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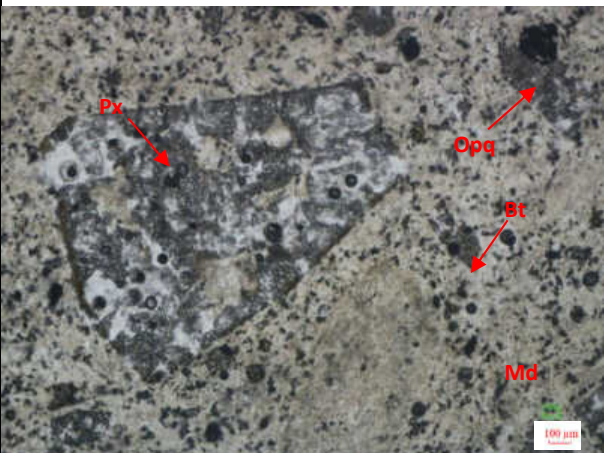
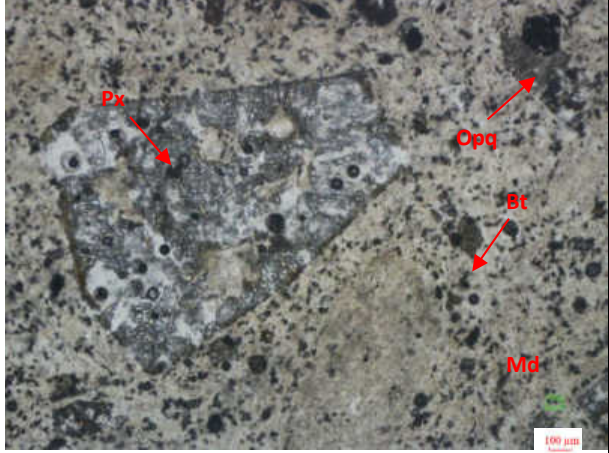
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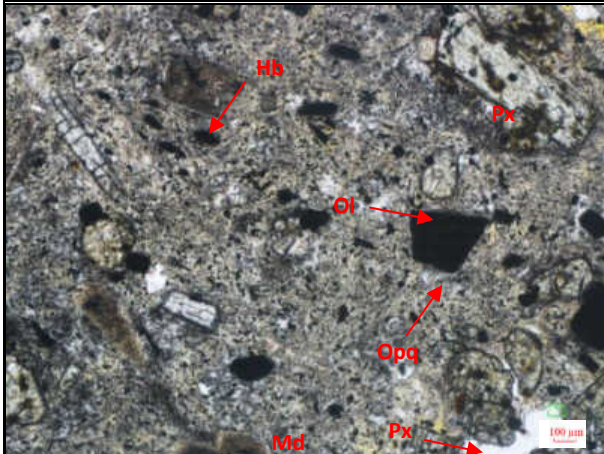
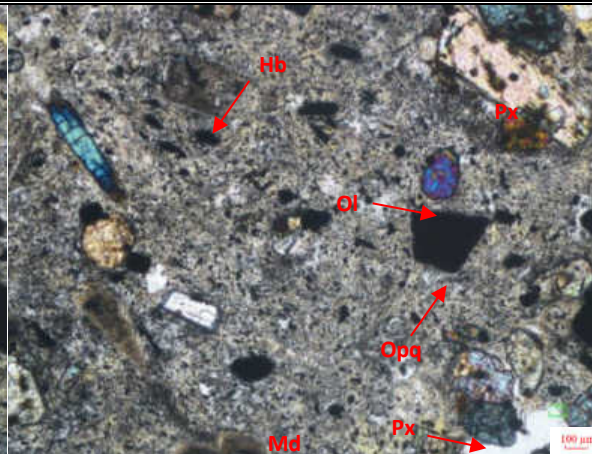
LAMPIRAN



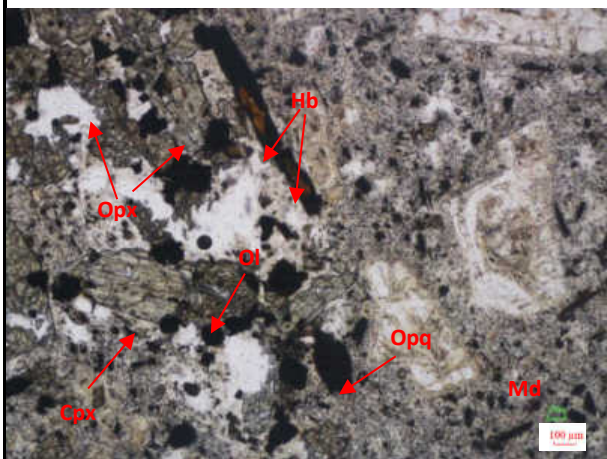
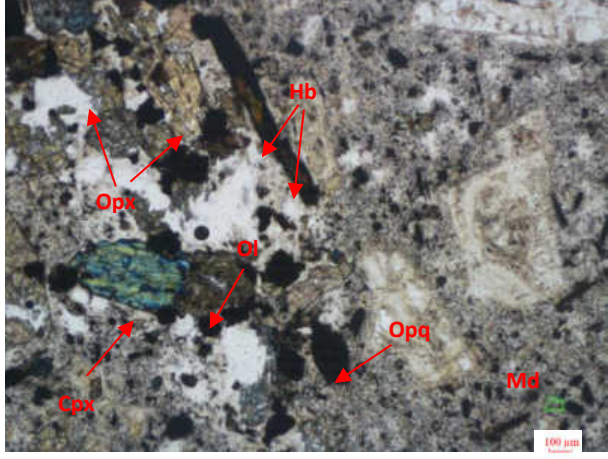
Lampiran 1. Deskripsi petrografi batuan vulkanik

Kode Sampel : 1B		Satuan : Basalt Porfiri	
Lokasi : Sinjai		Litologi : Basalt Porfiri	
			
// - Nikol		X - Nikol	
Tipe Batuan		Batuan Beku	
Tipe Struktur		Masif	
Klasifikasi		Travis (1955)	
Kenampakan Mikroskopis		Warna absorpsi kuning kecoklatan, warna interferensi coklat kehitaman. Tekstur terdiri dari kristanilitas hipokristalin, granularitas porfiri afanitik, bentuk mineral euhedral - anhedral. Relasi Inequigranular, Komposisi mineral terdiri dari mineral Piroksin, Biotit, Mineral Opaq, dan Massa Dasar. Ukuran mineral 0,05 m – 1 mm.	
Deskripsi Mineral			
Komposisi Mineral	(%)	Keterangan Optik Mineral	
Biotit (Bt)	20%	Warna Absorpsi Coklat, Bentuk subhedral - anhedral, Relief rendah, Pleokroisme Dwikroik, Ukuran 0,05 Mm. Warna Interferensi Coklat, Belahan Satu Arah, Indeks Bias $N_{min} > N_{cb}$, Sudut Gelapan 20°, Jenis Gelapan Miring	
Piroksin (Px)	30%	Warna Absorpsi <i>Colourless</i> , Bentuk Euhedral - Subhedral, Relief Tinggi, Pleokroisme Dwikroik, Ukuran 0,1 - 1 Mm, Warna Interferensi Abu-abu kehitaman, Belahan Satu Arah, Indeks Bias $N_{min} > N_{cb}$, Sudut Gelapan 23°, Jenis Gelapan Miring. Piroksen hadir sebagai fenokris.	
Mineral Opaq (Opg)	5%	Mineral Opaq Memiliki Warna Absorpsi Hitam, Warna Interferensi Hitam. Memiliki Relief Tinggi, Bentuk Anhedral, Intensitas Rendah, Ukuran 0,02 – 0,5 Mm.	
Massa Dasar (Md)	45%	Massa Dasar Terdiri Dari mikrokristalin. Massa Dasar Gelas Memiliki Warna Absorpsi Transparan/ <i>Colourless</i> , Warna Interferensi abu-abu kehitaman, Bentuk Anhedral, Ukuran Mineral < 0,02 Mm.	
Nama Batuan		Basalt Porfiri (Travis, 1955)	

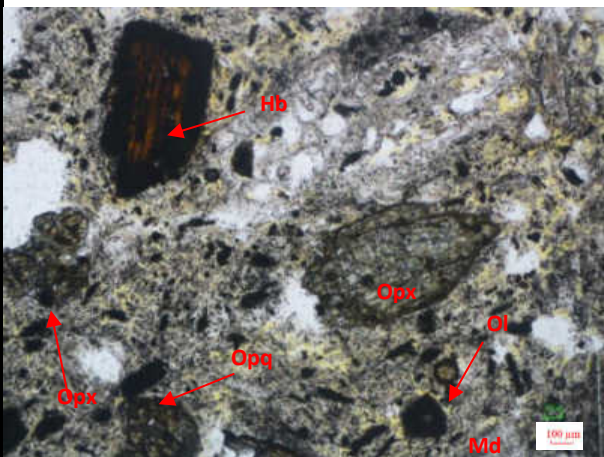
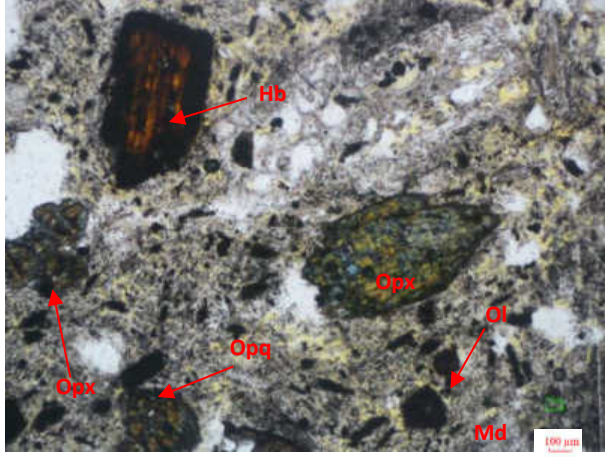


Kode Sampel : 2B		Satuan : Basalt Porfiri	
Lokasi : Sinjai		Litologi : Basalt Porfiri	
			
// - Nikol		X - Nikol	
Tipe Batuan	Batuan Beku		
Tipe Struktur	Masif		
Klasifikasi	Travis (1955)		
Kenampakan Mikroskopis	Warna absorpsi abu-abu kecoklatan, warna interferensi abu kehitaman. Tekstur terdiri dari kristanilitas hipokristalin, granularitas porfiro afanitik, bentuk mineral euhedral – anhedral, relasi inequigranular. Komposisi mineral terdiri dari mineral Olivin, Hornblend, Piroksin, Mineral Opaq, dan Massa Dasar. Ukuran mineral 0,2 m – 0,7 mm		
Deskripsi Mineral			
Komposisi Mineral	(%)	Keterangan Optik Mineral	
Olivin (Ol)	15%	Warna absorpsi Transparan/Colourless, intensitas sedang dengan relief tinggi, pecahan uneven, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna biru keunguan, pleokroisme Dwikroik, dengan sudut gelap 35°, jenis gelap miring. Ukuran mineral 0,3 mm dengan bentuk mineral euhedral sampai subhedral.	
Hornblend (Hb)	5%	Warna absorpsi coklat, intensitas rendah dengan relief tinggi, pecahan uneven, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna kuning kecoklatan, pleokroisme monokroik, dengan bentuk mineral euhedral-subhedral, sudut gelap 37°, jenis gelap miring dengan ukuran mineral 0,5 cm.	
Piroksin (Px)	25%	Warna absorpsi Transparan/Colourless, intensitas tinggi dengan relief tinggi, pecahan uneven dengan belahan satu arah, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna biru kemerahan, pleokroisme monokroik, dengan sudut gelap 51°, jenis gelap paralel. Ukuran mineral 0,3-0,7 mm dengan bentuk mineral euhedral sampai subhedral.	
Mineral Opaq (Opq)	15%	Mineral Opaq Memiliki Warna Absorpsi Hitam, Warna Interferensi Hitam. Memiliki Relief Tinggi, Bentuk Anhedral, Intensitas Rendah, Ukuran 0,02 – 0,6 Mm.	
Dasar (Md)	40%	Massa Dasar Terdiri Dari mikrokristalin. Massa Dasar Gelas Memiliki Warna Absorpsi Transparan/Colourless, Warna Interferensi abu-abu kehitaman, Bentuk Anhedral, Ukuran Mineral < 0,02 Mm.	
Nama Batuan		Basalt Porfiri (Travis, 1955)	

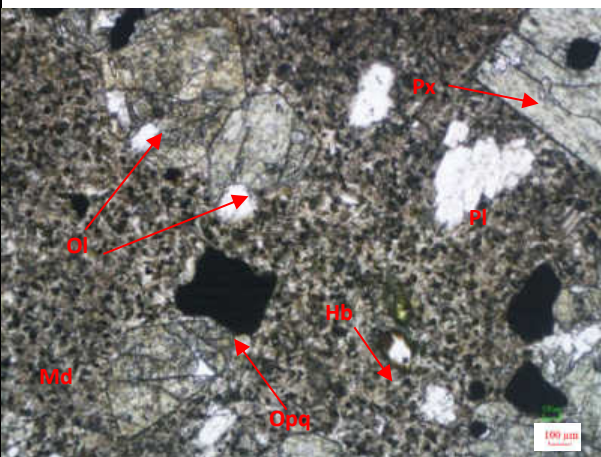
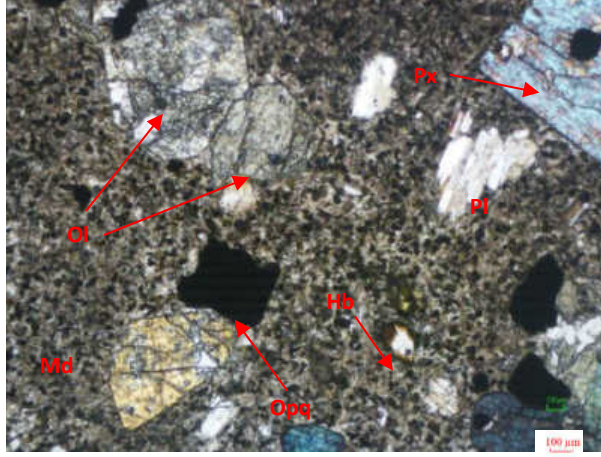


Kode Sampel : 3B		Satuan : Trakit Porfiri	
Lokasi : Sinjai		Litologi : Basalt Porfiri	
			
// - Nikol		X - Nikol	
Tipe Batuan		Batuan Beku	
Tipe Struktur		Masif	
Klasifikasi		Travis (1955)	
Kenampakan Mikroskopis		Warna absorpsi abu-abu kecoklatan, warna interferensi abu-abu kehitaman. Tekstur terdiri dari kristanilitas hipokristalin, granularitas porfiro afanitik, bentuk mineral euhedral – anhedral, relasi inequigranular. Komposisi mineral terdiri dari mineral <i>Orthopyroxene</i> , <i>Clinopyroxene</i> , Hornblend, Olivin, Mineral Opaq, dan Massa Dasar. Ukuran mineral 0,2 m – 0,8 mm.	
Deskripsi Mineral			
Komposisi Mineral	(%)	Keterangan Optik Mineral	
<i>Orthopyroxene</i> (Opx)	15%	Warna absorpsi Transparan/ <i>Colourless</i> , intensitas tinggi dengan relief tinggi, pecahan <i>uneven</i> dengan belahan dua arah, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna kuning kecoklatan, pleokroisme monokroik, dengan sudut gelapan 40°. Jenis gelapan miring, Ukuran mineral 0,5 mm dengan bentuk mineral euhedral sampai subhedral.	
<i>Clinopyroxene</i> (Cpx)	5%	Warna absorpsi Transparan/ <i>Colourless</i> , intensitas tinggi dengan relief tinggi, pecahan <i>uneven</i> dengan belahan dua arah, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna hijau kebiruan, pleokroisme monokroik, dengan sudut gelapan 48°. Jenis gelapan parallel, Ukuran mineral 0,3-0,7 mm dengan bentuk mineral euhedral sampai subhedral.	
Olivin (Ol)	20%	Warna absorpsi Transparan/ <i>Colourless</i> , intensitas sedang dengan relief tinggi, pecahan <i>uneven</i> , Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna kuning kecoklatan, pleokroisme Dwikroik, dengan sudut gelapan 40°, jenis gelapan miring, Ukuran mineral 0,3-0,5 mm dengan bentuk mineral euhedral sampai subhedral.	
Hornblend (Hb)	10%	Warna absorpsi coklat, intensitas sedang-tinggi dengan relief tinggi, pecahan <i>uneven</i> , Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna coklat kehitaman, pleokroisme monokroik, dengan bentuk mineral euhedral-subhedral, sudut gelapan 47°, jenis gelapan miring, ukuran mineral 0,2-0,8 cm.	
Mineral Opaq (Opq)	10%	Mineral Opaq Memiliki Warna Absorpsi Hitam, Warna Interferensi Hitam. Memiliki Relief Tinggi, Bentuk Anhedral, Ukuran 0,02 – 0,4 Mm	
Massa Dasar (Md)	40%	Massa Dasar Terdiri Dari mikrokristalin. Massa Dasar Gelas Memiliki Warna Absorpsi Transparan/ <i>Colourless</i> , Warna Interferensi abu-abu kehitaman, Bentuk Anhedral, Ukuran Mineral < 0,02 Mm.	
Nama Batuan		Basalt Porfiri (Travis, 1955)	

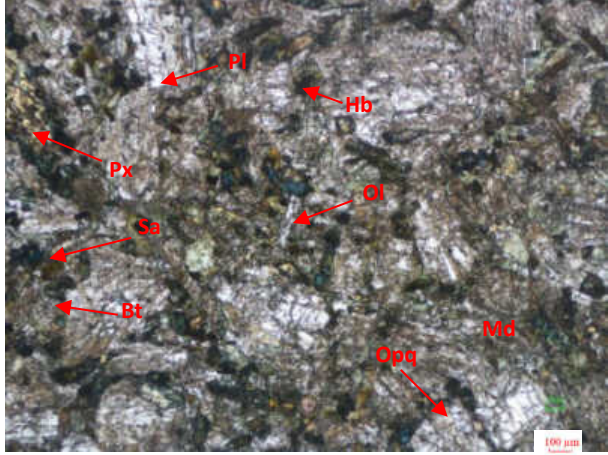


Kode Sampel : 4B		Satuan : Basalt Porfiri	
Lokasi : Sinjai		Litologi : Basalt Porfiri	
			
// - Nikol		X - Nikol	
Tipe Batuan		Batuan Beku	
Tipe Struktur		Vesicle	
Klasifikasi		Travis (1955)	
Kenampakan Mikroskopis		Warna absorpsi abu-abu kecoklatan, warna interferensi abu-abu kehitaman. Tekstur terdiri dari kristanilitas hipokristalin, granularitas Porfiro afanitik, bentuk mineral euhedral – anhedral, relasi inequigranular. Komposisi mineral terdiri dari mineral Hornblend, <i>Orthopyroxene</i> , Olivin, Mineral Opaq, dan Massa Dasar. Ukuran mineral 0,2 m – 1,2 mm.	
Deskripsi Mineral			
Komposisi Mineral	(%)	Keterangan Optik Mineral	
Hornblend (Hb)	15%	Warna Absorpsi Coklat, Bentuk Euhedral - Subhedral, Relief Tinggi, Pleokroisme Monokroik, Ukuran 1,2 Mm. Warna Interferensi Coklat, Belahan Satu Arah, Indeks Bias $N_{min} > N_{cb}$, Sudut Gelapan 20°, Jenis Gelapan Miring.	
<i>Orthopyroxene</i> (Opx)	30%	Warna Absorpsi Transparan/ <i>Colourless</i> , Bentuk Subhedral-anhedral, Relief Tinggi, Pleokroisme Dwikroik, Ukuran 0,5 – 1 Mm. Warna Interferensi kuning kecoklatan, Pecahan <i>Uneven</i> , Belahan Satu Arah, Indeks Bias $N_{min} < N_{cb}$, Sudut Gelapan 35°, Jenis Gelapan Miring	
Olivin (Ol)	5%	Warna Absorpsi Transparan/ <i>Colourless</i> , Bentuk Subhedral-anhedral, Relief Tinggi, Pleokroisme Dwikroik, Ukuran 0,3 Mm. Warna Interferensi biru keunguan, Pecahan <i>Uneven</i> , Indeks Bias $N_{min} > N_{cb}$, Sudut Gelapan 39°, Jenis Gelapan Miring	
Mineral Opaq (Opaq)	20%	Mineral Opaq Memiliki Warna Absorpsi Hitam, Warna Interferensi Hitam. Memiliki Relief Tinggi, Bentuk Anhedral, Ukuran 0,02 – 0,5 Mm	
Massa Dasar	30%	Massa Dasar Terdiri Dari mikrokristalin. Massa Dasar Gelas Memiliki Warna Absorpsi Transparan/ <i>Colourless</i> , Warna Interferensi abu-abu kehitaman, Bentuk Anhedral, Ukuran Mineral < 0,02 Mm.	
Nama Batuan		Basalt Porfiri (Travis, 1955)	



Kode Sampel : 6B		Satuan : Basalt Porfiri	
Lokasi : Maros		Litologi : Basalt Porfiri	
			
// - Nikol		X - Nikol	
Tipe Batuan		Batuan Beku	
Tipe Struktur		Masif	
Klasifikasi		Travis (1955)	
Kenampakan Mikroskopis		Warna absorpsi abu-abu kecoklatan, warna interferensi abu-abu kehitaman. Tekstur terdiri dari kristanilitas hipokristalin, granularitas Porfiro afanitik, bentuk mineral euhedral – anhedral, relasi inequigranular. Komposisi mineral terdiri dari mineral Olivin, Piroksen, Hornblend, Plagioklas, Mineral Opaq, dan Massa Dasar. Ukuran mineral 0,2 m – 1,5 mm.	
Deskripsi Mineral			
Komposisi Mineral	(%)	Keterangan Optik Mineral	
Olivin (Ol)	20%	Warna Absorpsi Transparan/Colourless, Bentuk Euhedral-subhedral, Relief Tinggi, Pleokroisme Dwikroik, Ukuran 0,8 Mm. Warna Interferensi abu-abu kehitaman, Pecahan Uneven, Indeks Bias $N_{min} > N_{cb}$, Sudut Gelapan 29°, Jenis Gelapan Miring	
Piroksen (Px)	15%	Warna absorpsi Transparan/Colourless , intensitas rendah sampai sedang dengan relief tinggi, pecahan uneven dengan belahan dua arah, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna biru, pleokroisme monokroik, dengan sudut gelapan 35°, jenis gelapan miring. Ukuran mineral 0,7-1,5 mm dengan bentuk mineral euhedral sampai subhedral.	
Hornblend (Hb)	5%	Warna absorpsi coklat, intensitas sedang dengan relief tinggi, pecahan uneven, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna kuning kecoklatan, pleokroisme monokroik, dengan bentuk mineral euhedral-subhedral, sudut gelapan miring 25°, jenis gelapan miring, ukuran mineral 0,3 cm.	
Plagioklas (Plg)	5%	Warna Absorpsi Transparan/Colourless, Warna Interferensi abu-abu kehitaman, Bentuk Subhedral - anhedral, Relief rendah, Pleokroisme monokroik, Ukuran 0,2-0,5 Mm, Belahan satu arah, Indeks Bias $N_{min} < N_{cb}$, Kembaran Albit, Sudut Gelapan 32° Jenis Gelapan Miring Sehingga Jenis Plagioklas adalah Mineral Labradorite.	
Mineral Opaq (Opq)	10%	Mineral Opaq Memiliki Warna Absorpsi Hitam, Warna Interferensi Hitam. Memiliki Relief Tinggi, Bentuk Anhedral, Ukuran 0,2-0,6 Mm	
Massa Dasar (Md)	45%	Massa Dasar Terdiri Dari mikrokristalin. Massa Dasar Gelas Memiliki Warna Absorpsi Transparan/Colourless, Warna Interferensi abu-abu kecoklatan, Bentuk Anhedral, Ukuran Mineral < 0,02 Mm.	
Nama Batuan		Basalt Porfiri (Travis, 1955)	

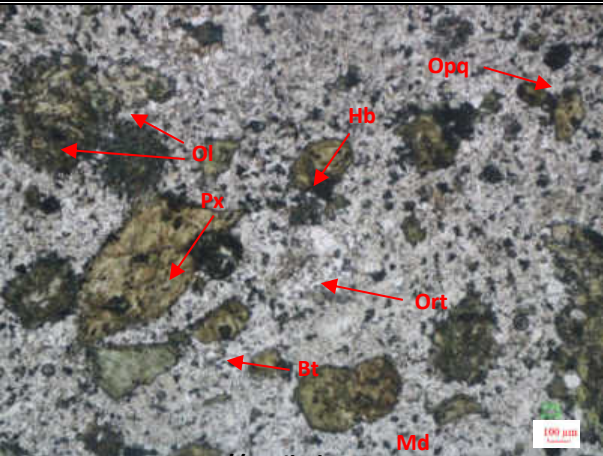
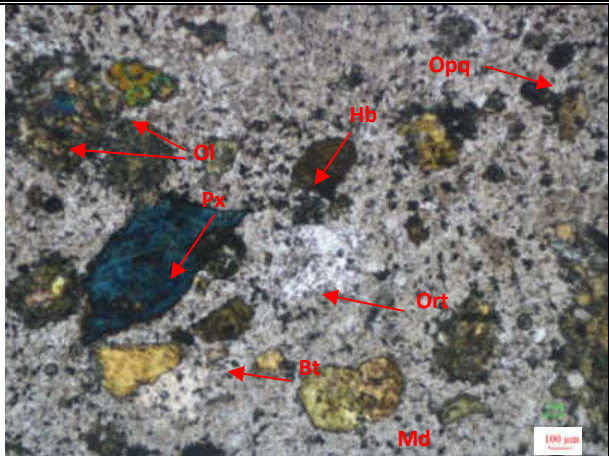


Kode Sampel : 7B ₁		Satuan : Basalt Porfiri	
Lokasi : Maros		Litologi : Basalt Porfiri	
			
// - Nikol		X - Nikol	
Tipe Batuan		Batuan Beku	
Tipe Struktur		Masif	
Klasifikasi		Travis (1955)	
Kenampakan Mikroskopis		Warna absorpsi abu-abu kecoklatan, warna interferensi abu-abu kehitaman. Tekstur terdiri dari kristanilitas hipokristalin, granularitas Porfiro afanitik, bentuk mineral euhedral – anhedral, relasi inequigranular. Komposisi mineral terdiri dari mineral Olivin, Piroksen, Hornblend, Biotit, Sanidin, Plagioklas, Mineral Opaq, dan Massa Dasar. Ukuran mineral 0,1 m – 0,7 mm.	
Deskripsi Mineral			
Komposisi Mineral	(%)	Keterangan Optik Mineral	
Olivin (Ol)	15%	Warna Absorpsi Transparan/Colourless, Bentuk Subhedral-Anhedral, Relief Tinggi, Pleokroisme Dwikroik, Ukuran 0,3 Mm. Warna Interferensi biru, Pecahan Uneven, Indeks Bias Nmin>Ncb, Sudut Gelapan 23°, Jenis Gelapan Miring	
Piroksen (Px)	5%	Warna absorpsi Transparan/Colourless, intensitas rendah sampai sedang dengan relief tinggi, pecahan uneven dengan belahan satu arah, Indeks bias Nmin < Ncb. Pada warna interferensi berwarna abu-abu kecoklatan, pleokroisme monokroik, dengan sudut gelap 35°, jenis gelap miring. Ukuran mineral 0,3-0,5 mm dengan bentuk mineral euhedral sampai subhedral.	
Hornblend (Hb)	10%	Warna absorpsi coklat, intensitas sedang dengan relief tinggi, pecahan uneven, Indeks bias Nmin < Ncb. Pada warna interferensi berwarna kuning kecoklatan, pleokroisme monokroik, dengan bentuk mineral subhedral-anhedral, sudut gelap miring 25°, jenis gelap miring, ukuran mineral 0,2 cm.	
Biotit (Bt)	20%	Warna absorpsi coklat, intensitas sedang dengan relief rendah, pecahan uneven, Indeks bias Nmin < Ncb. Pada warna interferensi berwarna kuning kecoklatan, pleokroisme monokroik, dengan bentuk mineral subhedral-anhedral, sudut gelap miring 28°, jenis gelap miring, ukuran mineral 0,15 cm.	
Sanidin (Sa)	5%	Warna Absorpsi Transparan/Colourless, Warna Interferensi abu-abu kehitaman, Bentuk Subhedral - anhedral, Relief rendah, Pleokroisme monokroik, Ukuran 0,2-0,5 Mm, Belahan satu arah, Indeks Bias Nmin<Ncb, Kembaran Carlsbad, Sudut Gelapan 21° Jenis Gelapan Miring	
Plagioklas (Plg)	5%	Warna Absorpsi Transparan/Colourless, Warna Interferensi abu-abu kehitaman, Bentuk Subhedral - anhedral, Relief rendah, Pleokroisme monokroik, Ukuran 0,2-0,7 Mm, Belahan satu arah, Indeks Bias	



		Nmin<Ncb, Kembaran Albit, Sudut Gelapan 38° Jenis Gelapan Miring Sehingga Jenis Plagioklas adalah Mineral Labradorite.
Mineral Opaq (Opq)	25%	Mineral Opaq Memiliki Warna Absorpsi Hitam, Warna Interferensi Hitam. Memiliki Relief Tinggi, Bentuk Anhedral, Ukuran 0,1-0,3 Mm
Massa Dasar (Md)	15%	Massa Dasar Terdiri Dari mikrokristalin. Massa Dasar Gelas Memiliki Warna Absorpsi Transparan/ <i>Colourless</i> , Warna Interferensi abu-abu kecoklatan, Bentuk Anhedral, Ukuran Mineral < 0,02 Mm.
Nama Batuan		Basalt Porfiri (Travis, 1955)

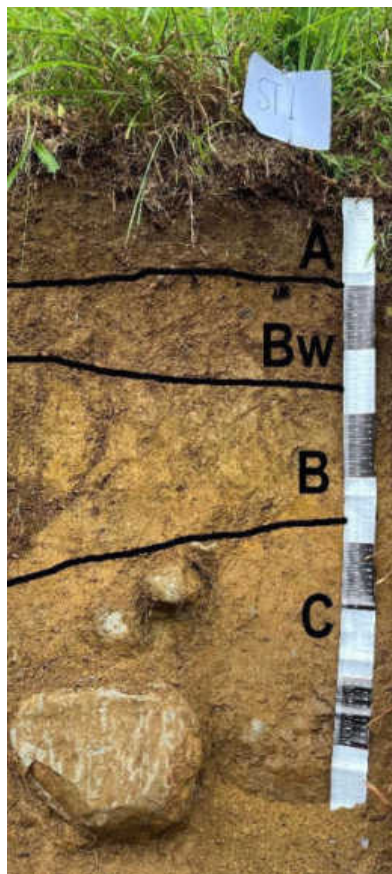


Kode Sampel : 7B2 Lokasi : Maros		Satuan : Basalt Porfiri Litologi : Basalt Porfiri	
			
// - Nikol		X - Nikol	
Tipe Batuan		Batuan Beku	
Tipe Struktur		Masif	
Klasifikasi		Travis (1955)	
Kenampakan Mikroskopis		Warna absorpsi abu-abu kecoklatan, warna interferensi abu-abu kehitaman. Tekstur terdiri dari kristanilitas hipokristalin, granularitas Porfiro afanitik, bentuk mineral euhedral – anhedral, relasi inequigranular. Komposisi mineral terdiri dari mineral Olivin, Piroksen, Hornblend, Biotit, Ortoklas, Mineral Opaq, dan Massa Dasar. Ukuran mineral 0,2 m – 0,7 mm.	
Deskripsi Mineral			
Komposisi Mineral	(%)	Keterangan Optik Mineral	
Olivin (Ol)	10%	Warna Absorpsi Transparan/Colourless, Bentuk Subhedral-Anhedral, Relief Tinggi, Pleokroisme Dwikroik, Ukuran 0,3-0,5 Mm. Warna Interferensi biru keunguan, Pecahan Uneven, Indeks Bias $N_{min} > N_{cb}$, Sudut Gelapan 33°, Jenis Gelapan Miring	
Piroksen (Px)	20%	Warna absorpsi Transparan/Colourless , intensitas rendah sampai sedang dengan relief tinggi, pecahan uneven dengan belahan satu arah, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna biru, pleokroisme monokroik, dengan sudut gelapan 37°, jenis gelapan miring. Ukuran mineral 0,3-0,7 mm dengan bentuk mineral euhedral sampai subhedral.	
Hornblend (Hb)	15%	Warna absorpsi coklat, intensitas sedang dengan relief tinggi, pecahan uneven, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna coklat kehitaman, pleokroisme monokroik, dengan bentuk mineral subhedral-anhedral, sudut gelapan miring 19°, jenis gelapan miring, ukuran mineral 0,2-0,4 cm.	
Biotit (Bt)	5%	Warna absorpsi coklat, intensitas sedang dengan relief rendah, pecahan uneven, Indeks bias $N_{min} < N_{cb}$. Pada warna interferensi berwarna kuning kecoklatan, pleokroisme monokroik, dengan bentuk mineral subhedral-anhedral, sudut gelapan miring 24°, jenis gelapan miring, ukuran mineral 0,2 cm.	
Ortoklas (Ort)	5%	Warna Absorpsi Transparan/Colourless, Warna Interferensi abu-abu kehitaman, Bentuk Subhedral - anhedral, Relief rendah, pecahan uneven Pleokroisme monokroik, Ukuran 0,2-0,4 Mm, Indeks Bias $N_{min} < N_{cb}$, Kembaran Carlsbad, Sudut Gelapan 21° Jenis Gelapan Miring	
Mineral Opaq (Opq)	10%	Mineral Opaq Memiliki Warna Absorpsi Hitam, Warna Interferensi Hitam. Memiliki Relief Tinggi, Bentuk Anhedral, Ukuran 0,1-0,2 Mm	
Massa Dasar (Md)	35%	Massa Dasar Terdiri Dari mikrokristalin. Massa Dasar Gelas Memiliki Warna Absorpsi Transparan/Colourless, Warna Interferensi abu-abu kecoklatan, Bentuk Anhedral, Ukuran Mineral < 0,02 Mm.	
Nama Batuan		Basalt Porfiri (Travis, 1955)	



Lampiran 2. Deskripsi profil *soil***Tabel 1.** Deskripsi profil stasiun pengamatan 1

Keterangan	Deskripsi
Kode profil	ST 1
Tanggal pengamatan	27 Mei 2023
Koordinat	5°17'4" LS dan 119°59'23" BT
Desa	Barania, Sinjai Barat
Altitude	1370 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	Pohon pinus, semak dan rumput
Kedalaman <i>soil</i>	93 cm
Horizon A (0-19 cm)	Tekstur liat berdebu
Horizon Bw (19-52 cm)	Tekstur liat
Horizon B (52-93 cm)	Tekstur liat berdebu



(a)



(b)

Gambar 1. Penampang profil *soil* (a) dan bentang lahan (b) ST 1

Tabel 2. Deskripsi profil stasiun pengamatan 2

Keterangan	Deskripsi
Kode profil	ST 2
Tanggal pengamatan	27 Mei 2023
Koordinat	5°17'8" LS dan 119°58'50" BT
Desa	Gunung Perak, Sinjai Barat
Altitude	1520 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	Pohon pinus, semak dan rumput
Kedalaman <i>soil</i>	14 cm
Horizon A (0-14 cm)	Tekstur liat



(a)



(b)

Gambar 2. Penampang profil *soil* (a) dan bentang lahan (b) ST 2

Tabel 3. Deskripsi profil stasiun pengamatan 3

Keterangan	Deskripsi
Kode profil	ST 3
Tanggal pengamatan	27 Mei 2023
Koordinat	5°17'25" LS dan 119°58'25" BT
Desa	Gunung Perak, Sinjai Barat
Altitude	1610 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	Pohon pinus, semak dan rumput
Kedalaman <i>soil</i>	30 cm
Horizon A (0-30 cm)	Tekstur lempung liat berdebu



(a)

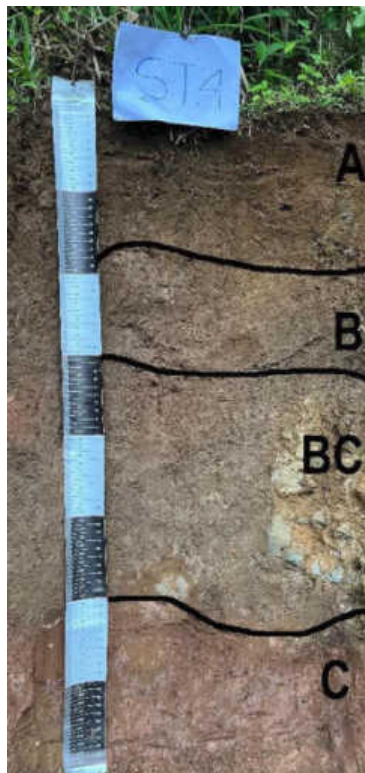


(b)

Gambar 3. Penampang profil *soil* (a) dan bentang lahan (b) ST 3

Tabel 4. Deskripsi profil stasiun pengamatan 4

Keterangan	Deskripsi
Kode profil	ST 4
Tanggal pengamatan	28 Mei 2023
Koordinat	5°15'51" LS dan 119°58'53" BT
Desa	Gunung Perak, Sinjai Barat
Altitude	1230 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	Pohon pinus, semak dan rumput
Kedalaman <i>soil</i>	120 cm
Horizon A (0-36 cm)	Tekstur liat berdebu
Horizon B (36-60 cm)	Tekstur liat berdebu
Horizon BC (60-120 cm)	Tekstur liat berdebu



(a)



(b)

Gambar 4. Penampang profil *soil* (a) dan bentang lahan (b) ST 4

Tabel 5. Deskripsi profil stasiun pengamatan 5

Keterangan	Deskripsi
Kode profil	ST 5
Tanggal pengamatan	28 Mei 2023
Koordinat	5°16'38" LS dan 119°58'33" BT
Desa	Gunung Perak, Sinjai Barat
Altitude	1430 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	Pohon pinus, semak dan rumput
Kedalaman <i>soil</i>	20 cm
Horizon A (0-20 cm)	Tekstur lempung liat berdebu



(a)



(b)

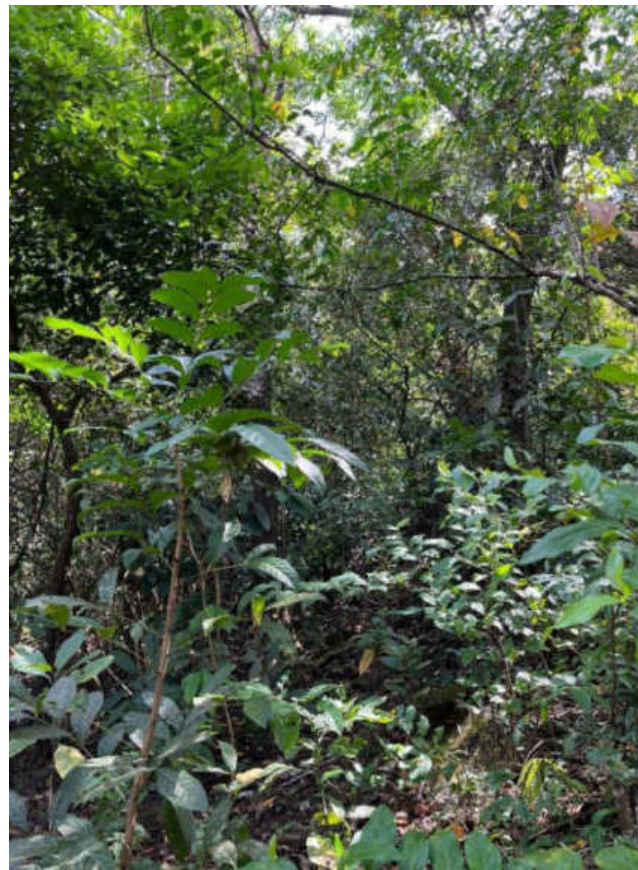
Gambar 5. Penampang profil *soil* (a) dan bentang lahan (b) ST 5

Tabel 6. Deskripsi profil stasiun pengamatan 6

Keterangan	Deskripsi
Kode profil	ST 6
Tanggal pengamatan	22 Juli 2023
Koordinat	5°08'27" LS dan 119°48'09" BT
Desa	Bonto Somba, Tompobulu
Altitude	538 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	
Kedalaman <i>soil</i>	105 cm
Horizon A (0-13 cm)	Tekstur lempung
Horizon Bt (13-105 cm)	Tekstur lempung liat



(a)

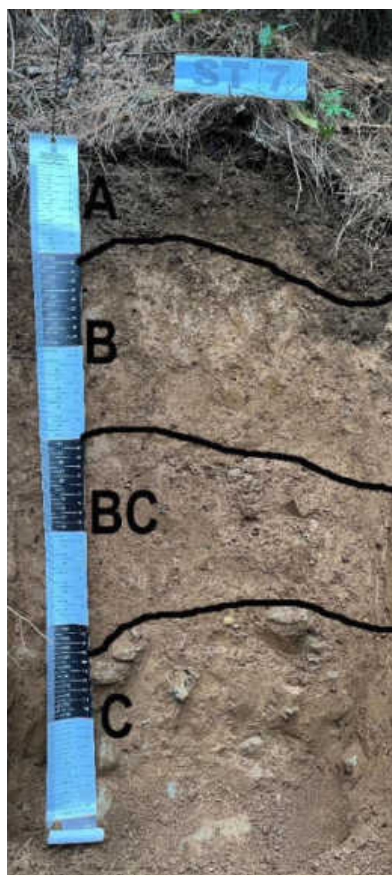


(b)

Gambar 6. Penampang profil *soil* (a) dan bentang lahan (b) ST 6

Tabel 7. Deskripsi profil stasiun pengamatan 7

Keterangan	Deskripsi
Kode profil	ST 7
Tanggal pengamatan	22 Juli 2023
Koordinat	5°09'32.3" LS dan 119°48'57.9" BT
Desa	Bonto Somba, Tompobulu
Altitude	565 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	Pohon pinus, pakis dan semak
Kedalaman <i>soil</i>	106 cm
Horizon A (0-22 cm)	Tekstur lempung
Horizon B (22-60 cm)	Tekstur lempung berdebu
Horizon BC (60-106 cm)	Tekstur lempung liat



(a)



(b)

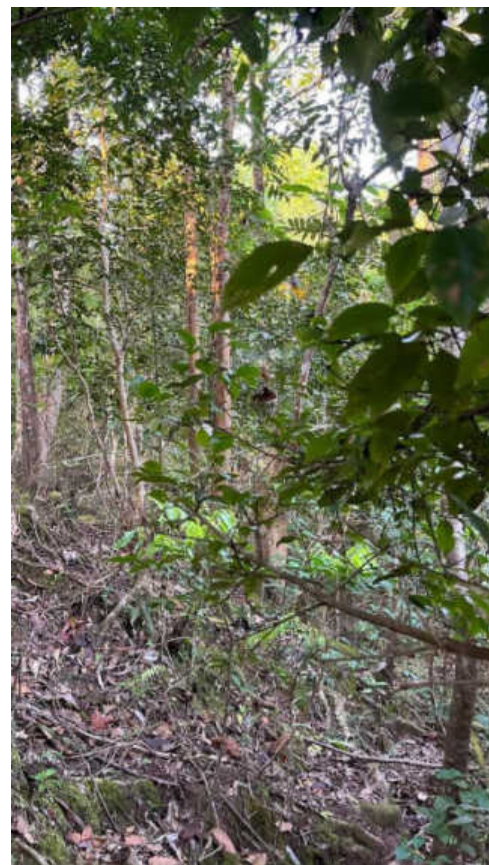
Gambar 7. Penampang profil *soil* (a) dan bentang lahan (b) ST 7

Tabel 8. Deskripsi profil stasiun pengamatan 8

Keterangan	Deskripsi
Kode profil	ST 8
Tanggal pengamatan	22 Juli 2023
Koordinat	5°09'12.4" LS dan 119°48'31.3" BT
Desa	Bonto Somba, Tompobulu
Altitude	503 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	
Kedalaman <i>soil</i>	52 cm
Horizon A (0-19 cm)	Tekstur lempung
Horizon B (19-52 cm)	Tekstur lempung berdebu



(a)

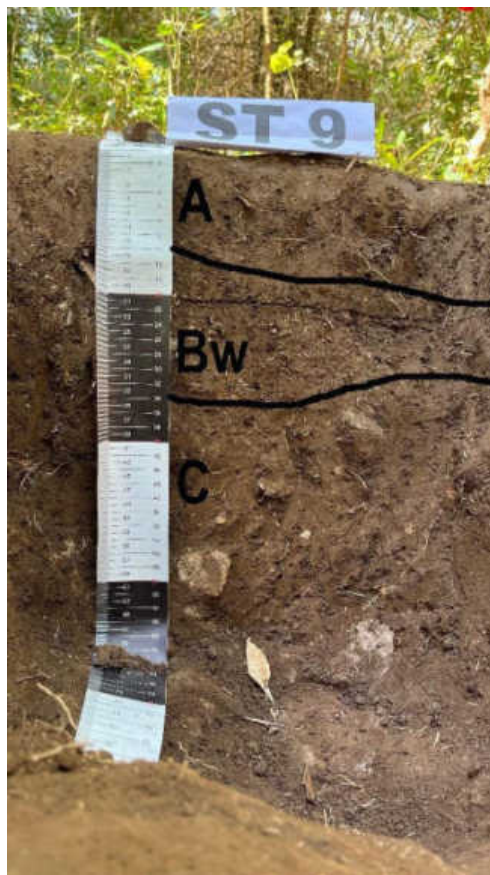


(b)

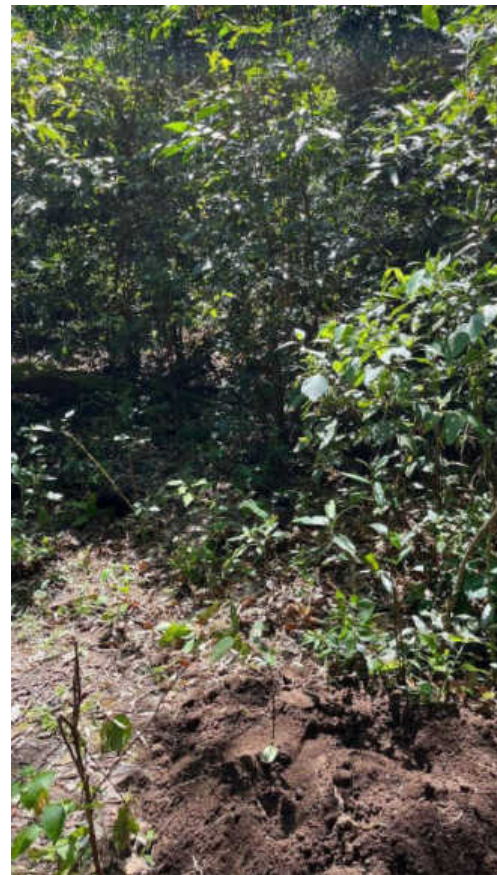
Gambar 8. Penampang profil *soil* (a) dan bentang lahan (b) ST 8

Tabel 9. Deskripsi profil stasiun pengamatan 9

Keterangan	Deskripsi
Kode profil	ST 9
Tanggal pengamatan	23 Juli 2023
Koordinat	5°08'33.2" LS dan 119°46'49" BT
Desa	Bonto Somba, Tompobulu
Altitude	533 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	
Kedalaman <i>soil</i>	34 cm
Horizon A (0-14 cm)	Tekstur lempung
Horizon Bw (14-34 cm)	Tekstur lempung



(a)



(b)

Gambar 9. Penampang profil *soil* (a) dan bentang lahan (b) ST 9

Tabel 10. Deskripsi profil stasiun pengamatan 10

Keterangan	Deskripsi
Kode profil	ST 10
Tanggal pengamatan	23 Juli 2023
Koordinat	5°08'41.1" LS dan 119°46'19.9" BT
Desa	Bonto Somba, Tompobulu
Altitude	452 mdpl
Penggunaan lahan	Hutan sekunder
Vegetasi	Semak, rumput, pinus
Kedalaman <i>soil</i>	10 cm
Horizon A (0-10 cm)	Tekstur lempung



(a)



(b)

Gambar 10. Penampang profil *soil* (a) dan bentang lahan (b) ST 10

Lampiran 3. Hasil analisis karakteristik sifat fisik *soil* lokasi penelitian

Fomasi Batuan	Profil	Horizon	Kedalaman cm	Tekstur			Kelas tekstur
				Pasir	Debu	Liat	
				----- % -----			
Batuan Gunungapi Lompobattang	ST 1	A	0-19	6	42	52	Liat berdebu
		Bw	19-52	5	39	56	Liat
		B	52-93	7	40	53	Liat berdebu
	ST 2	A	0-14	9	38	53	Liat
	ST 3	A	0-30	17	46	37	Lempung liat berdebu
	ST 4	A	0-36	6	52	42	Liat berdebu
		B	36-60	8	49	43	Liat berdebu
		BC	60-120	12	45	43	Liat berdebu
	ST 5	A	0-20	10	51	39	Lempung liat berdebu
	Batuan Gunungapi Baturape Cindakko	ST 6	A	0-13	42	42	16
Bt			13-105	23	41	36	Lempung liat
ST 7		A	0-22	39	36	25	Lempung
		B	22-60	29	50	21	Lempung berdebu
		BC	60-106	32	37	31	Lempung liat
ST 8		A	0-19	38	42	20	Lempung
		B	19-52	23	53	24	Lempung Berdebu
ST 9		A	0-14	26	47	26	Lempung
		Bw	14-34	39	45	16	Lempung
ST 10		A	0-10	39	43	18	Lempung



Lampiran 4. Hasil analisis karakteristik sifat kimia *soil* lokasi penelitian

Formasi Batuan	Profil	Horizon	Kedalaman	C-Organik	Basa-basa tukar				Jumlah
					Ca	Mg	K	Na	
					-----cmol kg ⁻¹ -----				
Batuan Gunungapi Lompobattang	ST 1	A	0-19	1,35	5,15	0,95	0,25	0,27	6,62
		AB	20-51	1,04	3,25	1,02	0,16	0,22	4,65
		B	52-95	0,65	1,95	0,85	0,11	0,19	3,1
	ST 2	A	0-14	2,12	8,25	1	0,41	0,32	9,98
	ST 3	A	0-30	1,33	5,45	1,74	0,38	0,3	7,87
	ST 4	A	0-32	1,25	3,25	2,01	0,35	0,25	5,86
		B1	33-60	0,82	2,95	1,65	0,25	0,11	4,96
		B2	61-119	1,02	1,25	0,73	0,19	0,11	2,28
	ST 5	A	0-36	2,81	8,05	0,85	0,25	0,41	9,56
	Batuan Gunungapi Baturape Cindakko	ST 6	A	0-13	2,67	2,64	3,24	0,38	0,53
B1			13-73	0,65	1,46	1,31	0,43	0,43	3,63
ST 7		A	0-22	2,37	1,41	1,5	0,68	0,62	4,21
		B1	22-66	0,7	0,32	0,18	0,25	0,47	1,22
		B2	66-142	0,49	0,73	0,5	0,29	0,36	1,88
ST 8		A	0-19	1,05	2,17	3,95	0,31	0,51	6,94
		AB	19-52	0,3	0,5	0,42	0,24	0,42	1,58
ST 9		A	0-20	1,92	5,65	6,85	0,98	0,41	64,74
		B	20-66	1,19	4,95	5,8	0,89	0,43	56,37
ST 10		A	0-10	2,65	1,99	1,43	0,56	0,53	4,51



Lampiran 5. Kriteria penilaian hasil analisis kimia *soil*

Sifat Tanah	Sangat Rendah	Rendah	Sedang	Tinggi	Sangat Tinggi
Karbon (%)	< 1,00	1,00-2,00	2,01-3,00	3,01-5,00	>5,00
Nitrogen(%)	< 0,10	0,10-0,20	0,21-0,50	0,51-0,75	>0,75
C/N	<5,0	5,0-7,9	8,0-12,0	12,1-17,0	>17
P ₂ O ₅ eks- HCl (%)	<0,021	0,021-0,039	0,040-0,060	0,061-0,100	>0,100
P-avl Bray-II (ppm)	<8,0	8,0-15	16-25	26-35	>35
P-avl Olsen (ppm)	<10	10-25	26-45	46-60	>60
K ₂ O eks-HCl (mg/100)	<0,03	0,03-0,06	0,07-0,11	0,12-0,20	>0,20
KTK/CEC (me/100)	<5	10-16	17-24	25-40	>40
Susunan Kation					
K-tukar (me/ 100)	<0,1	0,1-0,2	0,3-0,5	0,6-1,0	>1,0
Na-tukar (me/ 100)	<0,1	0,1-0,3	0,4-0,7	0,8-1,0	>1,0
Mg-tukar (me/ 100)	<0,4	0,4-1,0	1,1-2,0	2,1-8,0	>8,0
Ca-tukar (me/ 100)	<2,0	2-5	6-10	11-20	>20
Kejenuhan Basa (%)	<20	20-35	36-50	51-70	>70
Kejenuhan Al (%)	<10	10-20	21-30	31-60	>60

	Sangat Masam	Masam	Agak Masam	Netral	Agak Alkalis	Alkalis
pH (H ₂ O)	<4,5	4,5-5,5	5,6-6,5	6,6-7,5	7,6-8,5	>8,5
pH (KCl)	<2,5	2,5-4,0	-----	4,1-6,0	6,1-6,5	>6,5

Sumber : Hardjowigeno, S. 1995. Ilmu Tanah



Match! Phase Analysis Report

Sample: ST-1-L1 (5-70)

Sample Data

File name ST-1-L1.RAW
File path D:/S2 Geologi/data hasil/17 sampel/ST-1-L1
Data collected Jan 5, 2024 16:52:42
Data range 5.040° - 70.040°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction 0.04°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	34.0	Olivine	Mg2 O4 Si
B	24.4	Potassium aluminium silicate hydroxide * Muscovite 2M1	Al3 H K O12 Si3
C	24.3	Vermiculite	H2 Mg3 O12 Si4
D	9.7	Bytownite	Al7.76 Ca3.44 Na0.56 O32 Si8.24
E	3.3	Hornblende	Al2.42 Ca1.802 Fe1.846 H2 Mg2.224 Na0.628 O24 Si6.44 Ti0.07
F	2.7	Anorthite	Al Ca O4 Si
G	1.6	Quartz	O2 Si
H	0.0	Montmorillonite	Al2 Ca O12 Si4
	1.1	Unidentified peak area	

A: Olivine (34.0 %)

Formula sum Mg2 O4 Si
Entry number 96-900-6399
Figure-of-Merit (FoM) 0.662490
Total number of peaks 168
Peaks in range 168
Peaks matched 39
Intensity scale factor 0.58
Space group P b n m
Crystal system orthorhombic
Unit cell a= 4.7536 Å b= 10.2066 Å c= 5.9845 Å
I/Icor 0.83
Calc. density 3.217 g/cm³
Reference Muller-Sommer M, Hock R., Kirfel A., "Rietveld refinement study of the cation distribution in (Co, Mg)-olivine solid solution Sample: MR100", Physics and Chemistry of Minerals **24**, 17-23 (1997)

B: Potassium aluminium silicate hydroxide * Muscovite 2M1 (24.4 %)

Formula sum Al3 H K O12 Si3
Entry number 96-110-0014
Figure-of-Merit (FoM) 0.775981
Total number of peaks 299
Peaks in range 299
Peaks matched 111
Intensity scale factor 1.61
Space group C 1 2/c 1
Crystal system monoclinic
Unit cell a= 5.1800 Å b= 9.0200 Å c= 20.0400 Å β= 95.500 °
I/Icor 3.19
Meas. density 2.798 g/cm³
Calc. density 2.830 g/cm³
Reference Jackson WW, West J, "", Zeitschrift fuer Kristallographie, Kristallgeometrie, Kristallphysik, Kristallchemie (-144,1977) **76**, 211-227 (1931)

C: Vermiculite (24.3 %)

Formula sum H2 Mg3 O12 Si4
Entry number 96-900-0017
Figure-of-Merit (FoM) 0.727104
Total number of peaks 263
Peaks 2
Peaks 2
Intensity 0.78
Space group I c 1
Crystal system monoclinic
Unit cell a= 5.3300 Å b= 9.1800 Å c= 28.8500 Å β= 93.250 °
I/Icor 0.53
Calc. density 2.86 g/cm³
Reference Andriks S. B., Jefferson M. E., "Crystal structure of vermiculites and mixed vermiculite-chlorites", American Mineralogist **23**, 1-862 (1938)



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D: Bytownite (9.7 %)

Formula sum Al7.76 Ca3.44 Na0.56 O32 Si8.24
Entry number 96-901-1202
Figure-of-Merit (FoM) 0.665004

Figure-of-Merit (FoM)	0.685031
Total number of peaks	500
Peaks in range	500
Peaks matched	219
Intensity scale factor	0.14
Space group	P -1
Crystal system	triclinic (anorthic)
Unit cell	a= 8.1830 Å b= 12.8830 Å c= 14.1860 Å α= 93.380° β= 115.870° γ= 90.820°
I/lcor	0.60
Calc. density	2.729 g/cm ³
Reference	Facchinelli A, Bruno E., Chiari G., "The structure of bytownite quenched from 1723 K Locality: satellite dyke, Traversella stock, Sesia Lanzo zone, Western Alps, Italy Sample: BytQ, P-1 model", Acta Crystallographica, Section B 35 (1), 34-42 (1979)

E: Hornblende (3.3 %)

Formula sum	Al2.42 Ca1.802 Fe1.846 H2 Mg2.224 Na0.628 O24 Si6.44 Ti0.07
Entry number	96-900-1226
Figure-of-Merit (FoM)	0.716829
Total number of peaks	299
Peaks in range	299
Peaks matched	119
Intensity scale factor	0.07
Space group	C 1 2/m 1
Crystal system	monoclinic
Unit cell	a= 9.8130 Å b= 18.0550 Å c= 5.3210 Å β= 104.970°
I/lcor	1.09
Calc. density	3.205 g/cm ³
Reference	Phillips M W., Draheim J. E., Popp R. K., Clowe C. A., Pinkerton A A., "Effects of oxidation-dehydrogenation in tschermakitic hornblende sample H-1, natural", American Mineralogist 74 , 764-773 (1989)

F: Anorthite (2.7 %)

Formula sum	Al Ca O4 Si
Entry number	96-900-1173
Figure-of-Merit (FoM)	0.654532
Total number of peaks	248
Peaks in range	248
Peaks matched	131
Intensity scale factor	0.04
Space group	I -1
Crystal system	triclinic (anorthic)
Unit cell	a= 8.0820 Å b= 12.7670 Å c= 14.0320 Å α= 92.790° β= 115.770° γ= 91.680°
I/lcor	0.68
Calc. density	3.250 g/cm ³
Reference	Angel R. J., "High-pressure structure of anorthite Sample: P = 25 kbar", American Mineralogist 73 , 1114-1119 (1988)

G: Quartz (1.6 %)

Formula sum	O2 Si
Entry number	96-900-5025
Figure-of-Merit (FoM)	0.695380
Total number of peaks	35
Peaks in range	35
Peaks matched	11
Intensity scale factor	0.12
Space group	P 32 2 1
Crystal system	trigonal (hexagonal axes)
Unit cell	a= 4.9841 Å c= 5.4500 Å
I/lcor	3.65
Calc. density	2.552 g/cm ³
Reference	Kihara K., "An X-ray study of the temperature dependence of the quartz structure Sample: at T = 838 K", European Journal of Mineralogy 2 , 63-77 (1990)

H: Montmorillonite (0.0 %)

Formula sum	Al2 Ca O12 Si4
Entry number	96-110-1055
Figure-of-Merit (FoM)	0.671105
Total number of peaks	92
Peaks in range	91
Peaks matched	66
Intensity scale factor	0.00
Space group	P 1
Crystal system	triclinic (anorthic)
Unit cell	a= 5.1800 Å b= 8.9800 Å c= 15.0000 Å α= 90.000° β= 90.000° γ= 90.000°
I/lcor	20.53
Calc. density	1.800 g/cm ³

Name
Bytowr



Formula

Al7.76 Ca3.44 Na0.56 O32 Si8.2496-901-1202

Entry No.

FoM

0.6464

Candidates

Search-Match

Setting

Refered to: D-Inorg REV89244 2013.10.11

Autom

Minim

Parameter/influence 2theta 0.50

Parameter/influence intensities 0.50

Parameter multiple/single phase(s) 0.50

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Selection Criteria

Compound:

Name: Bytownite

Peak List

No.	2 θ [°]	d [Å]	I/I ₀	FWHM	Matched
1	6.68	13.2181	171.24	0.8800	C,D,H
2	11.63	7.6043	240.49	0.8800	D,E,F
3	11.93	7.4100	281.14	0.8800	H
4	12.13	7.2880	356.65	0.8800	C
5	12.49	7.0827	308.68	0.8800	
6	12.84	6.8890	165.03	0.8800	D,F
7	12.84	6.8890	155.55	0.8800	
8	12.84	6.8890	155.55	0.8800	
9	18.36	4.8271	258.80	0.8800	AB,C,D,E,F,H
10	19.43	4.5659	156.65	0.8800	B,C,D,E,F,H
11	20.11	4.4128	1000.00	0.8800	B,C,D,E,H
12	20.35	4.3607	897.75	0.8800	C,D
13	20.43	4.3444	899.14	0.8800	C,F
14	20.66	4.2957	746.56	0.8800	AB,D,G,H
15	20.66	4.2957	752.71	0.8800	
16	20.66	4.2957	746.56	0.8800	
17	20.90	4.2469	657.95	0.8800	C,D,F
18	20.90	4.2469	649.04	0.8800	
19	20.90	4.2469	649.04	0.8800	
20	21.17	4.1933	652.03	0.8800	B,E
21	21.34	4.1604	634.53	0.8800	D
22	21.34	4.1604	625.75	0.8800	
23	21.34	4.1604	625.75	0.8800	
24	21.50	4.1298	583.17	0.8800	C,D
25	21.50	4.1298	575.06	0.8800	
26	21.50	4.1298	575.06	0.8800	
27	21.68	4.0959	511.35	0.8800	B,C,F
28	21.68	4.0959	521.78	0.8800	
29	21.68	4.0959	511.35	0.8800	
30	21.98	4.0413	413.43	0.8800	D,E
31	22.28	3.9869	373.00	0.8800	E,F
32	22.28	3.9869	373.00	0.8800	
33	22.30	3.9834	377.24	0.8800	C,D,F
34	22.51	3.9470	334.64	0.8800	B,D,F
35	22.68	3.9183	333.58	0.8800	D,F
36	22.92	3.8765	292.36	0.8800	A,C,D,E,F
37	23.05	3.8558	290.63	0.8800	B,D,F,H
38	23.36	3.8053	260.69	0.8800	C,D,F
39	23.74	3.7442	289.00	0.8800	AB,D,F,H
40	24.23	3.6702	408.24	0.8800	C,D,F
41	24.42	3.6423	393.80	0.8800	D,F
42	24.77	3.5914	486.16	0.8800	B,C,D,F
43	25.09	3.5463	499.48	0.8800	D,F
44	25.32	3.5147	429.74	0.8800	C,D
45	25.38	3.5065	414.60	0.8800	AD,F
46	25.38	3.5065	414.60	0.8800	
47	25.71	3.4621	317.49	0.8800	AB,D
48	25.96	3.4295	229.43	0.8800	C,D,F
49	26.38	3.3752	254.32	0.8800	D,E,G
50	26.66	3.3413	285.46	0.8800	B,F,H
51	26.82	3.3213	275.55	0.8800	B,C
52	27.06	3.2925	183.64	0.8800	E
53	27.49	3.2421	194.36	0.8800	B,C,D,E,F
54	33.25	2.6924	249.94	0.8800	AB,C,D,E,F
55	35.11	2.5539	422.57	0.8800	AB,C,D,E,F,H
56	35.62	2.5183	650.51	0.8800	AC,D,F
57	35.92	2.4981	496.56	0.8800	B,C,D,F,H
58	35.92	2.4981	496.56	0.8800	
59	35.94	2.4968	485.67	0.8800	B,D,G,H
60	36.18	2.4810	373.06	0.8800	C,D,E,F
61	36.66	2.4494	396.49	0.8800	AB,C,D,E,F,H
62	36.82	2.4391	360.12	0.8800	B,C,D,E,F
63	36.82	2.4391	352.35	0.8800	
64			2.35	0.8800	
65			7.89	0.8800	D,E,F
66			3.71	0.8800	C,D,E
67			2.35	0.8800	AB,C,D,E,F
68			3.27	0.8800	D,F
69			3.86	0.8800	AC,F
70			3.82	0.8800	E
71			3.82	0.8800	
72			0.75	0.8800	C,D,F
73			6.33	0.8800	AB,D,E,F
74			0.76	0.8800	C,D,E,F,G,H
75			3.67	0.8800	C,D,E,F
76	39.44	2.2829	214.33	0.8800	C,D
77	39.44	2.2829	206.22	0.8800	
78	39.44	2.2829	206.22	0.8800	



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79	39.61	2.2733	198.13	0.8800	A,C,D,F,G
80	39.94	2.2554	164.80	0.8800	B,C,D,E
81	40.14	2.2445	199.51	0.8800	AB,F,H
82	40.38	2.2319	193.69	0.8800	B,C,D,E,F
83	40.75	2.2123	211.04	0.8800	B,C,D,E,F,H
84	41.02	2.1988	154.06	0.8800	B,C,D,E,F,H
85	41.74	2.1621	164.72	0.8800	AB,C,D,E,F,G,H
86	54.42	1.6847	302.11	0.8800	AB,C,E,G,H
87	54.74	1.6756	275.48	0.8800	B,C,E,G
88	54.92	1.6705	232.64	0.8800	AB,C,H
89	55.07	1.6663	221.95	0.8800	
90	55.32	1.6592	215.75	0.8800	B
91	55.57	1.6524	190.25	0.8800	B,C,E,H
92	56.10	1.6381	155.92	0.8800	AB,C,E,G
93	56.80	1.6196	163.56	0.8800	AB,C,E
94	57.12	1.6112	164.34	0.8800	AB,C,E,H
95	62.43	1.4864	421.49	0.8800	AB,C,E,H
96	62.58	1.4831	404.73	0.8800	B,E
97	62.88	1.4768	275.45	0.8800	AB,C,E,H
98	62.88	1.4768	289.45	0.8800	
99	62.88	1.4768	275.45	0.8800	
100	63.18	1.4705	171.50	0.8800	AB,C,E,G,H
101	63.18	1.4705	154.30	0.8800	
102	63.18	1.4705	154.30	0.8800	
103	63.86	1.4565	201.95	0.8800	B,E
104	64.08	1.4520	162.23	0.8800	AB,C,E,G,H

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

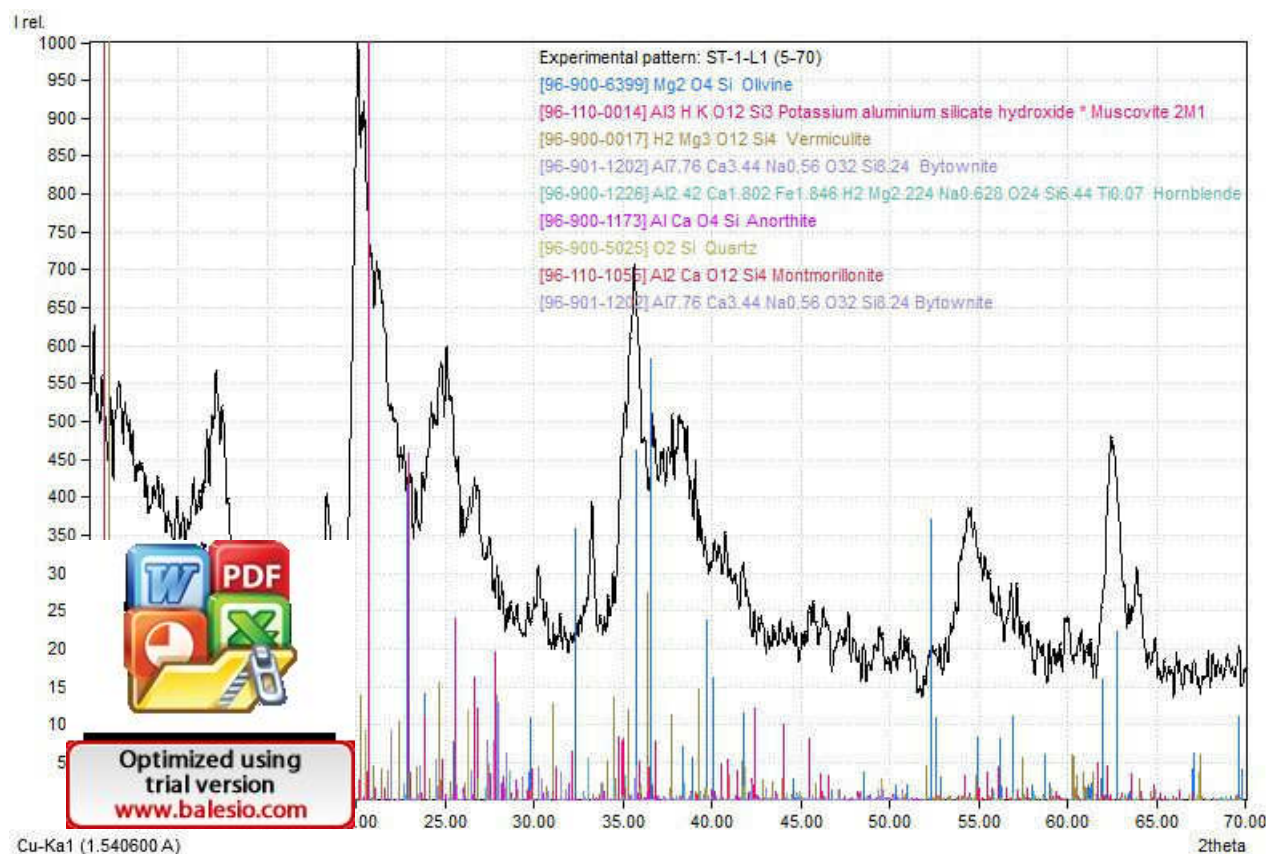
Based on calculated profile

Profile area	Counts	Amount
Overall diffraction profile	113962	100.00%
Background radiation	75249	66.03%
Diffraction peaks	38713	33.97%
Peak area belonging to selected phases	37411	32.83%
Unidentified peak area	1302	1.14%

Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	4940	100.00%
Peak intensity belonging to selected phases	1026	20.77%
Unidentified peak intensity	3914	79.23%

Diffraction Pattern Graphics





Match! Phase Analysis Report

Sample: ST-2-L1 (5-70)

Sample Data

File name ST-2-L1.RAW
File path D:/S2 Geologi/data hasil/17 sampel/ST-2-L1
Data collected Jan 5, 2024 16:52:43
Data range 5.030° - 70.030°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction 0.03°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	33.3	Anorthite	Al ₂ CaO ₈ Si ₂
B	23.2	Olivine	Fe _{0.2} Mg _{1.792} Mn _{0.003} Ni _{0.005} O ₄ Si
C	10.9	Quartz	O ₂ Si
D	10.5	Potassium aluminium silicate hydroxide * Muscovite 2M	Al ₃ H ₂ KO ₁₂ Si ₃
E	10.5	Hornblende	Al _{2.42} Ca _{1.83} Fe _{1.852} H _{1.34} Mg _{2.218} Na _{0.706} O ₂₄ Si _{6.44} Ti _{0.07}
F	7.0	Pyroxene	Al _{1.388} Ca _{0.742} Fe _{0.162} Mg _{0.016} O ₆ Si _{1.5}
G	3.1	Montmorillonite	Al _{0.86} Fe _{0.1} HLi _{0.08} Mg _{0.14} O ₁₀ Si _{3.9}
H	1.5	Vermiculite	Al _{0.57} H _{1.4} Mg _{1.705} O _{7.86} Si _{1.43}
	18.0	Unidentified peak area	

A: Anorthite (33.3 %)

Formula sum Al₂CaO₈Si₂
Entry number 96-900-0362
Figure-of-Merit (FoM) 0.770657
Total number of peaks 500
Peaks in range 500
Peaks matched 288
Intensity scale factor 0.31
Space group P -1
Crystal system triclinic (anorthic)
Unit cell a= 8.1940 Å b= 12.8970 Å c= 14.1900 Å α= 92.980° β= 115.820° γ= 91.150°
I/Cor 0.60
Calc. density 2.745 g/cm³
Reference Foit F. F., Peacor D. R., "The anorthite crystal structure at 410 and 830 C T= 410 C", American Mineralogist **58**, 665-675 (1973)

B: Olivine (23.2 %)

Formula sum Fe_{0.2}Mg_{1.792}Mn_{0.003}Ni_{0.005}O₄Si
Entry number 96-900-5917
Figure-of-Merit (FoM) 0.714859
Total number of peaks 168
Peaks in range 168
Peaks matched 42
Intensity scale factor 0.33
Space group P b n m
Crystal system orthorhombic
Unit cell a= 4.7631 Å b= 10.2230 Å c= 5.9920 Å
I/Cor 0.90
Calc. density 3.353 g/cm³
Reference Ottonello G., Princivalle F., Della Giusta A., "Temperature, composition, and fo₂ effects on intersite distribution of Mg and Fe²⁺ in olivines Sample: CC37a", Physics and Chemistry of Minerals **17**, 301-312 (1990)

C: Quartz (10.9 %)

Formula sum O₂Si
Entry number 96-901-2603
Figure-of-Merit (FoM) 0.702299
Total number of peaks 32
Peaks in range 32
Peaks matched 11
Intensity scale factor 0.43
Space group P 31 2 1
Crystal system trigonal (hexagonal axes)
Unit cell a= 4.7050 Å c= 5.2500 Å
I/Cor 0.73
Calc. density 2.654 g/cm³
Reference Finger R. M., Finger L. W., Hemley R. J., Mao H. K., "High-pressure crystal chemistry and amorphization of alpha-quartz Locality: Synthetic Sample: P = 5.1 GPa", Solid State Communications **72**, 507-511 (1989)



D: Pot

hydro
%)
Formu H₂KO₁₂Si₃
Entry number 96-101-1050
Figure-of-Merit (FoM) 0.824334
Total number of peaks 299
Peaks in range 299

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Peaks in range	299
Peaks matched	105
Intensity scale factor	0.53
Space group	C 1 2/c 1
Crystal system	monoclinic
Unit cell	a= 5.1800 Å b= 9.0200 Å c= 20.0400 Å β= 95.500 °
I/Icor	3.19
Meas. density	2.798 g/cm³
Calc. density	2.830 g/cm³
Reference	Jackson W W, West J, "The Crystal Structure of Muscovite - K Al ₃ Si ₃ O ₁₀ (OH) ₂ ", Zeitschrift fuer Kristallographie, Kristallgeometrie, Kristallphysik, Kristallchemie (-144,1977) 76 , 211-227 (1931)

E: Hornblende (10.5 %)

Formula sum	Al _{2.42} Ca _{1.83} Fe _{1.852} H _{1.34} Mg _{2.218} Na _{0.706} O ₂₄ Si _{6.44} Ti _{0.07}
Entry number	96-900-1228
Figure-of-Merit (FoM)	0.765631
Total number of peaks	299
Peaks in range	299
Peaks matched	118
Intensity scale factor	0.16
Space group	C 1 2/m 1
Crystal system	monoclinic
Unit cell	a= 9.7860 Å b= 18.0240 Å c= 5.3060 Å β= 105.090 °
I/Icor	0.97
Calc. density	3.241 g/cm³
Reference	Phillips M W, Draheim J. E., Popp R. K., Clowe C. A., Pinkerton A A, "Effects of oxidation-dehydrogenation in tschermakitic hornblende sample H-3, after annealing at 700 C", American Mineralogist 74 , 764-773 (1989)

F: Pyroxene (7.0 %)

Formula sum	Al _{1.388} Ca _{0.742} Fe _{0.162} Mg _{0.016} O ₆ Si _{1.5}
Entry number	96-900-6564
Figure-of-Merit (FoM)	0.764486
Total number of peaks	229
Peaks in range	229
Peaks matched	57
Intensity scale factor	0.13
Space group	C 1 2/c 1
Crystal system	monoclinic
Unit cell	a= 9.7190 Å b= 8.8140 Å c= 5.3050 Å β= 106.040 °
I/Icor	1.20
Calc. density	3.266 g/cm³
Reference	Okui M, Sawada H., Marumo F., "Structure refinement of a nonstoichiometric pyroxene synthesized under ambient pressure", Physics and Chemistry of Minerals 25 , 318-322 (1998)

G: Montmorillonite (3.1 %)

Formula sum	Al _{0.86} Fe _{0.1} H Li _{0.08} Mg _{0.14} O ₁₀ Si _{3.9}
Entry number	96-901-0959
Figure-of-Merit (FoM)	0.754715
Total number of peaks	251
Peaks in range	251
Peaks matched	62
Intensity scale factor	0.29
Space group	C 1 2/m 1
Crystal system	monoclinic
Unit cell	a= 5.1710 Å b= 8.9570 Å c= 9.7400 Å β= 96.100 °
I/Icor	5.89
Calc. density	2.245 g/cm³
Reference	Gournis D., Lappas A, Karakassides M A, Tobbens D., Moukarika A, "A neutron diffraction study of alkali cation migration in montmorillonites Sample: Li-mont-300", Physics and Chemistry of Minerals 35 , 49-58 (2008)

H: Vermiculite (1.5 %)

Formula sum	Al _{0.57} H _{1.4} Mg _{1.705} O _{7.86} Si _{1.43}
Entry number	96-900-0147
Figure-of-Merit (FoM)	0.775151
Total number of peaks	291
Peaks in range	290
Peaks matched	124
Intensity scale factor	0.22
Space group	C 1 2/c 1
Crystal system	monoclinic
Unit cell	a= 5.3490 Å b= 9.2550 Å c= 28.8900 Å β= 97.120 °
I/Icor	9.57
Calc. density	2.096 g/cm³
Reference	~irozu H., Bailey S. W., "Crystal structure of a two-layer Mg-vermiculite", American Mineralogist 51 , 1124-1143 (1966)

Name	Formula	Entry No.	FoM
Quartz	O2 Si	96-900-9667	0.6995
Silicor	O2 Si	96-500-0036	0.6946
Quartz	O2 Si	96-901-3322	0.6941
Quartz	O2 Si	96-901-2601	0.6937
Quartz	O2 Si	96-901-2605	0.6817
Quartz	O2 Si	96-900-8094	0.6793
Quartz	O2 Si	96-901-2602	0.6721
Quartz	O2 Si	96-900-8093	0.6699
Quartz	O2 Si	96-901-2604	0.6525
Quartz	O2 Si	96-900-5033	0.6521

Candidates



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Quartz	O2 Si	96-900-5034	0.6521
Quartz	O2 Si	96-900-5027	0.6321
Quartz	O2 Si	96-900-5026	0.6316
Quartz	O2 Si	96-900-5029	0.6315
Quartz	O2 Si	96-900-5028	0.6310
Quartz	O2 Si	96-900-5030	0.6309
Quartz	O2 Si	96-900-5031	0.6309
Quartz	O2 Si	96-900-5032	0.6299
Quartz	O2 Si	96-901-2603	0.6154

Search-Match

Settings

Reference database used	COD-Inorg REV89244 2013.10.11
Automatic zeropoint adaptation	Yes
Minimum figure-of-merit (FoM)	0.60
Parameter/influence 2theta	0.50
Parameter/influence intensities	0.50
Parameter multiple/single phase(s)	0.50

Selection Criteria

Compound:

Name: Quartz

Peak List

No.	2theta [°]	d [Å]	I/I0	FWHM	Matched
1	6.28	14.0648	159.14	0.2800	H
2	6.42	13.7642	178.65	0.2800	
3	6.79	13.0067	195.00	0.2800	
4	6.91	12.7820	185.33	0.2800	A
5	7.17	12.3197	169.36	0.2800	
6	7.33	12.0505	152.91	0.2800	
7	7.55	11.6951	150.06	0.2800	
8	7.75	11.3975	129.31	0.2800	
9	8.34	10.5902	138.70	0.2800	
10	8.84	9.9945	129.96	0.2800	D,G
11	10.58	8.3578	122.84	0.2800	AE
12	10.89	8.1144	121.34	0.2800	A
13	11.39	7.7628	141.37	0.2800	
14	11.65	7.5929	191.95	0.2800	
15	12.12	7.2962	231.23	0.2800	A
16	12.27	7.2077	211.19	0.2800	H
17	12.45	7.1039	160.25	0.2800	
18	12.63	7.0031	121.17	0.2800	A
19	18.32	4.8383	348.26	0.2800	AD,E,G
20	18.49	4.7947	272.66	0.2800	AE,F,H
21	20.08	4.4195	508.59	0.2800	AD,E,F,G,H
22	20.22	4.3887	525.35	0.2800	AD
23	20.35	4.3605	509.10	0.2800	
24	20.47	4.3352	491.15	0.2800	AB,H
25	20.65	4.2970	440.98	0.2800	AD
26	20.96	4.2351	324.89	0.2800	AD,H
27	21.30	4.1671	345.45	0.2800	AD,E,G,H
28	22.02	4.0337	1000.00	0.2800	AC,E,G,H
29	22.60	3.9318	208.94	0.2800	AD,E,G
30	22.83	3.8919	227.93	0.2800	AB,E,H
31	23.09	3.8487	213.28	0.2800	AD
32	23.35	3.8072	239.59	0.2800	AH
33	23.75	3.7434	398.52	0.2800	AB,D
34	23.99	3.7059	241.16	0.2800	AE
35	24.33	3.6558	293.42	0.2800	AF
36	24.69	3.6036	300.33	0.2800	AH
37	24.79	3.5887	304.99	0.2800	AD,H
38	25.05	3.5515	289.06	0.2800	A
39	25.55	3.4835	230.15	0.2800	AB,D,H
40	25.81	3.4494	268.07	0.2800	AG,H
41	26.34	3.3809	171.45	0.2800	AE
42	26.74	3.3315	257.99	0.2800	D,F,H
43	26.87	3.3154	244.87	0.2800	
44	27.06	3.2930	221.69	0.2800	AG
45			1.10	0.2800	AE,G,H
46			3.05	0.2800	AC,D,F
47			3.08	0.2800	AH
48			3.82	0.2800	AD,E,G
49			3.86	0.2800	AB,D,E,F,H
50			3.43	0.2800	A
51			1.36	0.2800	AE,F
52			1.55	0.2800	AD,F,H
53			1.61	0.2800	AD,E,G,H
54			3.85	0.2800	AB,D,E,F,G,H
55			3.11	0.2800	AB,E,F,G
56			3.58	0.2800	AD,E,H
57	36.35	2.4695	280.81	0.2800	AE
58	36.53	2.4577	286.31	0.2800	AB,D,E,G
59	36.86	2.4368	194.45	0.2800	AD,E,G,H



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60	37.34	2.4060	225.20	0.2800	AE,G
61	37.50	2.3965	210.78	0.2800	AE,H
62	37.73	2.3821	232.32	0.2800	AB,D,G,H
63	37.95	2.3690	179.43	0.2800	AE,F,H
64	38.16	2.3564	198.81	0.2800	E
65	38.27	2.3502	199.10	0.2800	AB,C
66	38.57	2.3324	150.60	0.2800	AE,F,G,H
67	38.93	2.3116	143.48	0.2800	AB,D,E,H
68	39.21	2.2960	163.92	0.2800	AE,F,H
69	39.39	2.2857	144.65	0.2800	AE
70	39.61	2.2735	118.18	0.2800	AB,H
71	40.25	2.2388	146.74	0.2800	AB,D,E,H
72	40.40	2.2307	145.48	0.2800	AD,E,F,G,H
73	40.80	2.2100	139.52	0.2800	AC,E,F,G,H
74	40.93	2.2032	135.86	0.2800	AD,F,H
75	41.09	2.1951	121.21	0.2800	A
76	41.22	2.1882	122.30	0.2800	AD,E,G
77	41.77	2.1609	196.95	0.2800	AB,C,D,E,F,H
78	44.38	2.0396	117.62	0.2800	AB,C,D,E,F,G,H
79	49.47	1.8409	164.99	0.2800	AB,D,E,F,G,H
80	49.81	1.8292	161.57	0.2800	AB,D,E,F,G,H
81	50.66	1.8004	149.38	0.2800	AB,E,F,H
82	51.01	1.7891	171.44	0.2800	AB,D,H
83	51.15	1.7844	155.45	0.2800	AD,F,G,H
84	52.33	1.7470	142.23	0.2800	B,C,E,G,H
85	52.46	1.7428	141.50	0.2800	B,D,H
86	53.13	1.7226	130.73	0.2800	D,E,F,G,H
87	53.65	1.7070	124.15	0.2800	B,D,E,H
88	54.11	1.6936	252.36	0.2800	D,E
89	54.22	1.6903	259.55	0.2800	D,E,G,H
90	54.49	1.6825	184.64	0.2800	D,E,G
91	54.93	1.6702	149.96	0.2800	B,D,E,F,H
92	55.10	1.6655	143.69	0.2800	F
93	55.23	1.6619	141.52	0.2800	D,H
94	55.41	1.6568	128.74	0.2800	D,G,H
95	55.92	1.6429	130.19	0.2800	D,E,G,H
96	56.21	1.6352	137.35	0.2800	B,D,E,F
97	56.37	1.6309	120.66	0.2800	D,E,G,H
98	57.11	1.6116	140.78	0.2800	B,C,D,E,F,G,H
99	57.30	1.6066	136.31	0.2800	B,C,D,E,F,H
100	62.10	1.4934	186.29	0.2800	B,D,E,F,G,H
101	62.44	1.4861	296.34	0.2800	D,E,F,G
102	62.68	1.4809	195.17	0.2800	B,D,H
103	62.85	1.4774	181.56	0.2800	C,E,G
104	63.00	1.4744	176.03	0.2800	D,E,F,G,H
105	63.26	1.4688	139.05	0.2800	B,D,E,F
106	63.41	1.4657	125.91	0.2800	B,D,E,G
107	63.96	1.4544	161.70	0.2800	D,E,F,G,H

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

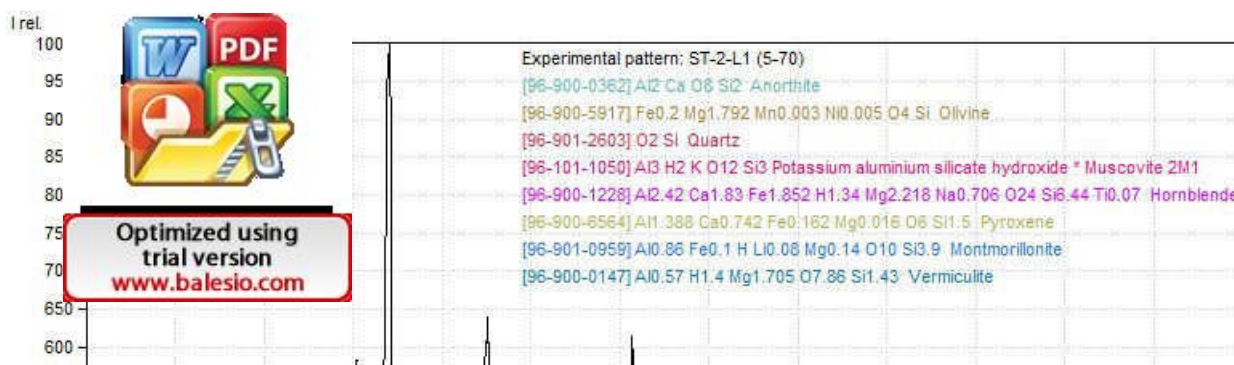
Based on calculated profile

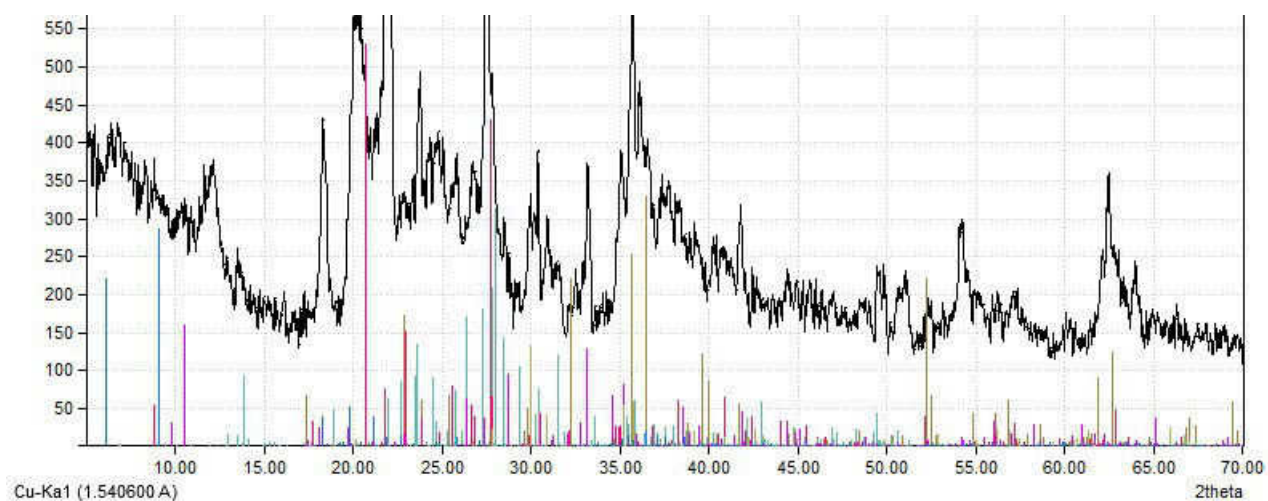
Profile area	Counts	Amount
Overall diffraction profile	123329	100.00%
Background radiation	75775	61.44%
Diffraction peaks	47554	38.56%
Peak area belonging to selected phases	25302	20.52%
Unidentified peak area	22252	18.04%

Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	1442	100.00%
Peak intensity belonging to selected phases	542	37.61%
Unidentified peak intensity	899	62.39%

Diffraction Pattern Graphics





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Match! Phase Analysis Report

Sample: ST-3-L1 (5-70)

Sample Data

File name ST-3-L1.RAW
File path D:/S2 Geologi/data hasil/17 sampel/ST-3-L1
Data collected Jan 5, 2024 16:52:43
Data range 5.010° - 70.010°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction 0.01°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	34.1	Muscovite	Al2.73 Ba0.01 H2.86 K0.465 N0.36 Na0.03 O11.236 Si3.128
B	30.4	Anorthite	Al2 Ca O8 Si2
C	22.3	Olivine	Co0.249 Mg1.749 O4 Si
D	8.3	Quartz	O2 Si
E	4.4	Montmorillonite	Al0.86 Fe0.1 H Li0.08 Mg0.14 O10 Si3.9
F	0.5	Vermiculite	H2 Mg3 O12 Si4
	11.1	Unidentified peak area	

A: Muscovite (34.1 %)

Formula sum Al2.73 Ba0.01 H2.86 K0.465 N0.36 Na0.03 O11.236 Si3.128
Entry number 96-901-5971
Figure-of-Merit (FoM) 0.759576
Total number of peaks 300
Peaks in range 300
Peaks matched 125
Intensity scale factor 0.46
Space group C 1 2/c 1
Crystal system monoclinic
Unit cell a= 9.0270 Å b= 5.1999 Å c= 20.6160 Å β= 100.113 °
I/Icor 0.77
Calc. density 2.576 g/cm³
Reference Mesto E., Scordari F., Lacalamita M., Schingaro E., "Tobelite and NH4+ -rich muscovite single crystals from Ordovician American sandstones (Brittany, France): Structure and crystal chemistry Note: sample musc_4", American Mineralogist **97**, 1460-1468 (2012)

B: Anorthite (30.4 %)

Formula sum Al2 Ca O8 Si2
Entry number 96-900-0362
Figure-of-Merit (FoM) 0.762263
Total number of peaks 500
Peaks in range 500
Peaks matched 328
Intensity scale factor 0.32
Space group P -1
Crystal system triclinic (anorthic)
Unit cell a= 8.1940 Å b= 12.8970 Å c= 14.1900 Å α= 92.980° β= 115.820° γ= 91.150 °
I/Icor 0.60
Calc. density 2.745 g/cm³
Reference Foit F. F., Peacor D. R., "The anorthite crystal structure at 410 and 830 C T = 410 C", American Mineralogist **58**, 665-675 (1973)

C: Olivine (22.3 %)

Formula sum Co0.249 Mg1.749 O4 Si
Entry number 96-900-6398
Figure-of-Merit (FoM) 0.705303
Total number of peaks 168
Peaks in range 168
Peaks matched 39
Intensity scale factor 0.40
Space group P b n m
Crystal system orthorhombic
Unit cell a= 4.7585 Å b= 10.2148 Å c= 5.9846 Å
I/Icor 0.2
Calc. density 3.607 g/cm³
Reference Müller-Sommer M., Hock R., Kirfel A., "Rietveld refinement study of the cation distribution in (Co, Mg)-olivine solid solution sample: M87", Physics and Chemistry of Minerals **24**, 17-23 (1997)

D: Quartz (8.3 %)

Formula sum Si
Entry number 96-900-5020
Figure-of-Merit (FoM) 0.7379
Total number of peaks 33
Peaks in range 33
Peaks matched 11
Intensity scale factor 0.45
Space group P 63/mc 2



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Space group	$P 32 2 1$
Crystal system	trigonal (hexagonal axes)
Unit cell	$a=4.9297 \text{ \AA}$ $c=5.4151 \text{ \AA}$
V/cor	3.07
Calc. density	2.625 g/cm^3
Reference	Kihara K., "An X-ray study of the temperature dependence of the quartz structure Sample: at $T = 498 \text{ K}$ ", European Journal of Mineralogy 2 , 63-77 (1990)

E: Montmorillonite (4.4 %)

Formula sum	$\text{Al}_0.86 \text{Fe}_{0.1} \text{H} \text{Li}_{0.08} \text{Mg}_{0.14} \text{O}_{10} \text{Si}_{3.9}$
Entry number	96-901-0959
Figure-of-Merit (FoM)	0.730056
Total number of peaks	251
Peaks in range	251
Peaks matched	60
Intensity scale factor	0.45
Space group	$C 1 2/m 1$
Crystal system	monoclinic
Unit cell	$a=5.1710 \text{ \AA}$ $b=8.9570 \text{ \AA}$ $c=9.7400 \text{ \AA}$ $\beta=96.100^\circ$
V/cor	5.89
Calc. density	2.245 g/cm^3
Reference	Gournis D., Lappas A., Karakassides M. A., Tobbens D., Moukarika A., "A neutron diffraction study of alkali cation migration in montmorillonites Sample: Li-mont-300", Physics and Chemistry of Minerals 35 , 49-58 (2008)

F: Vermiculite (0.5 %)

Formula sum	$\text{H}_2 \text{Mg}_3 \text{O}_{12} \text{Si}_4$
Entry number	96-900-0017
Figure-of-Merit (FoM)	0.732434
Total number of peaks	263
Peaks in range	262
Peaks matched	120
Intensity scale factor	0.20
Space group	$C 1 c 1$
Crystal system	monoclinic
Unit cell	$a=5.3300 \text{ \AA}$ $b=9.1800 \text{ \AA}$ $c=28.8500 \text{ \AA}$ $\beta=93.250^\circ$
V/cor	23.53
Calc. density	1.786 g/cm^3
Reference	Hendricks S. B., Jefferson M. E., "Crystal structure of vermiculites and mixed vermiculite-chlorites", American Mineralogist 23 , 851-862 (1938)

Candidates

Name	Formula	Entry No.	FoM
	$\text{Br}_2 \text{F}_{17} \text{Sb}_2$	96-431-1869	0.7895
	$\text{C F}_{20} \text{O}_4 \text{Te}_4$	96-430-8273	0.7872
Antigorite-T	$\text{Mg}_{48} \text{O}_{147} \text{Si}_{34}$	96-901-6234	0.7865
potassium hydroxopentafluoroarsenate	$\text{As F}_5 \text{H K O}$	96-200-3122	0.7863
Tridymite	$\text{O}_2 \text{Si}$	96-900-0521	0.7852
	$\text{Cs F}_{30} \text{O}_6 \text{Sb Te}_6$	96-432-7610	0.7837
Tridymite	$\text{O}_2 \text{Si}$	96-901-3492	0.7828
Kaolinite	$\text{Al}_2 \text{H}_4 \text{O}_9 \text{Si}_2$	96-900-9235	0.7824
Antigorite	$\text{H}_{79} \text{Mg}_{48} \text{O}_{147} \text{Si}_{34}$	96-900-3104	0.7799
Antigorite	$\text{H}_{62} \text{Mg}_{48} \text{O}_{147} \text{Si}_{34}$	96-900-4515	0.7791
Tridymite	$\text{O}_2 \text{Si}$	96-901-3493	0.7788
Kaolinite	$\text{Al}_2 \text{H}_4 \text{O}_9 \text{Si}_2$	96-900-9231	0.7764
Metaschoepite	$\text{H}_{34} \text{Na}_{0.47} \text{O}_{37.082} \text{U}_8$	96-901-0199	0.7748
Metaschoepite	$\text{H}_{34} \text{Na}_{0.48} \text{O}_{37.91} \text{U}_8$	96-901-0195	0.7745
meta-schoepite	$\text{H}_{16} \text{O}_{20} \text{U}_4$	96-210-2097	0.7742
Metaschoepite	$\text{H}_{34} \text{O}_{40} \text{U}_8$	96-901-1299	0.7742
Hexaammine cobalt(III) hydrogenarsenate tetrahydrate	$\text{As}_3 \text{Co}_2 \text{H}_{47} \text{N}_{12} \text{O}_{16}$	96-150-0006	0.7719
Metaschoepite	$\text{H}_{34} \text{Na}_{1.16} \text{O}_{37.9} \text{U}_8$	96-901-0198	0.7700
	$\text{Cl H}_{64} \text{K}_5 \text{Mo}_{12} \text{O}_{57} \text{P}_2 \text{S}_{1296}$	96-411-2785	0.7696
Metaschoepite	$\text{H}_{32} \text{Na}_{1.22} \text{O}_{39.09} \text{U}_8$	96-901-0197	0.7693
Potassium aluminium silicate hydroxide * (Muscovite 2M1)	$\text{Al}_3 \text{H}_2 \text{K O}_{12} \text{Si}_3$	96-101-1050	0.7683
Potassium aluminium silicate hydroxide * (Muscovite 2M1)	$\text{Al}_3 \text{H K O}_{12} \text{Si}_3$	96-110-0014	0.7683
Tetaborane	$\text{B}_4 \text{H}_{10}$	96-430-4422	0.7680
Metaschoepite	$\text{H}_{32} \text{Na}_{1.09} \text{O}_{38.328} \text{U}_8$	96-901-0196	0.7671
	$\text{C}_2 \text{F}_{22} \text{N}_2 \text{S}_6 \text{Sb}_4$	96-431-6375	0.7668
	$\text{C}_6 \text{Al}_3 \text{O}_{15} \text{P}_2$	96-700-6431	0.7668
Dimolybdenum(V) dioxide tetraphosphate	$\text{Mo}_2 \text{O}_{15} \text{P}_4$	96-100-1468	0.7665
	$\text{Mo O}_{7.5} \text{P}_2$	96-901-4374	0.7665
Tamarugite	$\text{Al H}_{12} \text{Na O}_{14} \text{S}_2$	96-900-0182	0.7662
	$\text{As}_2 \text{F}_{20} \text{Mg Xe}_4$	96-431-0410	0.7661
Khademite	$\text{Al F H}_{10} \text{O}_9 \text{S}$	96-900-9710	0.7655
	$\text{Cl F}_{17} \text{Sb}_2$	96-431-1868	0.7653
	$\text{C O}_8 \text{Fe}_2 \text{N}_6 \text{O}_8 \text{P}_2$	96-411-4284	0.7646
	$\text{Mo}_3 \text{O}_{22.5} \text{P}_6$	96-430-3254	0.7642
Antigorite	$\text{H}_{58} \text{Mg}_{45} \text{O}_{138} \text{Si}_{32}$	96-900-4000	0.7642
XeF ₂ F ₄	$\text{F}_{14} \text{O}_3 \text{Os Xe}_2$	96-431-5229	0.7639
lanthir	$\text{H}_{22.62} \text{O}_{27} \text{U}_6$	96-901-2907	0.7635
Alumir	$\text{Al}_2 \text{H}_4 \text{O}_9 \text{Si}_2$	96-101-1046	0.7634
Dickite	$\text{Al}_2 \text{H}_4 \text{O}_9 \text{Si}_2$	96-900-0123	0.7634
	$\text{O}_{45} \text{P}_9 \text{V}_{12}$	96-400-1392	0.7631
Zippeite	$\text{H}_{27} \text{Na}_5 \text{O}_{52} \text{S}_4 \text{U}_8$	96-900-4756	0.7628
Triam	$\text{H}_{12} \text{In N}_3 \text{O}_{12} \text{S}_3$	96-100-4054	0.7622
Bementite	$\text{Mn}_{6.683} \text{O}_{23} \text{Si}_6$	96-900-1585	0.7608
Tridymite	$\text{O}_2 \text{Si}$	96-900-6969	0.7595
	$\text{Mo}_2 \text{O}_{15} \text{P}_4$	96-430-3255	0.7592
	$\text{As}_2 \text{F}_{18} \text{Os}$	96-430-4075	0.7586



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Dickite	As ₂ F ₁₂ S ₈	96-432-1675	0.7586
Kaolinite	Al ₂ H ₄ O ₉ Si ₂	96-900-3082	0.7576
	Al ₂ O ₉ Si ₂	96-901-5000	0.7574
	C ₂ F ₁₁ N ₃ S ₃ Sb ₂	96-431-6377	0.7572
(Xe ₂ F ₁₁)(AuF ₆)	Mb ₃ O _{22.5} P ₆	96-430-3256	0.7570
Dialuminium phyllo-disilicate tetrahydroxide (Nacrite 6M)	AuF ₁₇ Xe ₂	96-433-1415	0.7569
and 150 others...	Al ₂ H ₄ O ₉ Si ₂	96-101-1063	0.7568

Search-Match

Settings

Reference database used	COD-Inorg REV89244 2013.10.11
Automatic zeropoint adaptation	Yes
Minimum figure-of-merit (FoM)	0.60
Parameter/influence 2theta	0.50
Parameter/influence intensities	0.50
Parameter multiple/single phase(s)	0.50

Peak List

No.	2theta [°]	d[Å]	I/I0	FWHM	Matched
1	6.85	12.8969	167.10	0.4800	B,F
2	8.76	10.0888	206.15	0.4800	A
3	9.00	9.8146	235.21	0.4800	B,E
4	11.91	7.4275	208.76	0.4800	B
5	12.15	7.2802	271.82	0.4800	
6	12.35	7.1612	241.51	0.4800	F
7	12.35	7.1612	241.51	0.4800	
8	12.52	7.0620	183.94	0.4800	B
9	18.36	4.8277	1000.00	0.4800	AB,E,F
10	20.29	4.3731	885.07	0.4800	AB,E,F
11	20.47	4.3352	833.58	0.4800	B,F
12	20.47	4.3352	833.58	0.4800	
13	20.71	4.2855	578.55	0.4800	AB,D
14	20.71	4.2855	578.55	0.4800	
15	20.88	4.2502	501.82	0.4800	AB,F
16	21.13	4.2012	445.41	0.4800	E
17	21.13	4.2012	445.41	0.4800	
18	21.47	4.1355	408.73	0.4800	B,F
19	22.06	4.0259	698.50	0.4800	B,E,F
20	22.45	3.9571	229.29	0.4800	AB,F
21	22.45	3.9571	229.29	0.4800	
22	22.62	3.9270	187.77	0.4800	B,E
23	22.81	3.8961	219.29	0.4800	AB,C,F
24	23.22	3.8275	213.14	0.4800	AB
25	23.46	3.7887	194.71	0.4800	B,F
26	23.69	3.7521	226.11	0.4800	B
27	23.88	3.7227	306.48	0.4800	B,C
28	24.27	3.6639	385.11	0.4800	AB,F
29	24.62	3.6128	373.00	0.4800	AB,F
30	24.86	3.5790	363.88	0.4800	B,F
31	25.13	3.5408	249.77	0.4800	AB
32	25.13	3.5408	249.77	0.4800	
33	25.47	3.4950	264.68	0.4800	B,C,F
34	25.91	3.4363	187.97	0.4800	AB,E,F
35	26.77	3.3277	637.38	0.4800	B,D,F
36	26.99	3.3009	494.33	0.4800	E
37	26.99	3.3009	494.33	0.4800	
38	27.29	3.2649	184.33	0.4800	B
39	27.57	3.2324	224.18	0.4800	AB,E,F
40	27.77	3.2103	212.42	0.4800	B
41	27.96	3.1884	169.81	0.4800	AB
42	28.74	3.1043	179.97	0.4800	AB,C,E,F
43	33.22	2.6948	425.75	0.4800	B,C,E,F
44	35.09	2.5551	185.60	0.4800	AB,C,E,F
45	35.33	2.5388	240.89	0.4800	AB,E,F
46	35.70	2.5127	547.36	0.4800	AB,C,E,F
47	36.11	2.4855	411.91	0.4800	AB,F
48	36.67	2.4487	477.46	0.4800	AB,C,D,E,F
49	37.09	2.4222	348.45	0.4800	AB,E
50	37.22	2.4140	361.90	0.4800	AB
51	37.67	2.3861	458.33	0.4800	AB,C,F
52	37.93	2.3702	322.03	0.4800	B,E
53	37.93	2.3702	322.03	0.4800	
54	41.73	2.1016	226.07	0.4800	AB,C,D,E,F
55	44.36	2.0404	165.44	0.4800	AB,C,E,F
56	45.49	1.9923	154.46	0.4800	B,F
57	45.64	1.9875	160.00	0.4800	AB,C,D,E,F



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67	45.61	1.9875	160.90	0.4800	AB,C,D,E,F
68	49.52	1.8391	162.24	0.4800	AB,C,D,E,F
69	54.21	1.6906	349.59	0.4800	A,C,E,F
70	54.49	1.6827	222.12	0.4800	A,E,F
71	54.87	1.6718	168.12	0.4800	A,C,D,F
72	55.31	1.6596	209.02	0.4800	AC,D,E,F
73	62.45	1.4859	319.47	0.4800	A,C,E,F
74	62.69	1.4808	245.56	0.4800	A,C,E,F
75	62.69	1.4808	245.56	0.4800	A,C,E,F
76	63.96	1.4544	284.41	0.4800	A,C,D,E,F

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

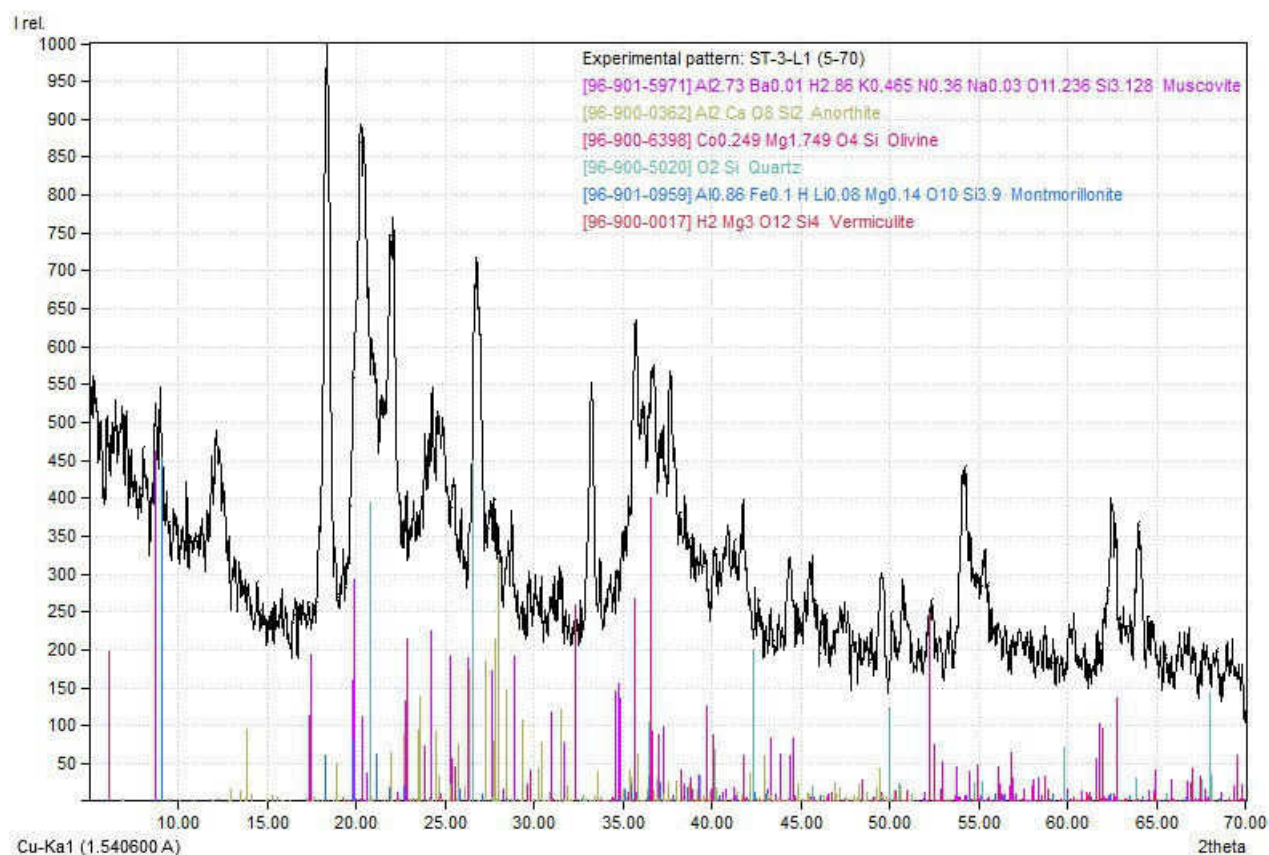
Based on calculated profile

Profile area	Counts	Amount
Overall diffraction profile	116218	100.00%
Background radiation	79054	68.02%
Diffraction peaks	37164	31.98%
Peak area belonging to selected phases	24262	20.88%
Unidentified peak area	12902	11.10%

Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	1700	100.00%
Peak intensity belonging to selected phases	749	44.07%
Unidentified peak intensity	951	55.93%

Diffraction Pattern Graphics



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Match! Phase Analysis Report

Sample: ST-4-L1 (5-70)

Sample Data

File name ST-4-L1.RAW
File path D:/S2 Geologi/data hasil/17 sampel/ST-4-L1
Data collected Jan 5, 2024 16:52:43
Data range 5.000° - 70.000°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	65.4	Muscovite	Al2.726 Ca0.011 Fe0.03 K0.776 Mg0.02 Na0.181 O11 Si3.15 Ti0.02
B	16.3	Olivine	Mg0.5 Ni1.5 O4 Si
C	10.5	Quartz	O2 Si
D	7.8	Anorthite	Al2 Ca O8 Si2
	8.6	Unidentified peak area	

A: Muscovite (65.4 %)

Formula sum Al2.726 Ca0.011 Fe0.03 K0.776 Mg0.02 Na0.181 O11 Si3.15 Ti0.02
Entry number 96-901-2887
Figure-of-Merit (FoM) 0.762865
Total number of peaks 144
Peaks in range 142
Peaks matched 110
Intensity scale factor 0.50
Space group C 1
Crystal system triclinic (anorthic)
Unit cell a= 5.2234 Å b= 9.1790 Å c= 20.2353 Å α= 90.000° β= 95.834° γ= 90.000°
I/Icor 0.77
Calc. density 2.588 g/cm³
Reference Tomita K., Shiraki K., Kawano M., "Crystal structure of dehydroxylated 2M1 sericite and its relationship with mixed-layer mica/smectite Locality: Goto mine, Nagasaki Prefecture, Japan Sample: heated", Clay Science **10**, 423-441 (1998)

B: Olivine (16.3 %)

Formula sum Mg0.5 Ni1.5 O4 Si
Entry number 96-900-1102
Figure-of-Merit (FoM) 0.735460
Total number of peaks 164
Peaks in range 164
Peaks matched 42
Intensity scale factor 0.30
Space group P b n m
Crystal system orthorhombic
Unit cell a= 4.7331 Å b= 10.1565 Å c= 5.9285 Å
I/Icor 1.83
Calc. density 4.480 g/cm³
Reference Bostrom D., "Single-crystal X-ray diffraction studies of synthetic Ni-Mg olivine solid solutions Sample: XNi2+ = .75", American Mineralogist **72**, 965-972 (1987)

C: Quartz (10.5 %)

Formula sum O2 Si
Entry number 96-900-7379
Figure-of-Merit (FoM) 0.677413
Total number of peaks 35
Peaks in range 35
Peaks matched 11
Intensity scale factor 0.31
Space group P 32 2 1
Crystal system trigonal (hexagonal axes)
Unit cell a= 4.9800 Å c= 5.4600 Å
I/Icor 2.99
Calc. density 2.552 g/cm³
Reference sa A. L., El-Barbary AA Heggie M. I., Briddon P. R., "Structural and thermodynamic properties of water related defects in alpha-quartz Note: Hypothetical structure derived using density-functional theory", Physics and Chemistry of Minerals **32**, 323-331 (2005)

D: An

Formu ! Ca O8 Si2
Entry r -900-0362
Figure '32741
Total n 0
Peaks 5
Peaks 7
Intensity scale factor 0.05
Space group P -1
Crystal system triclinic (anorthic)
Unit cell a= 5.4040 Å b= 4.0070 Å c= 11.4000 Å α= 90.000° β= 115.000° γ= 90.450°



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Unit cell a= 8.1940 Å b= 12.8970 Å c= 14.1900 Å α= 92.980° β= 115.820° γ= 91.150°
 I/cor 0.60
 Calc. density 2.745 g/cm³
 Reference Foit F. F., Peacor D. R., "The anorthite crystal structure at 410 and 830 C T = 410 C", American Mineralogist **58**, 665-675 (1973)

Candidates

Name	Formula	Entry No.	FoM
Antigorite	H62 Mg48 O147 Si34	96-900-4515	0.8044
Antigorite	H58 Mg45 O138 Si32	96-900-4000	0.8018
	O5 S Ti	96-210-0943	0.7994
	Cs F30 O6 Sb Te6	96-432-7610	0.7977
Antigorite	H79 Mg48 O147 Si34	96-900-3104	0.7970
Becquerelite	Ca O30 U6	96-900-1111	0.7966
aluminum phosphate	Al O4 P	96-201-0797	0.7949
Antigorite-T	Mg48 O147 Si34	96-901-6234	0.7942
	Br2 F17 Sb2	96-431-1869	0.7930
Becquerelite	Ca H22 O30 U6	96-901-2088	0.7925
Becquerelite	Ca H22 O30 U6	96-900-2701	0.7918
Billietite	Ba O23 U6	96-900-1112	0.7896
Olympite	Li Na5 O8 P2	96-901-2623	0.7891
Khademite	Al F H10 O9 S	96-900-9710	0.7879
	As2 F20 Mg Xe4	96-431-0410	0.7843
Ianthinite	H22.62 O27 U6	96-901-2907	0.7840
tetrafluorobromate undecafluorodiantimonate	Br F15 Sb2	96-431-8956	0.7823
	C2 F22 N2 S6 Sb4	96-431-6375	0.7814
K Au5 P2 S8	Au5 K P2 S8	96-402-7827	0.7806
Billietite	Ba H6 O30 U6	96-901-0600	0.7798
Dialuminium phyllo-disilicate tetrahydroxide * (Nacrite 4M)	Al2 H4 O9 Si2	96-101-1081	0.7796
Kaolinite	Al2 H4 O9 Si2	96-900-9235	0.7786
	F14 Sb2 Xe	96-431-6547	0.7784
Triammonium indium sulfate	H12 In N3 O12 S3	96-100-4054	0.7780
Thallium phosphate hydrate *	H23.5 O36.5 P8 Ti3.5	96-100-8946	0.7770
	Al1.75 Na1.75 O4 Si0.25	96-200-2894	0.7769
Nacrite	Al2 O9 Si2	96-901-4064	0.7757
Schoepite	H18 O21 U4	96-900-4445	0.7752
	Cs2 Ga6 H2 O26 P6	96-400-1539	0.7751
Magnesium thiosulfate hexahydrate	H12 Mg O9 S2	96-110-0073	0.7741
	Mg O9 S2	96-901-5179	0.7741
Brianroulstonite	B5 Ca3 Cl2 H23 O21	96-900-4452	0.7734
Tridymite	O2 Si	96-901-3492	0.7729
potassium hydroxopentafluoroarsenate	As F5 H K O	96-200-3122	0.7728
(Xe2 F11) (Au F6)	Au F17 Xe2	96-433-1415	0.7724
Metaschoepite	H34 Na1.16 O37.9 U8	96-901-0198	0.7719
	C0 F8 Fe2 N6 O8 P2	96-411-4284	0.7717
	C12 Co N8	96-433-0306	0.7715
	Al Cs2 O10 P3	96-700-9288	0.7714
	Mo N2 O9 P2	96-901-4933	0.7714
Vandendriesscheite	H33 O48 Pb2 U10	96-900-1892	0.7709
	As2 H16 Mn O20 U2	96-900-4896	0.7708
	Cs4 O18 P3 Te	96-901-4415	0.7705
	C14 Cl2 F8	96-402-4047	0.7699
Kaolinite	Al2 H4 O9 Si2	96-900-9231	0.7698
	C18 B3 F15 N18	96-701-0277	0.7697
	C F20 O4 Te4	96-430-8273	0.7695
1,10-octachlorophenanthroline	C12 Cl8 N2	96-432-3462	0.7695
	O36.5 P8 Ti3.5	96-901-5355	0.7695
	B3 H3 N12	96-432-2188	0.7692
	Cs2 Ga O10 P3	96-700-9289	0.7692
	C16 Br2 N8 O2 Pd	96-900-7878	0.7692

and 151 others...

Search-Match

Settings

Reference database used COD-Inorg REV89244 2013.10.11
 Automatic zeropoint adaptation Yes
 Minimum figure-of-merit (FoM) 0.60
 Parameter/influence 2theta 0.50
 Parameter/influence intensities 0.50
 Parameter multiple/single phase(s) 0.50

Peak List

No.	I/I0	FWHM	Matched
1	1.17	0.8800	D
2	0.95	0.8800	
3	0.24	0.8800	
4	0.12	0.8800	D
5	0.91	0.8800	
6	0.64	0.8800	AD
7	0.00	0.8800	AD
8	0.48	0.8800	AD
9	0.74	0.8800	AB,C,D
10	0.33	0.8800	D
11	21.24 4.1800 651.54	0.8800	A
12	21.50 4.1298 542.64	0.8800	D
13	21.77 4.0795 545.44	0.8800	
14	21.88 4.0627 536.28	0.8800	AD



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14	21.99	4.0387	536.39	0.8800	A,D
15	22.28	3.9864	354.03	0.8800	A,D
16	22.62	3.9274	286.95	0.8800	A,D
17	22.84	3.8901	233.65	0.8800	D
18	23.11	3.8453	272.91	0.8800	B,D
19	23.33	3.8098	280.37	0.8800	A,D
20	23.71	3.7500	264.83	0.8800	D
21	24.09	3.6918	389.41	0.8800	B,D
22	24.22	3.6723	390.66	0.8800	D
23	24.87	3.5773	496.23	0.8800	A,D
24	25.18	3.5339	425.71	0.8800	A,D
25	25.44	3.4988	384.89	0.8800	D
26	25.61	3.4750	341.27	0.8800	B,D
27	25.95	3.4304	290.41	0.8800	D
28	26.27	3.3896	239.25	0.8800	A,C
29	26.46	3.3658	204.93	0.8800	A,D
30	26.66	3.3406	194.73	0.8800	
31	27.04	3.2952	189.39	0.8800	
32	27.37	3.2562	191.36	0.8800	A,D
33	30.23	2.9540	148.07	0.8800	A,B,D
34	33.18	2.6977	223.20	0.8800	A,B,D
35	35.04	2.5585	434.90	0.8800	A,B,D
36	35.67	2.5151	771.73	0.8800	A,D
37	35.98	2.4941	453.24	0.8800	A,B,C,D
38	36.14	2.4834	386.32	0.8800	D
39	36.32	2.4715	334.37	0.8800	D
40	36.60	2.4530	308.63	0.8800	A,D
41	36.76	2.4429	284.16	0.8800	A,B,D
42	36.96	2.4300	281.22	0.8800	D
43	37.16	2.4175	301.35	0.8800	D
44	37.33	2.4070	320.59	0.8800	A,D
45	37.48	2.3976	298.15	0.8800	D
46	37.64	2.3878	280.08	0.8800	A,D
47	37.87	2.3736	302.91	0.8800	B,D
48	38.28	2.3496	364.89	0.8800	A,D
49	38.62	2.3292	280.00	0.8800	A,B,D
50	38.80	2.3191	288.11	0.8800	D
51	39.08	2.3029	245.44	0.8800	A,B,C,D
52	39.58	2.2752	176.93	0.8800	A,C,D
53	40.36	2.2330	154.32	0.8800	A,B,D
54	40.72	2.2140	163.67	0.8800	A,D
55	40.94	2.2026	149.44	0.8800	A,D
56	41.17	2.1908	163.73	0.8800	A,B,C,D
57	53.54	1.7103	152.90	0.8800	A,B
58	54.12	1.6932	261.71	0.8800	A,B,C
59	54.45	1.6837	207.19	0.8800	A
60	54.68	1.6772	235.36	0.8800	A,C
61	54.96	1.6693	182.73	0.8800	
62	55.08	1.6660	176.72	0.8800	A
63	55.30	1.6599	181.17	0.8800	A,B
64	55.54	1.6534	171.54	0.8800	A
65	57.20	1.6093	141.18	0.8800	A,B,C
66	62.05	1.4946	238.95	0.8800	B
67	62.43	1.4863	391.58	0.8800	
68	62.64	1.4819	342.23	0.8800	B
69	62.80	1.4785	237.49	0.8800	
70	63.01	1.4741	208.32	0.8800	B,C
71	63.72	1.4593	174.66	0.8800	B
72	63.88	1.4561	166.85	0.8800	
73	64.06	1.4524	150.84	0.8800	B,C

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

Based on calculated profile

Profile area	Counts	Amount
Overall diffraction profile	103549	100.00%
Background radiation	68273	65.93%
Diffraction peaks	35277	34.07%
Peak area belonging to selected phases	26383	25.48%
Unidentified	8893	8.59%

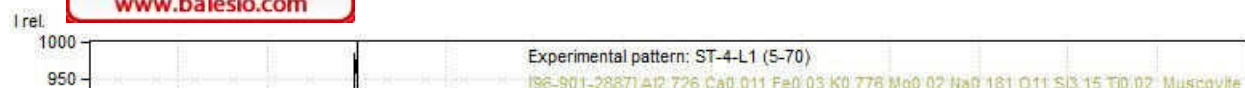
Peak Residuals

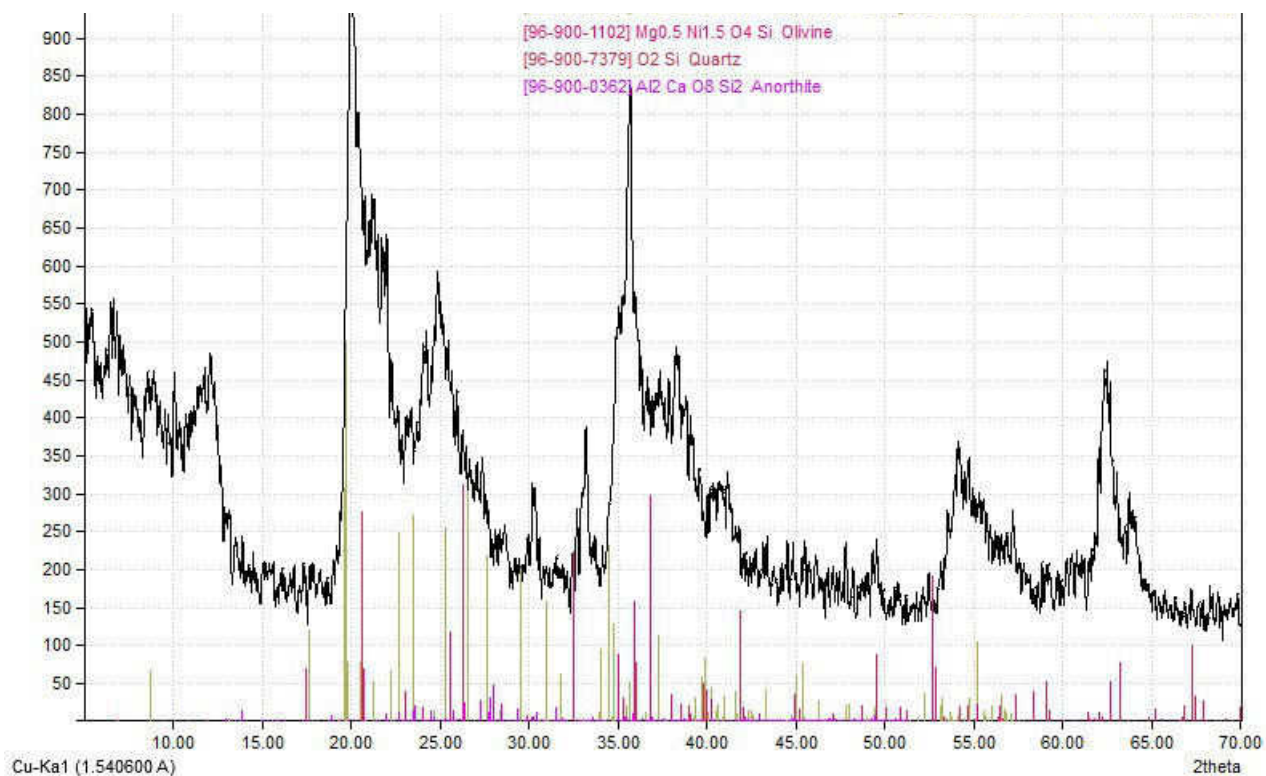
Peak	Counts	Amount
Overall	3038	100.00%
Peak identified	985	32.43%
Unidentified	2052	67.57%



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Diffraction Pattern Graphics





Cu-Ka1 (1.540600 Å)

2theta

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Match! Phase Analysis Report

Sample: ST-5-L1 (5-70)

Sample Data

File name ST-5-L1.RAW
File path D:/S2 Geologi/data hasil/17 sampel/ST-5-L1
Data collected Jan 5, 2024 16:52:43
Data range 5.130° - 70.130°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction 0.13°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	32.9	Muscovite	Al _{2.8} Ba _{0.01} Fe _{0.08} H ₂ K _{0.9} Mg _{0.04} Na _{0.07} O ₁₂ Si _{3.04} Ti _{0.04}
B	21.8	Anorthite	Al ₂ Ca O ₈ Si ₂
C	12.6	Olivine	Mg ₂ O ₄ Si
D	12.5	Albite	Al _{1.005} Na _{0.986} O ₈ Si _{2.995}
E	8.0	Pyroxene	Fe _{0.086} Mg _{0.914} O ₃ Si
F	6.1	Labradorite	Al _{0.824} Ca _{0.325} Na _{0.175} O ₄ Si _{1.174}
G	6.1	Quartz	O ₂ Si
	9.6	Unidentified peak area	

A: Muscovite (32.9 %)

Formula sum Al_{2.8} Ba_{0.01} Fe_{0.08} H₂ K_{0.9} Mg_{0.04} Na_{0.07} O₁₂ Si_{3.04} Ti_{0.04}
Entry number 96-900-6328
Figure-of-Merit (FoM) 0.704421
Total number of peaks 300
Peaks in range 300
Peaks matched 76
Intensity scale factor 0.32
Space group C 1 2/c 1
Crystal system monoclinic
Unit cell a= 5.1510 Å b= 8.9310 Å c= 19.3990 Å β= 95.800 °
I/I_{cor} 0.47
Calc. density 2.995 g/cm³
Reference Comodi P., Zanazzi P. F., "High-pressure structural study of muscovite Sample: K-Ms, P = 2.80 GPa", Physics and Chemistry of Minerals **22**, 170-177 (1995)

B: Anorthite (21.8 %)

Formula sum Al₂ Ca O₈ Si₂
Entry number 96-900-1259
Figure-of-Merit (FoM) 0.796398
Total number of peaks 500
Peaks in range 500
Peaks matched 201
Intensity scale factor 0.29
Space group P -1
Crystal system triclinic (anorthic)
Unit cell a= 8.1750 Å b= 12.8730 Å c= 14.1700 Å α= 93.110° β= 115.890° γ= 91.280°
I/I_{cor} 0.63
Calc. density 2.762 g/cm³
Reference Angel R. J., Carpenter M. A., Finger L. W., "Structural variation associated with compositional variation and order-disorder behavior in anorthite-rich feldspars sample from Val Pasmada", American Mineralogist **75**, 150-162 (1990)

C: Olivine (12.6 %)

Formula sum Mg₂ O₄ Si
Entry number 96-900-1097
Figure-of-Merit (FoM) 0.656816
Total number of peaks 167
Peaks in range 167
Peaks matched 17
Intensity scale factor 0.20
Space group Rh n m
Crystal system orthorhombic
Unit cell a= 4.7490 Å b= 10.1985 Å c= 5.9792 Å
I/I_{cor} 7
Calc. density 3.25 g/cm³
Reference Strom D., "Single-crystal X-ray diffraction studies of synthetic Ni-Mg olivine solid solutions Sample: XNi²⁺ = 0.00", American Mineralogist **72**, 965-972 (1987)

D: Albite

Formula Al_{1.005} Na_{0.986} O₈ Si_{2.995}
Entry number 96-900-0784
Figure-of-Merit (FoM) 47406
Total number of peaks 250
Peaks in range 250
Peaks matched 87
Intensity scale factor 0.20



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intensity scale factor	0.22
Space group	C -1
Crystal system	triclinic (anorthic)
Unit cell	a= 8.1420 Åb= 12.7850 Åc= 7.1590 Åα= 94.190° β= 116.610 ° γ= 87.680 °
I/Icor	0.84
Calc. density	2.616 g/cm³
Reference	Harlow G. E., Brown G. E., "Low albite: An X-Ray and neutron diffraction study Sample: X-rays single Na atom Note: this sample of feldspar is from Amelia, Virginia", American Mineralogist 65 , 986-995 (1980)

E: Pyroxene (8.0 %)	
Formula sum	Fe0.086 Mg0.914 O3 Si
Entry number	96-900-1220
Figure-of-Merit (FoM)	0.711879
Total number of peaks	430
Peaks in range	430
Peaks matched	63
Intensity scale factor	0.10
Space group	P b c a
Crystal system	orthorhombic
Unit cell	a= 18.2430 Åb= 8.8340 Åc= 5.1900 Å
I/Icor	0.60
Calc. density	3.274 g/cm³
Reference	Mblin G. M., "Crystal-chemical study of cation disordering in Al-rich and Al-poor orthopyroxenes from spinel lherzolite xenoliths sample Le9A, 1150 C, 10 min", American Mineralogist 74 , 593-598 (1989)

F: Labradorite (6.1 %)	
Formula sum	Al0.824 Ca0.325 Na0.175 O4 Si1.174
Entry number	96-900-0748
Figure-of-Merit (FoM)	0.744135
Total number of peaks	251
Peaks in range	251
Peaks matched	84
Intensity scale factor	0.10
Space group	C -1
Crystal system	triclinic (anorthic)
Unit cell	a= 8.1747 Åb= 12.8706 Åc= 7.1014 Åα= 93.461° β= 116.086 ° γ= 90.514 °
I/Icor	0.75
Calc. density	2.704 g/cm³
Reference	Wenk H. R., Joswig W., Tagai T., Korekawa M., Smith B. K., "The average structure of An 62-66 labradorite Lake County neutron feldspar", American Mineralogist 65 , 81-95 (1980)

G: Quartz (6.1 %)	
Formula sum	O2 Si
Entry number	96-901-2604
Figure-of-Merit (FoM)	0.677968
Total number of peaks	31
Peaks in range	31
Peaks matched	5
Intensity scale factor	0.43
Space group	P 31 2 1
Crystal system	trigonal (hexagonal axes)
Unit cell	a= 4.6250 Åc= 5.2160 Å
I/Icor	3.59
Calc. density	3.096 g/cm³
Reference	Hazen R. M., Finger L. W., Hemley R. J., Mao H. K., "High-pressure crystal chemistry and amorphization of alpha-quartz Locality: synthetic Sample: P = 8.0 GPa", Solid State Communications 72 , 507-511 (1989)

Candidates

Name	Formula	Entry No.	FoM
Quartz	O2 Si	96-901-3322	0.6670
Silicon oxide (Quartz)	O2 Si	96-500-0036	0.6668
Quartz	O2 Si	96-901-2604	0.6588
Quartz	O2 Si	96-901-2603	0.6118

Search-Match

Settings	
Reference database used	COD-Inorg REV89244 2013.10.11
Automatic zeropoint adaptation	Yes
Minimum figure-of-merit (FoM)	0.60
Parameter/influence 2theta	0.50
Parameter/influence intensities	0.50
Parameter/influence ...	0.50



Selection Criteria

Compound Name	Quartz
---------------	--------

Peak List

No.	Optimized using trial version www.balesio.com		I/I0	FWHM	Matched
1			0.07	0.2800	B
2	7.04	12.5478	241.68	0.2800	
3	7.14	12.3662	237.79	0.2800	
4	19.90	4.4582	257.76	0.2800	A B E

5	20.27	4.3771	281.16	0.2800	AD,E
6	20.37	4.3562	261.44	0.2800	AB,E,F
7	20.37	4.3562	261.44	0.2800	
8	20.54	4.3205	249.58	0.2800	B,C
9	20.76	4.2757	231.01	0.2800	B
10	20.98	4.2302	306.18	0.2800	AB
11	21.14	4.1992	290.88	0.2800	
12	21.31	4.1662	262.70	0.2800	AB
13	21.31	4.1662	262.70	0.2800	
14	21.78	4.0769	295.16	0.2800	B
15	22.00	4.0369	333.37	0.2800	AB,D,E,F
16	22.30	3.9829	294.73	0.2800	B,E,G
17	22.53	3.9432	256.83	0.2800	B
18	22.53	3.9432	256.83	0.2800	
19	22.84	3.8908	272.17	0.2800	AB,C,F
20	23.06	3.8535	261.31	0.2800	B,D
21	23.32	3.8112	326.33	0.2800	AB,D
22	23.77	3.7404	324.55	0.2800	B,F
23	23.94	3.7136	374.83	0.2800	B,C,D
24	24.36	3.6515	783.83	0.2800	AB,D,F
25	24.80	3.5871	329.19	0.2800	B
26	25.16	3.5363	298.69	0.2800	AB,D
27	25.72	3.4612	410.75	0.2800	B,C,D,F
28	26.01	3.4226	378.96	0.2800	AB,D,F
29	26.40	3.3734	310.94	0.2800	B,D,E,F
30	26.88	3.3140	535.17	0.2800	D,E
31	27.23	3.2729	384.90	0.2800	A
32	27.41	3.2513	328.79	0.2800	B,F
33	27.41	3.2513	328.79	0.2800	
34	27.92	3.1925	562.05	0.2800	AB,D,E,F
35	28.11	3.1719	379.73	0.2800	B,D,E,F,G
36	28.11	3.1719	379.73	0.2800	
37	28.29	3.1516	344.14	0.2800	B,D,E
38	28.57	3.1218	222.38	0.2800	AB,F
39	28.57	3.1218	222.38	0.2800	
40	28.84	3.0931	217.98	0.2800	B
41	29.06	3.0700	214.25	0.2800	AB
42	29.51	3.0249	257.93	0.2800	B,C,E,F
43	30.11	2.9655	479.41	0.2800	B,C,D,F
44	30.44	2.9345	356.82	0.2800	B,D,E,F
45	30.62	2.9177	359.00	0.2800	AB,D,F
46	30.93	2.8887	219.67	0.2800	B
47	31.11	2.8729	234.09	0.2800	B,D,E,F
48	33.39	2.6815	964.12	0.2800	AB,D,E,F
49	35.80	2.5060	1000.00	0.2800	AB,C,D,E,F
50	36.21	2.4788	228.04	0.2800	AB,C,D,E,F
51	36.21	2.4788	228.04	0.2800	
52	39.57	2.2759	214.66	0.2800	AB,C,D,E,F
53	41.07	2.1958	388.92	0.2800	AB,D,E,F,G
54	42.70	2.1159	230.42	0.2800	AB,D,E,F,G
55	49.64	1.8349	372.55	0.2800	AB,C,D,E,F
56	54.26	1.6891	408.48	0.2800	AC,D,E,F
57	62.61	1.4825	289.63	0.2800	AC,D,E,F
58	64.15	1.4505	250.37	0.2800	AD,E,F,G

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

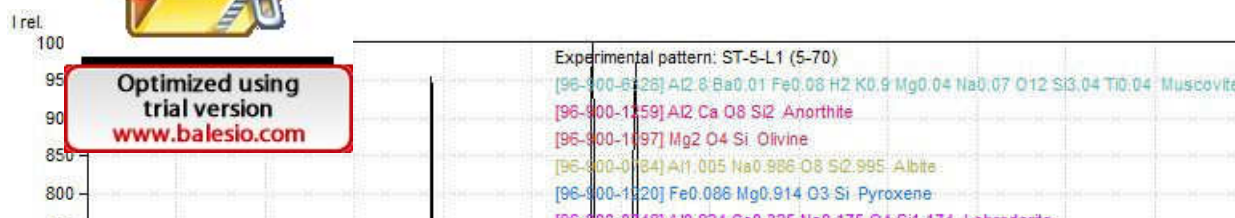
Based on calculated profile

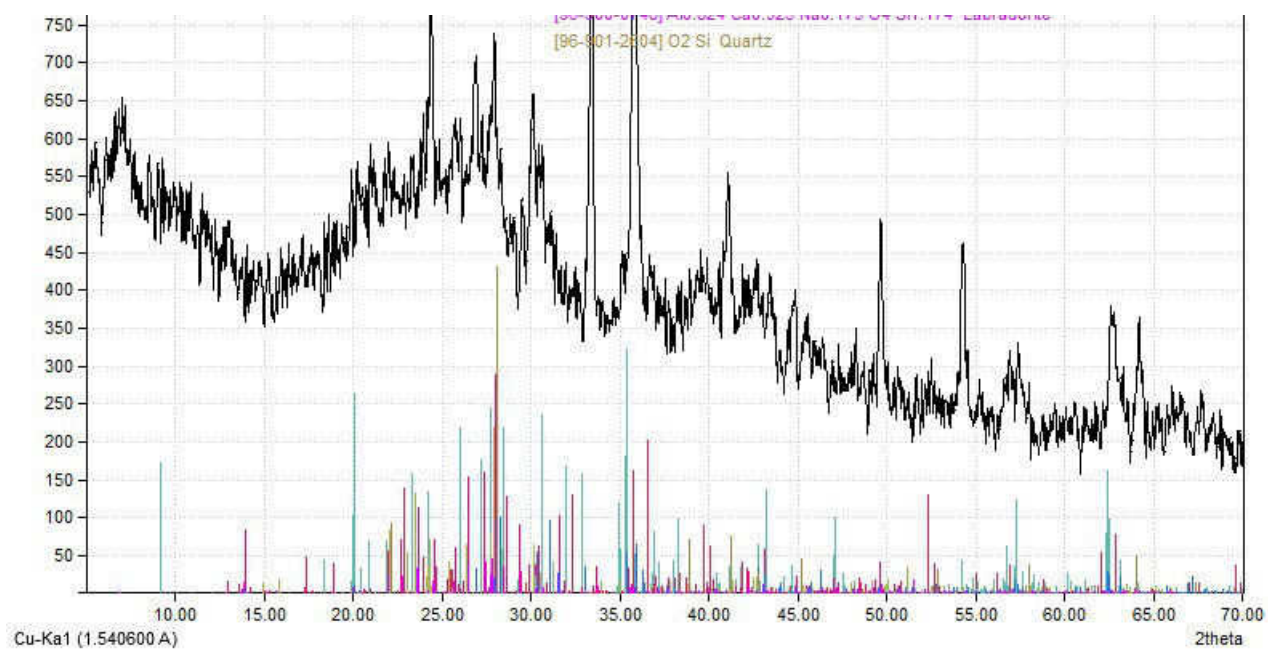
Profile area	Counts	Amount
Overall diffraction profile	172122	100.00%
Background radiation	136953	79.57%
Diffraction peaks	35170	20.43%
Peak area belonging to selected phases	18637	10.83%
Unidentified peak area	16533	9.61%

Peak Residuals

Peak data	Counts	Amount
Overall	791	100.00%
Peak identified	367	46.34%
Unidentified	425	53.66%

Diffraction Pattern Graphics





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Match! Phase Analysis Report

Sample: ST-6-L1 (5-70)

Sample Data

File name ST-6-L1.RAW
File path D:/S2 Geologi/data hasil/XRD Bu Ulva/ST-6-L1
Data collected Jan 5, 2024 16:52:42
Data range 5.080° - 70.080°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction 0.08°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	54.8	Anorthite	Al ₂ Ca O ₈ Si ₂
B	15.1	Olivine	Fe Mg O ₄ Si
C	13.8	Quartz	O ₂ Si
D	6.7	Vermiculite	C ₁₂ Mg ₃ N ₄ O ₁₂ Si ₄
E	5.4	Montmorillonite	Al _{0.86} Fe _{0.1} H Li _{0.08} Mg _{0.14} O ₁₀ Si _{3.9}
F	4.2	Pyroxene	Fe _{0.44} Mg _{0.56} O ₃ Si
	1.7	Unidentified peak area	

A: Anorthite (54.8 %)

Formula sum Al₂ Ca O₈ Si₂
Entry number 96-900-0362
Figure-of-Merit (FoM) 0.817616
Total number of peaks 500
Peaks in range 500
Peaks matched 405
Intensity scale factor 0.60
Space group P -1
Crystal system triclinic (anorthic)
Unit cell a= 8.1940 Å b= 12.8970 Å c= 14.1900 Å α= 92.980° β= 115.820° γ= 91.150°
I/Icor 0.60
Calc. density 2.745 g/cm³
Reference Foit F. F., Peacor D. R., "The anorthite crystal structure at 410 and 830 C T = 410 C", American Mineralogist **58**, 665-675 (1973)

B: Olivine (15.1 %)

Formula sum Fe Mg O₄ Si
Entry number 96-900-6878
Figure-of-Merit (FoM) 0.732500
Total number of peaks 173
Peaks in range 173
Peaks matched 60
Intensity scale factor 0.35
Space group P b n m
Crystal system orthorhombic
Unit cell a= 4.8106 Å b= 10.3863 Å c= 6.0682 Å
I/Icor 1.27
Calc. density 3.773 g/cm³
Reference Redfern S. A. T., Artioli G., Rinaldi R., Henderson C. M. B., Knight K. S., Wood B. J., "Octahedral cation ordering in olivine at high temperature. II: an in situ neutron powder diffraction study on synthetic MgFeSiO₄ (Fa50) Sample: T = 300 C", Physics and Chemistry of Minerals **27**, 630-637 (2000)

C: Quartz (13.8 %)

Formula sum O₂ Si
Entry number 96-900-5022
Figure-of-Merit (FoM) 0.668722
Total number of peaks 35
Peaks in range 35
Peaks matched 14
Intensity scale factor 0.61
Space group P 3 2 1
Crystal system trigonal (hexagonal axes)
Unit cell a= 4.9509 Å c= 5.4285 Å
I/Icor 0.4
Calc. density 2.65 g/cm³
Reference Kura K., "An X-ray study of the temperature dependence of the quartz structure Sample: at T = 697 K", European Journal of Mineralogy **2**, 63-77 (1990)

D: Vermiculite

Formula sum C₁₂ Mg₃ N₄ O₁₂ Si₄
Entry number 96-900-0119
Figure-of-Merit (FoM) 0.64982
Total number of peaks 305
Peaks in range 305
Peaks matched 156
Intensity scale factor 0.43
Space group C 2/c



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Space group	C 1 2 1
Crystal system	monoclinic
Unit cell	a= 5.3300 Å b= 9.1800 Å c= 17.1200 Å β= 97.000 °
I/lcor	3.46
Calc. density	2.306 g/cm³
Reference	Haase D. J., Weiss E. J., Steinfink H., "The crystal structure of a hexamethylene-diamine-vermiculite complex", American Mineralogist 48 , 261-270 (1963)

E: Montmorillonite (5.4 %)

Formula sum	Al0.86 Fe0.1 H Li0.08 Mg0.14 O10 Si3.9
Entry number	96-901-0959
Figure-of-Merit (FoM)	0.703046
Total number of peaks	251
Peaks in range	251
Peaks matched	85
Intensity scale factor	0.58
Space group	C 1 2/m 1
Crystal system	monoclinic
Unit cell	a= 5.1710 Å b= 8.9570 Å c= 9.7400 Å β= 96.100 °
I/lcor	5.89
Calc. density	2.245 g/cm³
Reference	Gournis D., Lappas A., Karakassides M. A., Tobbens D., Moukarika A., "A neutron diffraction study of alkali cation migration in montmorillonites Sample: Li-mont-300", Physics and Chemistry of Minerals 35 , 49-58 (2008)

F: Pyroxene (4.2 %)

Formula sum	Fe0.44 Mg0.56 O3 Si
Entry number	96-900-1578
Figure-of-Merit (FoM)	0.739670
Total number of peaks	437
Peaks in range	437
Peaks matched	139
Intensity scale factor	0.05
Space group	P b c a
Crystal system	orthorhombic
Unit cell	a= 18.3120 Å b= 8.9170 Å c= 5.2170 Å
I/lcor	0.59
Calc. density	3.564 g/cm³
Reference	Hazen R. M., Finger L. W., Ko J., "Effects of pressure on Mg-Fe ordering in orthopyroxene synthesized at 11.3 GPa and 1600 C", American Mineralogist 78 , 1336-1339 (1993)

Candidates

Name	Formula	Entry No.	FoM
Quartz	O2 Si	96-901-2606	0.6688
Quartz	O2 Si	96-901-1495	0.6674
Quartz	O2 Si	96-900-5022	0.6670
Quartz	O2 Si	96-900-5023	0.6668
Quartz	O2 Si	96-900-0778	0.6634
Quartz	O2 Si	96-900-5020	0.6567
Quartz	O2 Si	96-901-2603	0.6523
Quartz	O2 Si	96-900-5024	0.6506
Quartz	O2 Si	96-900-5021	0.6505
Quartz	O2 Si	96-901-0145	0.6501
Quartz	O2 Si	96-900-5019	0.6449
Quartz	O2 Si	96-901-2604	0.6441
Quartz	O2 Si	96-901-1494	0.6426
Quartz	O2 Si	96-901-2605	0.6420
Quartz	O2 Si	96-900-0777	0.6371
Quartz	O2 Si	96-900-0780	0.6334
Quartz	O2 Si	96-901-0146	0.6332
Quartz	O2 Si	96-901-0147	0.6315
Quartz	O2 Si	96-900-0779	0.6256
Quartz	O2 Si	96-900-5025	0.6255
Quartz	O2 Si	96-900-0776	0.6228
Quartz	O2 Si	96-900-5018	0.6218
Quartz	O2 Si	96-901-5023	0.6142
Quartz	O2 Si	96-900-7379	0.6091

Search-Match

Settings	
Reference database used	COD-Inorg REV89244 2013.10.11
Automatic zeropoint adaptation	Yes
Minimum figure of merit (FoM)	0.60
Parametric search	no
Parametric search	no
Parametric search	no



Selection Criteria

Compound Name	Quartz
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Peak List

No.	2theta [°]	d[Å]	I/I0	FWHM	Matched
1	6.70	13.1878	184.00	1.2400	AD

2	7.05	12.5349	179.70	1.2400	AE
3	11.44	7.7311	210.74	1.2400	AD,F
4	11.73	7.5395	262.85	1.2400	
5	12.02	7.3565	301.96	1.2400	A
6	12.32	7.1773	329.62	1.2400	
7	12.58	7.0308	205.01	1.2400	AF
8	12.58	7.0308	205.01	1.2400	
9	19.48	4.5540	265.49	1.2400	AB,D,E,F
10	20.03	4.4287	951.41	1.2400	AD,E,F
11	20.34	4.3626	764.61	1.2400	AB,F
12	20.34	4.3626	764.61	1.2400	
13	20.52	4.3247	675.54	1.2400	AC,D
14	20.52	4.3247	675.54	1.2400	
15	21.03	4.2211	878.36	1.2400	AD
16	21.19	4.1891	877.03	1.2400	E
17	21.55	4.1195	554.26	1.2400	A
18	21.74	4.0847	507.08	1.2400	F
19	21.74	4.0847	507.08	1.2400	
20	21.93	4.0499	403.82	1.2400	AD,E,F
21	22.18	4.0044	426.27	1.2400	AF
22	22.58	3.9344	384.21	1.2400	AB
23	22.75	3.9053	318.53	1.2400	AE
24	22.98	3.8663	304.86	1.2400	AD
25	23.14	3.8405	321.04	1.2400	A
26	23.62	3.7640	692.62	1.2400	AB,D
27	23.98	3.7080	422.79	1.2400	A
28	23.98	3.7080	422.79	1.2400	
29	24.35	3.6526	385.79	1.2400	A
30	24.58	3.6183	484.06	1.2400	AF
31	25.01	3.5579	587.49	1.2400	AB,D
32	25.33	3.5140	595.75	1.2400	AB
33	25.74	3.4584	679.46	1.2400	AE
34	26.23	3.3946	334.67	1.2400	AC,D,F
35	26.77	3.3277	843.99	1.2400	F
36	27.10	3.2873	990.94	1.2400	AD,E
37	27.60	3.2290	1000.00	1.2400	AE,F
38	28.25	3.1569	178.07	1.2400	AD,E,F
39	29.95	2.9811	370.96	1.2400	AB,F
40	30.21	2.9556	198.01	1.2400	AD,F
41	30.88	2.8936	391.77	1.2400	AB,D,F
42	34.15	2.6236	183.21	1.2400	AB,D,E,F
43	34.39	2.6058	288.40	1.2400	AB,D,E,F
44	35.07	2.5566	798.23	1.2400	AD,E,F
45	35.26	2.5433	662.03	1.2400	AB,D
46	35.26	2.5433	662.03	1.2400	
47	35.64	2.5174	547.51	1.2400	AD,E,F
48	35.84	2.5035	492.25	1.2400	AE,F
49	35.84	2.5035	492.25	1.2400	
50	36.18	2.4808	358.07	1.2400	AB,C,D,F
51	36.18	2.4808	358.07	1.2400	
52	36.38	2.4678	320.92	1.2400	AE
53	36.65	2.4503	305.71	1.2400	AD
54	36.83	2.4384	288.91	1.2400	AD
55	37.22	2.4140	386.09	1.2400	AB,D,E,F
56	37.73	2.3825	384.15	1.2400	AB,D,E,F
57	37.90	2.3720	369.74	1.2400	A
58	37.90	2.3720	369.74	1.2400	
59	38.03	2.3642	364.77	1.2400	A
60	38.35	2.3451	462.49	1.2400	AB,D
61	38.56	2.3328	440.48	1.2400	AE
62	38.90	2.3133	294.05	1.2400	AD,F
63	38.90	2.3133	294.05	1.2400	
64	39.13	2.3002	246.83	1.2400	AB,D
65	39.26	2.2928	236.49	1.2400	AC,D,F
66	39.51	2.2787	203.67	1.2400	AB,D,F
67	39.99	2.2525	206.14	1.2400	AC,D,F
68	40.36	2.2332	230.59	1.2400	AD,E,F
69	40.63	2.2188	195.51	1.2400	AD,E
70	41.02	2.1983	196.34	1.2400	AD,F
71	41.21	2.1889	186.71	1.2400	AB,D
72	41.36	2.1812	178.70	1.2400	AB,E
73	41.36	2.1812	178.70	1.2400	
74	41.84	2.1574	436.73	1.2400	AC,D,F
75			.60	1.2400	AB,D,E,F
76			.10	1.2400	AB,C,D,E,F
77			.47	1.2400	AB,D,E,F
78			.88	1.2400	AB,C,D,E,F
79			.80	1.2400	B,D,E,F
80			.13	1.2400	C,D,E,F
81			.80	1.2400	C,D,F
82			.69	1.2400	B,D,F
83			.69	1.2400	
84			.24	1.2400	B,E
85			.08	1.2400	D,F
86			.08	1.2400	
87	55.79	1.6465	179.08	1.2400	B,C,D,E,F
88	60.51	1.5288	239.62	1.2400	B,C,D,E,F
89	61.68	1.5027	229.58	1.2400	B,D,E,F



90	62.24	1.4905	403.15	1.2400	D,E,F
91	62.53	1.4843	317.51	1.2400	B,C,D,E,F

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

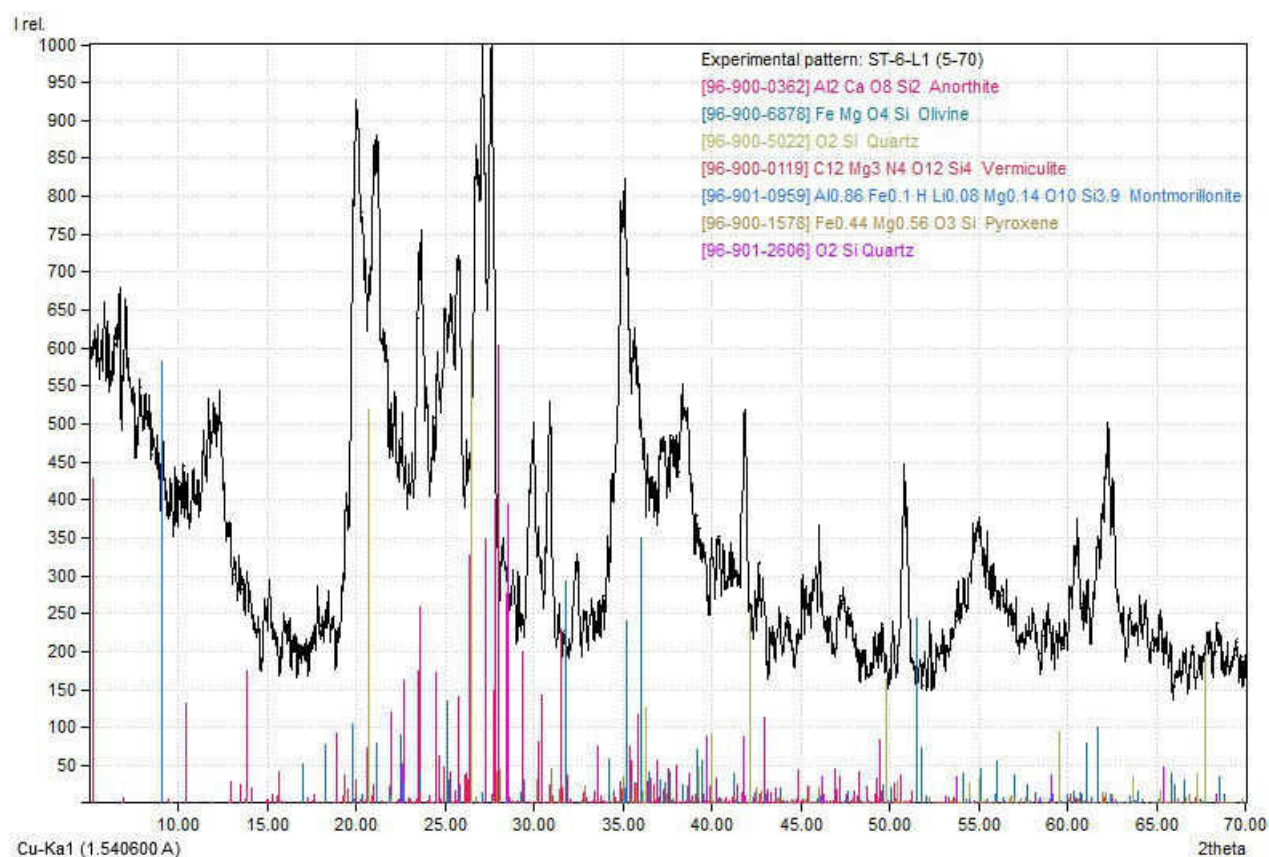
Based on calculated profile

Profile area	Counts	Amount
Overall diffraction profile	114888	100.00%
Background radiation	67436	58.70%
Diffraction peaks	47452	41.30%
Peak area belonging to selected phases	45492	39.60%
Unidentified peak area	1960	1.71%

Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	5837	100.00%
Peak intensity belonging to selected phases	1959	33.55%
Unidentified peak intensity	3879	66.45%

Diffraction Pattern Graphics



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Match! Phase Analysis Report

Sample: ST-7-L1 (5-70)

Sample Data

File name ST-7-L1.RAW
File path D:/S2 Geologi/data hasil/XRD Bu Ulva/ST-7-L1
Data collected Jan 5, 2024 16:52:42
Data range 4.940° - 69.940°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction -0.06°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	26.0	Muscovite	Al2.038 Ba0.01 Ca0.02 Fe0.14 H2 K0.87 Mg0.5 Na0.07 O12 Si3.392 Ti0.01
B	25.6	Anorthite	Al2 Ca O8 Si2
C	16.2	Biotite	Al1.322 Fe0.864 K Mg1.638 O12 Si2.84 Ti0.336
D	12.7	Olivine	Fe Mg O4 Si
E	8.4	Quartz	O2 Si
F	5.3	Pyroxene	Fe0.44 Mg0.56 O3 Si
G	5.3	Vermiculite	C12 Mg3 N4 O12 Si4
H	0.5	Montmorillonite	Al2 Ca O12 Si4
	16.5	Unidentified peak area	

A: Muscovite (26.0 %)

Formula sum Al2.038 Ba0.01 Ca0.02 Fe0.14 H2 K0.87 Mg0.5 Na0.07 O12 Si3.392 Ti0.01
Entry number 96-901-6013
Figure-of-Merit (FoM) 0.854215
Total number of peaks 300
Peaks in range 172
Peaks matched 136
Intensity scale factor 0.36
Space group C 1 2/c 1
Crystal system monoclinic
Unit cell a= 5.2112 Å b= 9.0383 Å c= 19.9473 Å β= 95.769 °
I/Icor 0.75
Calc. density 2.858 g/cm³
Reference Guven N., "The crystal structures of 2M1 phengite and 2M1 muscovite Note: sample 2M1 phengite", Zeitschrift für Kristallographie **134**, 196-212 (1971)

B: Anorthite (25.6 %)

Formula sum Al2 Ca O8 Si2
Entry number 96-900-0363
Figure-of-Merit (FoM) 0.837021
Total number of peaks 500
Peaks in range 500
Peaks matched 340
Intensity scale factor 0.28
Space group P -1
Crystal system triclinic (anorthic)
Unit cell a= 8.2230 Å b= 12.9150 Å c= 14.2040 Å α= 92.750° β= 115.800° γ= 91.020 °
I/Icor 0.60
Calc. density 2.725 g/cm³
Reference Foit F. F., Peacor D. R., "The anorthite crystal structure at 410 and 830 C T = 830 C", American Mineralogist **58**, 665-675 (1973)

C: Biotite (16.2 %)

Formula sum Al1.322 Fe0.864 K Mg1.638 O12 Si2.84 Ti0.336
Entry number 96-900-0844
Figure-of-Merit (FoM) 0.827107
Total number of peaks 297
Peaks in range 177
Peaks matched 119
Intensity scale factor 0.46
Space group C 1 2/c 1
Crystal system monoclinic
Unit cell a= 5.3175 Å b= 9.2120 Å c= 19.9760 Å β= 95.090 °
I/Icor 0.4
Calc. density 3.71 g/cm³
Reference Takeda H., Takeda H., Takeuchi Y., "Mica polytypism: Similarities in the crystal structures of coexisting 1M and 2M(1) oxybiotite sample: in the 2M1 setting", American Mineralogist **67**, 298-310 (1982)

D: Olivine

Formula Mg O4 Si
Entry number 96-900-6881
Figure-of-Merit (FoM) 0.37179
Total number of peaks 174
Peaks in range 63
Peaks matched 42
Intensity scale factor 0.28



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Intensity scale factor	0.29
Space group	P b n m
Crystal system	orthorhombic
Unit cell	a= 4.8168 Å b= 10.4029 Å c= 6.0800 Å
I/I _{cor}	1.24
Calc. density	3.754 g/cm ³
Reference	Redfern S. A. T., Artioli G., Rinaldi R., Henderson C. M. B., Knight K. S., Wood B. J., "Octahedral cation ordering in olivine at high temperature. II: an in situ neutron powder diffraction study on synthetic MgFeSiO ₄ (Fa50) Sample: T = 450 °C", Physics and Chemistry of Minerals 27 , 630-637 (2000)

E: Quartz (8.4 %)

Formula sum	O ₂ Si
Entry number	96-901-3322
Figure-of-Merit (FoM)	0.760559
Total number of peaks	35
Peaks in range	14
Peaks matched	14
Intensity scale factor	0.69
Space group	P 3 2 1
Crystal system	trigonal (hexagonal axes)
Unit cell	a= 4.9134 Å c= 5.4051 Å
I/I _{cor}	4.74
Calc. density	2.649 g/cm ³
Reference	Antao S. M., Hassan I., Wang J., Lee P. L., Toby B. H., "State-of-the-art high-resolution powder x-ray diffraction (HRPXRD) illustrated with Rietveld structure refinement of quartz, sodalite, tremolite, and meionite Locality: not specified", The Canadian Mineralogist 46 , 1501-1509 (2008)

F: Pyroxene (5.3 %)

Formula sum	Fe _{0.44} Mg _{0.56} O ₃ Si
Entry number	96-900-1578
Figure-of-Merit (FoM)	0.777974
Total number of peaks	437
Peaks in range	148
Peaks matched	114
Intensity scale factor	0.06
Space group	P b c a
Crystal system	orthorhombic
Unit cell	a= 18.3120 Å b= 8.9170 Å c= 5.2170 Å
I/I _{cor}	0.59
Calc. density	3.564 g/cm ³
Reference	Hazen R. M., Finger L. W., Ko J., "Effects of pressure on Mg-Fe ordering in orthopyroxene synthesized at 11.3 GPa and 1600 °C", American Mineralogist 78 , 1336-1339 (1993)

G: Vermiculite (5.3 %)

Formula sum	C ₁₂ Mg ₃ N ₄ O ₁₂ Si ₄
Entry number	96-900-0119
Figure-of-Merit (FoM)	0.789110
Total number of peaks	306
Peaks in range	167
Peaks matched	126
Intensity scale factor	0.33
Space group	C 1 2 1
Crystal system	monoclinic
Unit cell	a= 5.3300 Å b= 9.1800 Å c= 17.1200 Å β= 97.000 °
I/I _{cor}	3.46
Calc. density	2.306 g/cm ³
Reference	Haase D. J., Weiss E. J., Steinfink H., "The crystal structure of a hexamethylene-diamine-vermiculite complex", American Mineralogist 48 , 261-270 (1963)

H: Montmorillonite (0.5 %)

Formula sum	Al ₂ Ca O ₁₂ Si ₄
Entry number	96-110-1055
Figure-of-Merit (FoM)	0.716328
Total number of peaks	92
Peaks in range	92
Peaks matched	79
Intensity scale factor	0.20
Space group	P 1
Crystal system	triclinic (anorthic)
Unit cell	a= 5.1800 Å b= 8.9800 Å c= 15.0000 Å α= 90.000° β= 90.000° γ= 90.000°
I/I _{cor}	20.53
Calc. density	1.800 g/cm ³

Candidates

Name	Formula	Entry No.	FoM
Quartz	O ₂ Si	96-901-2602	0.7248
Quartz	O ₂ Si	96-900-5023	0.7047
Quartz	O ₂ Si	96-900-0781	0.6840
Quartz	O ₂ Si	96-901-3322	0.6828
Quartz	O ₂ Si	96-901-2601	0.6819
Silicor	O ₂ Si	96-500-0036	0.6814
Quartz	O ₂ Si	96-900-5021	0.6785
Quartz	O ₂ Si	96-900-5024	0.6781
Quartz	O ₂ Si	96-900-5025	0.6772
Quartz	O ₂ Si	96-900-5022	0.6757
Quartz	O ₂ Si	96-900-7379	0.6715
Quartz	O ₂ Si	96-901-1496	0.6269



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Search-Match

Settings

Reference database used COD-Inorg REV89244 2013.10.11
Automatic zeropoint adaptation Yes
Minimum figure-of-merit (FoM) 0.60
Parameter/influence 2theta 0.50
Parameter/influence intensities 0.50
Parameter multiple/single phase(s) 0.50

Selection Criteria

Compound:

Name: quartz

Peak List

No.	2theta [°]	d[Å]	I/I0	FWHM	Matched
1	5.22	16.9043	135.20	0.2800	G
2	5.64	15.6527	180.11	0.2800	
3	5.80	15.2254	138.97	0.2800	
4	5.92	14.9275	133.10	0.2800	H
5	6.09	14.5068	155.91	0.2800	
6	6.62	13.3369	228.53	0.2800	
7	6.77	13.0433	288.58	0.2800	B
8	6.98	12.6450	218.41	0.2800	B
9	7.27	12.1553	222.83	0.2800	
10	7.56	11.6845	180.16	0.2800	
11	7.93	11.1395	178.73	0.2800	
12	8.16	10.8256	194.55	0.2800	
13	8.35	10.5842	199.85	0.2800	
14	8.54	10.3443	190.84	0.2800	
15	8.79	10.0521	174.28	0.2800	
16	8.97	9.8537	188.78	0.2800	A,C
17	9.38	9.4180	155.98	0.2800	B
18	9.65	9.1556	141.54	0.2800	F
19	10.26	8.6167	138.88	0.2800	B,G
20	10.56	8.3687	138.17	0.2800	
21	10.81	8.1795	134.98	0.2800	B
22	10.97	8.0571	149.13	0.2800	
23	11.49	7.6980	211.63	0.2800	
24	12.02	7.3558	253.56	0.2800	B,H
25	12.16	7.2727	238.16	0.2800	
26	12.40	7.1333	222.38	0.2800	
27	12.58	7.0308	161.52	0.2800	B
28	13.36	6.6207	140.50	0.2800	
29	13.48	6.5657	152.37	0.2800	B,F
30	17.62	5.0294	153.28	0.2800	AB,C,H
31	19.35	4.5833	204.43	0.2800	AB,C,F,G
32	19.97	4.4424	489.04	0.2800	AB,C,F,G,H
33	20.14	4.4055	394.70	0.2800	AB,C
34	20.33	4.3650	396.48	0.2800	AB,D,F,G
35	20.92	4.2439	723.75	0.2800	AB,C,E,G,H
36	21.12	4.2032	614.85	0.2800	C
37	21.40	4.1489	317.87	0.2800	B
38	21.67	4.0980	285.36	0.2800	AB
39	21.82	4.0698	257.16	0.2800	B,C,F,G
40	22.18	4.0038	250.99	0.2800	B,F
41	22.41	3.9634	310.84	0.2800	AB,D
42	22.80	3.8970	191.83	0.2800	AB,C
43	23.13	3.8426	250.13	0.2800	B,G,H
44	23.56	3.7736	573.78	0.2800	B,C,D,G,H
45	23.89	3.7213	245.00	0.2800	AB
46	24.55	3.6229	339.38	0.2800	B,C,F
47	24.93	3.5686	351.76	0.2800	AB,G
48	25.12	3.5429	375.62	0.2800	B,C,D
49	25.65	3.4701	494.39	0.2800	AB
50	26.16	3.4033	275.11	0.2800	B,C,F,G
51	26.67	3.3398	675.53	0.2800	AB,E,F,H
52	26.94	3.3065	700.38	0.2800	A,C,G
53	27.16	3.2806	596.01	0.2800	B
54			1.00	0.2800	B,C
55			3.75	0.2800	AB,F
56			1.07	0.2800	AB,C,F,G
57			1.44	0.2800	B,C,D,F,H
58			7.77	0.2800	AB,F,G
59			1.78	0.2800	AB,C,F,G,H
60			3.15	0.2800	AB,C,D,G
61			1.29	0.2800	AB,C,F,G
62			3.66	0.2800	AB,C,D,F,G
63			3.65	0.2800	A,C,D,F,G,H
64			5.59	0.2800	AB,G
65			2.51	0.2800	AB,D,F,H
66	35.26	2.5433	379.91	0.2800	B
67	35.47	2.5287	342.57	0.2800	B,F,G
68	35.59	2.5202	346.23	0.2800	B,F,G



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69	35.86	2.5022	250.70	0.2800	AB,C,F,H
70	36.02	2.4914	221.59	0.2800	AB,C,D,F,G,H
71	36.22	2.4781	184.01	0.2800	AB,F
72	36.41	2.4656	193.35	0.2800	AB
73	36.62	2.4520	209.77	0.2800	AB,E,G,H
74	37.08	2.4223	292.97	0.2800	B,C,F,G
75	37.41	2.4019	340.21	0.2800	AB,D,F,G
76	37.76	2.3805	235.90	0.2800	AB,F,G
77	37.97	2.3675	252.08	0.2800	B
78	38.21	2.3535	236.78	0.2800	A
79	38.38	2.3435	207.99	0.2800	B,C,D,G
80	38.58	2.3317	251.40	0.2800	B
81	38.90	2.3133	203.00	0.2800	AB,F,G
82	39.12	2.3010	181.60	0.2800	B,C,D,G,H
83	39.39	2.2855	141.63	0.2800	B,C,D,E,F,G
84	39.87	2.2590	132.03	0.2800	AB,C,F,G,H
85	40.44	2.2285	139.74	0.2800	AB,C,E,F,G,H
86	41.26	2.1862	147.58	0.2800	AB,C,D,F,G,H
87	41.71	2.1635	449.37	0.2800	AB,C,F,G,H
88	42.52	2.1242	139.66	0.2800	AB,C,E,F,G,H
89	44.93	2.0159	172.26	0.2800	AB,F,G
90	45.17	2.0059	169.30	0.2800	B,F,G
91	45.34	1.9986	138.58	0.2800	AB,F,G
92	45.52	1.9910	156.73	0.2800	C
93	45.82	1.9788	140.43	0.2800	AB,D,E,F,G
94	46.12	1.9666	144.68	0.2800	AB,C,F,H
95	47.06	1.9294	157.85	0.2800	AB,F,G,H
96	47.22	1.9231	152.00	0.2800	B,C,F,H
97	47.51	1.9121	150.04	0.2800	AB,C,D,F,G
98	50.68	1.7998	309.83	0.2800	AB,C,E,F,G,H
99	53.64	1.7074	135.36	0.2800	AC,F,G
100	53.84	1.7015	174.30	0.2800	AF,G
101	54.03	1.6960	159.03	0.2800	AC,D,F,G,H
102	54.15	1.6922	179.02	0.2800	AC,F
103	54.33	1.6872	178.78	0.2800	AC,F,G,H
104	54.79	1.6741	239.37	0.2800	AC,E,F,G,H
105	55.15	1.6639	236.90	0.2800	AC,D,F,G
106	55.42	1.6567	147.83	0.2800	AC,D,E,F,G,H
107	55.55	1.6530	152.71	0.2800	AG,H
108	55.75	1.6476	159.96	0.2800	AD,F,G
109	56.05	1.6395	187.79	0.2800	AC,G
110	56.49	1.6277	166.39	0.2800	AC,D,F
111	56.73	1.6214	134.27	0.2800	AC,D,E,F,G,H
112	59.94	1.5419	149.28	0.2800	AC,D,E,F,G,H
113	60.24	1.5350	252.59	0.2800	AC,D,F,G
114	60.80	1.5222	138.44	0.2800	AC,D,F,G,H
115	61.53	1.5060	157.94	0.2800	AC,D,F,G
116	62.01	1.4954	266.59	0.2800	AF,G,H
117	62.15	1.4923	289.14	0.2800	AF
118	62.42	1.4866	216.67	0.2800	AC,D,F,G,H
119	62.68	1.4810	146.90	0.2800	AC,F,G,H
120	63.48	1.4643	132.77	0.2800	AC,D,F,G,H
121	64.09	1.4518	173.67	0.2800	AC,D,E,F,G
122	64.79	1.4377	131.13	0.2800	AD,F,G,H
123	65.13	1.4311	135.01	0.2800	F,H

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

Based on calculated profile

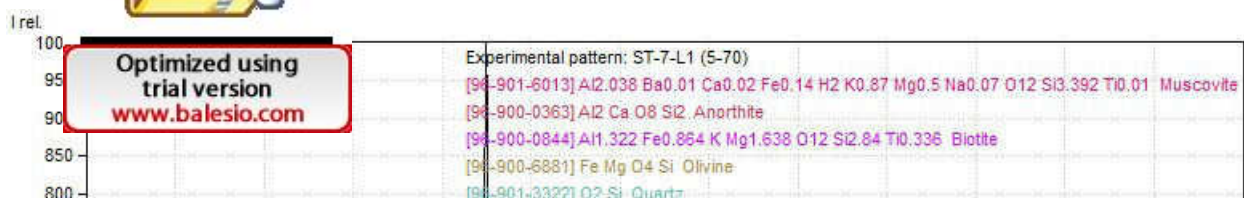
Profile area	Counts	Amount
Overall diffraction profile	105586	100.00%
Background radiation	65865	62.38%
Diffraction peaks	39721	37.62%
Peak area belonging to selected phases	22305	21.13%
Unidentified peak area	17416	16.49%

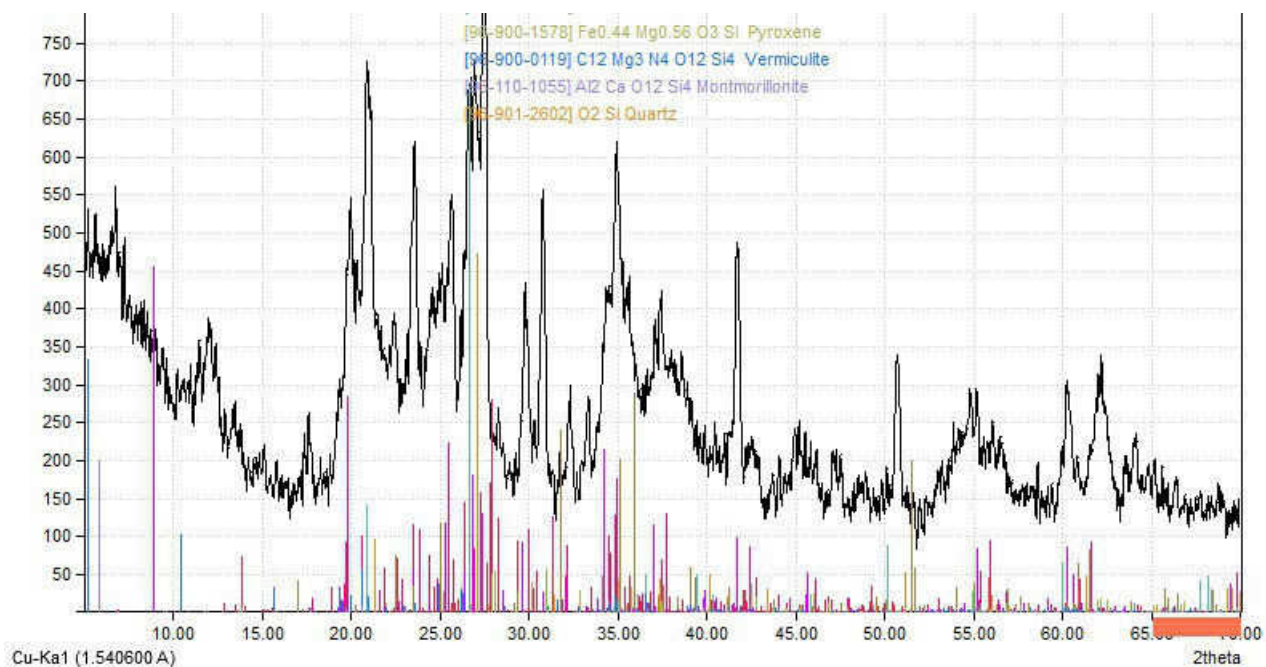
Peak Residuals

Peak	Counts	Amount
Overall	1469	100.00%
Peak i	567	38.62%
Unide	902	61.38%



Diffraction Pattern Graphics





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Match! Phase Analysis Report

Sample: ST-8-L1 (5-70)

Sample Data

File name ST-8-L1.RAW
File path D:/S2 Geologi/data hasil/XRD Bu Ulva/ST-8-L1
Data collected Jan 5, 2024 16:52:42
Data range 5.020° - 70.020°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction 0.02°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	44.7	Muscovite	Al2.908 H0.12 K O10 Si3.092
B	26.8	Anorthite	Al2 Ca O8 Si2
C	12.6	Tridymite	O2 Si
D	11.0	Olivine	Fe Mg O4 Si
E	4.6	Hornblende	Al2.42 Ca1.83 Fe1.852 H1.34 Mg2.218 Na0.706 O24 Si6.44 Ti0.07
F	0.4	Vermiculite	Al0.721 Fe0.24 H3 Mg1.338 O9 Si1.36
	3.9	Unidentified peak area	

A: Muscovite (44.7 %)

Formula sum Al2.908 H0.12 K O10 Si3.092
Entry number 96-900-1059
Figure-of-Merit (FoM) 0.757165
Total number of peaks 300
Peaks in range 300
Peaks matched 162
Intensity scale factor 0.44
Space group C 1 2/c 1
Crystal system monoclinic
Unit cell a= 5.1890 Å b= 9.0040 Å c= 20.2560 Å β= 95.740 °
I/Icor 0.52
Calc. density 2.571 g/cm³
Reference Guggenheim S., Chang Y. H., Koster van Groos A. F., "Muscovite dehydroxylation: High-temperature studies Sample: T = 650 deg C, from Panasqueira, Portugal", American Mineralogist **72**, 537-550 (1987)

B: Anorthite (26.8 %)

Formula sum Al2 Ca O8 Si2
Entry number 96-900-0363
Figure-of-Merit (FoM) 0.772052
Total number of peaks 500
Peaks in range 498
Peaks matched 390
Intensity scale factor 0.30
Space group P -1
Crystal system triclinic (anorthic)
Unit cell a= 8.2230 Å b= 12.9150 Å c= 14.2040 Å α= 92.750° β= 115.800° γ= 91.020 °
I/Icor 0.60
Calc. density 2.725 g/cm³
Reference Foit F. F., Peacor D. R., "The anorthite crystal structure at 410 and 830 C T = 830 C", American Mineralogist **58**, 665-675 (1973)

C: Tridymite (12.6 %)

Formula sum O2 Si
Entry number 96-900-6969
Figure-of-Merit (FoM) 0.820244
Total number of peaks 500
Peaks in range 499
Peaks matched 207
Intensity scale factor 0.37
Space group P 21 21 21
Crystal system orthorhombic
Unit cell a= 26.1753 Å b= 4.9844 Å c= 8.2006 Å
I/Icor 0.8
Calc. density 2.37 g/cm³
Reference Jaesch H., "X-ray powder diffraction study on the modulated high temperature forms of SiO2 tridymite between 110 and 220 C sample: superstructure, T = 115 C Locality: synthetic", Physics and Chemistry of Minerals **28**, 313-321 (2001)

D: Olivine

Formula sum Mg O4 Si
Entry number 96-900-6878
Figure-of-Merit (FoM) 0.25974
Total number of peaks 3
Peaks matched 3
Intensity scale factor 0.26
Space group P b n m
Crystal system orthorhombic



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Crystal system	orthorhombic
Unit cell	a= 4.8106 Å b= 10.3863 Å c= 6.0682 Å
I/Cor	1.27
Calc. density	3.773 g/cm ³
Reference	Redfern S. A. T., Artioli G., Rinaldi R., Henderson C. M. B., Knight K. S., Wood B. J., "Octahedral cation ordering in olivine at high temperature. II: an in situ neutron powder diffraction study on synthetic MgFeSiO ₄ (Fa50) Sample: T = 300 C", Physics and Chemistry of Minerals 27 , 630-637 (2000)

E: Hornblende (4.6 %)

Formula sum	Al _{2.42} Ca _{1.83} Fe _{1.852} H _{1.34} Mg _{2.218} Na _{0.706} O ₂₄ Si _{6.44} Ti _{0.07}
Entry number	96-900-1228
Figure-of-Merit (FoM)	0.760294
Total number of peaks	299
Peaks in range	299
Peaks matched	164
Intensity scale factor	0.08
Space group	C 1 2/m 1
Crystal system	monoclinic
Unit cell	a= 9.7860 Å b= 18.0240 Å c= 5.3060 Å β= 105.090 °
I/Cor	0.97
Calc. density	3.241 g/cm ³
Reference	Phillips M. W., Draheim J. E., Popp R. K., Clowe C. A., Pinkerton A. A., "Effects of oxidation-dehydrogenation in tschermakitic hornblende sample H-3, after annealing at 700 C", American Mineralogist 74 , 764-773 (1989)

F: Vermiculite (0.4 %)

Formula sum	Al _{0.721} Fe _{0.24} H ₃ Mg _{1.338} O ₉ Si _{1.36}
Entry number	96-900-0061
Figure-of-Merit (FoM)	0.742667
Total number of peaks	292
Peaks in range	291
Peaks matched	171
Intensity scale factor	0.15
Space group	C 1 c 1
Crystal system	monoclinic
Unit cell	a= 5.3300 Å b= 9.1800 Å c= 28.9000 Å β= 97.000 °
I/Cor	19.70
Calc. density	1.185 g/cm ³
Reference	Mathieson A. M., Walker G. F., "Crystal structure of magnesium-vermiculite", American Mineralogist 39 , 231-255 (1954)

Candidates

Name	Formula	Entry No.	FoM
Antigorite-T	Mg ₄₈ O ₁₄₇ Si ₃₄	96-901-6234	0.7958
meta-schoepite	H ₁₆ O ₂₀ U ₄	96-210-2097	0.7956
Metaschoepite	H ₃₄ O ₄₀ U ₈	96-901-1299	0.7956
Metaschoepite	H ₃₂ Na _{1.09} O _{38.328} U ₈	96-901-0196	0.7916
Metaschoepite	H ₃₄ Na _{0.48} O _{37.91} U ₈	96-901-0195	0.7872
Metaschoepite	H ₃₄ Na _{0.47} O _{37.082} U ₈	96-901-0199	0.7871
Metaschoepite	H ₃₄ Na _{1.16} O _{37.9} U ₈	96-901-0198	0.7814
Metaschoepite	Cs F ₃₀ O ₆ Sb Te ₆	96-432-7610	0.7808
Porphyrazinealuminiumchloride	C ₁₆ Al Cl N ₁₆ S ₄	96-430-9965	0.7793
Metaschoepite	H ₃₂ Na _{1.22} O _{39.09} U ₈	96-901-0197	0.7792
Becquerelite	Ca H ₂₂ O ₃₀ U ₆	96-900-2701	0.7785
Vandendriesscheite	H ₃₃ O ₄₈ Pb ₂ U ₁₀	96-900-1892	0.7774
Becquerelite	Ca O ₃₀ U ₆	96-900-1111	0.7762
Billietite	Ba O ₂₃ U ₆	96-900-1112	0.7739
Bikitaite	Al H ₂ Li O _{6.385} Si ₂	96-900-3102	0.7725
potassium hydroxypentafluoroarsenate	As F ₅ H K O	96-200-3122	0.7718
Schoepite	H ₁₈ O ₂₁ U ₄	96-900-4445	0.7701
Billietite	Ba H ₆ O ₃₀ U ₆	96-901-0600	0.7692
Becquerelite	Ca H ₂₂ O ₃₀ U ₆	96-901-2088	0.7688
Fourmarierite	H ₁₀ O ₁₉ Pb U ₄	96-900-4592	0.7683
Porphyrazinegalliumchloride	C ₁₆ Cl Ga N ₁₆ S ₄	96-430-9966	0.7681
Dickite	Al ₂ H ₄ O ₉ Si ₂	96-900-0123	0.7681
Sidwillite	H ₄ Mo O ₅	96-901-1123	0.7664
	Cl F ₁₇ Sb ₂	96-431-1868	0.7652
	Br ₂ F ₁₇ Sb ₂	96-431-1869	0.7649
CuNO blue	C ₁₂ Cu N ₆ O ₇	96-710-1658	0.7649
Khademite	Al F H ₁₀ O ₉ S	96-900-9710	0.7649
Tricaesium cyclo-triphosphate telluric acid hydrate	Cs ₃ H ₈ O ₁₆ P ₃ Te	96-100-8369	0.7644
Potassium aluminium silicate hydroxide * (Muscovite 2M1)	Al ₃ H ₂ K O ₁₂ Si ₃	96-101-1050	0.7641
Potassium aluminium silicate hydroxide * (Muscovite 2M1)	Al ₃ H K O ₁₂ Si ₃	96-110-0014	0.7641
Dialuminium phyllo-disilicate tetrahydroxide (Nacrite 6M)	Al ₂ H ₄ O ₉ Si ₂	96-101-1063	0.7640
Gemirite	As Cu H ₃ O ₅	96-901-6163	0.7635
Gobbite	Al _{2.8} Ca _{0.47} H ₁₂ Na ₂ O ₂₂ Si _{5.296}	96-900-9491	0.7629
	B ₃ H ₃ N ₁₂	96-432-2188	0.7624
Fourm	H ₁₀ O ₁₉ Pb U ₄	96-900-4591	0.7624
Nacrite	Al ₂ O ₉ Si ₂	96-901-4064	0.7620
Dialun	Al ₂ H ₄ O ₉ Si ₂	96-101-1081	0.7614
Dickite	Al ₂ H ₄ O ₉ Si ₂	96-901-3920	0.7612
Cobalt	As Co H ₃ O ₅	96-901-2337	0.7610
lanthir	H _{22.62} O ₂₇ U ₆	96-901-2907	0.7601
ammo	B ₃ H ₈ N O _{9.5} P	96-221-6154	0.7595
Fourm	H ₁₀ O ₁₉ Pb U ₄	96-900-4588	0.7595
	H ₁₅ Na N ₄ O ₂₂ P ₄	96-432-3133	0.7592
Magnesium hydrogensulfate hydrate	H ₄ Mg O ₉ S ₂	96-110-0087	0.7591
Monoclinic titanium metaphosphate	O ₂₇ P ₉ Ti ₃	96-201-0156	0.7591
Fourmarierite	H ₁₀ O ₁₉ Pb U ₄	96-900-4584	0.7590
Fourmarierite	H ₁₀ O ₁₉ Pb U ₄	96-900-4588	0.7588



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Fourmarierite

Antigorite

Fourmarierite

Bikitaite

Fourmarierite

and 153 others...

H10 O19 Pb U4

Cr2 H8 Mg O16 U2

H58 Mg45 O138 Si32

H10 O19 Pb U4

Al Li O6 Si2

H10 O19 Pb U4

96-900-4590

96-400-1609

96-900-4000

96-900-4581

96-900-3103

96-900-4589

0.7589

0.7587

0.7587

0.7583

0.7580

0.7579

Search-Match

Settings

Reference database used COD-Inorg REV89244 2013.10.11

Automatic zeropoint adaptation Yes

Minimum figure-of-merit (FoM) 0.60

Parameter/influence 2theta 0.50

Parameter/influence intensities 0.50

Parameter multiple/single phase(s) 0.50

Peak List

No.	2theta [°]	d [Å]	I/I0	FWHM	Matched
1	7.33	12.0447	147.58	0.6800	B,C,F
2	11.47	7.7109	158.16	0.6800	B,E
3	11.87	7.4476	215.09	0.6800	B
4	12.17	7.2647	264.72	0.6800	
5	12.36	7.1554	214.53	0.6800	F
6	12.50	7.0756	156.15	0.6800	B,C
7	19.92	4.4530	1000.00	0.6800	AB,C,E,F
8	20.30	4.3708	703.74	0.6800	AB,C,D,F
9	20.56	4.3164	549.47	0.6800	AB,C,F
10	20.81	4.2653	619.08	0.6800	B,C
11	20.98	4.2308	600.47	0.6800	AF
12	21.18	4.1914	485.07	0.6800	C,E
13	21.38	4.1521	412.80	0.6800	B,F
14	21.55	4.1197	351.00	0.6800	AB
15	21.72	4.0884	321.08	0.6800	C
16	21.96	4.0443	291.75	0.6800	B,C,E,F
17	22.14	4.0118	265.62	0.6800	B
18	22.46	3.9557	273.36	0.6800	AB,C,D,E,F
19	22.66	3.9211	266.66	0.6800	B,C
20	22.86	3.8871	223.54	0.6800	AE
21	23.08	3.8509	256.61	0.6800	B,C,F
22	23.56	3.7735	358.50	0.6800	B,D,F
23	23.76	3.7418	288.10	0.6800	AB
24	24.33	3.6553	423.37	0.6800	B,E,F
25	24.59	3.6180	401.59	0.6800	B,C
26	24.76	3.5925	388.79	0.6800	AB,F
27	25.05	3.5513	403.59	0.6800	B,C,D
28	25.40	3.5041	408.59	0.6800	AB,F
29	25.62	3.4743	454.40	0.6800	B,C
30	25.82	3.4478	344.76	0.6800	B
31	26.00	3.4237	274.46	0.6800	B,C,E,F
32	26.73	3.3325	946.49	0.6800	AB,E,F
33	27.16	3.2806	376.87	0.6800	B,C,E
34	27.50	3.2409	413.60	0.6800	B
35	27.70	3.2179	303.25	0.6800	AB,C,F
36	27.86	3.1998	211.01	0.6800	B
37	28.24	3.1572	165.17	0.6800	B,C,F
38	28.59	3.1200	148.06	0.6800	AB,E
39	28.98	3.0786	163.55	0.6800	B,C,E
40	29.21	3.0553	169.17	0.6800	B,C,D,F
41	29.92	2.9843	167.83	0.6800	AB,C,E,F
42	33.18	2.6981	284.88	0.6800	AB,C,D,E,F
43	34.96	2.5643	864.73	0.6800	AB,C,D,E,F
44	35.28	2.5419	541.42	0.6800	B,D,E,F
45	35.58	2.5212	569.23	0.6800	AB,C
46	35.86	2.5022	378.14	0.6800	B
47	36.00	2.4927	316.70	0.6800	B,C,D,F
48	36.16	2.4821	255.73	0.6800	AB,C,E
49	36.69	2.4471	286.39	0.6800	AB,E,F
50	37.01	2.4270	259.23	0.6800	AB,C,E,F
51	37.28	2.4100	238.96	0.6800	B,C,D,E
52	37.61	2.3895	324.68	0.6800	AB,C,D,E,F
53	37.78	2.3795	311.67	0.6800	AB,C,E,F
54	37.98	2.3695	298.66	0.6800	AB,C,E
55	38.18	2.3595	285.65	0.6800	B,C,D
56	38.38	2.3495	272.64	0.6800	AB,E,F
57	38.58	2.3395	259.63	0.6800	B,C,E
58	38.78	2.3295	246.62	0.6800	B,C,D,E,F
59	38.98	2.3195	233.61	0.6800	AB,C,D,E,F
60	39.18	2.3095	220.60	0.6800	AB,C,E,F
61	39.38	2.2995	207.59	0.6800	AB,C,D,F
62	39.58	2.2895	194.58	0.6800	AB,C,E,F
63	39.78	2.2795	181.57	0.6800	AB,C,E,F
64	39.98	2.2695	168.56	0.6800	AB,C,D,E,F
65	40.18	2.2595	155.55	0.6800	AB,C,D,E,F
66	40.38	2.2495	142.54	0.6800	AB,C,D,E,F
67	40.58	2.2395	129.53	0.6800	AB,C,D,E,F
68	40.78	2.2295	116.52	0.6800	AB,C,D,E,F
69	40.98	2.2195	103.51	0.6800	AB,C,D,E,F
70	41.18	2.2095	90.50	0.6800	AB,C,D,E,F
71	41.38	2.1995	77.49	0.6800	AB,C,D,E,F
72	41.58	2.1895	64.48	0.6800	AB,C,D,E,F
73	41.78	2.1795	51.47	0.6800	AB,C,D,E,F
74	41.98	2.1695	38.46	0.6800	AB,C,D,E,F
75	42.18	2.1595	25.45	0.6800	AB,C,D,E,F
76	42.38	2.1495	12.44	0.6800	AB,C,D,E,F
77	42.58	2.1395	0.43	0.6800	AB,C,D,E,F



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68	54.10	1.6939	319.52	0.6800	AC,D,E,F
69	54.40	1.6852	197.72	0.6800	AC,E,F
70	54.90	1.6709	280.24	0.6800	AC,E,F
71	55.02	1.6677	271.24	0.6800	AF
72	55.19	1.6630	265.72	0.6800	AD
73	55.37	1.6580	279.05	0.6800	AC,D,F
74	55.74	1.6479	221.26	0.6800	AC,E,F
75	56.01	1.6404	221.14	0.6800	AC,D,E,F
76	56.27	1.6336	206.01	0.6800	AC,E
77	56.46	1.6285	174.82	0.6800	AC,D,E
78	56.61	1.6244	176.17	0.6800	C,F
79	56.82	1.6190	144.25	0.6800	AC,D,E,F
80	57.22	1.6087	148.85	0.6800	AC,D,E,F
81	60.10	1.5383	160.33	0.6800	AC,D,E,F
82	62.02	1.4951	431.24	0.6800	AC,D,E,F
83	62.40	1.4870	492.18	0.6800	AC,D,E,F
84	62.62	1.4823	339.46	0.6800	AC,D,E,F
85	62.74	1.4797	269.32	0.6800	AC,E,F
86	64.06	1.4524	177.10	0.6800	AC,D,E,F

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

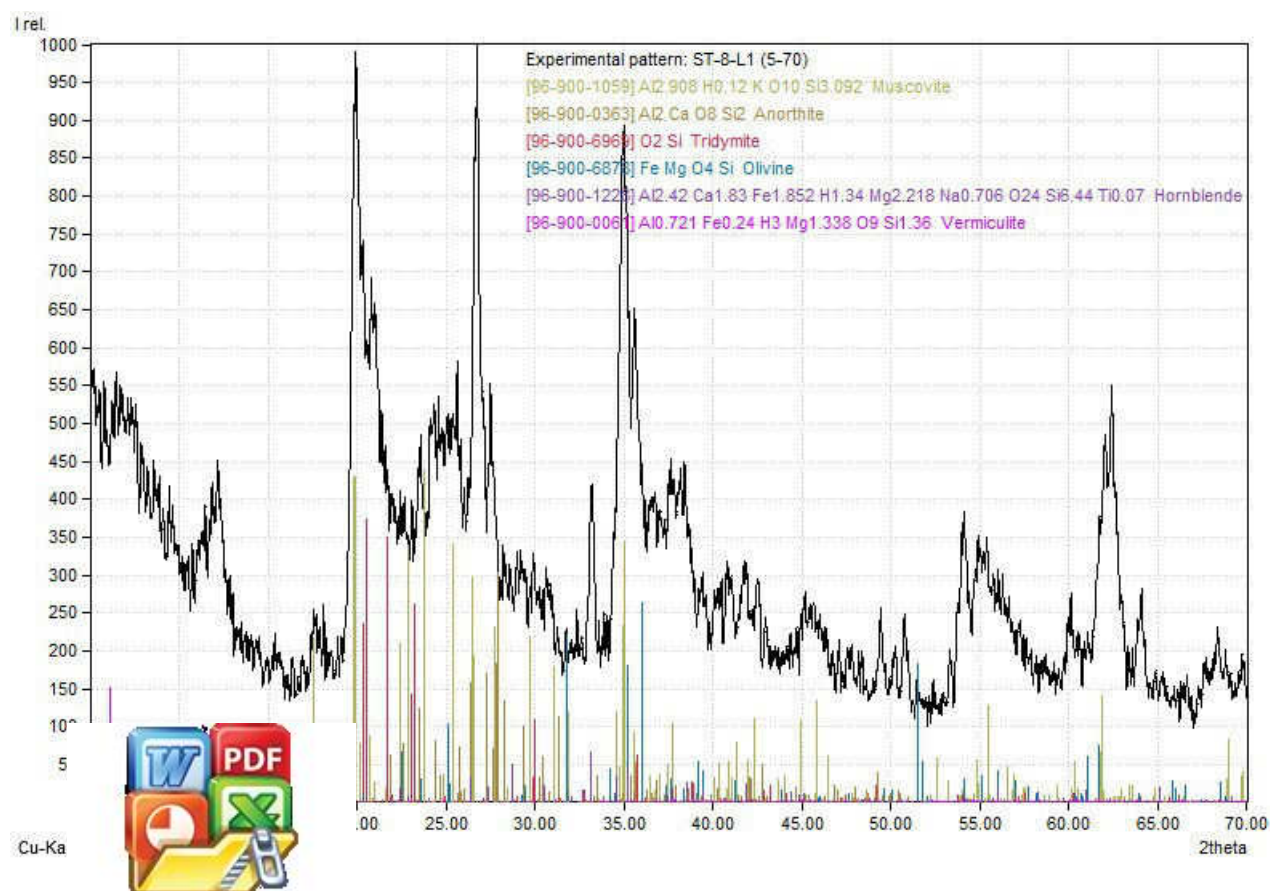
Based on calculated profile

Profile area	Counts	Amount
Overall diffraction profile	113138	100.00%
Background radiation	69300	61.25%
Diffraction peaks	43837	38.75%
Peak area belonging to selected phases	39447	34.87%
Unidentified peak area	4390	3.88%

Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	2950	100.00%
Peak intensity belonging to selected phases	1621	54.95%
Unidentified peak intensity	1329	45.05%

Diffraction Pattern Graphics



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Match! Phase Analysis Report

Sample: ST-9-L1 (5-70)

Sample Data

File name ST-9-L1.RAW
File path D:/S2 Geologi/data hasil/XRD Bu Ulva/ST-9-L1
Data collected Jan 5, 2024 16:52:42
Data range 5.110° - 70.110°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction 0.11°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	43.2	Anorthite	Al1.911 Ca0.716 Mn0.196 Na0.045 O8 Si2.089
B	42.8	Muscovite	Al3 K O12 Si3
C	5.9	Olivine	Fe Mg O4 Si
D	4.3	Vermiculite	C12 Mg3 N4 O12 Si4
E	2.0	Quartz	O2 Si
F	1.6	Montmorillonite	Al0.86 Fe0.1 H Li0.08 Mg0.14 O10 Si3.9
	0.9	Unidentified peak area	

A: Anorthite (43.2 %)

Formula sum Al1.911 Ca0.716 Mn0.196 Na0.045 O8 Si2.089
Entry number 96-900-5310
Figure-of-Merit (FoM) 0.815525
Total number of peaks 251
Peaks in range 251
Peaks matched 202
Intensity scale factor 0.55
Space group C -1
Crystal system triclinic (anorthic)
Unit cell a= 8.1310 Å b= 12.8470 Å c= 7.0690 Å α= 94.030° β= 115.900° γ= 90.720°
I/Icor 0.62
Calc. density 2.797 g/cm³
Reference Matsui T., Kimata M., "Crystal chemistry of synthetic Mn-bearing anorthite: Incorporation of MnAl2Si2O8 end-member into feldspar Note: U(1,1) for Obo altered to reproduce Uiso", European Journal of Mineralogy **9**, 333-344 (1997)

B: Muscovite (42.8 %)

Formula sum Al3 K O12 Si3
Entry number 96-901-5624
Figure-of-Merit (FoM) 0.739222
Total number of peaks 300
Peaks in range 300
Peaks matched 155
Intensity scale factor 0.37
Space group C 1 2/c 1
Crystal system monoclinic
Unit cell a= 5.1890 Å b= 8.9950 Å c= 20.0970 Å β= 95.180°
I/Icor 0.42
Calc. density 2.818 g/cm³
Reference Gatineau L., "Localisation des remplacements isomorphiques dans la muscovite Note: polytype Muscovite 2M1 _cod_database_code 1000042", Comptes Rendus Hebdomadaires des Seances de l **256**, 4648-4649 (1963)

C: Olivine (5.9 %)

Formula sum Fe Mg O4 Si
Entry number 96-900-6881
Figure-of-Merit (FoM) 0.707344
Total number of peaks 174
Peaks in range 174
Peaks matched 55
Intensity scale factor 0.15
Space group P b n m
Crystal system orthorhombic
Unit cell a= 4.8168 Å b= 10.4029 Å c= 6.0800 Å
I/Icor 0.4
Calc. density 3.54 g/cm³
Reference S. A. T., Artioli G., Rinaldi R., Henderson C. M. B., Knight K. S., Wood B. J., "Octahedral cation ordering in olivine at high temperature. II: an in situ neutron powder diffraction study on synthetic MgFeSiO4 (Fa50) Sample: T = 450 C", Physics and Chemistry of Minerals **27**, 630-637 (2000)

D: Vermiculite (4.3 %)

Formula sum C12 Mg3 N4 O12 Si4
Entry number 96-900-0119
Figure-of-Merit (FoM) 0.66380
Total number of peaks 306
Peaks in range 305
Peaks matched 145
Intensity scale factor 0.24



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intensity scale factor 0.51
 Space group C 1 2 1
 Crystal system monoclinic
 Unit cell a= 5.3300 Å b= 9.1800 Å c= 17.1200 Å β= 97.000 °
 I/Cor 3.46
 Calc. density 2.306 g/cm³
 Reference Haase D. J., Weiss E. J., Steinfink H., "The crystal structure of a hexamethylene-diamine-vermiculite complex", American Mineralogist **48**, 261-270 (1963)

E: Quartz (2.0 %)

Formula sum O2 Si
 Entry number 96-901-5023
 Figure-of-Merit (FoM) 0.687666
 Total number of peaks 35
 Peaks in range 35
 Peaks matched 13
 Intensity scale factor 0.12
 Space group P 32 2 1
 Crystal system trigonal (hexagonal axes)
 Unit cell a= 4.9290 Å c= 5.3190 Å
 I/Cor 2.83
 Calc. density 2.673 g/cm³
 Reference Gibbs G. V., Boisen M. B., Downs R. T., Lasaga A. C., "Mathematical Modeling of the structures and bulk moduli of TX2 quartz and cristobalite structure types, T = C, Si, Ge and X = O, S", Materials Research Society Symposia Proceedings **121**, 155-165 (1988)

F: Montmorillonite (1.6 %)

Formula sum Al0.86 Fe0.1 H Li0.08 Mg0.14 O10 Si3.9
 Entry number 96-901-0959
 Figure-of-Merit (FoM) 0.729550
 Total number of peaks 251
 Peaks in range 251
 Peaks matched 82
 Intensity scale factor 0.19
 Space group C 1 2/m 1
 Crystal system monoclinic
 Unit cell a= 5.1710 Å b= 8.9570 Å c= 9.7400 Å β= 96.100 °
 I/Cor 5.89
 Calc. density 2.245 g/cm³
 Reference Gournis D., Lappas A., Karakassides M. A., Tobbens D., Moukarika A., "A neutron diffraction study of alkali cation migration in montmorillonites Sample: Li-mont-300", Physics and Chemistry of Minerals **35**, 49-58 (2008)

Candidates

Name	Formula	Entry No.	FoM
Quartz	O2 Si	96-900-7379	0.6595
Quartz	O2 Si	96-901-5023	0.6564
Quartz	O2 Si	96-901-2603	0.6563
Quartz	O2 Si	96-900-5025	0.6540
Quartz	O2 Si	96-900-5022	0.6521
Quartz	O2 Si	96-900-5023	0.6518
Quartz	O2 Si	96-900-5021	0.6463
Quartz	O2 Si	96-900-5024	0.6462
Quartz	O2 Si	96-901-0145	0.6384
Quartz	O2 Si	96-901-1494	0.6384
Quartz	O2 Si	96-900-5020	0.6336
Quartz	O2 Si	96-901-2604	0.6330
Quartz	O2 Si	96-900-0777	0.6325
Quartz	O2 Si	96-900-5018	0.6317
Quartz	O2 Si	96-900-5019	0.6317
Quartz	O2 Si	96-900-0776	0.6314
Quartz	O2 Si	96-901-0147	0.6309
Quartz	O2 Si	96-901-0146	0.6283
Quartz	O2 Si	96-901-2605	0.6169
Quartz	O2 Si	96-901-2606	0.6124

Search-Match

Settings

Reference database used COD-Inorg REV89244 2013.10.11
 Automatic zeropoint adaptation Yes
 Minimum figure-of-merit (FoM) 0.60
 Parameter/influence 2theta 0.50
 Parameter/influence 2theta 0.50
 Parameter/influence 2theta 0



Compi

Name Quartz

Selection Criteria

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No.		I/I0	FWHM	Matched	
1	6.08 14.5268	177.08	1.2400	D	
2	6.21 14.2267	188.93	1.2400		
3	6.74 13.1056	213.63	1.2400		

Peak List

4	7.07	12.4931	255.81	1.2400	B,F
5	12.34	7.1649	170.69	1.2400	AD
6	19.91	4.4558	929.05	1.2400	AB,D,F
7	20.11	4.4120	800.77	1.2400	D
8	20.27	4.3775	715.78	1.2400	B,C
9	20.46	4.3373	614.99	1.2400	A
10	20.65	4.2978	554.74	1.2400	B,D,E
11	21.07	4.2126	711.02	1.2400	D,F
12	21.27	4.1739	628.86	1.2400	
13	21.45	4.1393	536.86	1.2400	
14	21.61	4.1090	471.35	1.2400	B
15	21.92	4.0515	373.35	1.2400	D,F
16	22.16	4.0077	328.83	1.2400	A
17	22.37	3.9706	309.02	1.2400	B
18	22.57	3.9366	375.13	1.2400	C
19	22.76	3.9038	370.39	1.2400	AF
20	23.15	3.8386	289.77	1.2400	B,D
21	23.64	3.7605	767.99	1.2400	AC,D
22	23.89	3.7217	433.06	1.2400	AB
23	24.08	3.6930	326.45	1.2400	
24	24.35	3.6532	436.50	1.2400	A
25	24.70	3.6018	436.58	1.2400	AB
26	24.98	3.5612	354.42	1.2400	C,D
27	25.43	3.5002	530.88	1.2400	AB
28	25.79	3.4522	586.60	1.2400	AF
29	26.23	3.3943	329.45	1.2400	AD
30	26.96	3.3042	949.53	1.2400	AB,D,E,F
31	27.48	3.2432	1000.00	1.2400	A
32	27.65	3.2236	957.25	1.2400	B,F
33	27.83	3.2031	759.14	1.2400	A
34	28.44	3.1356	234.99	1.2400	AF
35	28.70	3.1077	206.26	1.2400	AB,D
36	29.02	3.0741	257.11	1.2400	
37	29.17	3.0590	193.97	1.2400	AC
38	29.95	2.9814	466.45	1.2400	B,D
39	30.36	2.9414	239.83	1.2400	A
40	30.92	2.8901	301.36	1.2400	A
41	31.11	2.8725	176.71	1.2400	AB,C,D,F
42	34.95	2.5653	892.57	1.2400	AB,C,D,F
43	35.19	2.5483	977.11	1.2400	AB,C,D,F
44	35.66	2.5154	576.14	1.2400	AD,F
45	35.97	2.4948	367.97	1.2400	AB,C
46	36.15	2.4830	334.55	1.2400	AB,D
47	36.35	2.4695	269.24	1.2400	AE
48	36.55	2.4565	247.29	1.2400	AB,D,F
49	36.95	2.4307	278.37	1.2400	AD
50	37.23	2.4133	286.22	1.2400	AB,C,D,F
51	37.43	2.4008	340.95	1.2400	
52	37.72	2.3828	375.55	1.2400	AB,C,D,F
53	37.94	2.3699	337.79	1.2400	A
54	38.32	2.3473	276.30	1.2400	AB,C,D
55	38.84	2.3166	245.37	1.2400	AB,C,D,F
56	39.85	2.2606	165.73	1.2400	AB,D,E
57	40.88	2.2055	182.31	1.2400	AB,D,F
58	41.05	2.1971	182.14	1.2400	B,C,D,F
59	41.81	2.1590	337.03	1.2400	AB,C,D,E,F
60	50.94	1.7912	251.80	1.2400	AB,C,D,E,F
61	54.07	1.6947	246.21	1.2400	AB,C,D,F
62	54.38	1.6858	185.36	1.2400	AB,D,F
63	55.11	1.6653	273.20	1.2400	AB,C,D,E
64	55.32	1.6592	244.88	1.2400	AB,C,D,F
65	55.57	1.6524	204.16	1.2400	B,D
66	55.83	1.6453	252.25	1.2400	B,C,D,F
67	56.00	1.6406	258.16	1.2400	AB,D,E
68	56.29	1.6330	198.61	1.2400	AB,C,D,F
69	56.61	1.6246	219.65	1.2400	AB,D
70	56.83	1.6187	198.85	1.2400	AB,C,D,E,F
71	60.56	1.5278	168.21	1.2400	AB,C,D,E,F
72	60.75	1.5233	172.61	1.2400	AB,C,D,F
73	61.16	1.5141	208.19	1.2400	AB,C,D,F
74	61.58	1.5048	342.12	1.2400	AC,D
75	61.76	1.5009	341.86	1.2400	AB
76	61.94	1.4969	362.23	1.2400	AB,D
77			1.25	1.2400	AB,D,F
78			1.45	1.2400	AB,F
79			1.76	1.2400	AC,D,F
80			1.90	1.2400	AB,D
81			1.54	1.2400	AB,D,F
82			1.50	1.2400	AB,C,D,E,F



Calculated using FullProf

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Rietveld Refinement using FullProf

Integrated Profile Areas

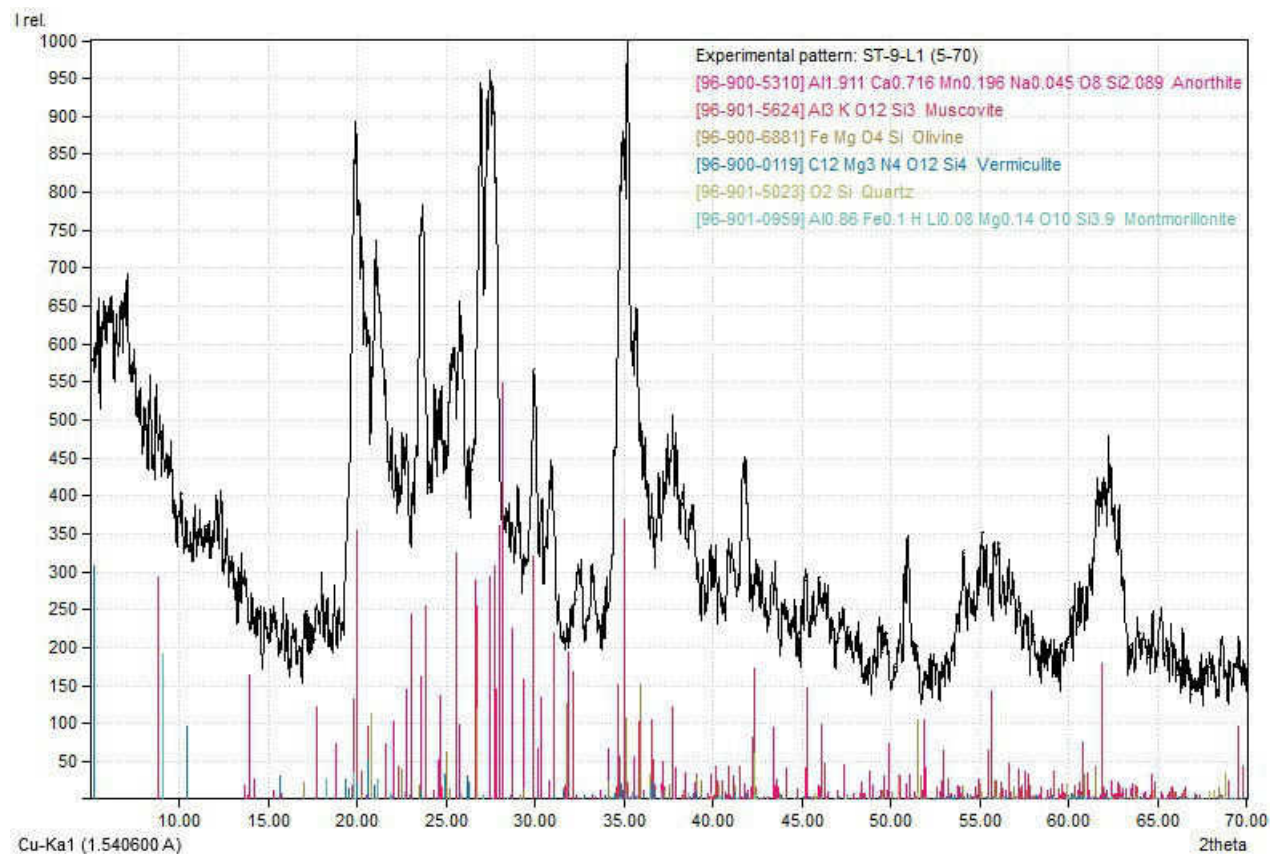
Based on calculated profile

Profile area	Counts	Amount
Overall diffraction profile	163805	100.00%
Background radiation	100675	61.46%
Diffraction peaks	63130	38.54%
Peak area belonging to selected phases	61658	37.64%
Unidentified peak area	1472	0.90%

Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	6312	100.00%
Peak intensity belonging to selected phases	3143	49.80%
Unidentified peak intensity	3169	50.20%

Diffraction Pattern Graphics



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Match! Phase Analysis Report

Sample: ST-10-L1 (5-70)

Sample Data

File name ST-10-L1.RAW
File path D:/S2 Geologi/data hasil/XRD Bu Ulva/ST-10-L1
Data collected Jan 5, 2024 16:52:42
Data range 4.950° - 69.950°
Number of points 3251
Step size 0.020
Rietveld refinement converged No
Alpha2 subtracted No
Background subtr. No
Data smoothed Yes
2theta correction -0.05°
Radiation X-rays
Wavelength 1.540600 Å

Matched Phases

Index	Amount (%)	Name	Formula sum
A	39.8	Muscovite	Al3 H2 K O12 Si3
B	33.6	Anorthite	Al2 Ca O8 Si2
C	10.9	Olivine	Fe Mg O4 Si
D	7.7	Quartz	O2 Si
E	6.9	Biotite	Al Fe H2 K Mg2 O12 Si3
F	0.6	Montmorillonite	Al2 Ca O12 Si4
G	0.6	Vermiculite	H2 Mg3 O12 Si4
	5.9	Unidentified peak area	

A: Muscovite (39.8 %)

Formula sum Al3 H2 K O12 Si3
Entry number 96-900-0838
Figure-of-Merit (FoM) 0.801323
Total number of peaks 300
Peaks in range 300
Peaks matched 168
Intensity scale factor 0.38
Space group C 1 2/c 1
Crystal system monoclinic
Unit cell a= 5.1988 Å b= 9.0266 Å c= 20.1058 Å β= 95.782 °
I/lor 0.41
Calc. density 2.818 g/cm³
Reference Richardson S. M., Richardson J. W., "Crystal structure of a pink muscovite from Archer's Post, Kenya: Implications for reverse pleochroism in dioctahedral micas", American Mineralogist **67**, 69-75 (1982)

B: Anorthite (33.6 %)

Formula sum Al2 Ca O8 Si2
Entry number 96-900-0362
Figure-of-Merit (FoM) 0.856007
Total number of peaks 500
Peaks in range 500
Peaks matched 422
Intensity scale factor 0.48
Space group P -1
Crystal system triclinic (anorthic)
Unit cell a= 8.1940 Å b= 12.8970 Å c= 14.1900 Å α= 92.980° β= 115.820° γ= 91.150°
I/lor 0.60
Calc. density 2.745 g/cm³
Reference Foit F. F., Peacor D. R., "The anorthite crystal structure at 410 and 830 C T = 410 C", American Mineralogist **58**, 665-675 (1973)

C: Olivine (10.9 %)

Formula sum Fe Mg O4 Si
Entry number 96-900-6893
Figure-of-Merit (FoM) 0.758596
Total number of peaks 177
Peaks in range 177
Peaks matched 61
Intensity scale factor 0.28
Space group P b n m
Crystal system orthorhombic
Unit cell a= 4.8497 Å b= 10.5034 Å c= 6.1418 Å
I/lor 0.8
Calc. density 3.57 g/cm³
Reference S. A. T., Artioli G., Rinaldi R., Henderson C. M. B., Knight K. S., Wood B. J., "Octahedral cation ordering in olivine at high temperature. II: an in situ neutron powder diffraction study on synthetic MgFeSiO4 (Fa50) Sample: T = 1250 C", Physics and Chemistry of Minerals **27**, 630-637 (2000)

D: Quartz

Formula Si
Entry number 96-900-0778
Figure-of-Merit (FoM) 10301
Total number of peaks 34
Peaks in range 34
Peaks matched 13
Intensity scale factor 0.40



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Intensity scale factor	0.49
Space group	P 32 2 1
Crystal system	trigonal (hexagonal axes)
Unit cell	a= 4.7736 Å c= 5.3010 Å
I/lcor	2.70
Calc. density	2.861 g/cm ³
Reference	Levien L., Prewitt C. T., Weidner D. J., "Structure and elastic properties of quartz at pressure P = 37.6 kbar", American Mineralogist 65 , 920-930 (1980)

E: Biotite (6.9 %)

Formula sum	Al Fe H2 K Mg2 O12 Si3
Entry number	96-900-1266
Figure-of-Merit (FoM)	0.767219
Total number of peaks	279
Peaks in range	279
Peaks matched	95
Intensity scale factor	0.28
Space group	C 1 2/m 1
Crystal system	monoclinic
Unit cell	a= 5.3460 Å b= 9.2520 Å c= 10.2380 Å β= 100.020 °
I/lcor	1.74
Calc. density	2.987 g/cm ³
Reference	Brigatti M. F., Davoli P., "Crystal-structure refinement of 1M plutonic biotites sample M32 from a syenite in the Valle del Cervo pluton", American Mineralogist 75 , 305-313 (1990)

F: Montmorillonite (0.6 %)

Formula sum	Al2 Ca O12 Si4
Entry number	96-110-1055
Figure-of-Merit (FoM)	0.731583
Total number of peaks	92
Peaks in range	92
Peaks matched	87
Intensity scale factor	0.28
Space group	P 1
Crystal system	triclinic (anorthic)
Unit cell	a= 5.1800 Å b= 8.9800 Å c= 15.0000 Å α= 90.000° β= 90.000° γ= 90.000°
I/lcor	20.53
Calc. density	1.800 g/cm ³

G: Vermiculite (0.6 %)

Formula sum	H2 Mg3 O12 Si4
Entry number	96-900-0010
Figure-of-Merit (FoM)	0.780436
Total number of peaks	262
Peaks in range	262
Peaks matched	160
Intensity scale factor	0.31
Space group	C 1 c 1
Crystal system	monoclinic
Unit cell	a= 5.3100 Å b= 9.2000 Å c= 28.4600 Å β= 97.140 °
I/lcor	23.48
Calc. density	1.826 g/cm ³
Reference	Gruner J. W., "The structures of vermiculites and their collapse by dehydration Locality: Structure results from data of many samples", American Mineralogist 19 , 557-575 (1934)

Candidates

Name	Formula	Entry No.	FoM
Quartz	O2 Si	96-900-0776	0.7137
Quartz	O2 Si	96-900-0777	0.6433
Quartz	O2 Si	96-900-0778	0.6698
Quartz	O2 Si	96-900-0779	0.7095
Quartz	O2 Si	96-900-0780	0.7092
Quartz	O2 Si	96-900-0781	0.6925
Quartz	O2 Si	96-900-5018	0.7126
Quartz	O2 Si	96-900-5019	0.6984
Quartz	O2 Si	96-900-5020	0.6946
Quartz	O2 Si	96-900-5021	0.6942
Quartz	O2 Si	96-900-5022	0.6785
Quartz	O2 Si	96-900-5023	0.7058
Quartz	O2 Si	96-900-5024	0.6983
Quartz	O2 Si	96-900-5025	0.7118
Quartz	O2 Si	96-900-7379	0.7217
Quartz	O2 Si	96-900-9667	0.7081
Quartz	O2 Si	96-901-0145	0.7090
Quartz	O2 Si	96-901-0146	0.7008
Quartz	O2 Si	96-901-0147	0.6971
Quartz	O2 Si	96-901-1494	0.6865
Quartz	O2 Si	96-901-1495	0.6770
Quartz	O2 Si	96-901-1496	0.6754
Quartz	O2 Si	96-901-1497	0.6109
Quartz	O2 Si	96-901-2601	0.7126
Quartz	O2 Si	96-901-2602	0.6917
Quartz	O2 Si	96-901-2603	0.6680
Quartz	O2 Si	96-901-2604	0.6741
Quartz	O2 Si	96-901-2605	0.6456
Quartz	O2 Si	96-901-2606	0.6306
Quartz	O2 Si	96-901-3322	0.7129



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Quartz	O2 Si	96-901-5023	0.7414
Silicon oxide (Quartz)	O2 Si	96-500-0036	0.7131

Search-Match

Settings

Reference database used	COD-Inorg REV89244 2013.10.11
Automatic zeropoint adaptation	Yes
Minimum figure-of-merit (FoM)	0.60
Parameter/influence 2theta	0.50
Parameter/influence intensities	0.50
Parameter multiple/single phase(s)	0.50

Selection Criteria

Compound:

Name: quartz

Peak List

No.	2theta [°]	d [Å]	I/I0	FWHM	Matched
1	5.07	17.4229	168.21	0.5600	
2	5.28	16.7106	242.33	0.5600	
3	5.56	15.8915	195.80	0.5600	
4	5.70	15.4986	181.26	0.5600	F
5	6.12	14.4295	242.31	0.5600	G
6	6.39	13.8205	277.01	0.5600	
7	6.54	13.4958	247.34	0.5600	
8	6.84	12.9213	271.54	0.5600	B
9	7.19	12.2803	242.87	0.5600	
10	7.33	12.0505	241.40	0.5600	
11	7.50	11.7778	242.37	0.5600	
12	7.85	11.2467	222.37	0.5600	
13	8.12	10.8831	190.10	0.5600	
14	8.53	10.3627	190.95	0.5600	
15	8.65	10.2171	190.46	0.5600	
16	8.88	9.9539	201.22	0.5600	AE
17	9.28	9.5260	182.00	0.5600	B
18	10.59	8.3497	198.38	0.5600	B
19	10.78	8.1973	243.44	0.5600	B
20	11.19	7.9013	190.20	0.5600	
21	11.64	7.5958	283.01	0.5600	
22	11.85	7.4607	334.04	0.5600	F
23	12.06	7.3320	367.92	0.5600	B
24	12.27	7.2077	293.10	0.5600	
25	12.46	7.1006	269.91	0.5600	G
26	12.69	6.9701	212.05	0.5600	B
27	13.65	6.4807	274.97	0.5600	B
28	17.59	5.0385	169.30	0.5600	B,C,E
29	17.77	4.9865	161.65	0.5600	AB,F
30	18.17	4.8797	189.15	0.5600	B
31	18.30	4.8453	190.62	0.5600	
32	18.46	4.8037	184.49	0.5600	B,G
33	19.80	4.4804	1000.00	0.5600	AB,E,F,G
34	19.93	4.4505	971.94	0.5600	AB,E
35	20.26	4.3789	875.26	0.5600	AB,C,G
36	20.53	4.3226	670.72	0.5600	AB,F,G
37	20.84	4.2585	708.64	0.5600	B
38	21.05	4.2166	675.52	0.5600	AE,G
39	21.28	4.1714	644.71	0.5600	B
40	21.51	4.1279	541.26	0.5600	AB,D,G
41	21.88	4.0588	478.39	0.5600	B
42	22.05	4.0277	438.65	0.5600	B,G
43	22.30	3.9839	393.72	0.5600	AB,C
44	22.57	3.9358	493.30	0.5600	B,E,G
45	22.78	3.8998	365.43	0.5600	B
46	22.97	3.8682	356.15	0.5600	AB,F,G
47	23.22	3.8277	403.56	0.5600	B,C
48	23.57	3.7709	474.25	0.5600	B,G
49	23.81	3.7341	360.86	0.5600	AF
50	24.05	3.6974	419.27	0.5600	B,E,G
51	24.60	3.6159	586.58	0.5600	B
52			3.35	0.5600	AB,C
53			3.77	0.5600	B,G
54			2.84	0.5600	AB,G
55			2.11	0.5600	B
56			3.19	0.5600	
57			3.42	0.5600	B,E,G
58			3.12	0.5600	AE,F
59			1.33	0.5600	G
60			2.14	0.5600	B
61			3.78	0.5600	B,D
62			3.21	0.5600	B
63			1.04	0.5600	AB,G
64	28.18	3.1640	225.06	0.5600	B,E
65	28.59	3.1194	200.82	0.5600	AB,G
66	28.73	3.1048	186.57	0.5600	B



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67	28.95	3.0818	222.84	0.5600	B,C
68	29.11	3.0651	186.01	0.5600	B,C
69	29.45	3.0310	193.75	0.5600	B,G
70	29.58	3.0176	179.40	0.5600	
71	29.77	2.9988	272.14	0.5600	AB,F
72	29.95	2.9812	210.11	0.5600	B
73	30.16	2.9607	157.21	0.5600	B,G
74	30.54	2.9249	185.58	0.5600	B,E
75	30.78	2.9027	244.37	0.5600	AB,C,F,G
76	32.92	2.7183	156.97	0.5600	AB,E,G
77	33.26	2.6919	183.95	0.5600	AB,C,E,G
78	34.97	2.5636	806.87	0.5600	AB,C,F,G
79	35.17	2.5496	603.99	0.5600	AB,F
80	35.35	2.5368	586.72	0.5600	B,E,G
81	35.63	2.5176	595.54	0.5600	AB,C,E
82	36.04	2.4901	488.55	0.5600	AB,F
83	36.43	2.4643	343.97	0.5600	AB,G
84	36.62	2.4520	309.31	0.5600	B,F
85	36.84	2.4380	362.24	0.5600	AB,E,G
86	37.15	2.4180	432.63	0.5600	AB,C,G
87	37.60	2.3902	424.85	0.5600	AB,D
88	37.89	2.3728	408.82	0.5600	B,G
89	38.07	2.3620	396.20	0.5600	AB,C
90	38.25	2.3511	384.02	0.5600	B,G
91	38.61	2.3302	421.41	0.5600	B,C
92	38.79	2.3196	320.20	0.5600	AB
93	38.99	2.3080	334.87	0.5600	B,C,E
94	39.20	2.2963	239.90	0.5600	B,F,G
95	39.40	2.2851	198.85	0.5600	B,E
96	39.59	2.2746	175.28	0.5600	B,E,G
97	39.73	2.2668	175.81	0.5600	AB
98	39.92	2.2564	172.48	0.5600	AB,E,G
99	40.27	2.2378	222.66	0.5600	AB,F,G
100	40.49	2.2258	222.75	0.5600	AB,D
101	40.67	2.2165	204.81	0.5600	E,F,G
102	40.88	2.2057	249.87	0.5600	AB,C,G
103	41.13	2.1931	201.48	0.5600	AB,E,F,G
104	41.65	2.1665	416.97	0.5600	B,D,E,G
105	41.97	2.1508	177.58	0.5600	AB,F,G
106	42.24	2.1379	196.55	0.5600	AB,F
107	42.42	2.1293	185.20	0.5600	AB,F,G
108	42.71	2.1153	170.47	0.5600	AB,C,E,G
109	44.23	2.0461	169.01	0.5600	AB,C,D,E,F,G
110	44.56	2.0316	177.97	0.5600	AB
111	44.71	2.0253	197.05	0.5600	B
112	45.02	2.0121	198.72	0.5600	B,E,G
113	45.17	2.0056	213.18	0.5600	AB,E,G
114	45.37	1.9974	185.22	0.5600	AB,C,G
115	45.81	1.9793	214.70	0.5600	AB,E,G
116	46.09	1.9679	187.11	0.5600	AB
117	46.26	1.9611	163.15	0.5600	B,E,F,G
118	46.52	1.9506	179.49	0.5600	AB,G
119	46.69	1.9440	175.02	0.5600	AB
120	47.09	1.9285	174.68	0.5600	AB,C,D,F,G
121	47.35	1.9183	157.42	0.5600	B,E
122	47.55	1.9107	186.13	0.5600	B
123	47.78	1.9022	157.40	0.5600	AB,C,E,F,G
124	49.29	1.8473	163.23	0.5600	AB,C,E,G
125	50.71	1.7990	179.19	0.5600	B,F,G
126	50.88	1.7932	188.69	0.5600	AB,C,D,E,G
127	53.28	1.7179	165.00	0.5600	AC,E,F,G
128	53.45	1.7130	163.97	0.5600	AC,G
129	53.74	1.7044	253.39	0.5600	AE,G
130	53.94	1.6986	280.24	0.5600	AF
131	54.10	1.6938	290.29	0.5600	AF,G
132	54.41	1.6848	305.52	0.5600	AC,E,F,G
133	54.77	1.6746	294.43	0.5600	AG
134	54.93	1.6701	303.26	0.5600	C,F,G
135	55.16	1.6639	315.99	0.5600	AC,E
136	55.45	1.6558	266.46	0.5600	AE,F,G
137	55.70	1.6489	266.09	0.5600	AF
138	55.87	1.6443	213.86	0.5600	AC,G
139	56.13	1.6374	238.76	0.5600	AE
140			3.95	0.5600	AC,D,G
141			2.44	0.5600	AD,G
142			3.85	0.5600	AE,G
143			0.75	0.5600	C
144			0.82	0.5600	AF,G
145			3.11	0.5600	AC,E,F,G
146			0.69	0.5600	AC,D,E,F,G
147			1.09	0.5600	AC
148			7.58	0.5600	AC,E,G
149			5.95	0.5600	AC,E,F,G
150			2.83	0.5600	AE
151			5.60	0.5600	AC,D,F,G
152	62.17	1.4919	501.98	0.5600	G
153	62.37	1.4877	399.06	0.5600	AF
154	62.63	1.4821	265.26	0.5600	AE,G



155	62.80	1.4785	252.14	0.5600	A,E,F,G
156	63.14	1.4713	195.47	0.5600	A,C,E,F,G
157	64.08	1.4519	158.60	0.5600	A,C,E,F,G

Rietveld Refinement using FullProf

Calculation was not run or did not converge.

Integrated Profile Areas

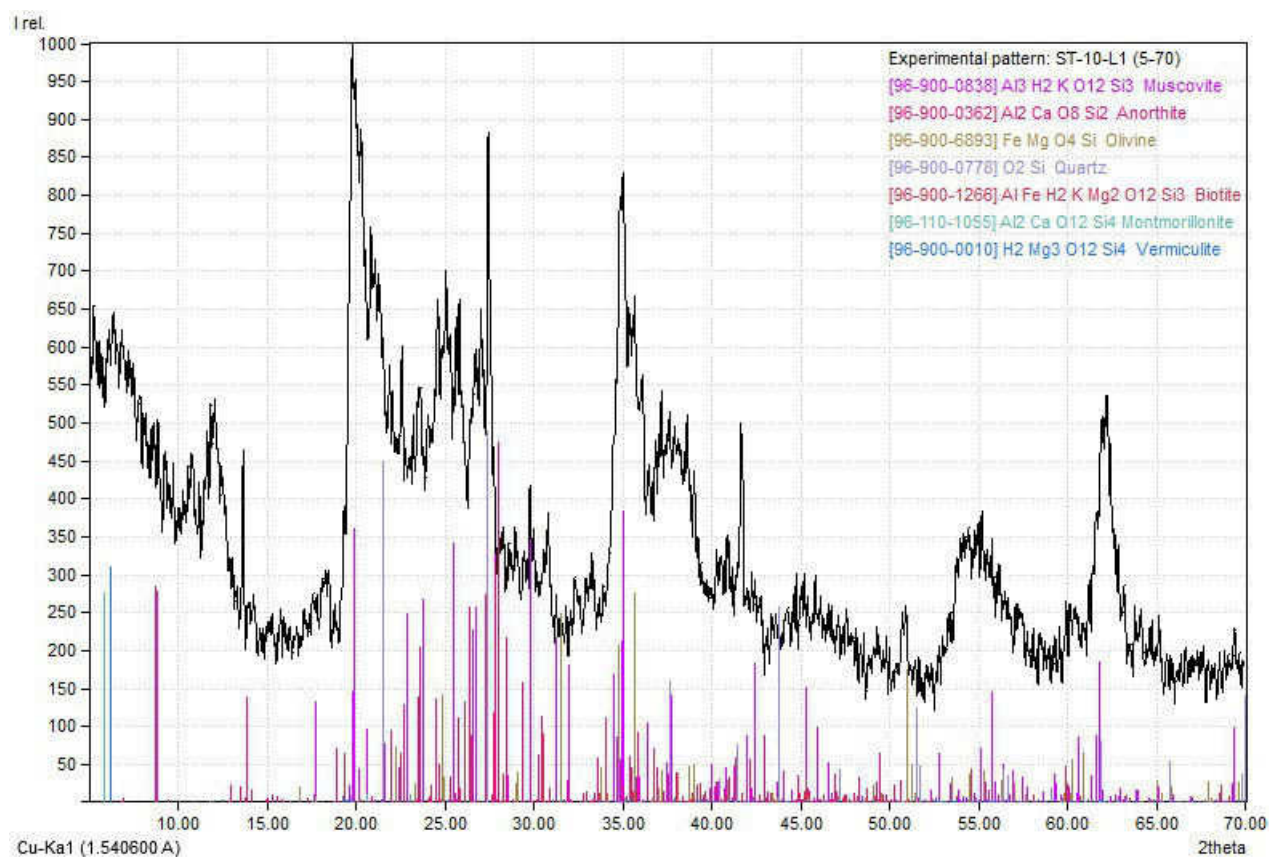
Based on calculated profile

Profile area	Counts	Amount
Overall diffraction profile	124378	100.00%
Background radiation	79458	63.88%
Diffraction peaks	44920	36.12%
Peak area belonging to selected phases	37575	30.21%
Unidentified peak area	7345	5.91%

Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	4256	100.00%
Peak intensity belonging to selected phases	1658	38.96%
Unidentified peak intensity	2598	61.04%

Diffraction Pattern Graphics



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FINAL REPORT

Ref : F5501_Rev01

Job : 232320

SAMPLE	Al2O3	CaO	Cr2O3	Fe2O3	K2O	MgO	MnO	Na2O	P2O5
ST1L1	26.84	0.22	0.04	13.96	0.19	0.71	0.12	0.15	0.482
ST1L5	30.55	0.06	0.02	13.67	0.16	0.79	0.12	0.12	0.517
ST2L1	25.44	1.37	0.01	10.05	0.95	1.82	0.11	0.68	0.520
ST2L2	30.74	0.27	0.01	11.53	0.17	0.71	0.08	0.16	0.499
ST3L1	27.05	0.63	0.03	13.68	0.42	1.19	0.09	0.24	0.535
ST3L3	31.44	0.98	0.02	14.61	0.43	1.85	0.11	0.23	0.659
ST4L1	28.18	0.10	0.02	13.04	0.55	1.27	0.11	0.28	0.408
ST4L4	29.81	0.16	<0.01	11.93	0.82	1.10	0.21	0.33	0.260
ST5L1	17.80	1.14	0.01	9.19	0.31	1.60	0.11	0.25	0.649
ST5L3	27.41	1.59	<0.01	12.14	0.41	3.41	0.16	0.42	0.747
UNI	%	%	%	%	%	%	%	%	%
DET	0.01	0.01	0.01	0.01	0.01	0.01	0.01	0.01	0.002
SCH	XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250



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FINAL REPORT

Ref : F5501_Rev01

Job : 232320

UNI			%	%	%	%	%			
DET			0.01	0.01	0.002	0	0.01			
SCH			XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	LOI			

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Job Number: 232320
Client Ref: F5501_Rev01

FINAL REPORT

Ref : F5588

Job : 232624

SAMPLE	Al2O3	CaO	Cr2O3	Fe2O3	K2O	MgO	MnO	Na2O	P2O5
ST6L1	24.34	0.16	0.01	11.15	3.79	1.30	0.31	0.15	0.201
ST6L3	26.47	0.03	<0.01	9.08	3.59	1.31	0.24	0.16	0.116
ST7L1	21.78	1.90	0.01	13.11	4.49	1.93	0.36	0.26	0.279
ST7L3	22.99	1.77	0.02	13.45	4.74	2.22	0.37	0.25	0.279
ST8L1	28.15	0.30	<0.01	9.56	3.49	0.87	0.33	0.27	0.213
ST8L3	26.24	0.05	<0.01	13.27	1.60	1.32	0.88	0.09	0.218
ST9L1	23.68	1.32	<0.01	9.08	4.65	1.62	0.38	0.30	0.263
ST9L2	24.58	1.24	0.02	9.27	4.47	1.70	0.38	0.31	0.286
ST10L1	25.96	0.31	<0.01	11.66	2.64	1.13	0.40	0.22	0.234
ST10L3	29.92	0.18	0.02	13.02	1.63	1.11	0.43	0.07	0.209
UNI	%	%	%	%	%	%	%	%	%
DET	0.01	0.01	0.01	0.01	0.01	0.01	0.01	0.01	0.002
SCH	XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250



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Ref : F5588

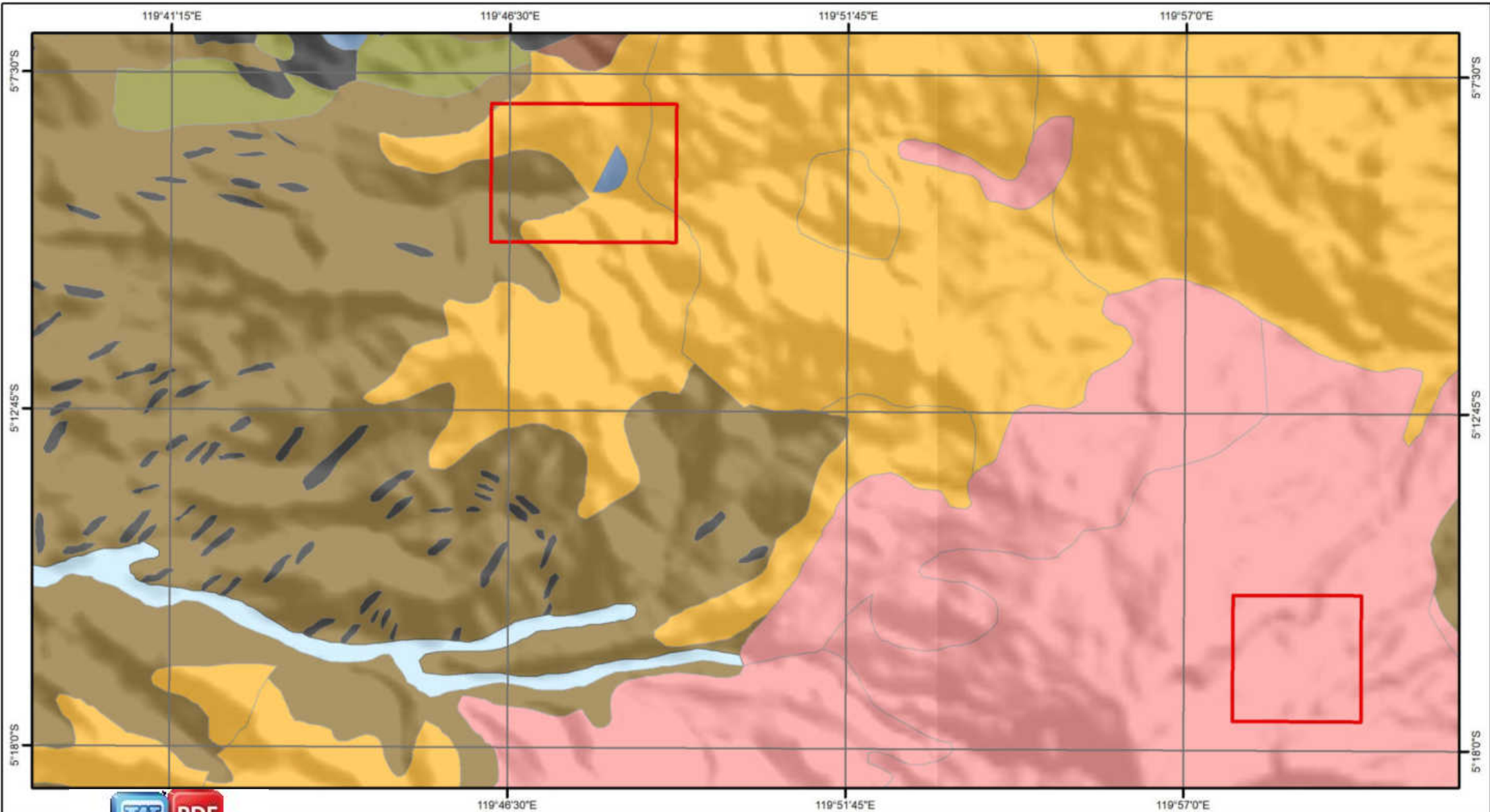
Job : 232624

SAMPLE	SiO2	TiO2	S	Total	LOI				
ST6L1	42.82	1.07	0.005	99.6	14.24				
ST6L3	48.29	1.03	<0.002	100.0	9.85				
ST7L1	43.28	1.16	0.008	100.0	11.50				
ST7L3	44.63	1.22	<0.002	100.0	8.24				
ST8L1	45.32	1.03	<0.002	100.0	10.59				
ST8L3	44.73	1.24	<0.002	100.0	10.56				
ST9L1	46.02	1.00	0.004	99.4	11.11				
ST9L2	46.55	1.04	<0.002	100.0	10.06				
ST10L1	41.80	1.12	<0.002	100.0	14.51				
ST10L3	38.81	1.13	<0.002	100.0	13.69				
UNI	%	%	%	%	%				
DET	0.01	0.01	0.002	0	0.01				
SCH	XRF250	FB1/XRF250	FB1/XRF250	FB1/XRF250	LOI				



This

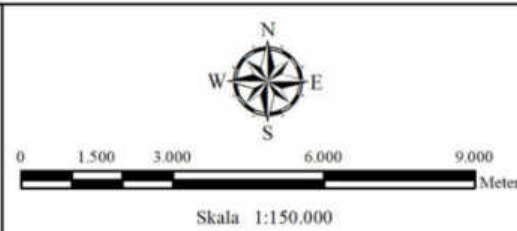
reproduced or presented in full.



**GIONAL
AN**

SIR

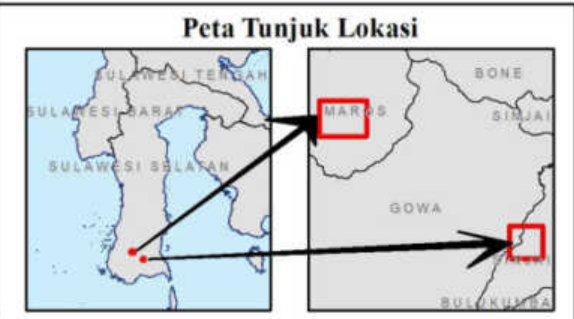
GOWA
2024

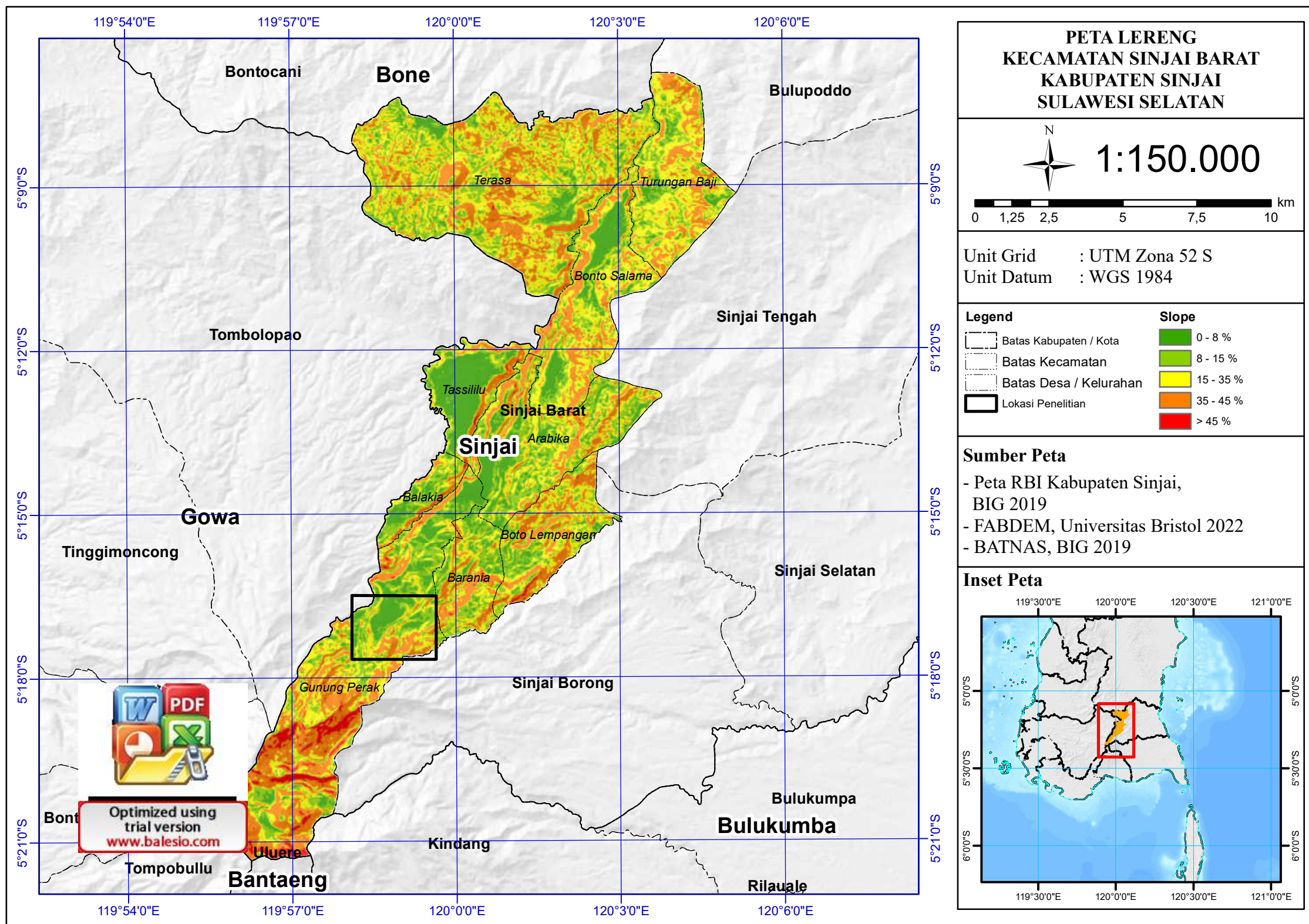


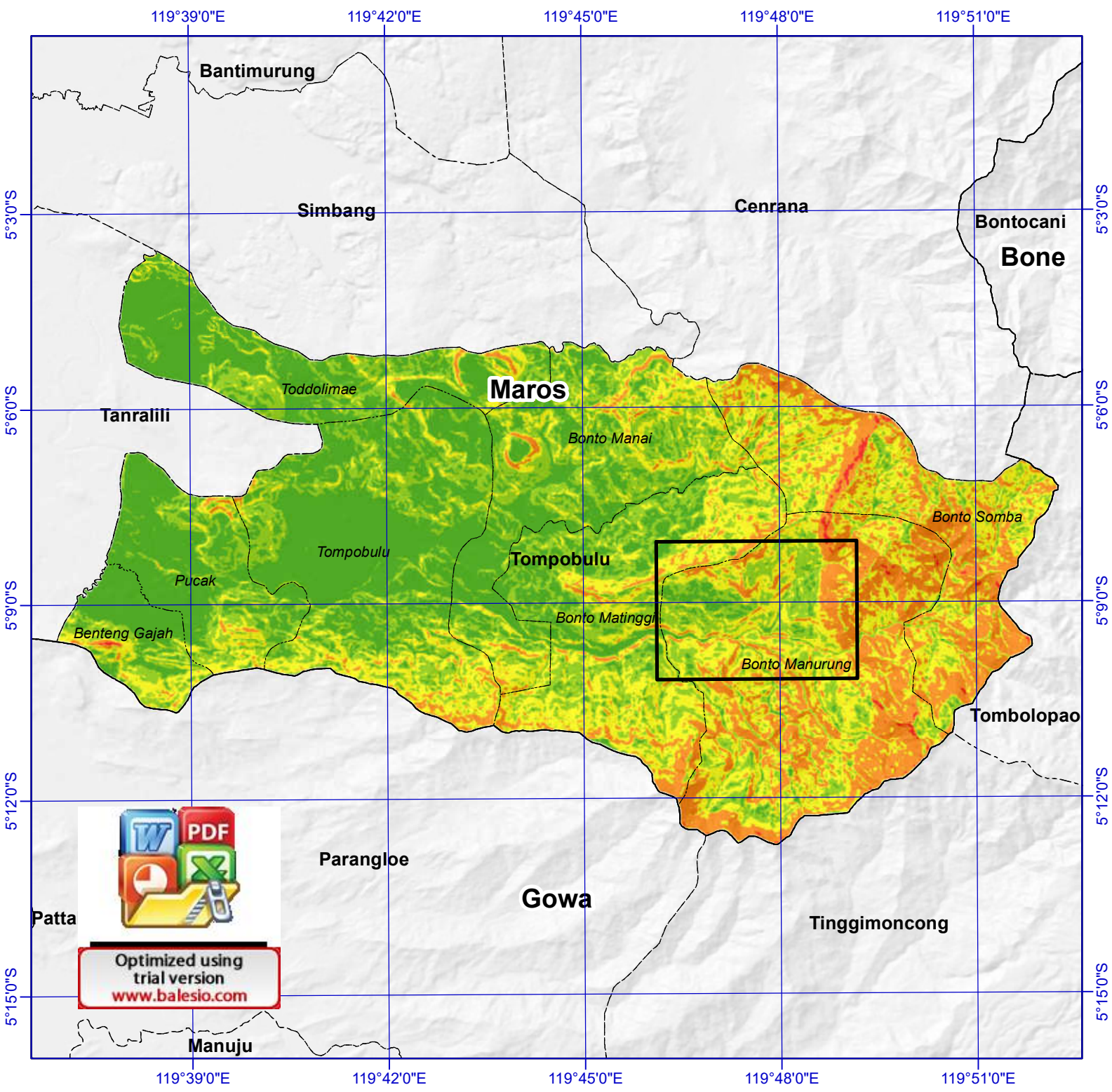
Proyeksi : Universal Traver Mercator
Datum : Datum Geodetik Nasional (DGN) 1995
Koordinat : Geografis

- Legenda :**
- | | |
|-----------------------------------|-----------------------|
| Formasi Camba | Endapan Aluvium |
| Formasi Tonasa | Basal dan Retas Basal |
| Batuan Gunungapi Lompobatang | |
| Batuan Gunungapi Baturape-Cindako | |
| Diorit | |
| Batas Wilayah Penelitian | |

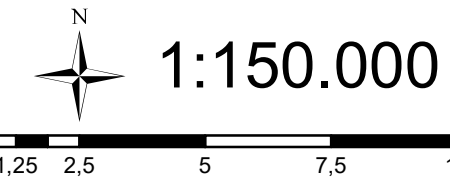
Sumber :
Peta Geologi Lembang, Ujungpandang, Benteng dan Simai, Sulawesi, Rabi Sukanto dan Sam Soeparto (1982)







**PETA LERENG
KECAMATAN TOMPOBULU
KABUPATEN MAROS
SULAWESI SELATAN**

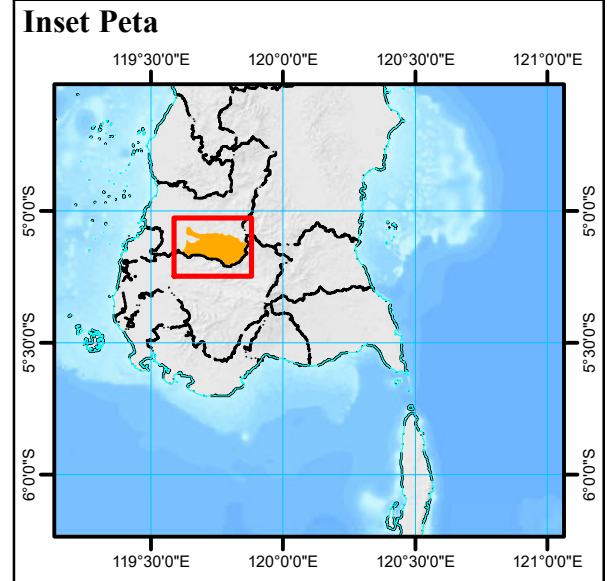


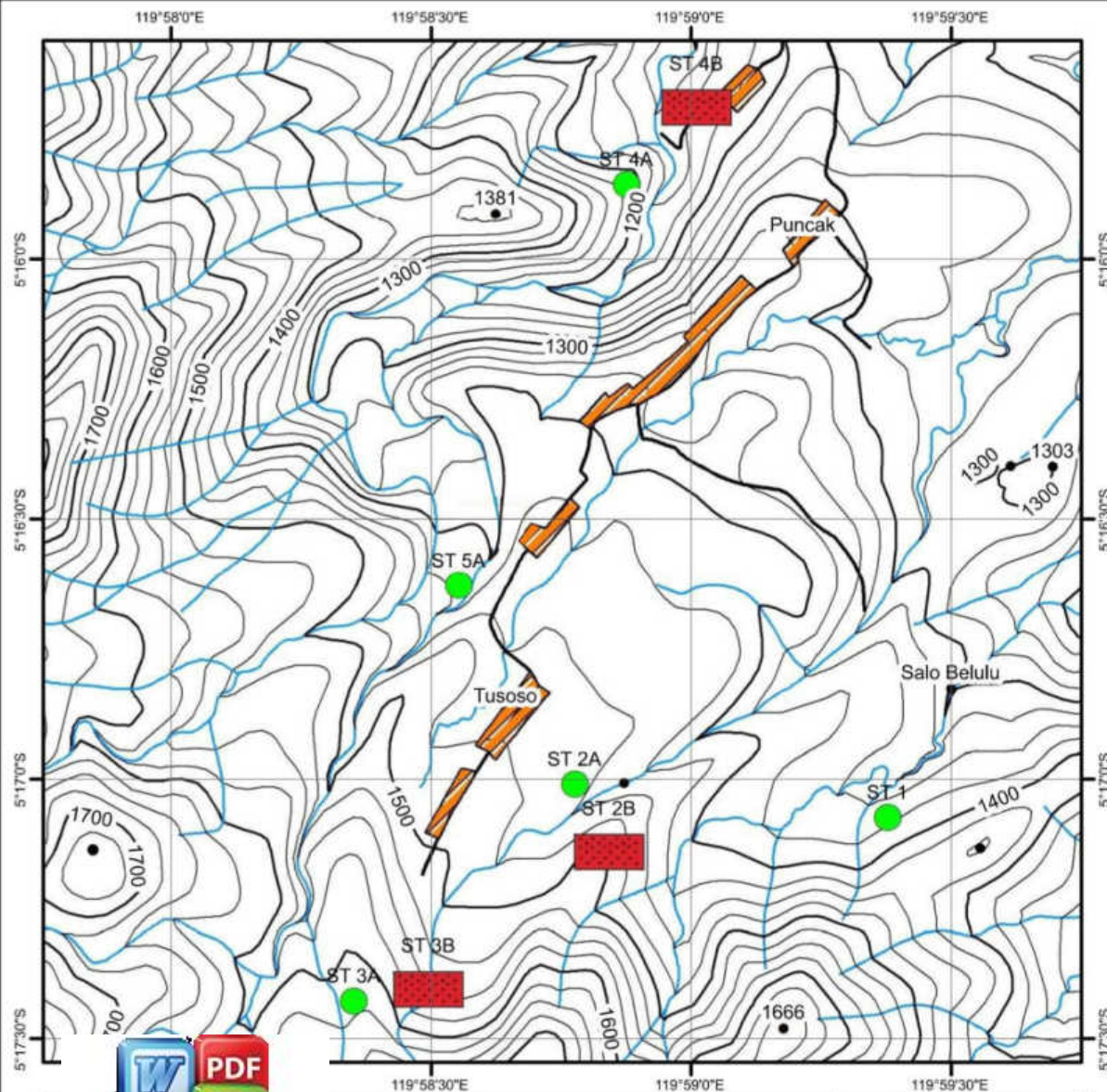
Unit Grid : UTM Zona 52 S
Unit Datum : WGS 1984

Legend	Slope
Batas Kabupaten / Kota	0 - 8 %
Batas Kecamatan	8 - 15 %
Batas Desa / Kelurahan	15 - 35 %
Lokasi Penelitian	35 - 45 %
	> 45 %

Sumber Peta

- Peta RBI Kabupaten Maros, BIG 2019
- FABDEM, Universitas Bristol 2022
- BATNAS, BIG 2019





KEMENTERIAN PENDIDIKAN KEBUDAYAAN RISET DAN TEKNOLOGI
UNIVERSITAS HASANUDDIN
FAKULTAS TEKNIK
DEPARTEMEN TEKNIK GEOLOGI
PROGRAM STUDI MAGISTER TEKNIK GEOLOGI

PETA LOKASI PENGAMBILAN SAMPEL

DAERAH GUNUNG PERAK KECAMATAN SINJAI BARAT
KABUPATEN SINJAI PROVINSI SULAWESI SELATAN



0 250 500 1,000 1,500 Meters

SKALA 1:25.000
IK 25 METER

OLEH:
KHAERUNNISA NASIR
D062211002

GOWA
2024

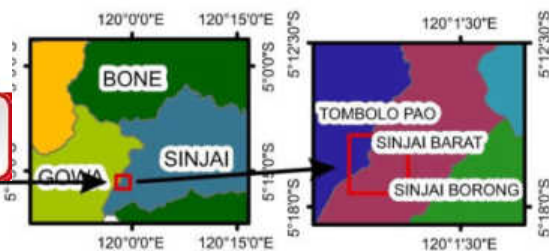
KETERANGAN

-  Basalt
-  Sampel Tanah
-  Titik Ketinggian
-  Kontur
-  Anak Sungai
-  Jalan
-  Pemukiman
-  Nama Daerah



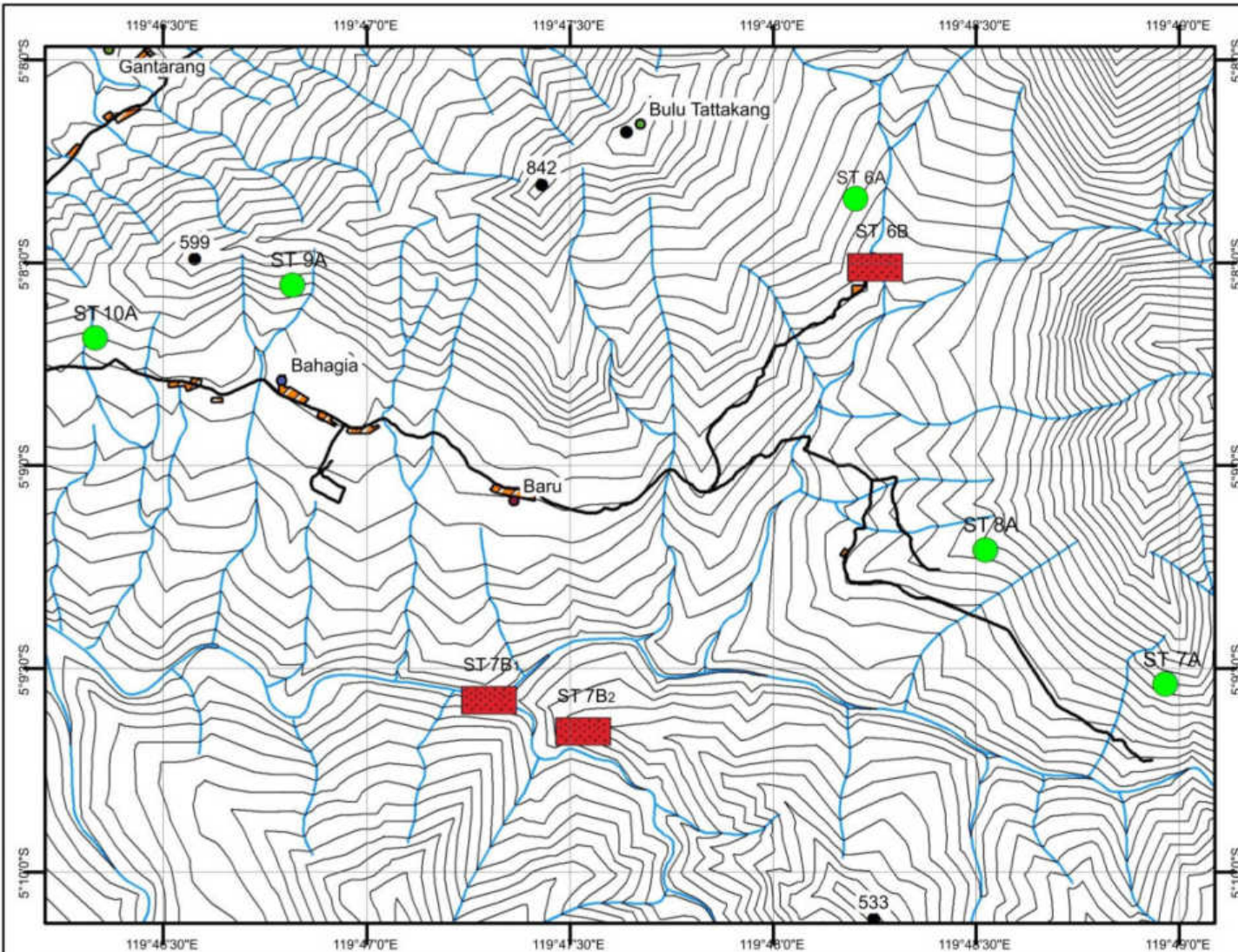
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PROGRAM STUDI MAGISTER TEKNIK GEOLOGI

PETA LOKASI PENGAMBILAN SAMPEL

DAERAH BONTOSOMBA KECAMATAN TOMPOBULU
KABUPATEN MAROS PROVINSI SULAWESI SELATAN



0 300 600 1,200 1,800 Meters

SKALA 1:25.000
IK 25 METER

OLEH:
KHAERUNNISA NASIR
D062211002

GOWA
2024

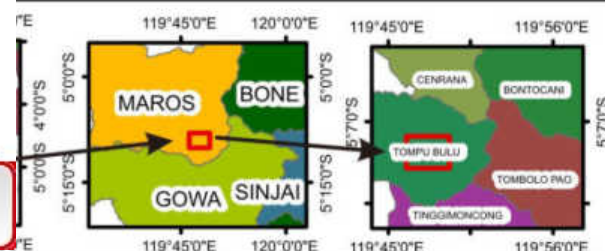
KETERANGAN

- Basalt
- Sampel Tanah
- Titik Ketinggian
- Kontur
- Anak Sungai
- Jalan
- Pemukiman
- Bontomanai
- Nama Daerah



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