





Hydrolysis and Photolysis Kinetics, and Identification of Degradation Products of the Novel Bactericide 2-(4-Fluorobenzyl)-5-(Methylsulfonyl)-1,3,4-Oxadiazole in Water

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¹H NMR, ¹³C NMR, ¹⁹F NMR spectra of FBEZF's degradation product P₁ (2-(4-fluorobenzyl)-5methoxy-1,3,4-oxadiazole) were recorded on a JEOL ECX 500 NMR spectrometer operated at room temperature and 500 MHz using DMSO-d6 as a solvent and TMS as an internal standard. MS spectra were recorded on LC/MSD Trap VL.



Figure S1. ¹H NMR spectrum of degradation product P₁.



Figure S3. ¹⁹F NMR spectrum of degradation product P₁.



Figure S4. Mass Spectrum of degradation product P1.

Table S1. Hydrolysis parameters of FBEZF in water with different pH value

Initial concentration	Temperature	pН	Kinetic	Κ	Half-life	R^2
(mg L ⁻¹)	(°C)		equation	(d-1)	(d)	
5.0	25	5	$C_t = 5.0386e^{-0.048t}$	0.048	14.44	0.9832
5.0	25	7	$C_t = 4.9574 e^{-0.432t}$	0.432	1.60	0.9972
5.0	25	9	/	/	/	/

 Table S2. Hydrolysis parameters of FBEZF with different initial concentrations in water.

Initial concentration	Temperature	pН	Kinetic	Κ	Half-life	D 2
(mg L ⁻¹)	(°C)		equation	(h-1)	(h)	K^2
1.0	25	7	$C_t = 0.7419e^{-0.019t}$	0.019	36.48	0.9869
5.0	25	7	$C_t = 4.9754e^{-0.018t}$	0.018	38.51	0.9972
10.0	25	7	$C_t = 11.765e^{-0.022t}$	0.022	31.51	0.9979

 Table S3. Hydrolysis parameters of FBEZF in water under different temperature.

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	Initial concentration	Temperature	ъH	Kinetic	Κ	Half-life	R 2
	(mg L ⁻¹)	(°C)	pm	equation	(h-1)	(h)	Λ^{-}
	5.0	15	7	$C_t = 4.9709 e^{-0.009t}$	0.009	77.02	0.9949
	5.0	25	7	$C_t = 4.9754e^{-0.018t}$	0.018	38.51	0.9972
	5.0	35	7	$C_t = 4.8602e^{-0.035t}$	0.035	19.80	0.9967
	5.0	45	7	$C_t = 5.5072 e^{-0.231t}$	0.231	3.00	0.9948

Initial concentration $(mg L^{-1})$	Temperature (°C)	pН	Kinetic equation	K (h ⁻¹)	Half-life (h)	R^2
1.0	25	7	$C_t = 0.9655 e^{-0.079t}$	0.079	8.77	0.9892
5.0	25	7	$C_t = 4.2420e^{-0.083t}$	0.061	8.35	0.9909
10.0	25	7	$C_t = 9.7266e^{-0.080t}$	0.080	8.66	0.9955