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Supplementary Table 1: Analytical results obtained by using Langmuir, D-R, Freundlich and Tempkin isotherm

	Langmuir			Dubinin-Radushkevich			Freundlich			Tempkin		
	q _e	b		Xm	Е		K _f			A _T		
_	(mg/g)	(l/mol)	χ^2	(mg/g)	(kJ/mol)	χ^2	(mmol/g)	n	χ^2	(L/mg)	b	χ^2
Am ³⁺	67	0.03	0.999	71	11.3	0.987	66	11	0.968	15	37	0.955
Eu ³⁺	93	0.04	0.999	89	16.8	0.992	98	10	0.966	13	39	0.932

Supplementary Table 2: Analytical results for sorption kinetics using different models

	Lagergren first order kinetics			Intra partic	usion	Pseudo 2 nd order			
	q _e	k _{ads}	χ^2	k _p (mg g ⁻¹ min ⁻¹)	С	χ^2	q _e (mg g ⁻¹)	$k_2 (mg g^{-1} min^{-1})$	χ^2
Am ³⁺	15	0.03	0.904	31	19	0.955	71	6.1E-04	0.999
Eu^{3+}	19	0.04	0.935	39	31	0.943	97	8.6E-04	0.999

	Ti K edge			Eu L	.3 edge
Path	Parameters	Eu:DPA	Path	Parameters	Eu:DPA
Ti-O	R (Å)	1.88 ± 0.02	Eu-O	R (Å)	2.19±0.01
	Ν	4.03±0.24		Ν	2.26±0.27
	σ^2	0.0036±0.0010		σ^2	0.0015±0.0010
Ti-Ti	R (Å)	2.83±0.03	Eu-O	R (Å)	2.39±0.02
	Ν	3.16±0.24		Ν	7.91±0.98
	σ^2	0.0014±0.0011		σ^2	0.0048±0.0017
Ti-Ti	R (Å)	3.58±0.03	Eu-N	R (Å)	3.01±0.04
	Ν	1.05±0.12		Ν	4.52±0.56
	σ^2	0.0014±0.0011		σ^2	0.0072±0.0013
			Eu-C	R (Å)	3.88±0.05
				Ν	9.04±1.36
				σ^2	0.0047±0.0018

Supplementary Table 3: Bond length, coordination number and disorder factor obtain by EXAFS fitting at Ti K-edge and Eu L3 edge.



Supplementary Fig. 1: Fourier transformed EXAFS spectra at Ti K-edge of Anatase TiO₂



Supplementary Fig. 2: Optimized structures of TiO₂-APTES-DPA molecule