

Table S1. Raman modes vibrations and their corresponding assignments.

PtMo/Nb₂O₅

Raman shift (cm ⁻¹)	30-140 106 115 125 165 195 230 262 306 367										625 464 545 673 733				500-850	992 850-1000
	150-400															
Assignations	Vibrations of the Nb octahedrons and tetrahedrons Cations occupying centers of octahedrons and tetrahedrons Bending of Nb–O–Nb bonds [23,25] Stretching modes of long Nb–O bond [21-24] Me-Me vibrations in which (Me=Mo or Nb) [21]										Oxygen framework vibrations bending of Nb-O-Nb at about 548 cm ⁻¹ [23,25] Stretching vibrations of the oxygen bonds Stretch of Nb-O-Nb bridges in octahedron and tetrahedron chains [21-24]				Vibrations of the Nb-O bond O-Nb=O bonds in highly distorted structures Mo-O bonds vibrations[10,21]	

NiMo/Nb₂O₅

Raman shift (cm ⁻¹)	107	113	127	163	195	230	262	308	368	467	549	622	669	29	500-850	991				
	30-140					150-400										850-1000				
Assignations	Vibrations of the Nb octahedrons and tetrahedrons Me-Me vibrations in which (Me=Mo or Nb) [21]									Cations occupying centers of octahedrons and tetrahedrons Bending of Nb–O–Nb bonds [23,25] Stretching modes of long Nb–O bond [21-24]						Oxygen framework vibrations bending of Nb-O-Nb at about 548 cm ⁻¹ [23,25] Stretching vibrations of the oxygen bonds Stretch of Nb-O-Nb bridges in octahedron and tetrahedron chains [21-24]			Vibrations of the Nb-O bond O-Nb=O bonds in highly distorted structures Mo-O bonds vibrations	

CoMo/Nb ₂ O ₅																
Raman shift (cm ⁻¹)	107	114	125	133	195	233	264	307	483	500-850	554	621	638	728	990	
	30-140		150-400												850-1000	
Assignations	Vibrations of the Nb octahedrons and tetrahedrons			Cations occupying centers of octahedrons and tetrahedrons Bending of Nb–O–Nb bonds [23,25] Stretching modes of long Nb–O bond [21-24]						Oxygen framework vibrations vibrations bending of Nb-O-Nb at about 548 cm ⁻¹ [23,25] Stretching vibrations of the oxygen bonds Stretch of Nb-O-Nb bridges in octahedron and tetrahedron chains [21-24]					Vibrations of the Nb-O bond O-Nb=O bonds in highly distorted structures Mo-O bonds vibrations	
	Me-Me vibrations in which (Me=Mo or Nb)[21]															