

化学

## 氨基羟基脒与Pu(IV)的还原动力学研究

肖松涛; 叶国安; 刘协春; 罗方祥; 兰天; 李峰峰

中国原子能科学研究院 放射化学研究所, 北京102413

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**摘要** 研究了氨基羟基脒(HSC)与Pu(IV)的还原反应动力学, 其动力学方程式为:  $-d_c(\text{Pu(IV)})/d_t = kc(\text{Pu(IV)})c^{1.06}(\text{HSC})c^{-0.43}(\text{H}^+)c^{-0.58}(\text{NO}_3^-)$ , 在22.1 °C时反应速率常数 $k=(11.8\pm 1.1)$  (mol/L)<sup>-0.046</sup> s<sup>-1</sup>, 活化能为(71.0±1.0) kJ/mol。研究了氨基羟基脒浓度、H<sup>+</sup>浓度、硝酸根浓度、Fe<sup>3+</sup>浓度、UO<sub>2</sub><sup>2+</sup>浓度对氨基羟基脒与Pu(IV)还原反应速率的影响, 增加氨基羟基脒浓度, 降低H<sup>+</sup>浓度、硝酸根浓度, Pu(IV)还原速度增加; UO<sub>2</sub><sup>2+</sup>浓度和Fe<sup>3+</sup>浓度对Pu(IV)还原速度基本无影响。

**关键词** [Pu\(IV\)](#) [氨基羟基脒](#) [还原反应](#) [反应速率](#)

分类号

## Kinetics of Reaction Between Pu(IV) and Hydroxysemicarbazide in Nitric Acid Solution

XIAO Song-tao; YE Guo-an; LIU Xi e-chun; LUO Fang-xi ang; LAN Ti an; LI Feng-f eng

China Institute of Atomic Energy, P. O. Box 27526, Beijing 102413, C h i n a

**Abstract** The kinetics of reaction between Pu(IV) and hydroxysemicarbazide (HSC) in nitric acid solution was studied. The rate equation is found to be:  $-d_c(\text{Pu(IV)})/d_t = kc(\text{Pu(IV)})c^{1.06}(\text{HSC})c^{-0.43}(\text{H}^+)c^{-0.58}(\text{NO}_3^-)$ , where  $k=(11.8\pm 1.1)$  (mol/L)<sup>-0.046</sup> s<sup>-1</sup> at 22.1 °C and the activation energy  $E_a=(71.0\pm 1.0)$  kJ/mol. Effects of  $c(\text{HSC})$ ,  $c(\text{H}^+)$ ,  $c(\text{Fe}^{3+})$ ,  $c(\text{UO}_2^{2+})$ , ionic strength and temperature on reduction rate of Pu(IV) were investigated. The results show that Pu(IV) can be rapidly reduced to Pu(III) by HSC under normal conditions. The reaction rate can be accelerated by increasing concentration of hydroxysemicarbazide or temperature, and decreasing concentration of HNO<sub>3</sub> or ionic strength. The influence of UO<sub>2</sub><sup>2+</sup> and Fe<sup>3+</sup> on reaction rate is negligible.

**Key words** [Pu\(IV\)](#) [hydroxysemicarbazide](#) [reduction](#) [reaction](#) [reaction](#) [rate](#)

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