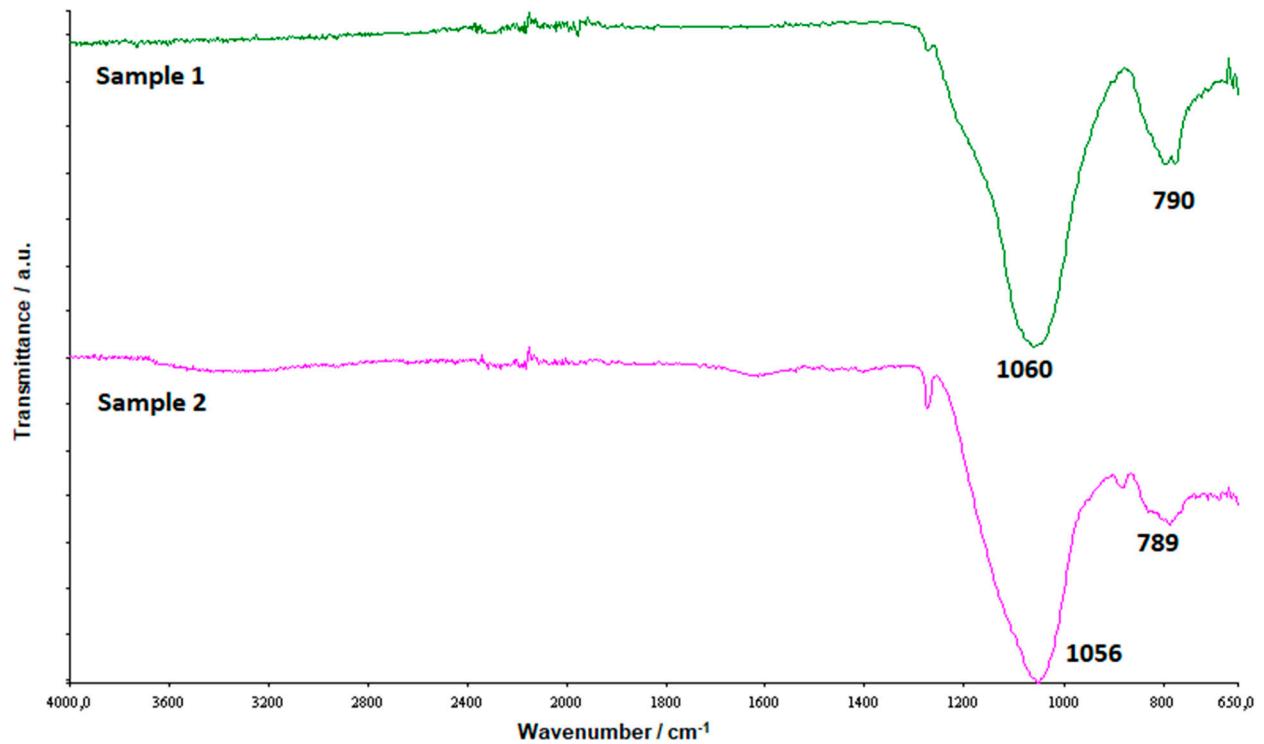


Figure SM1. FTIR spectra of the solid residues at 700 °C for the samples **1** and **2**.

Table SM1. Regression coefficients and apparent activation energies (E_a) of degradation by the FWO integral isoconversional method for **PS** in static air.

α	a	$b \cdot 10^{-3} /K$	r	$E_a /(\text{kJ}\cdot\text{mol}^{-1})$
0.1	24.8 (± 0.7)	13.9 (± 0.4)	0.9968	116 (± 3)
0.2	25.5 (± 0.6)	14.8 (± 0.4)	0.9979	123 (± 3)
0.3	25.4 (± 0.6)	15.1 (± 0.4)	0.9977	125 (± 3)
0.4	25.1 (± 0.6)	15.1 (± 0.7)	0.9931	125 (± 6)
0.5	24.5 (± 0.8)	14.8 (± 0.6)	0.9956	123 (± 5)
0.6	24.4 (± 0.8)	15.0 (± 0.5)	0.9964	125 (± 4)
0.7	24.2 (± 0.8)	14.9 (± 0.6)	0.9957	124 (± 5)
0.8	23.9 (± 0.7)	14.9 (± 0.5)	0.9967	124 (± 4)
0.9	22.8 (± 0.7)	14.4 (± 0.5)	0.9960	120 (± 4)

Table SM2. Regression coefficients and apparent activation energies (E_a) of degradation by the FWO integral isoconversional method for **PS** in flowing nitrogen.

α	a	$b \cdot 10^{-3} /K$	r	$E_a /(\text{kJ}\cdot\text{mol}^{-1})$
0.1	33.4 (± 0.7)	20.4 (± 0.4)	0.9986	170 (± 3)
0.2	33.3 (± 0.6)	20.8 (± 0.4)	0.9989	173 (± 3)
0.3	35.1 (± 0.6)	22.2 (± 0.4)	0.9990	185 (± 3)
0.4	34.8 (± 0.8)	22.3 (± 0.5)	0.9982	185 (± 4)
0.5	34.8 (± 0.7)	22.5 (± 0.5)	0.9986	187 (± 4)
0.6	34.6 (± 1.1)	22.5 (± 0.8)	0.9966	187 (± 7)
0.7	34.7 (± 0.8)	22.7 (± 0.6)	0.9981	189 (± 5)
0.8	34.2 (± 0.8)	22.6 (± 0.6)	0.9979	188 (± 5)
0.9	33.7 (± 0.7)	22.6 (± 0.5)	0.9986	188 (± 4)

Table SM3. Regression coefficients and apparent activation energies (E_a) of degradation by the FWO integral isoconversional method for sample **1** in static air.

α	a	$b \cdot 10^{-3} /K$	r	$E_a /(\text{kJ}\cdot\text{mol}^{-1})$
0.1	30.1 (± 1.4)	17.2 (± 0.9)	0.9923	143 (± 7)
0.2	29.9 (± 0.8)	17.5 (± 0.5)	0.9977	145 (± 4)
0.3	29.3 (± 1.1)	17.6 (± 0.7)	0.9950	146 (± 6)
0.4	29.1 (± 1.6)	17.7 (± 1.0)	0.9898	147 (± 8)
0.5	28.4 (± 1.4)	17.4 (± 0.9)	0.9915	145 (± 7)
0.6	28.4 (± 1.2)	17.5 (± 0.8)	0.9939	145 (± 7)
0.7	28.6 (± 1.5)	17.8 (± 1.0)	0.9907	148 (± 8)
0.8	28.3 (± 1.2)	17.7 (± 0.8)	0.9932	147 (± 7)
0.9	27.9 (± 1.4)	17.6 (± 1.0)	0.9911	146 (± 8)

Table SM4. Regression coefficients and apparent activation energies (E_a) of degradation by the FWO integral isoconversional method for sample **1** in flowing nitrogen.

α	a	$b \cdot 10^{-3} /K$	r	$E_a /(\text{kJ}\cdot\text{mol}^{-1})$
0.1	37.2 (± 1.3)	23.0 (± 0.9)	0.9958	191 (± 7)
0.2	36.8 (± 0.5)	23.2 (± 0.3)	0.9994	193 (± 2)
0.3	37.1 (± 0.5)	23.6 (± 0.4)	0.9993	196 (± 3)
0.4	37.1 (± 0.5)	23.8 (± 0.4)	0.9993	198 (± 3)
0.5	37.1 (± 0.6)	24.0 (± 0.4)	0.9990	199 (± 3)
0.6	37.1 (± 0.5)	24.2 (± 0.3)	0.9994	201 (± 2)
0.7	36.9 (± 0.7)	24.2 (± 0.5)	0.9989	201 (± 4)
0.8	36.1 (± 0.9)	24.0 (± 0.6)	0.9980	199 (± 5)
0.9	35.5 (± 1.6)	23.9 (± 1.2)	0.9930	199 (± 10)

Table SM5. Regression coefficients and apparent activation energies (E_a) of degradation by the FWO integral isoconversional method for sample **2** in static air.

α	a	$b \cdot 10^{-3} /K$	r	$E_a /(\text{kJ}\cdot\text{mol}^{-1})$
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0.1	31.1 (\pm 0.6)	18.2 (\pm 0.3)	0.9989	151 (\pm 2)
0.2	30.7 (\pm 1.5)	18.4 (\pm 0.9)	0.9918	153 (\pm 7)
0.3	30.7 (\pm 0.6)	18.4 (\pm 0.4)	0.9988	153 (\pm 3)
0.4	30.4 (\pm 1.5)	18.7 (\pm 1.0)	0.9915	155 (\pm 8)
0.5	30.1 (\pm 0.9)	18.7 (\pm 0.6)	0.9970	155 (\pm 5)
0.6	30.2 (\pm 1.2)	18.9 (\pm 0.8)	0.9948	157 (\pm 7)
0.7	29.4 (\pm 0.9)	18.7 (\pm 0.6)	0.9965	155 (\pm 5)
0.8	29.2 (\pm 0.9)	18.6 (\pm 0.6)	0.9965	155 (\pm 5)
0.9	28.8 (\pm 0.7)	18.5 (\pm 0.5)	0.9981	154 (\pm 4)

Table SM6. Regression coefficients and apparent activation energies (E_a) of degradation by the FWO integral isoconversional method for sample **2** in flowing nitrogen.

α	a	$b \cdot 10^{-3} / \text{K}$	r	$E_a / (\text{kJ} \cdot \text{mol}^{-1})$
0.1	37.5 (\pm 1.7)	23.3 (\pm 1.1)	0.9932	194 (\pm 9)
0.2	37.1 (\pm 1.1)	23.4 (\pm 0.7)	0.9970	194 (\pm 6)
0.3	36.0 (\pm 1.0)	23.6 (\pm 0.7)	0.9974	196 (\pm 6)
0.4	36.2 (\pm 0.9)	23.3 (\pm 0.6)	0.9977	194 (\pm 5)
0.5	36.2 (\pm 1.6)	23.5 (\pm 1.1)	0.9933	195 (\pm 9)
0.6	36.1 (\pm 1.9)	23.6 (\pm 1.3)	0.9909	196 (\pm 11)
0.7	36.2 (\pm 1.6)	23.9 (\pm 1.1)	0.9937	199 (\pm 9)
0.8	35.7 (\pm 0.9)	23.8 (\pm 0.6)	0.9980	198 (\pm 5)
0.9	34.9 (\pm 0.7)	23.6 (\pm 0.5)	0.9985	196 (\pm 4)