

Electronic Supporting Information

Nanometer-Sized Nickel and Cobalt Doped Forsterite Synthesis for Investigating Critical Element Recovery from Mafic and Ultramafic Rocks

Kelly A. Peterson^{a,*}, Mark E. Bowden^a, Bavan P. Rajan^a, Tenley E. Webb^a, Bridgette N. Carven^a, Libor Kovarik^a, Zsombor Molnár^a, Mark H. Engelhard^b, Sandra D. Taylor^a, Elsa A. Cordova^b, Thomas W. Wietsma^c, Sebastian T. Mergelsberg^a, Christopher J. Thompson^b, Sebastien N. Kerisit^a, and John S. Loring^{a,*}

^a *Physical and Computational Sciences Directorate, Pacific Northwest National Laboratory, Richland, WA 99352, United States*

^b *Energy and Environment Directorate, Pacific Northwest National Laboratory, Richland, WA 99352, United States*

^c *Earth and Biological Sciences Directorate, Pacific Northwest National Laboratory, Richland, WA 99352, United States*

*Corresponding Authors: kelly.peterson@pnnl.gov and john.loring@pnnl.gov

This file contains 7 Pages, 5 Figures, and 1 Table.

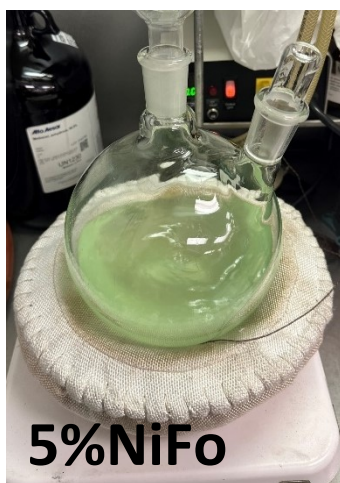
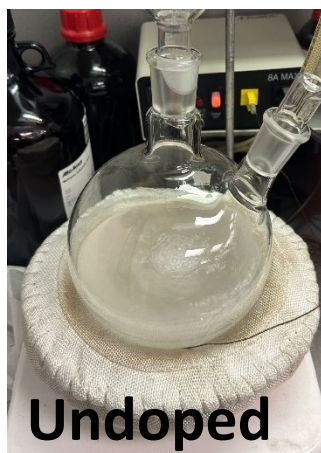


Figure S1. Pictures showing the reflux step for each of the undoped and metal doped forsterites after water additions.

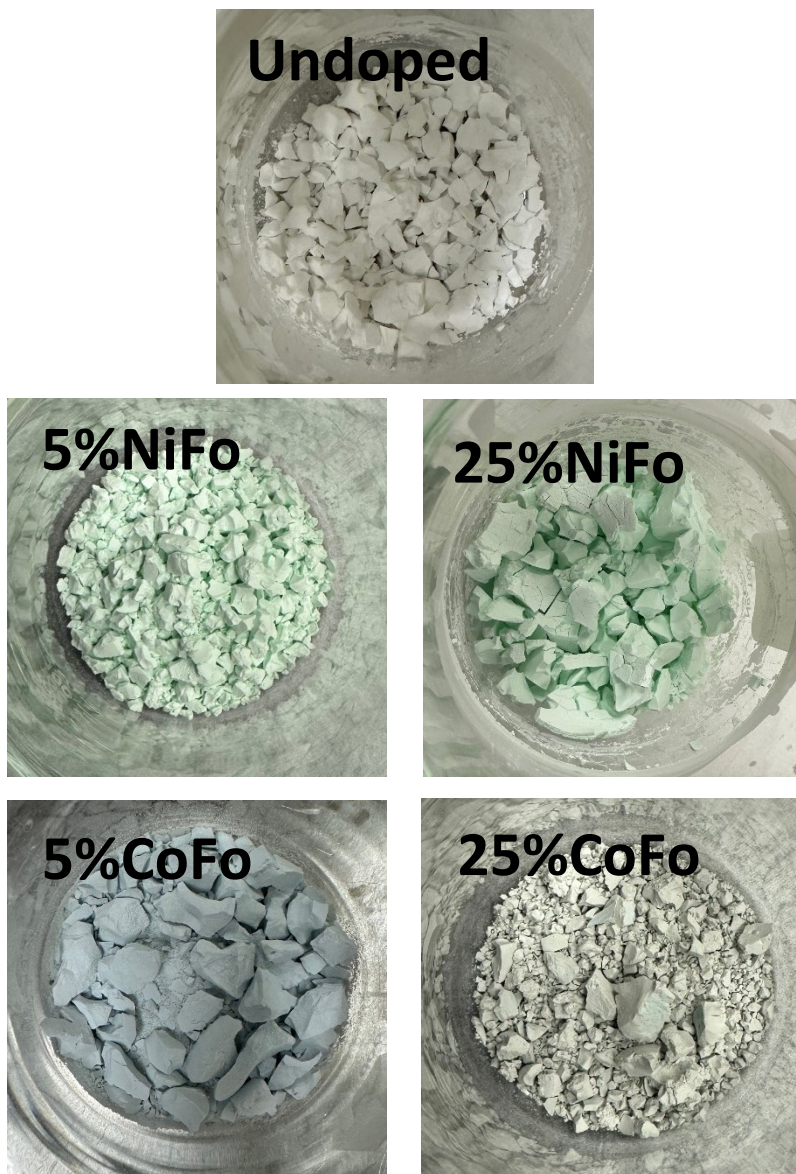


Figure S2. Pictures showing the dried gel after the washing step of the synthesis of undoped and metal doped forsterites.

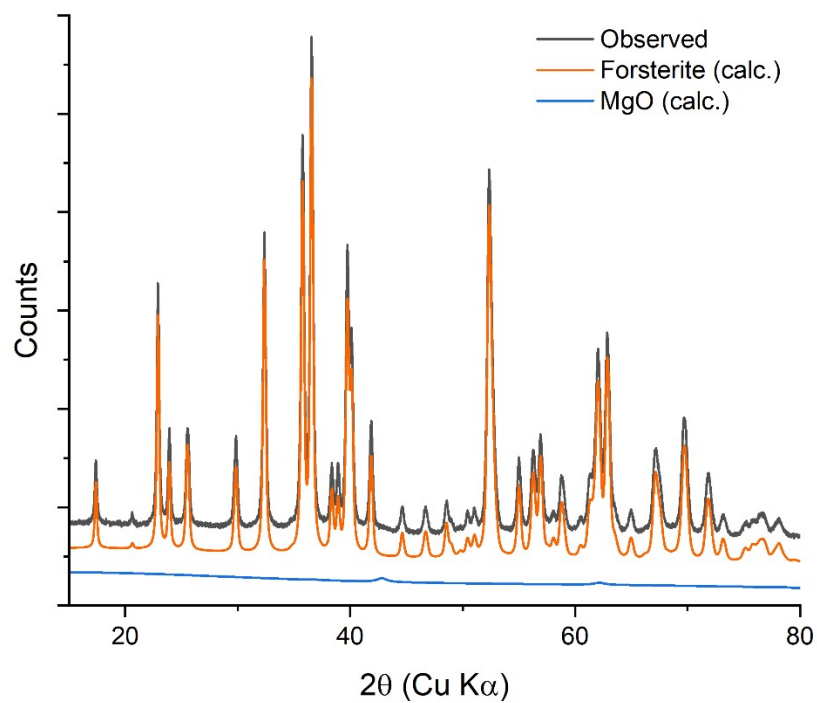


Figure S3. Experimental X-Ray diffraction pattern of undoped forsterite (black trace) shown with Rietveld refinement calculations of forsterite (orange trace) and MgO (blue trace) contributions.

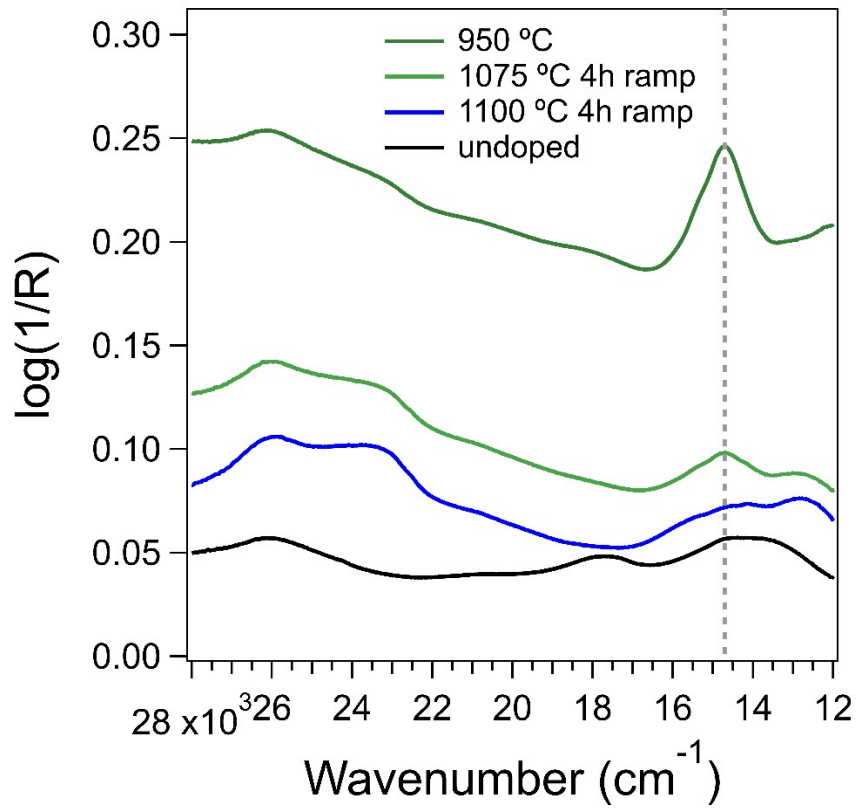


Figure S4. UV-Vis absorbance spectra of 5% Ni-doped forsterites synthesized under different calcination conditions. Absorbance band at $14,700 \text{ cm}^{-1}$ (marked with gray dashed line) was assigned to (Mg,Ni)O secondary phase.

Table S1. Mid-IR and far IR peak locations in wavenumber (cm^{-1}) of undoped and doped forsterite FTIR-ATR spectra. Peaks assigned based on assignments by Hofmeister and Pitman (*Phys. Chem. Miner.*, 2007).

Assignment	undoped	5%CoFo	5%NiFo	25%CoFo	25%NiFo
	1103.364	1102.239	1106.351		
	1005.193	1001.749	1002.509	1003.453	1004.626
$\nu_3(3)$	987.072	983.32	986.867	973.281	979.803
$\nu_3(2)$		956.132	956.433		
$\nu_3(1)$	871.848	870.062	874.602	870.668	873.465
ν_1	840.278	838.567	842.274	838.195	840.448
$\nu_4(3)$	613.672	611.732	613.855	604.505	603.128
		547.11	547.691		
$\nu_4(2)$	524.564	524.155	524.45	520.22	521.491
$\nu_4(3)$	504.343	501.582	504.756	492.806	496.032
$\nu_2(3)$	471.011	468.073	471.418	457.024	462.014
LO or ν_2	415.511	413.61	414.638		
R	396.887	394.042	397.326	400.094	405.104
M(7)	381.463	379.651	381.711		381.332
M(6)	353.526	347.219	354.124	339.634	345.091
M(4)	319.645				
M(5)				295.779	304.181
M(3)+LO	287.41	281.844	286.816		
M(1)	271.624	272.555	272.196	263.69	267.526
T(2)	200.996	197.094	200.469		
T(1)	142.125	137.87	139.228	128.935	131.301

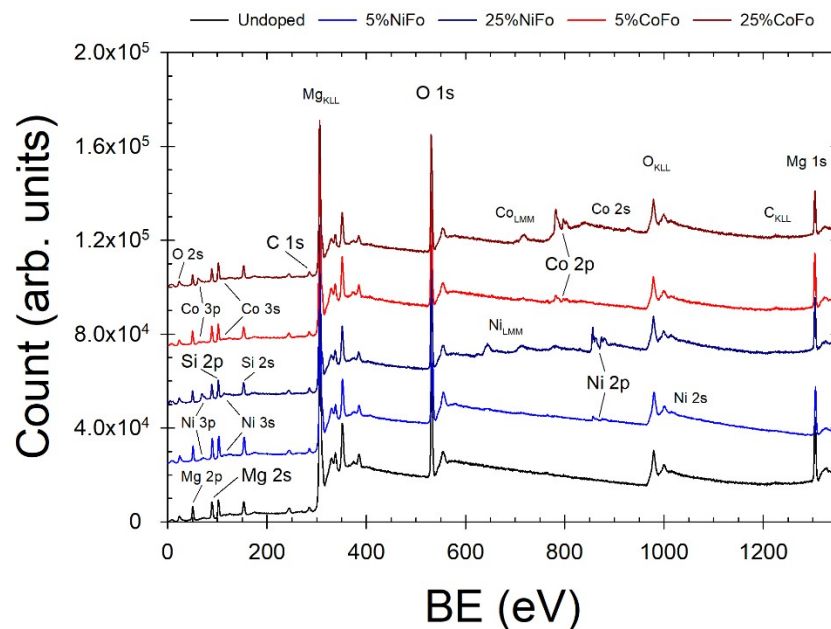


Figure S5. XPS spectra (survey scans) of undoped, 5% Ni-doped, 25% Ni-doped, 5% Co-doped, and 25% Co-doped forsterites. C, O, Mg, Si, Ni, and Co photoelectron lines are labeled. Larger font size is used to indicate the photoelectron lines used in the regional scans.