

Supporting information:

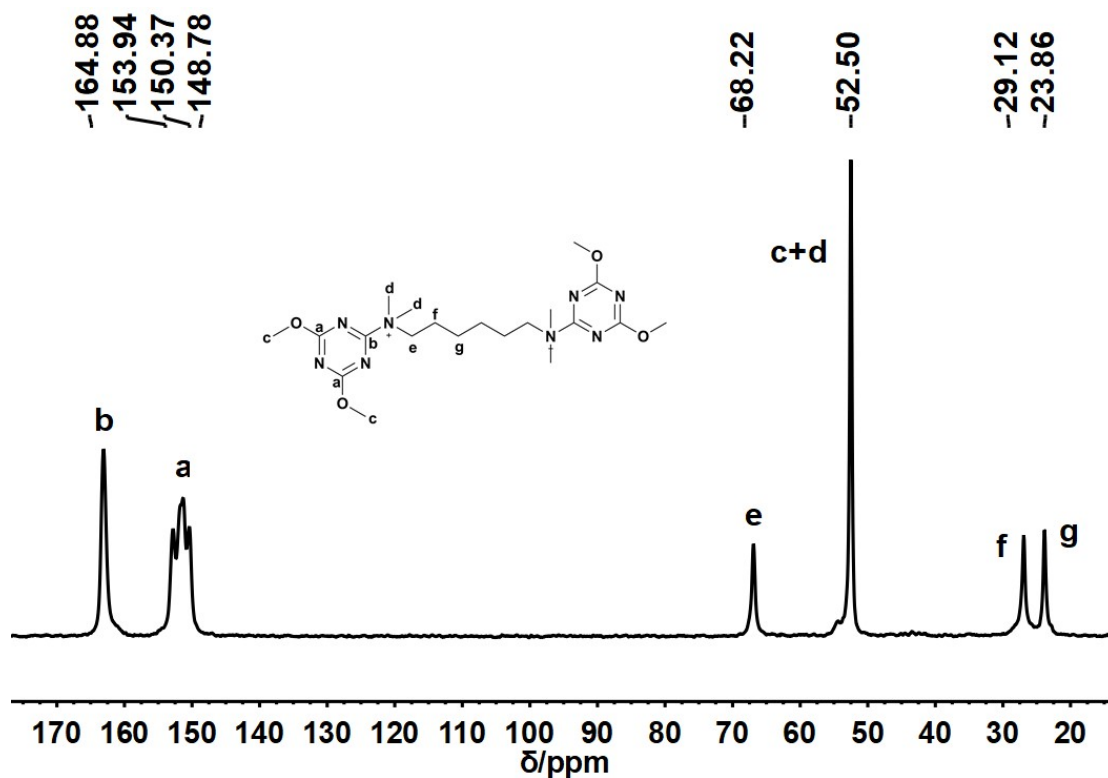


Fig. S1. The solid-state ^{13}C CP/MAS NMR spectrum of the TBQA-1.

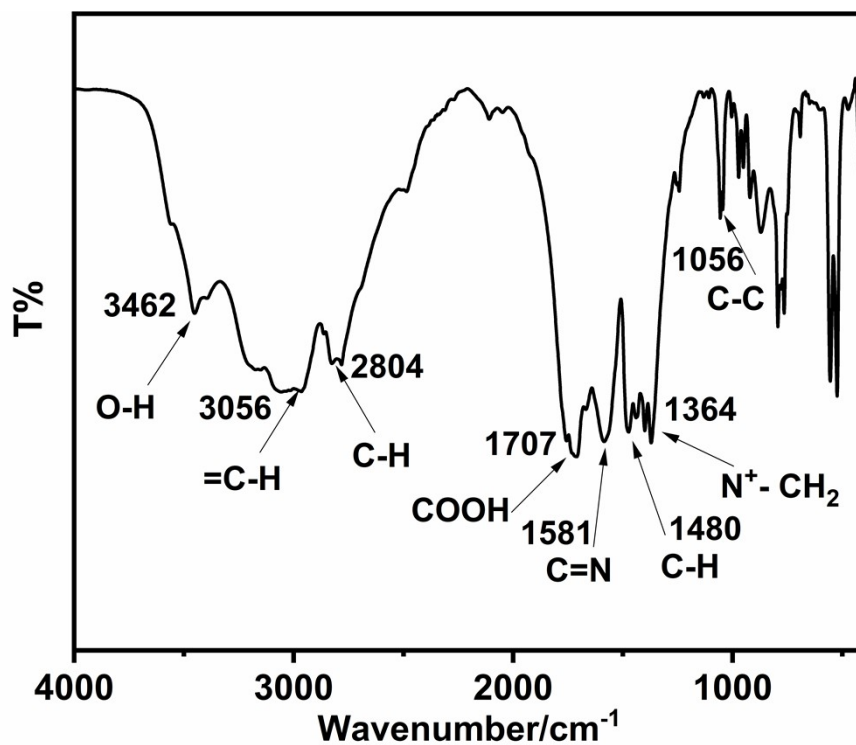


Fig. S2. The FT-IR spectrum of the TBQA-1. The absorption peak of $-\text{COOH}$ in FT-IR spectrum was probably attributed to the hydrolysis of residual acetonitrile, which was serving as a solvent.

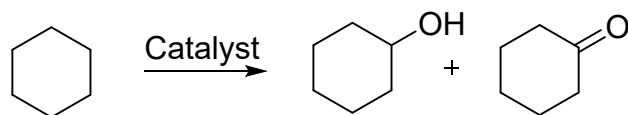


Fig. S3. Selective oxidation of cyclohexane. Similar reaction was performed with Pd loaded NSH-ZSM-48-300, SBA-15 and conventional Silicalite-1.

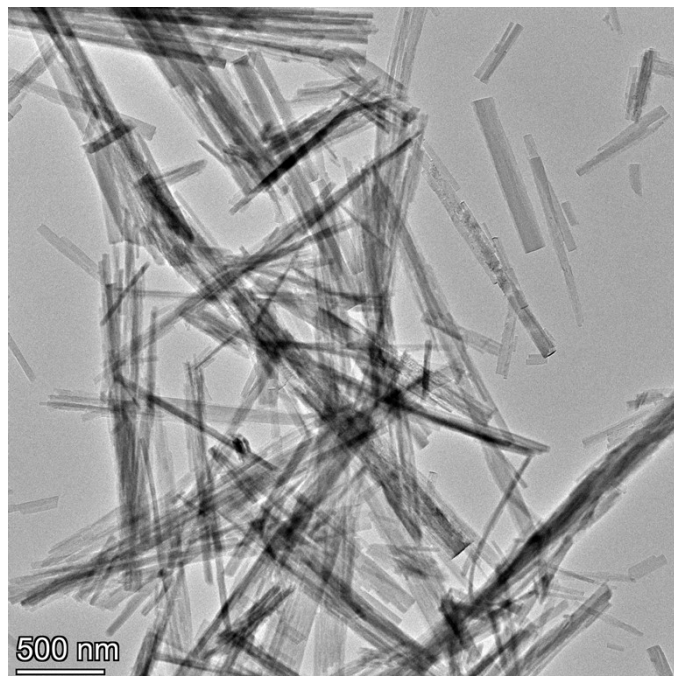


Fig. S4. The Low-magnification TEM images of NSH-ZSM-48-300.

	$S_{\text{BET}}^{\text{a/}}$	$S_{\text{micro}}^{\text{b/}}$	$S_{\text{meso}}^{\text{c/}}$	$V_{\text{total}}^{\text{d/}}$	$V_{\text{micro}}^{\text{b/}}$	$V_{\text{meso}}^{\text{c/}}$
	m^2/g	m^2/g	m^2/g	cm^3/g	cm^3/g	cm^3/g
Conventional SBA-15	858	117	714	1.292	0.065	1.227
Conventional Silicalite-1	275	220	57	0.185	0.116	0.069

Table S1. Porosity of the calcined samples.

^a calculated by applying the BET equation using the linear part ($0.05 < P/P_0 < 0.30$) of the adsorption isotherm. ^b calculated by t-plot method. ^c calculated by BJH adsorption model. ^d $V_{\text{total}} = V_{\text{micro}} + V_{\text{meso}}$.