

Supporting Information

Hydrothermal Synthesis, Crystal Structures, and X-ray Photoelectron Spectroscopy of Lead Tellurium(IV) and Tellurium(VI) Oxycompounds: $\text{Ba}_3\text{PbTe}_6\text{O}_{16}$ and $\text{Na}_2\text{Pb}_9(\mu_6\text{-O})_2(\text{Te}_2\text{O}_{10})_2$

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Figure S1. Energy dispersive X-ray spectroscopy analysis on a crystal of $\text{Ba}_3\text{PbTe}_6\text{O}_{16}$.

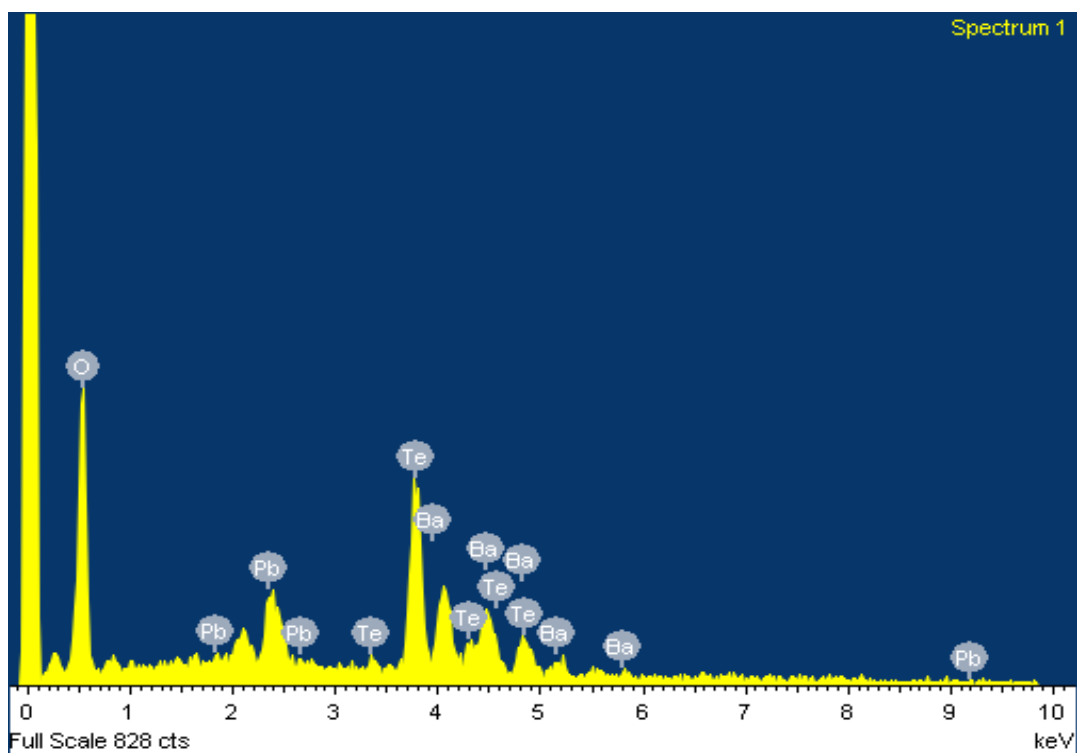
Figure S2. Pawley fit of the powder diffraction data to the structure of $\text{Ba}_3\text{PbTe}_6\text{O}_{16}$.

Figure S3. Energy dispersive X-ray spectroscopy analysis on a crystal of $\text{Na}_2\text{Pb}_9(\mu_6\text{-O})_2(\text{Te}_2\text{O}_{10})_2$.

Figure S4. Pawley fit of the powder diffraction data to the structure of $\text{Na}_2\text{Pb}_9(\mu_6\text{-O})_2(\text{Te}_2\text{O}_{10})_2$.

Figure S5. The infrared spectrum of $\text{Ba}_3\text{PbTe}_6\text{O}_{16}$ (KBr method).

Figure S6. The infrared spectrum of $\text{Na}_2\text{Pb}_9(\mu_6\text{-O})_2(\text{Te}_2\text{O}_{10})_2$ (KBr method).



EDS

Spectrum	O	Ba	Te	Pb
Spectrum 1	70.96	8.56	17.38	3.1
Spectrum 2	69.76	8.86	18.23	3.14
Mean	70.36	8.71	17.805	3.12

Atomic%

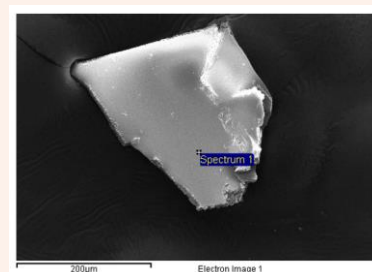
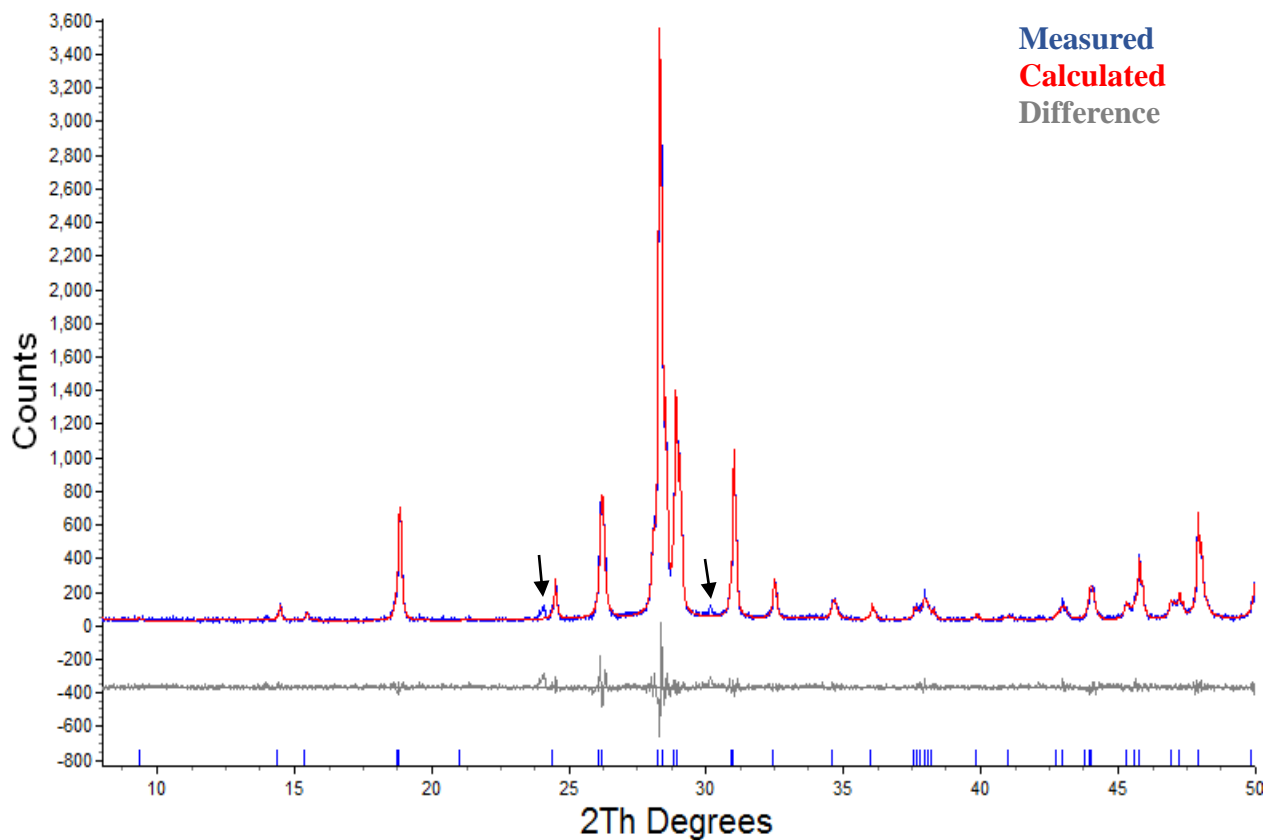
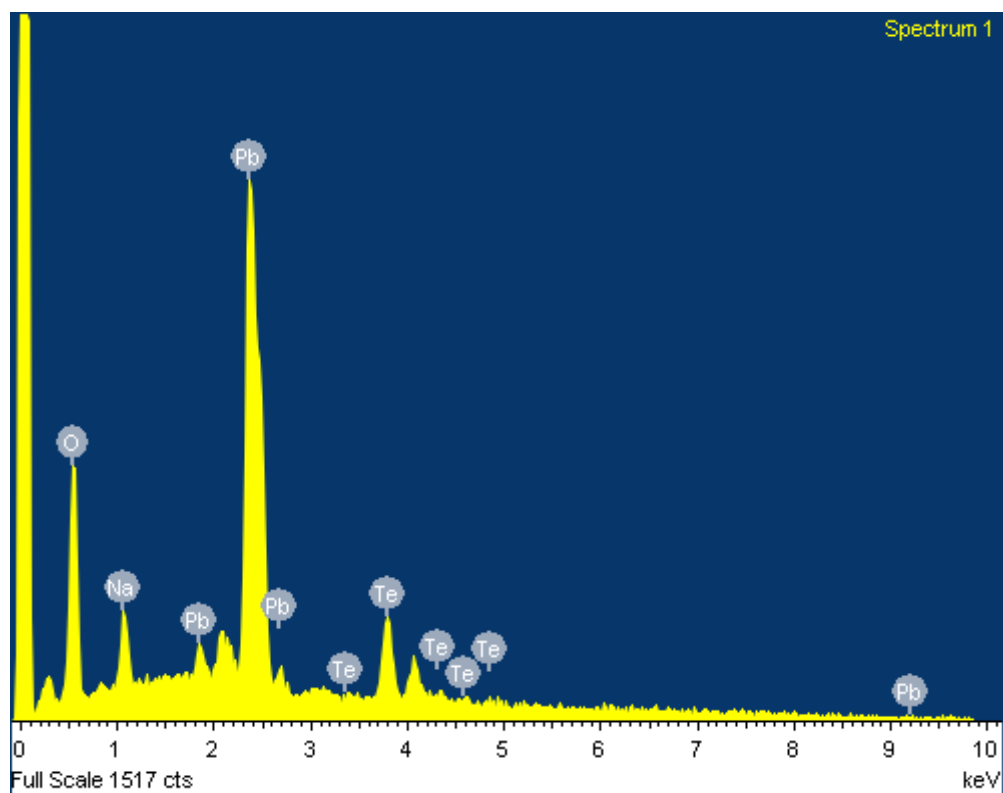


Figure S1. Energy dispersive X-ray spectroscopy analysis on a crystal of $Ba_3PbTe_6O_{16}$.



	Value
GOF	1.42
Rexp	9.64
Rwp	13.70
Rp	9.62
Rexp-dash	1.10
Rwp-dash	1.56
Rp-dash	1.86
Weighted Durbin Watson	1.28

Figure S2. Pawley fit of the powder diffraction data to the structure of $\text{Ba}_3\text{PbTe}_6\text{O}_{16}$. The small peaks marked with an arrow correspond to the impurity $\text{Ba}_3\text{Te}_4\text{O}_{11}$.



EDS

Spectrum	O	Na	Te	Pb
Spectrum 1	70.53	7.96	6.68	14.82
Spectrum 2	74.59	8.22	5.77	11.41
Mean	72.56	8.09	6.23	13.12

Atomic%

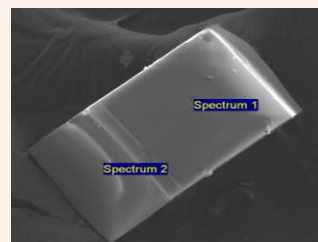
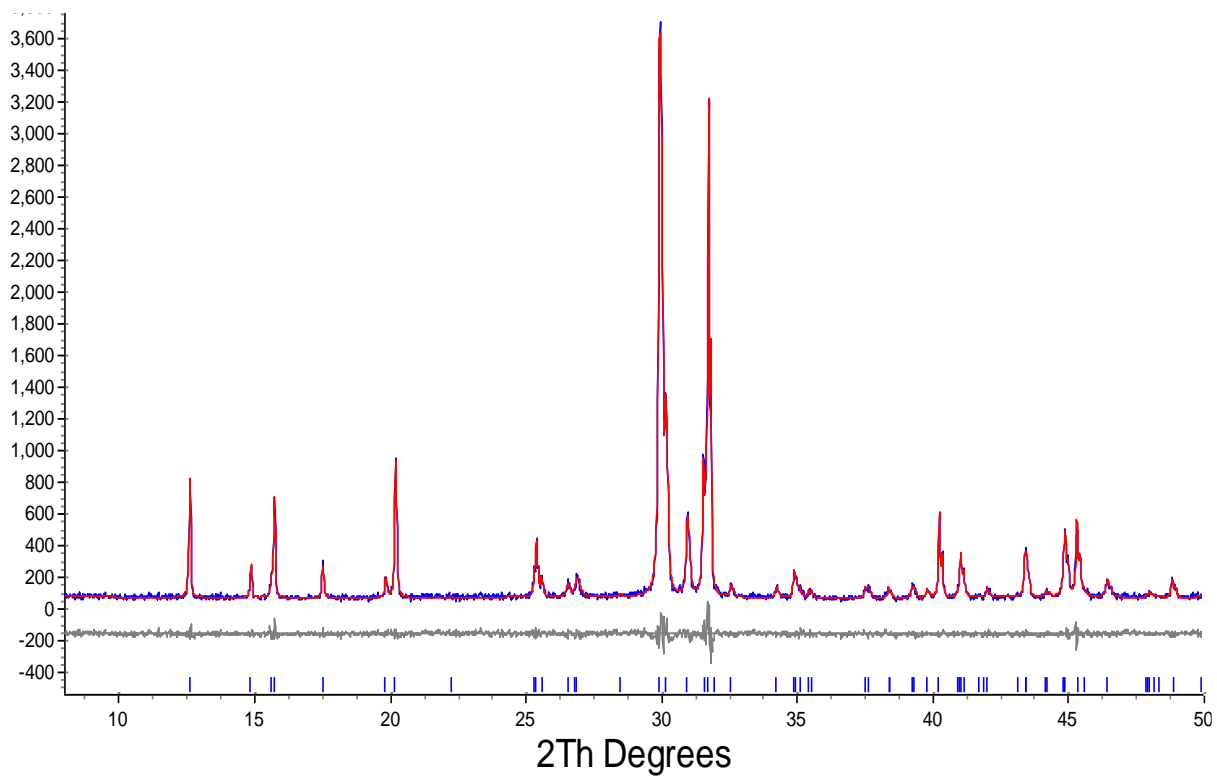


Figure S3. Energy dispersive X-ray spectroscopy analysis on a crystal of $\text{Na}_2\text{Pb}_9(\mu_6\text{-O})_2(\text{Te}_2\text{O}_{10})_2$



All range dependent		Rwps	Path	Display	Rpt/Text
		Value			
GOF		1.23			
Rexp		8.37			
Rwp		10.25			
Rp		7.66			
Rexp-dash		9.49			
Rwp-dash		11.62			
Rp-dash		9.02			
Weighted Durbin Watson		1.63			

Figure S4. Pawley fit of the powder diffraction data to the structure of $\text{Na}_2\text{Pb}_9(\mu_6\text{-O})_2(\text{Te}_2\text{O}_{10})_2$.

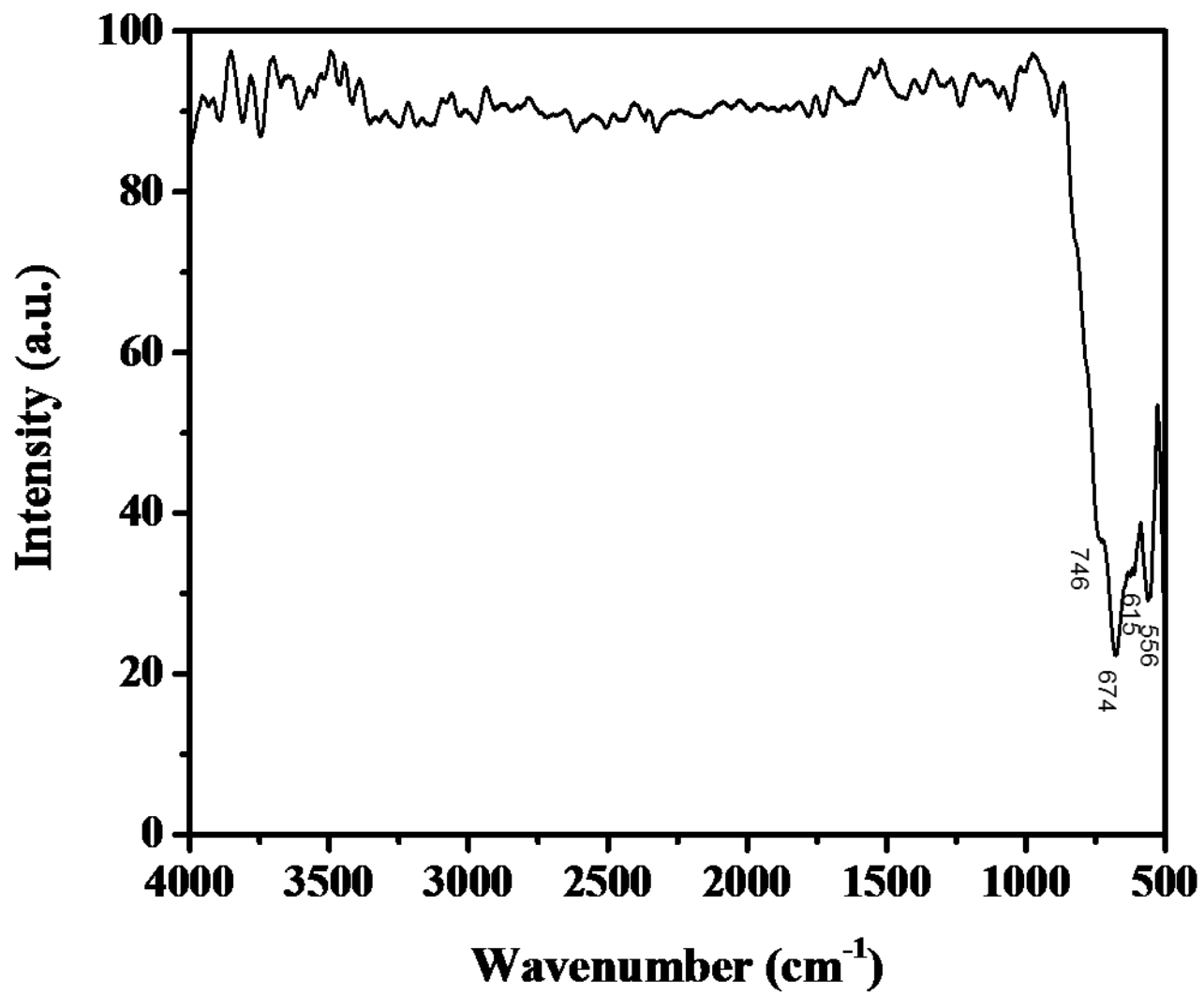


Figure S5. The infrared spectrum of Ba₃PbTe₆O₁₆ (KBr method).

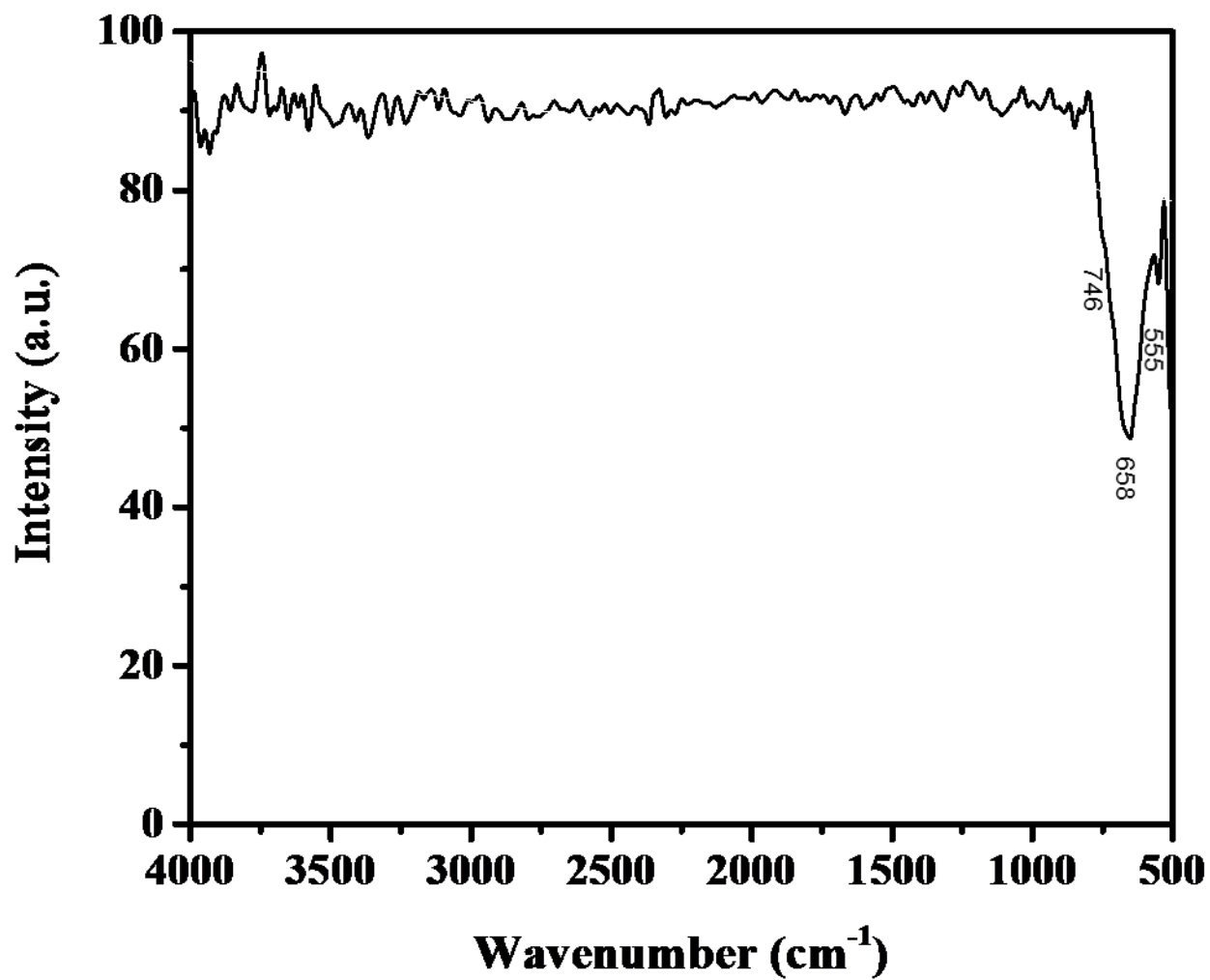


Figure S6. The infrared spectrum of $\text{Na}_2\text{Pb}_9(\mu_6\text{-O})_2(\text{Te}_2\text{O}_{10})_2$ (KBr method).