

## Supplementary Materials

**Table S1.** The Raman shifts and OH bands by FTIR ( $\nu_i$ ), temperature dependence of  $\nu_i$  ( $\partial\nu_i/\partial T$ )<sub>P</sub>, isobaric mode Grüneisen parameters, as well as intrinsic anharmonic parameters ( $a_i$ ) for hydrous forsterite.

| $\nu_i$ (cm <sup>-1</sup> ) | $(\partial\nu_i/\partial T)_P$ (cm <sup>-1</sup> K <sup>-1</sup> ) | $i_P$  | $i_T^b$ | $a_i$ (10 <sup>-5</sup> K <sup>-1</sup> ) |
|-----------------------------|--------------------------------------------------------------------|--------|---------|-------------------------------------------|
| 3612                        | -0.009 <sup>a</sup>                                                | 0.065  | -0.014  | -0.304                                    |
| 3578                        | -0.016 <sup>a</sup>                                                | 0.114  | -0.047  | -0.616                                    |
| 3566                        | -0.017 <sup>a</sup>                                                | 0.124  | -0.022  | -0.554                                    |
| 3551                        | -0.056 <sup>a</sup>                                                | 0.417  | -0.012  | -1.635                                    |
| 3477                        | 0.017 <sup>a</sup>                                                 | -0.129 |         |                                           |
| 968                         | -0.033 <sup>c</sup>                                                | 0.895  | 0.554   | -1.297                                    |
| 919                         | -0.034 <sup>c</sup>                                                | 0.971  |         |                                           |
| 880                         |                                                                    |        | 0.439   |                                           |
| 858                         | -0.023 <sup>c</sup>                                                | 0.704  | 0.403   | -1.146                                    |
| 825                         | -0.022 <sup>c</sup>                                                | 0.700  | 0.550   | -0.570                                    |
| 610                         | -0.012 <sup>c</sup>                                                | 0.516  | 0.692   | 0.671                                     |
| 590                         | -0.016 <sup>c</sup>                                                | 0.712  | 0.454   | -0.981                                    |
| 545                         |                                                                    |        | 0.442   |                                           |
| 437                         | -0.026 <sup>c</sup>                                                | 1.562  | 1.427   | -0.511                                    |
| 418                         | -0.027 <sup>c</sup>                                                | 1.695  |         |                                           |
| 336                         |                                                                    |        | 1.663   |                                           |
| 320                         |                                                                    |        | 1.416   |                                           |
| 305                         | -0.031 <sup>c</sup>                                                | 2.668  | 1.310   | -5.173                                    |
| 227                         | -0.016 <sup>c</sup>                                                | 1.850  | 0.652   | -4.566                                    |

*a*: mean temperature pressure dependences of OH band in this study; *b*: measured from Hushur et al. [24]; *c*: measured from Yang et al. [57]