



Oxidative degradation of the antineoplastic drugs 5-fluorouracil and cytarabine in aqueous solution by potassium permanganate

Pingping Cheng^a, Fangyuan Sun^a, Wei Wang^a, Jingwei Feng^a, Zhen-Hu Hu^a,
Shoujun Yuan^{a,b,*}, Qiquan Wang^{c,*}

^aDepartment of Municipal Engineering, School of Civil and Hydraulic Engineering, Hefei University of Technology, Hefei 230009, China, Tel. 86-551-62904144; emails: sjyuan@hfut.edu.cn (S. Yuan), chengpingping2010@126.com (P. Cheng), 1328095157@qq.com (F. Sun), dwhit@126.com (W. Wang), fengjingwei200@163.com (J. Feng), zhhu@hfut.edu.cn (Z.-H. Hu)

^bState Key Laboratory of Pollution Control and Resource Reuse, Nanjing University, Nanjing 210046, China

^cChemistry Department, Delaware State University, Dover, DE 19901, USA, Tel. 302-857-6547; email: qwang@desu.edu

Received 18 May 2016; Accepted 26 September 2016

ABSTRACT

5-Fluorouracil (5-FU) and cytarabine (Ara-C) are two commonly used antineoplastic drugs, which are ecotoxic and genotoxic and frequently detected in hospital wastewater. Potassium permanganate (KMnO₄) has been widely used to remove organic pollutants from wastewater. The oxidative degradation of 5-FU and Ara-C in aqueous solution by KMnO₄ has not been reported. In this study, the oxidative degradation of 5-FU and Ara-C with KMnO₄ in aqueous solution was investigated. 5-FU was oxidized rapidly by KMnO₄ due to the olefinic group in the central heterocyclic ring. The degradation kinetics of 5-FU was well described by the equation of $-d[5\text{-FU}]/dt = k_{\text{app}}[5\text{-FU}]_t^{0.77}[\text{KMnO}_4]_t^{0.86}$. The degradation kinetics of Ara-C followed a generalized second-order rate law. The apparent rate constant (k_{app}) of these two target compounds was strongly depended on the reaction temperature, pH, humic acid (HA) and water quality. The apparent activation energies (E_a) of 5-FU and Ara-C were 32 and 40 kJ·mol⁻¹, respectively, indicating that 5-FU was more easily decomposed by KMnO₄ as compared with Ara-C. The maximum values of k_{app} for 5-FU and Ara-C were $1.73 \times 10^{-4} \mu\text{M}^{-0.63} \cdot \text{s}^{-1}$ at pH 7.4 and $1.52 \text{ M}^{-1} \cdot \text{s}^{-1}$ at pH 3.5, respectively. The oxidative degradation of 5-FU was inhibited by HA, while the oxidation of Ara-C was enhanced by HA. Reaction of both 5-FU and Ara-C was influenced by water quality. Thus, this study provides new insights into the degradation of antineoplastic drugs.

Keywords: 5-Fluorouracil; Cytarabine; Potassium permanganate; Kinetics; Oxidative degradation

* Corresponding author.