

# Supplementary Materials: Integrated Utilization of Sewage Sludge and Coal Gangue for Cement Clinker Products: Promoting Tricalcium Silicate Formation and Trace Elements Immobilization

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## 1. The Equations of Compositional Parameters in Cement Chemistry:

Lime saturation coefficient:

$$KH = (\text{CaO} - 1.65 \text{Al}_2\text{O}_3 - 0.34 \text{Fe}_2\text{O}_3) / 2.8 \text{SiO}_2 \quad (\text{S1})$$

$$0.902 < KH < 0.915.$$

Silica Modulus:

$$SM = \text{SiO}_2 / (\text{Al}_2\text{O}_3 + \text{Fe}_2\text{O}_3) \quad (\text{S2})$$

$$2.25 < SM < 2.35.$$

Iron modulus:

$$IM = \text{Al}_2\text{O}_3 / \text{Fe}_2\text{O}_3 \quad (\text{S3})$$

$$1.3 < IM < 1.4.$$

Lime saturation:

$$\text{L.S.F} = \text{CaO} / (2.8 \text{SiO}_2 + 1.18 \text{Al}_2\text{O}_3 + 0.65 \text{Fe}_2\text{O}_3) \quad (\text{S4})$$

$$0.94 < \text{L.S.F} < 0.96.$$

## 2. Prepared Method

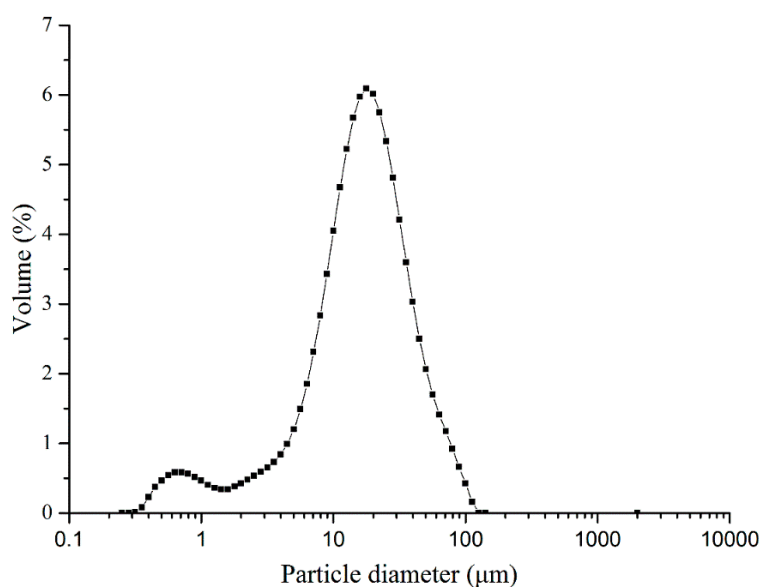
All mixtures were pressed to  $\phi 20 \text{ mm} \times 5 \text{ mm}$  slices by applying a pressure of 10 Mpa and calcined in the programmable electrically heated tube furnace. The furnace temperature was raised at the rate of  $10 \text{ }^\circ\text{C}/\text{min}$  from room temperature to  $1450 \text{ }^\circ\text{C}$ . The temperature was maintained at  $950 \text{ }^\circ\text{C}$  for 30 min to ensure the complete decomposition of  $\text{CaCO}_3$  and held at  $1450 \text{ }^\circ\text{C}$  for 2 h.

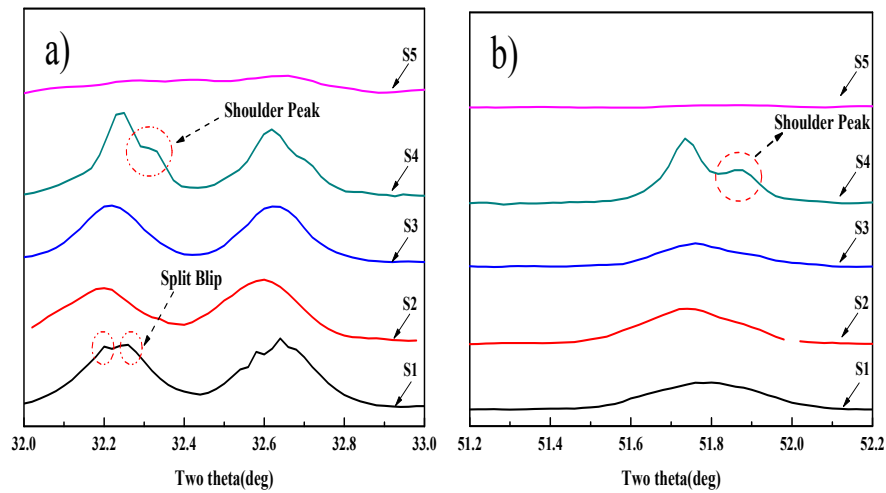
**Table S1.** Characteristic bands of FTIR spectra.

| Band Center (cm <sup>-1</sup> ) | Spectral Feature | Probable Assignment                                         |
|---------------------------------|------------------|-------------------------------------------------------------|
| 3694, 3669, 3653                | Band             | Outer OH stretching                                         |
| 3404                            | Band             | O–H stretching                                              |
| 2923, 2854                      | Shoulder         | Symmetric C–H stretching                                    |
| 2353                            | Band             | O–C=O stretching                                            |
| 1652                            | Band             | C=O stretching                                              |
| 1439                            | Peak             | C=O symmetrical stretching                                  |
| 1165                            | Band             | Trace of quartz                                             |
| 1115, 1110                      | Shoulder         | Apical Si–O stretching                                      |
| 1062                            | Shoulder         | (PO <sub>4</sub> ) groups                                   |
| 1034, 1009                      | Band             | Si–O stretching of Si–O–Si and Si–O–Al                      |
| 988                             | Shoulder         | $\nu_3(\text{SiO}_4)$ in C <sub>2</sub> S                   |
| 937                             | Band             | $\nu_3(\text{SiO}_4)$ in C <sub>3</sub> S                   |
| 934                             | Shoulder         | Al–OH libration (inner, outer)                              |
| 918                             | Shoulder         | $\nu_3(\text{SiO}_4)$ in C <sub>2</sub> S                   |
| 891                             | Band             | $\nu_3(\text{SiO}_4)$ in C <sub>3</sub> S                   |
| 840                             | Band             | $\nu_1(\text{SiO}_4)$ in C <sub>2</sub> S                   |
| 816                             | Shoulder         | $\nu_1(\text{SiO}_4)$ in C <sub>3</sub> S                   |
| 797, 780                        | Shoulder         | Trace of quartz                                             |
| 792, 754, 695                   | Shoulder         | OH translation                                              |
| 741                             | Shoulder         | Al–O vibrations in ferrite and tricalcium aluminate         |
| 522                             | Band             | SiO <sub>4</sub> tetrahedra out-of-plane bending vibrations |
| 538                             | Band             | Si–O–Al <sup>VI</sup> deformation                           |
| 470, 431                        | Band             | Si–O bending                                                |
| 454                             | Band             | Al–O vibration in C <sub>3</sub> A                          |

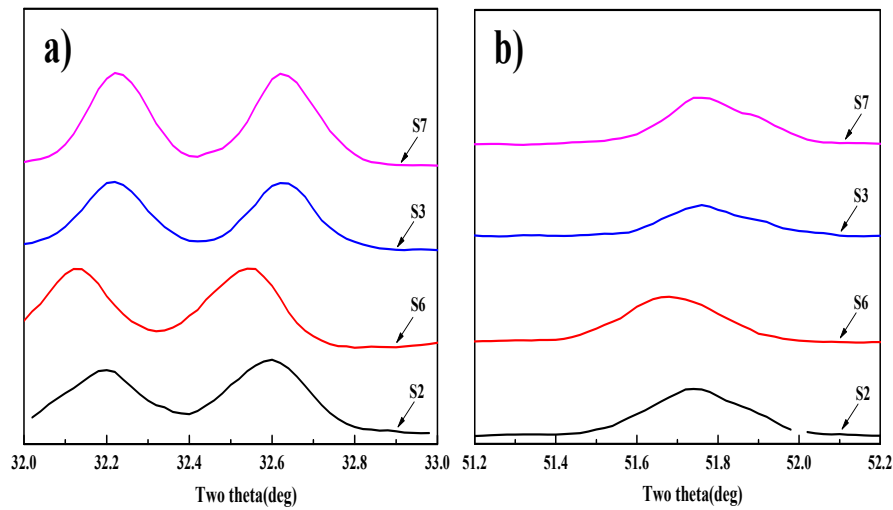
**Table S2.** The chemical analysis of the raw materials in the reference.

| Raw Meals                      | SiO <sub>2</sub> | Al <sub>2</sub> O <sub>3</sub> | Fe <sub>2</sub> O <sub>3</sub> | CaO  | LOI  |
|--------------------------------|------------------|--------------------------------|--------------------------------|------|------|
| Limestone                      | 5                | 2.4                            | 0.74                           | 50   | 40.7 |
| Sand                           | 88.6             | 2.9                            | 2.67                           | 2.2. | 2.61 |
| Clay                           | 37.1             | 10.6                           | 4.23                           | 18.7 | 20.6 |
| Fe <sub>2</sub> O <sub>3</sub> | 0                | 0                              | 96                             | 0    | 0.5  |

**Figure S1.** Particle size distribution of raw meals.



**Figure S2.** Detail of XRD peaks appearing between 32° and 33°, and between 51° and 52° with different amount of SS addition: (a) XRD peak between 32° and 33°; and (b) XRD peak between 51° and 52°.



**Figure S3.** The comparison detail of XRD peaks appearing between 32° and 33°, and between 51° and 52° with different amount of SS and CG addition: (a) XRD peak between 32° and 33°; and (b) XRD peak between 51° and 52°.