

Fig. S1 Structures of MEQ, OLA and QCT.

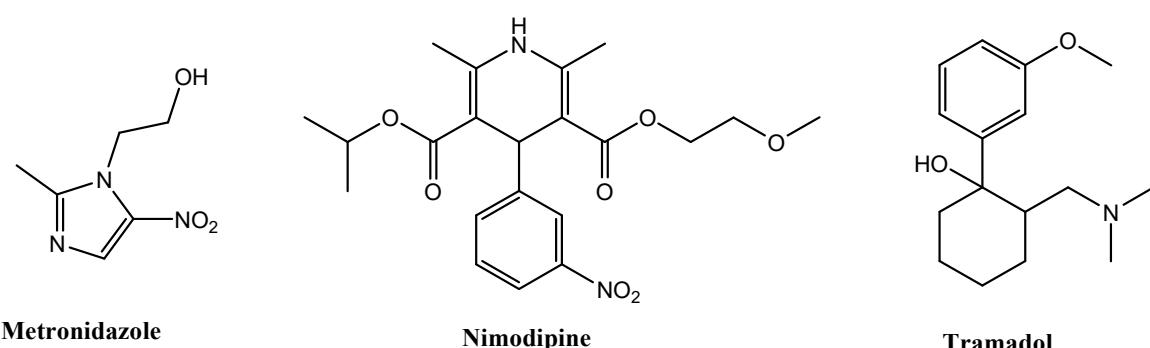


Fig. S2 Structures of metronidazole, nimodipine and tramadol.

Table S1 MS parameters for the QdNOs.

| Analytes | Precursor ions | Product ions | Collision energy |
|----------|----------------|--------------|------------------|
| OLA | 264.00 | 212.00 | 22 |
| MEQ | 219.00 | 143.00 | 23 |
| QCT | 307.00 | 273.00 | 20 |

Table S2 Scatchard equation, dissociation constant (K_d), correlation coefficient (R^2) and maximum adsorption amount (Q_{\max}) of $\text{Fe}_3\text{O}_4@\text{COOH}-\text{MIPs}$.

| Analytes | Scatchard Equation | R^2 | K_d ($\mu\text{g mL}^{-1}$) | Q_{\max} ($\mu\text{g g}^{-1}$) |
|----------|-------------------------|--------|---------------------------------|-------------------------------------|
| MEQ | $y = -0.0155x + 33.834$ | 0.9747 | 64.5 | 2182.3 |
| OLA | $y = -0.0982x + 61.498$ | 0.9887 | 10.2 | 627.3 |
| QCT | $y = -0.049x + 44.641$ | 0.9144 | 20.4 | 910.7 |

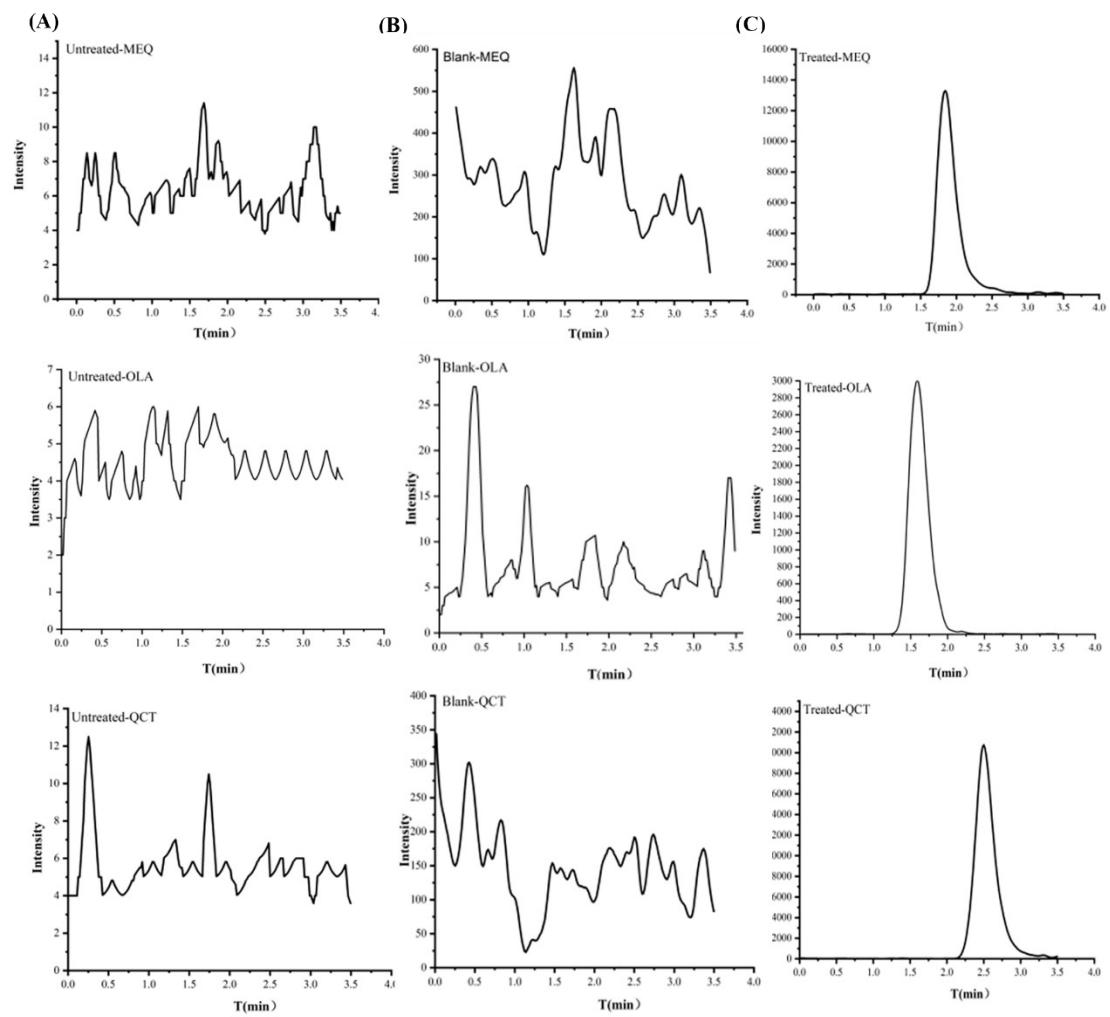


Fig. S3 Typical chromatograms of QdNOs in Hunhe Water sample. (A) un-spiked water sample before MSPE, (B) un-spiked water sample after MSPE, and (C) spiked river sample after MSPE.