

Supplementary information

Structural investigations and acidic properties of high surface area pyrochlore aluminium hydroxyfluoride

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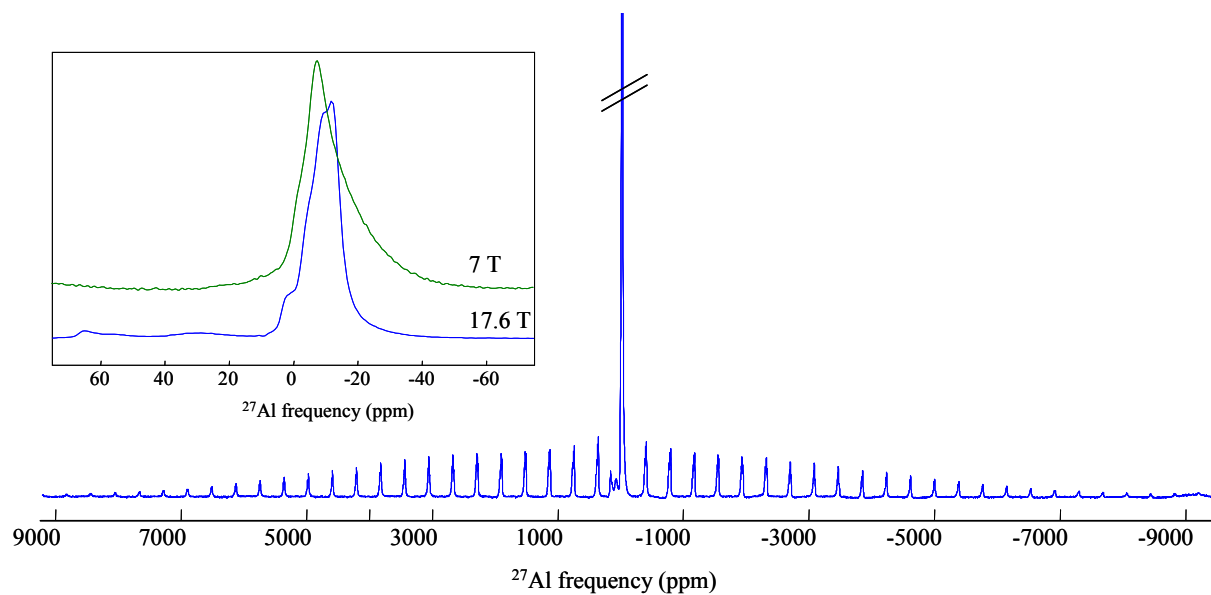


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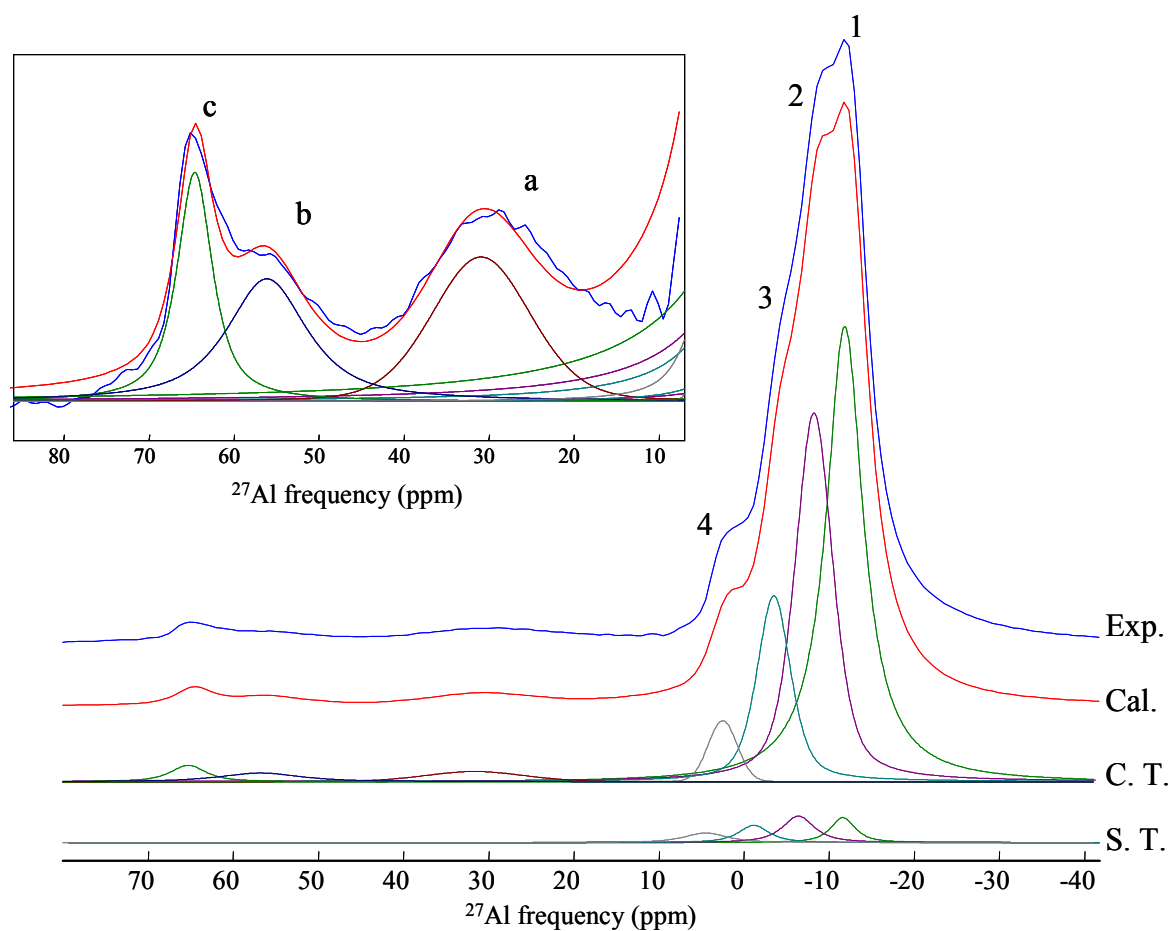


Figure S2. Experimental (blue) and calculated (red) central line of the ^{27}Al MAS NMR spectrum of the pyrochlore $\text{AlF}_{1.8}(\text{OH})_{1.2} \cdot 0.3\text{H}_2\text{O}$. The deconvolution, achieved using four contributions, takes into account the $N = 0$ band of both the satellite transitions $\langle 3/2 \rangle$ (S. T.) and the central transition $\langle 1/2 \rangle$ (C. T.). Resonances a, b and c were reconstructed using Gaussian lines to extract isotropic chemical shifts.

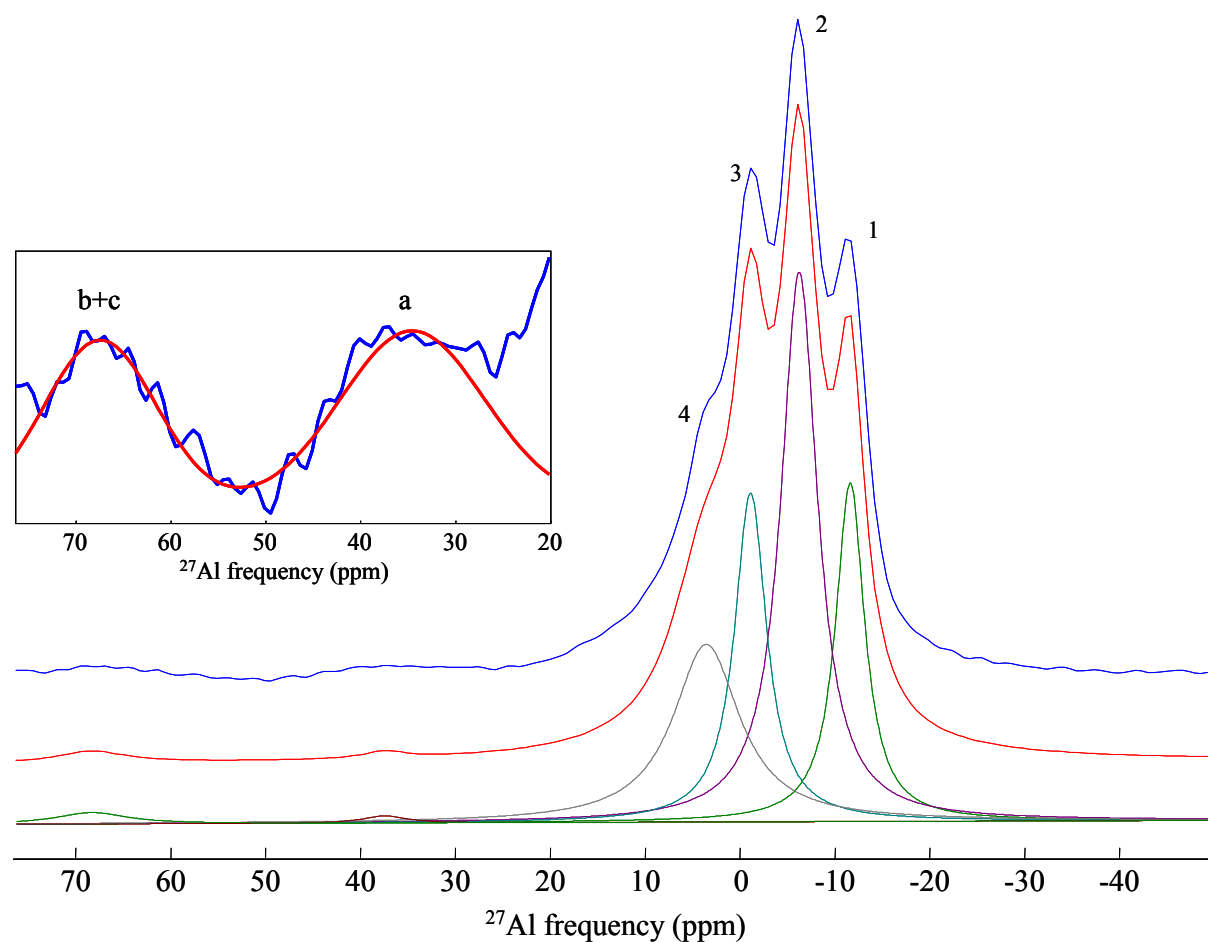


Figure S3. Experimental (blue) and calculated (red) full projection of the TOP reconstruction of the ^{27}Al MAS NMR spectrum of the pyrochlore $\text{AlF}_{1.8}(\text{OH})_{1.2} \cdot 0.3\text{H}_2\text{O}$. The deconvolution is achieved using Gaussian contributions.