

Lampiran 1

Data-data kristal sebagai *file input* dalam Rietveld.

- **Fasa Zirkonium *Baddeleyite***

Fasa	: Zirkonium <i>Baddeleyite</i> (ZrO ₂)		
Symmetri	: <i>monoclinic</i>		
Space group	: P 1 21/c 1		
Parameter cell	a = 5.145 (Å)	α = 90.0°	
	b = 5.207 (Å)	β = 99.2°	
	c = 5.311 (Å)	γ = 90.0°	
Total atom per cell	: 14		

Tabel 1.1 Data fasa Zirkonium *Baddeleyite*.

Atom	Wyck	x	y	z	Occ
Zr	4e	0.2758	0.0411	0.2082	1
O	4e	0.0703	0.3359	0.3406	1
O	4e	0.4423	0.7549	0.4789	1

Referensi : Smith, D. K and Newkirk, H.K. 1965. Crystal Structure of Baddeleyite (Monoclinic ZrO_2) and it's Relation to The Polymorphism of ZrO_2 . *Acta Crystallography*. Vol.18. Page. 982.

- **Fasa *Tenorite***

Fasa	: <i>Tenorite</i> (CuO)	
<i>Symmetri</i>	: <i>monoclinic</i>	
<i>Space group</i>	: C2/c	
<i>Parameter cell</i>	: a = 4.6530 (Å)	$\alpha = 90.0^\circ$
	b = 3.4100 (Å)	$\beta = 99.480^\circ$
	c = 5.1080 (Å)	$\gamma = 90.0^\circ$
Total atom per <i>cell</i>	: 8	

Tabel 1.2 Data fasa *Tenorite*.

Atom	Wyck	x	y	z	Occ
Cu	4c	0.2500	0.2500	0.0000	0.5
O	4c	0.0000	0.9160	0.2500	0.5

Referensi : Tunell, E.G., Ponsjak, C.J. and Ksanda, Z. 1935. Identification of The Copper Ore Minerals by Means of X-Rays Powder Diffraction Pattern. *Zeit Krist.* Vol.90. Page.138-139.