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LAMPIRAN



Lampiran 1
Hasil Analisis *X-Ray Diffraction* (XRD)



HASIL ANALISIS *X-RAY DIFFRACTION* (XRD) SAMPEL AWAL







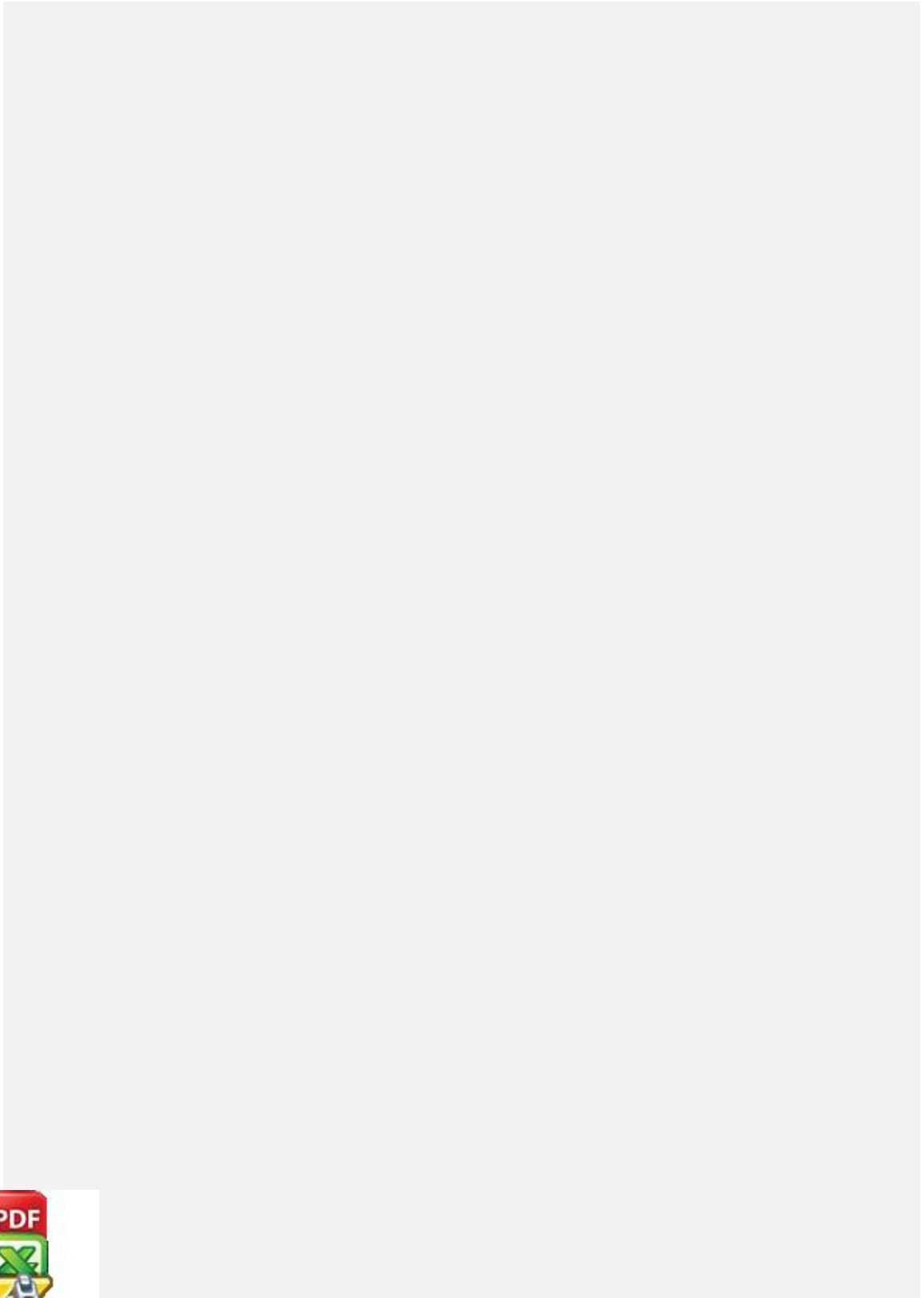


HASIL ANALISIS X-RAY *DIFFRACTION* (XRD) SAMPEL P -177+149







HASIL ANALISIS X-RAY *DIFFRACTION* (XRD) SAMPEL P -149+105





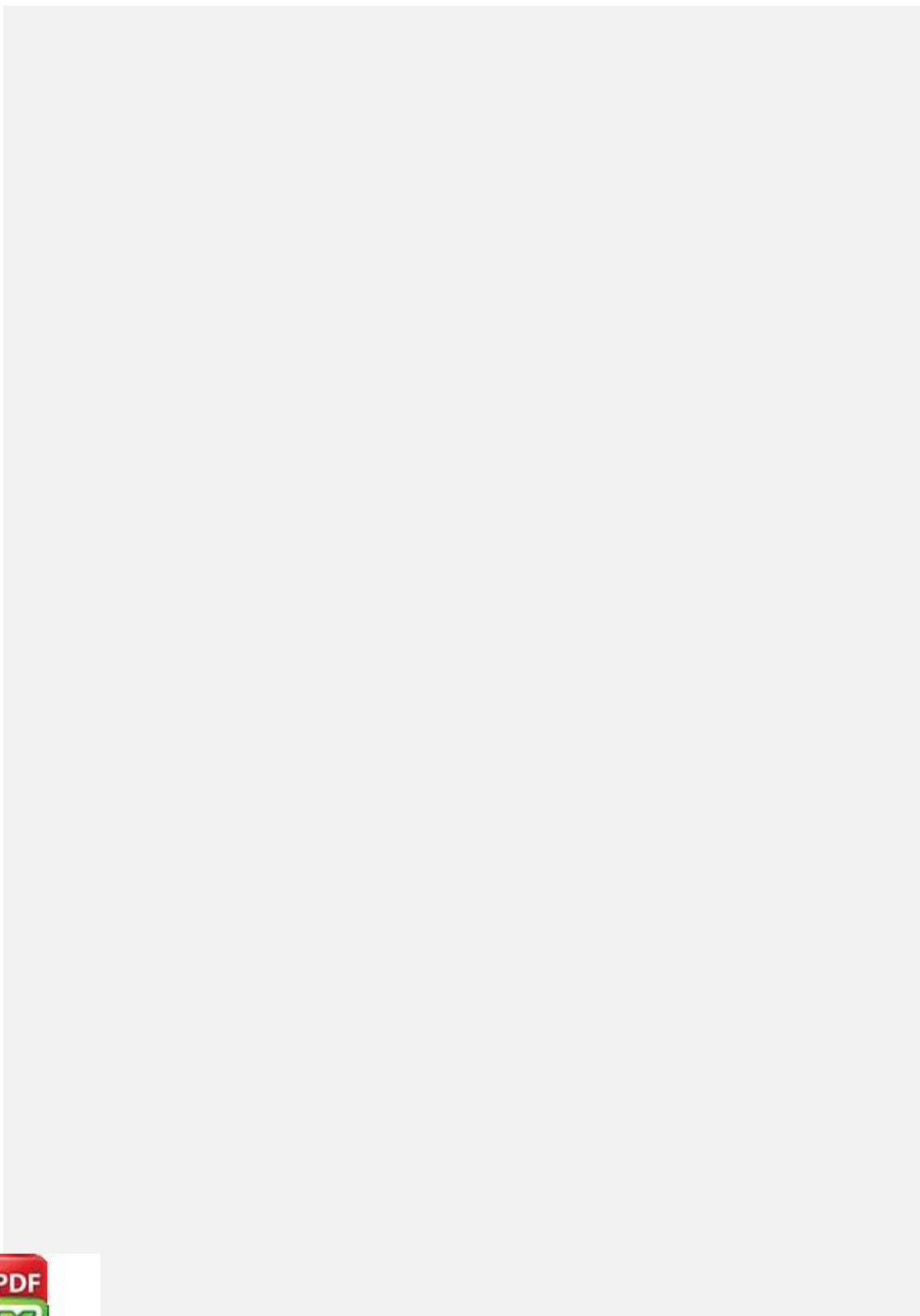


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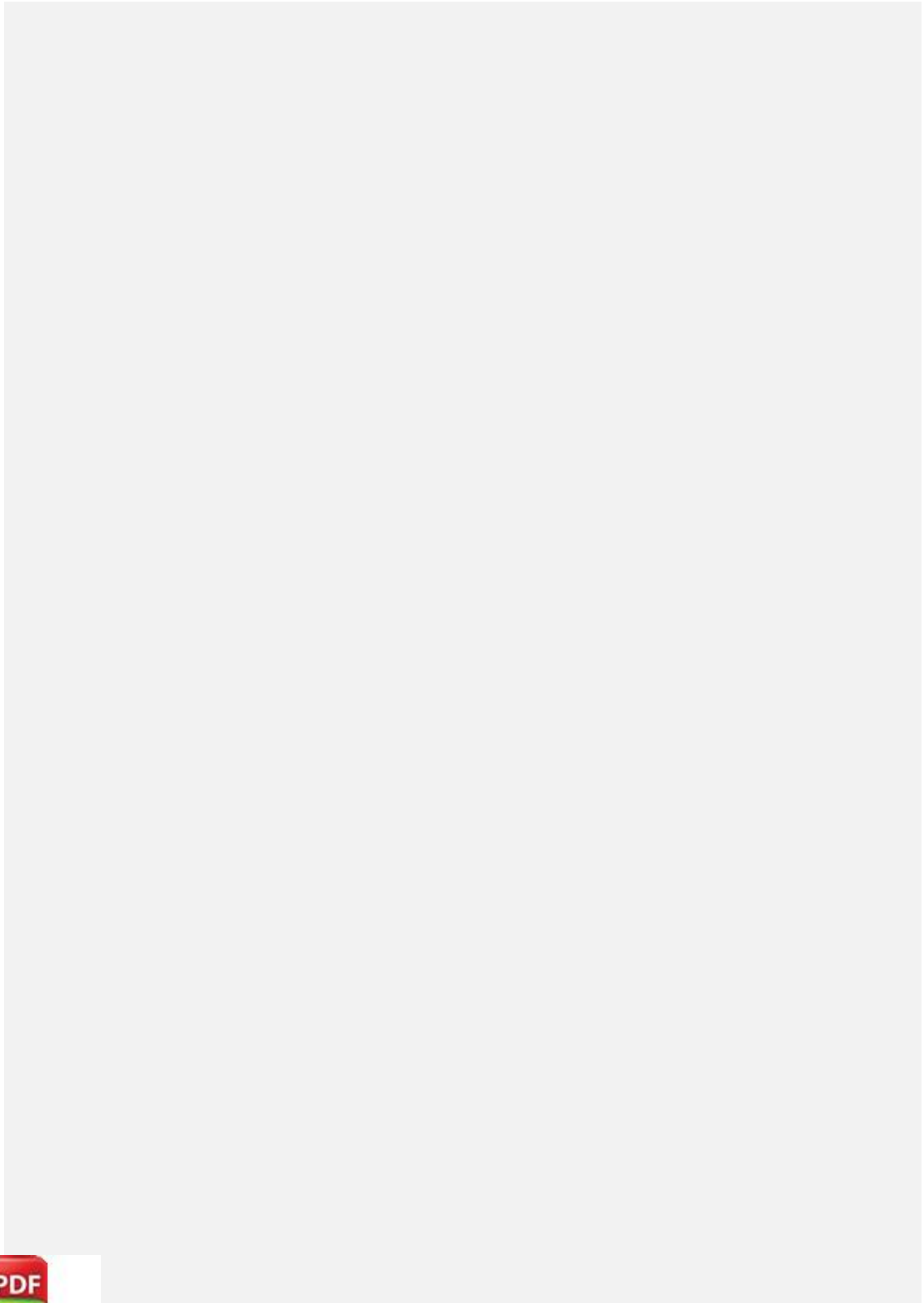


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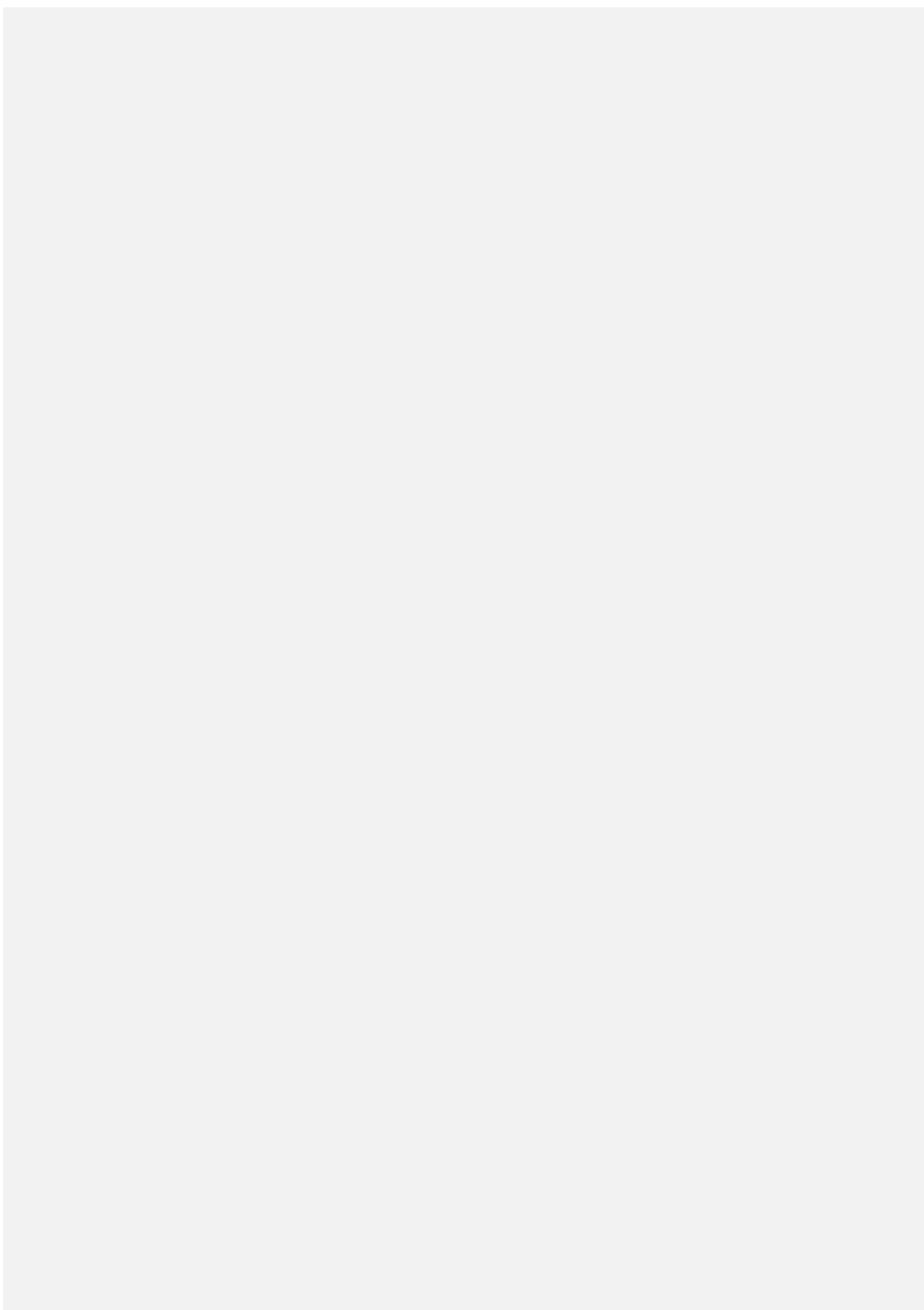






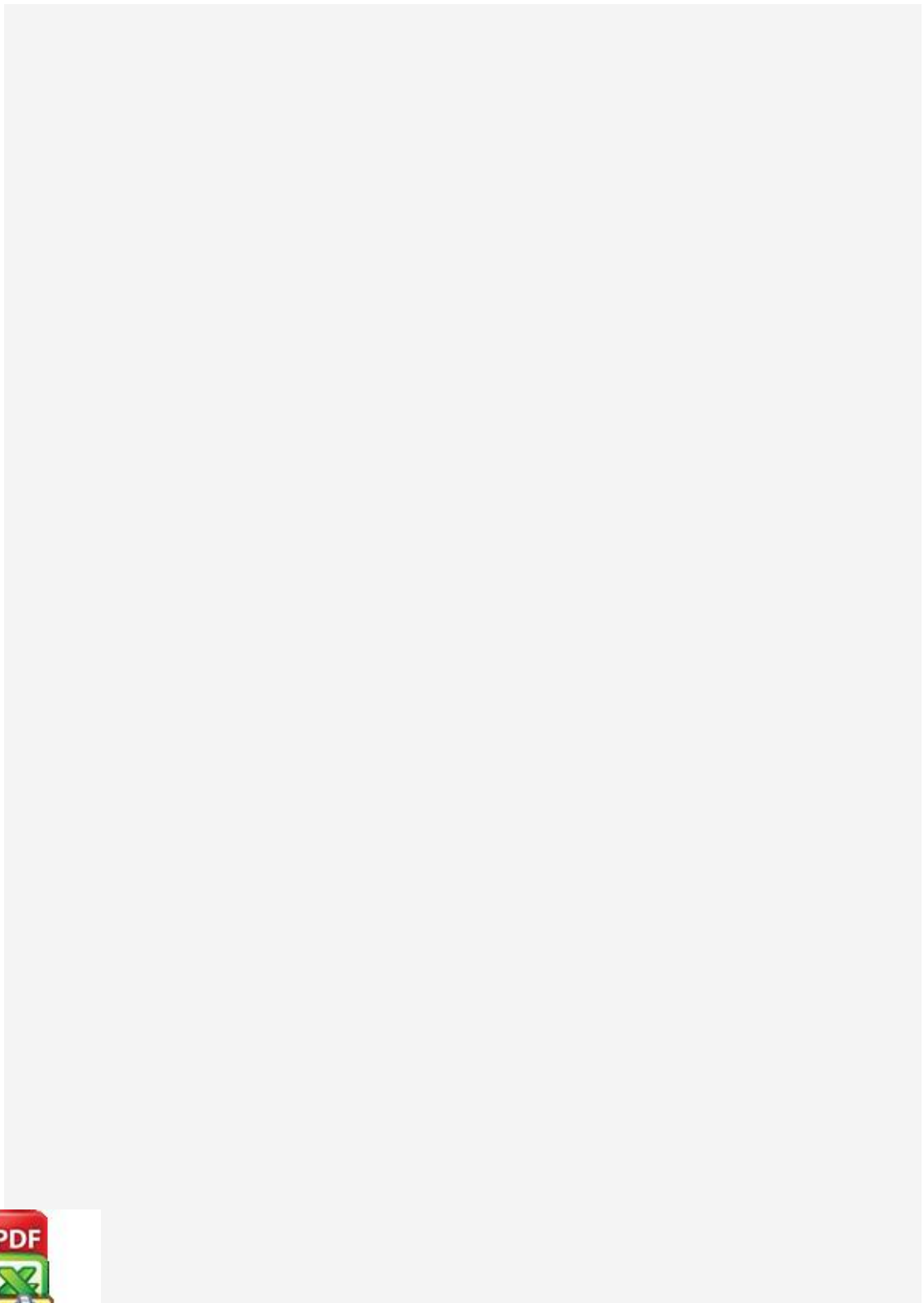
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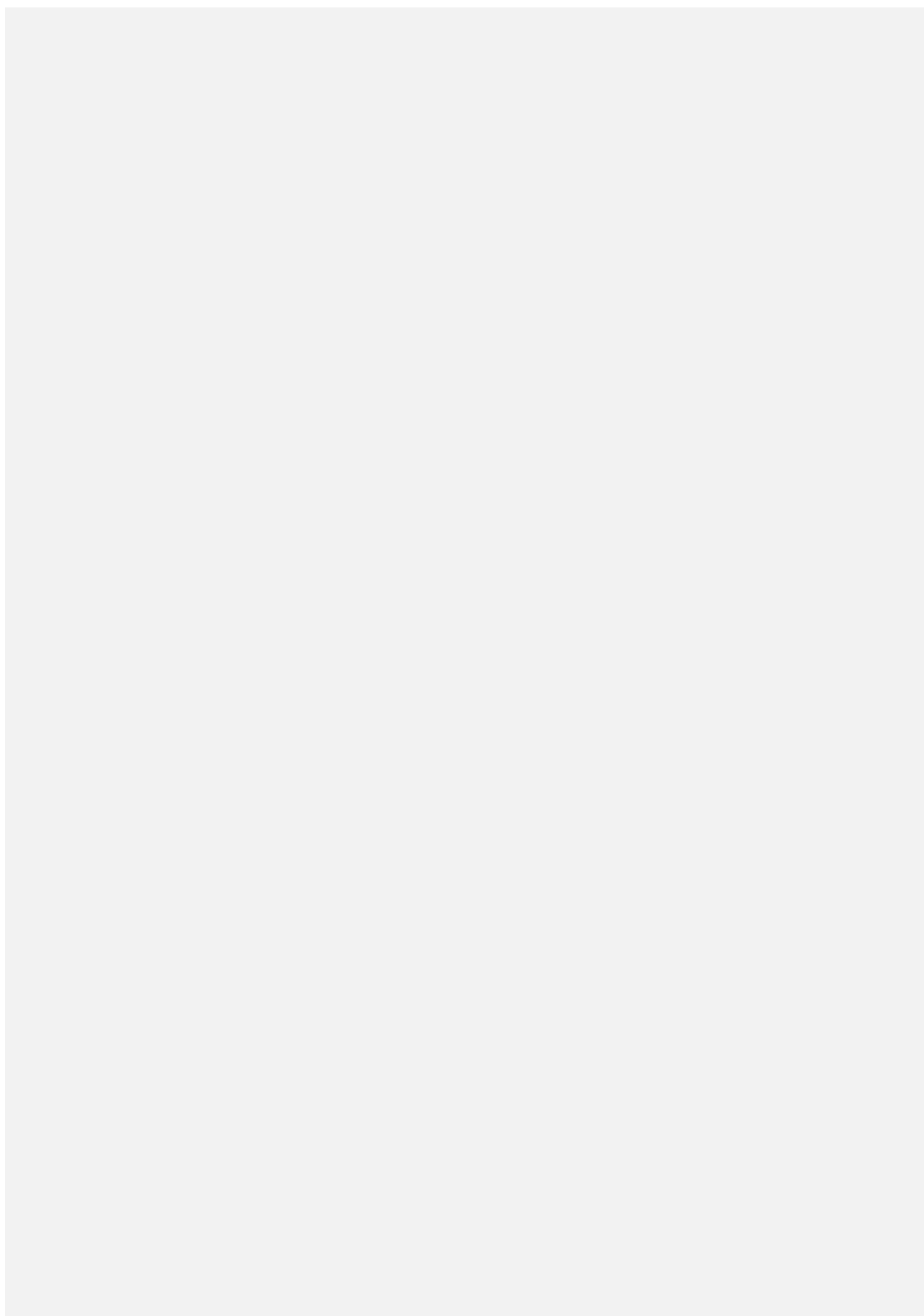






HASIL ANALISIS *X-RAY DIFFRACTION* (XRD) SAMPEL TP -149+105









HASIL ANALISIS X-RAY DIFFRACTION (XRD) SAMPEL TP -105+74

Match! Phase Analysis Report

Sample: NMD-TP-140 (5-70)

Sample Data

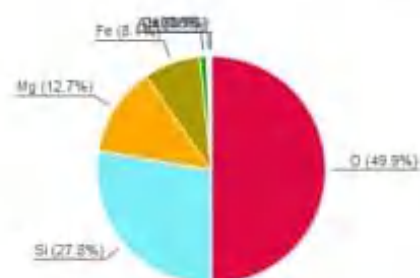
File name	NMD-TP-140.RAW
File path	D:\naufal\UNHAS\Skrpsi\Skrpsi\Data XRD Sampel Akhir\Naufal\NMD-TP-140
Data collected	Feb 2, 2024 13:13:21
Data range	5.000° - 70.000°
Original data range	5.000° - 70.000°
Number of points	3251
Step size	0.020
Rietveld refinement converged	No
Alpha2 subtracted	No
Background subtr.	No
Data smoothed	Yes
Radiation	X-rays
Wavelength	1.540600 Å

Analysis Results

Phase composition (Weight %)



Elemental composition (Weight %)



Index	Amount (%)	Name	Formula sum	Element	Amount (weight %)
A	30.6	Lizardite	Al _{0.22} Fe _{0.15} H ₄ Mg _{2.79} O ₉ Si _{1.84}	O	49.9% (*)
B	28.8	Silicon oxide Quartz low	O ₂ Si	Mg	12.7%
C	27.8	Talc	H ₂ Mg ₃ O ₁₂ Si ₄	Fe	8.1%
D	11.3	Goethite	Fe O ₂	Al	0.9%
E	1.5	Montmorillonite	Al ₂ Ca _{0.5} O ₁₂ Si ₄	Ca	0.1%
	12.0	Unidentified peak area		*LE (sum)	50.5%

Amounts calculated by RIR (Reference Intensity Ratio) method

Details of identified phases

A: Lizardite (30.6 %)

Formula sum	Al _{0.22} Fe _{0.15} H ₄ Mg _{2.79} O ₉ Si _{1.84}
Entry number	96-900-4995
Figure-of-Merit (FoM)	0.000000
Total number of peaks	111
Peaks in range	27
Peaks matched	6
Intensity scale factor	0.47
Space group	P 3 1 m
Crystal system	trigonal (hexagonal axes)
Unit cell	a = 5.3160 Å c = 7.1500 Å
V/c	1.38
Calc. density	2.674 g/cm ³
Reference	Mellini M., Zanazzi P. F., "Effects of pressure on the structure of lizardite-1T Sample: at P = 12.5 kbar", European Journal of Mineralogy 1 , 13-19 (1989)



B: Silicon oxide Quartz**low (28.8 %)**

Formula sum	O2 Si
Entry number	96-101-1160
Figure-of-Merit (FoM)	0.789206 [*]
Total number of peaks	70
Peaks in range	18
Peaks matched	10
Intensity scale factor	0.96
2theta correction	-0.016°
Space group	P 32 2 1 S
Crystal system	trigonal (hexagonal axes)
Unit cell	a= 4.9100 Å c= 5.4000 Å
V/c	3.01
Meas. density	2.660 g/cm ³
Calc. density	2.654 g/cm ³
Reference	Machatschki F, "Kristallstruktur von Tiefquarz", Fortschritte der Mineralogie 20 , 45-47 (1936)

C: Talc (27.6 %)

Formula sum	H2 Mg3 O12 Si4
Entry number	96-900-8298
Figure-of-Merit (FoM)	0.000000
Total number of peaks	251
Peaks in range	183
Peaks matched	65
Intensity scale factor	0.32
Space group	C -1
Crystal system	triclinic (anorthic)
Unit cell	a= 5.2900 Å b= 9.1730 Å c= 9.4600 Å α= 90.460° β= 98.680° γ= 90.090°
V/c	1.03
Calc. density	2.778 g/cm ³
Reference	Perdikatsis B., Burzlaff H., "Strukturverfeinerung am talc Mg3[(OH)2Si4O10]", Zeitschrift für Kristallographie 156 , 177-186 (1981)

D: Goethite (11.3 %)

Formula sum	Fe O2
Entry number	96-901-5697
Figure-of-Merit (FoM)	0.654345 [*]
Total number of peaks	187
Peaks in range	37
Peaks matched	14
Intensity scale factor	0.40
2theta correction	0.058°
Space group	P b n m
Crystal system	orthorhombic
Unit cell	a= 4.6188 Å b= 9.9528 Å c= 3.0236 Å
V/c	3.17
Calc. density	4.198 g/cm ³
Reference	Hazemann J.-L., Bézar J. F., Manceau A., "Rietveld studies of the aluminium-iron substitution in synthetic goethite", Materials Science Forum 79-82 , 821-826 (1991)

E: Montmorillonite (1.8 %)

Formula sum	Al2 Ca0.5 O12 Si4
Entry number	96-900-2780
Figure-of-Merit (FoM)	0.000000
Total number of peaks	92
Peaks in range	92
Peaks matched	30
Intensity scale factor	0.36
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°
V/c	17.33
Calc. density	1.801 g/cm ³
Reference	Viani A., Gualtieri A., Artioli G., "The nature of disorder in montmorillonite by simulation of X-ray powder patterns. Note: Structural simulation model", American Mineralogist 87 , 966-975 (2002)

^(*) 2theta values have been shifted internally for the calculation of the amounts, the intensity scaling factors as well as the figure-of-merit (FoM), due to the active search-match option 'Automatic zero point adaption'.

Candidates

Name	Formula	Entry No.	FoM
MOL-3CdS	C158 H138 Cd17 N4 S32	96-411-1893	0.7016
	C12 Ga O5	96-770-3817	0.6938
	Cs4 H62 K7 Na5 Nb23 O108 V5	96-712-7393	0.6888
	C10 H10 Co3 F2 N2 O8 P2	96-431-0851	0.6865
	Sc	96-151-2540	0.6801
	Be F2	96-153-1932	0.6786
	C114 Mn N7 O74 Si2 W18	96-711-6609	0.6714



Portlandite (deuterated)	Ca O2 O2	96-901-0908	0.6677
Lithium Cobalt Oxide	Co Li0.58 O2	96-155-0394	0.6672
	C116 Cl4 Fe N7 O72 Si2 W18	96-711-6607	0.6667
Li0.16 (Ni0.7 Fe0.15 Co0.15)1.03 O2	Co0.1545 Fe0.1545 Li0.16 Ni0.721 O2	96-152-6297	0.6631
Sodium tetrapropyl uranyl peroxide	C1.66 H3.6 N0.14 Na0.17 O1.68 U0.31	96-410-5386	0.6628
Si O2	O2 Si	96-153-8065	0.6588
Rb3GaF6	C8 H136 Fe6 N4 Na14 O184 Se6 W34	96-711-5857	0.6565
	F6 Ga Rb3	96-771-3195	0.6562
	O45 P9 V12	96-400-1392	0.6536
	C4 N5 S3	96-411-6670	0.6521
Lithium Cobalt Oxide	Co Li0.5 O2	96-155-0393	0.6476
Birnessite	H4 Mn Na0.29 O2.691	96-901-0156	0.6434
Li0.08 (Ni0.7 Fe0.15 Co0.15)1.03 O2	Co0.1545 Fe0.1545 Li0.08 Ni0.721 O2	96-152-6299	0.6430
	C116 Ag2 Cr Lu2 Mn4 N14 O78 Si2 W18	96-712-0260	0.6427
Zippelite	H27 Na5 O52 S4 U8	96-900-4756	0.6403
Na0.364 Mn O2 (H2 O)0.544	H1.088 Mn Na0.364 O2.544	96-153-1680	0.6400
Brucite	H2 Mg O2	96-900-2352	0.6370
	Cs4 H74 K8 Na5 Nb23 O115 V6	96-712-7394	0.6364
COV-3CdS-BPy	C166 H138 Cd17 N2 S30	96-410-0573	0.6346
	Hg Zr	96-152-2700	0.6335
beryllium bis(hypophosphite)	Be H4 O4 P2	96-201-4099	0.6321
(C (N H2)3)18 (Na (U O2)3 (O H) (H2 O)8 (As4 W40 O140 (W O))) (H2 O)25	C12.05 As4 N35.95 Na O179 U3 W41	96-153-6613	0.6312
Rb2 Sb8 S13 (H2 O)3.28	H6.56 O3.28 Rb2 S13 Sb8	96-152-6921	0.6303
	Ba Pb Zn	96-152-7803	0.6298
	Cu F H4 O5 P	96-154-5244	0.6265
	As2 Cd O6	96-220-0092	0.6254
	As2 Cd O6	96-900-7894	0.6254
Rb (Hg F3)	F3 Hg Rb	96-152-7535	0.6249
(Eu Yb)	Eu Yb	96-152-4623	0.6234
cis-Diaqua-(2,2-bipyridyl-N,N')-copper sulfate	C10 H12 Cu N2 O6 S	96-210-2510	0.6213
	Cs4 Se4 Si	96-810-1314	0.6209
Ho Ni C2	C2 Ho Ni	96-154-0779	0.6206
(P Cl4)3 (Ti Cl6) (P Cl6)	Cl24 P4 Ti	96-153-0749	0.6201
Ba Ga Sn D	Ba D Ga Sn	96-412-4017	0.6199
K (I O3)	I K O3	96-154-1797	0.6188
Zippelite	H2 N4 O21 S2 U4	96-900-4757	0.6181
Volborthite	Cu3 H6 O11 V2	96-901-4305	0.6173
	Nb Rh	96-231-0203	0.6115
Rb Ca F3	Ca F3 Rb	96-153-0113	0.6102
	C128 N8 O74 Si2 W20 Zn4	96-431-6561	0.6099
	Ca O2	96-153-0293	0.6095
Lithium Cobalt Oxide	Co Li0.42 O2	96-155-0391	0.6084
Pb0.855 ((U O2)4 O2.710 (O H)4.290) (H2 O)4	H12.29 O19 Pb0.855 U4	96-152-1509	0.6074
(Sn Cl (H2 O)2) (Sn Cl3) (H2 O)	Cl4 H8 O3 Sn2	96-153-4642	0.6052
Thallium tridecafluorotetraantimonate(III)	F13 Sb4 Ti	96-156-7935	0.6050
and 22 others...			

Search-Match

Settings

Reference database used	COD-Inorg 2023.12.05
Automatic zeropoint adaptation	Yes
Downgrade entries with low scaling factors	Yes
Minimum figure-of-merit (FoM)	0.60
2theta window for peak corr.	0.30 deg.
Minimum rel. int. for peak corr.	0
Parameter/influence 2theta	0.50
Parameter/influence intensities	0.50
Parameter multiple/single phase(s)	0.50

Criteria for entries added by user

Reference:

Entry number: 96-901-5697;96-900-4995;96-900-8298;96-900-2780

Peak List

No.	2theta [°]	d [Å]	I/I0 (peak height)	Counts (peak area)	FWHM	Matched
1	5.66	15.6017	349.51	32.85	0.6400	E
2	6.70	13.1821	103.83	4.27	0.2800	
3	9.32	9.4815	66.79	6.37	0.6493	C
4	12.38	7.1439	368.27	19.47	0.3600	A
5	18.66	4.7514	286.95	20.23	0.4800	
6	19.70	4.5029	246.43	27.50	0.7600	C,E
	0.98	4.2309	461.87	67.82	1.0000	B,D
	4.92	3.5702	376.28	24.31	0.4400	A



12	31.14	2.8698	262.91	15.44	0.4000	E
13	33.16	2.6995	187.96	8.83	0.3200	D
14	34.74	2.5802	253.08	29.73	0.8000	C,D,E
15	35.68	2.5144	339.53	55.84	1.1200	C,D,E
16	36.58	2.4545	485.36	45.61	0.6400	B,C,D,E
17	37.36	2.4051	102.53	10.84	0.7200	C
18	39.56	2.2762	130.96	13.08	0.6800	B,C
19	40.38	2.2319	117.48	15.87	0.9200	B,C,E
20	42.44	2.1282	84.01	4.38	0.3554	A,B,C,E
21	45.04	2.0112	103.72	3.66	0.2400	D
22	50.22	1.8152	162.44	7.63	0.3200	B,C
23	52.94	1.7282	176.04	8.27	0.3200	C,E
24	53.34	1.7162	198.67	10.50	0.3800	C,D
25	53.62	1.7079	120.15	29.64	1.6800	C
26	58.90	1.5667	119.44	19.64	1.1200	A,C,D,E
27	60.00	1.5406	104.93	7.47	0.4847	A,B,C,E
28	60.90	1.5200	135.01	11.10	0.5600	C,D,E
29	61.28	1.5115	135.04	22.21	1.1200	C,D
30	63.36	1.4688	111.24	5.23	0.3200	C,D,E
31	68.30	1.3722	101.65	8.36	0.5600	B,C,D

Integrated Profile Areas

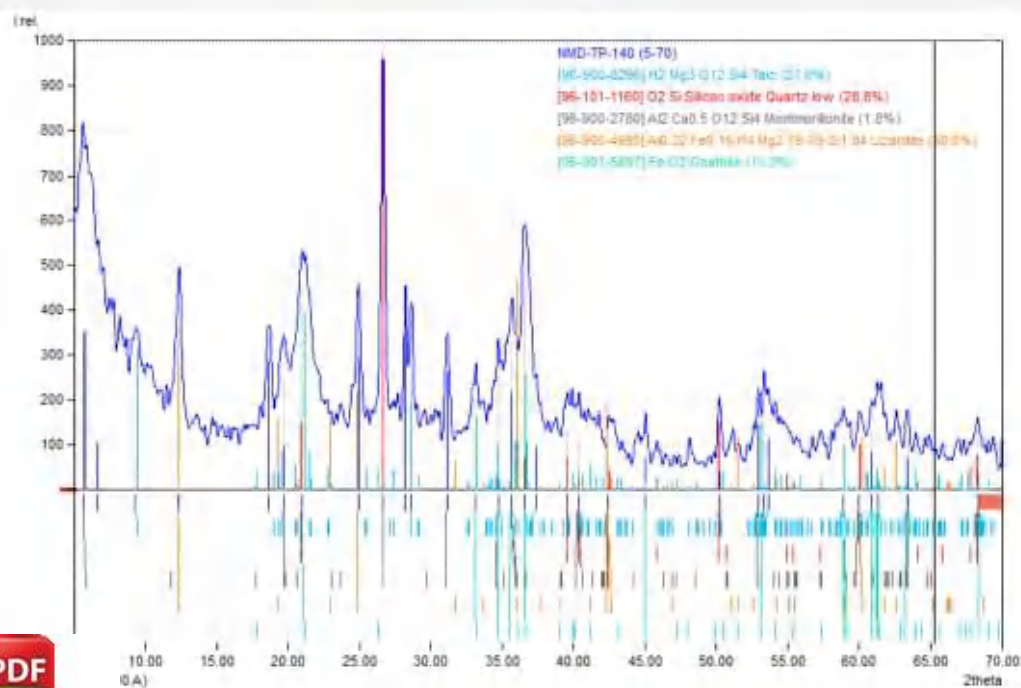
Based on calculated profile

Profile area	Counts	Amount
Overall diffraction profile	84625	100.00%
Background radiation	52647	62.21%
Diffraction peaks	31978	37.79%
Peak area belonging to selected phases	21798	25.76%
Peak area of phase A (Talc)	5989	7.08%
Peak area of phase B (Silicon oxide Quartz low)	4447	5.26%
Peak area of phase C (Montmorillonite)	1518	1.79%
Peak area of phase D (Lizardite)	4592	5.43%
Peak area of phase E (Goethite)	5251	6.21%
Unidentified peak area	10180	12.03%

Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	616	100.00%
Peak intensity belonging to selected phases	458	74.33%
Unidentified peak intensity	158	25.67%

Diffraction Pattern Graphics



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trial version
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HASIL ANALISIS X-RAY DIFFRACTION (XRD) SAMPEL TP -74

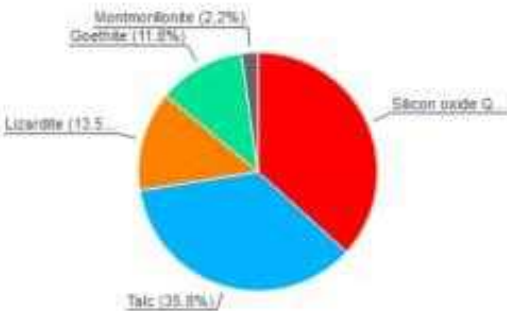
Match! Phase Analysis Report

Sample: NMD-TP-200 (5-70)

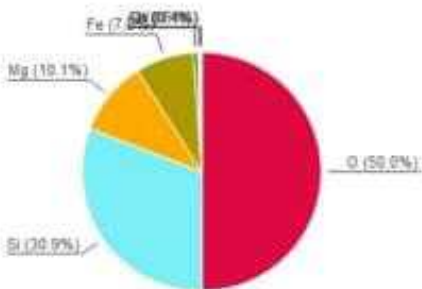
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Data collected	Feb 2, 2024 13:13:21
Data range	5.190° - 70.190°
Original data range	5.000° - 70.000°
Number of points	3251
Step size	0.020
Rietveld refinement converged	No
Alpha2 subtracted	No
Background subtr.	No
Data smoothed	Yes
2theta correction	0.19°
Radiation	X-rays
Wavelength	1.540600 Å

Analysis Results

Phase composition (Weight %)



Elemental composition (Weight %)



Index	Amount (%)	Name	Formula sum
A	36.7	Silicon oxide Quartz low Q2 Si	Q2 Si
B	35.8	Talc	H2 Mg3 O12 Si4
C	13.5	Lizardite	Al2 Si2 Fe0.15 H4 Mg2.75 O9 Si1.84
D	11.8	Goethite	Fe O3
E	2.2	Montmorillonite	Al2 Ca0.5 O12 Si4
	15.0	Unidentified peak area	

Element	Amount (weight %)
O	50.0% (*)
Si	20.9%
Mg	10.1%
Fe	7.0%
Al	0.6%
Ca	0.1%
*LE (sum)	50.4%

Amounts calculated by RIR (Reference Intensity Ratio) method

Details of identified phases

A: Silicon oxide Quartz low (36.7 %)	
Formula sum	Q2 Si
Entry number	96-101-1160
Figure-of-Merit (FoM)	0.757052
Total number of peaks	70
Peaks in range	18
Peaks matched	10
Intensity scale factor	0.96
2theta correction	-0.188°
Space group	P 32 2 1 S
	trigonal (hexagonal axes)
	a= 4.9100 Å c= 5.4000 Å
	3.01



Meas. density	2.660 g/cm ³
Calc. density	2.654 g/cm ³
Reference	Machatschki F, "Kristallstruktur von Tiefquarz", Fortschritte der Mineralogie 20 , 45-47 (1936)
B: Talc (35.8 %)	
Formula sum	H ₂ Mg ₃ O ₁₂ Si ₄
Entry number	96-900-8298
Figure-of-Merit (FoM)	0.000000
Total number of peaks	251
Peaks in range	184
Peaks matched	68
Intensity scale factor	0.32
Space group	C -1
Crystal system	triclinic (anorthic)
Unit cell	a= 5.2900 Å b= 9.1730 Å c= 9.4600 Å α= 90.460° β= 98.680° γ= 90.090°
I/c	1.03
Calc. density	2.776 g/cm ³
Reference	Perdikatsis B., Burzlaff H., "Strukturverfeinerung am talk Mg ₃ [(OH)2Si ₄ O ₁₀]", Zeitschrift für Kristallographie 156 , 177-186 (1981)
C: Lizardite (13.5 %)*	
Formula sum	Al _{0.22} Fe _{0.15} H ₄ Mg _{2.79} O ₈ Si _{1.84}
Entry number	96-900-4995
Figure-of-Merit (FoM)	0.586259*
Total number of peaks	111
Peaks in range	28
Peaks matched	9
Intensity scale factor	0.16
2theta correction	-0.051°
Space group	P 3 1 m
Crystal system	trigonal (hexagonal axes)
Unit cell	a= 5.3160 Å c= 7.1500 Å
I/c	1.38
Calc. density	2.674 g/cm ³
Reference	Mellini M., Zanazzi P. F., "Effects of pressure on the structure of lizardite-1T Sample: at P = 12.5 kbar", European Journal of Mineralogy 1 , 13-19 (1989)
D: Goethite (11.8 %)*	
Formula sum	Fe O ₂
Entry number	96-901-5697
Figure-of-Merit (FoM)	0.688008*
Total number of peaks	187
Peaks in range	37
Peaks matched	16
Intensity scale factor	0.32
2theta correction	-0.043°
Space group	P b n m
Crystal system	orthorhombic
Unit cell	a= 4.6188 Å b= 9.9528 Å c= 3.0236 Å
I/c	3.17
Calc. density	4.198 g/cm ³
Reference	Hazemann J.-L., Béjar J. F., Manceau A., "Rietveld studies of the aluminium-iron substitution in synthetic goethite", Materials Science Forum 79-82 , 821-826 (1991)
E: Montmorillonite (2.2 %)	
Formula sum	Al ₂ Ca _{0.5} O ₁₂ Si ₄
Entry number	96-900-2780
Figure-of-Merit (FoM)	0.000000
Total number of peaks	92
Peaks in range	92
Peaks matched	31
Intensity scale factor	0.33
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/c	17.33
Calc. density	1.801 g/cm ³
Reference	Viani A., Gualtieri A., Artioli G., "The nature of disorder in montmorillonite by simulation of X-ray powder patterns Note: Structural simulation model", American Mineralogist 87 , 966-975 (2002)

(*2theta values have been shifted internally for the calculation of the amounts, the intensity scaling factors as well as the figure-of-merit (FoM), due to the active search-match option 'Automatic zero point adaption'.

Candidates

Name	Formula	Entry No.	FoM
aluminum phosphate	Ga ₁₆ Se ₃₃ Zn ₄	96-410-6294	0.7196
	Al O ₄ P	96-201-0796	0.7145
	H _{78.6} Na _{12.6} O ₈₆	96-700-2723	0.7121
i) ((C ₅ H ₉ N H) ₂ (C H ₂) ₃)	P ₁₀ Pd _{15.4}		
	Cd ₄ In ₁₆ S ₃₃	96-412-3975	0.7104



	C28 H183 Cu8 K4 N28 Na7 Nb54 O184	96-712-9691	0.7077
	C76 H179 N10 O60 Sb W18	96-450-0754	0.7075
	C60 Bi10 Br40 O25	96-713-0614	0.7052
	C32 H19 Ag12 F50 O29.5	96-411-5643	0.7020
	C132 H42 F80 Fe Mn12 O50	96-432-0646	0.6961
	Cl2 Cs3 Fe H36 K4 O95 Si2 Ta6 W18	96-722-8912	0.6942
Tridymite	O2 Si	96-901-3394	0.6910
	C120 Cl8 Mn4 N7 Ni O70 Si2 W18	96-703-5027	0.6891
	C18 Cl17 Cs3 N6 O61 Ti12	96-412-6915	0.6866
	C102 H227 Mo17 N9 O62 V3	96-700-3207	0.6866
(Zn4 In16 S33) ((C5 H9 N H)2 (C H2)3)	In16 S33 Zn4	96-412-3976	0.6877
(Co4 In16 S33) ((C5 H9 N H)2 (C H2)3)	Co4 In16 S33	96-412-3977	0.6873
	C120 Cl8 Cu Mn4 N7 O70 Si2 W18	96-703-5028	0.6872
	C12 O7 P2 Zn	96-431-1830	0.6870
	C120 Cl8 Co Mn4 N7 O70 Si2 W18	96-703-5026	0.6861
	Al1.55 Na1.55 O4 Si0.45	96-200-2896	0.6834
MOL-3CdS	C158 H138 Cd17 N4 S32	96-411-1893	0.6821
	C48 H84 Br24 Cr12 O110 Si W12	96-400-1632	0.6808
	C52 B6 Co F3 N7 O2	96-723-4123	0.6784
	C116 Ag2 Cr Lu2 Mn4 N14 O78 Si2 W18	96-712-0260	0.6778
	C118 Cl6 Mn5 N7 O70 Si2 W18	96-703-5024	0.6773
	C77 H190.2 Cu5 Ge0 In30 N11 O7.1 S53	96-722-8736	0.6750
	C258 H579 N17 O156 Si4 Ti8 W40	96-110-1220	0.6748
	C48 H96 Al21 Na21 O216 Si75	96-150-7214	0.6731
	C20 H9 Ag8 F18 N3 O12	96-411-5641	0.6716
	C124 Cl12 Cu4 Fe N7 O72 Si2 W18	96-434-6914	0.6710
COV-3CdS-BPy	C166 H138 Cd17 N2 S30	96-410-0573	0.6709
	C12 H94 Cu3 Ge K2 N12 Na4 Nb12 O65 V2	96-451-0028	0.6674
	C156 H134 Br4 Cd17 N2 S28	96-723-3663	0.6638
	Cs4 H62 K7 Na5 Nb23 O108 V5	96-712-7393	0.6636
	C324 H270 Cd33 S60 C3 Cs2 K0.5 O88 P2 W18	96-410-9637	0.6611
	C10 H10 Co3 F2 N2 O8 P2	96-156-1589	0.6609
	C120 As24 O99 U10 C8 H136 Fe6 N4	96-431-0851	0.6608
	Na14 O184 Se6 W34 C35 H11 Ag11 F45 N3 O19	96-433-5132	0.6596
	C12 Ga O5 C347.64 H811.16 In60 N57.94 Se118.97	96-711-5857	0.6593
LiNaU20P12	Li6 Na18 O188 P12 S U20	96-411-5644	0.6581
LiNaU20P6	Li Na18 O162 P6 S U20	96-770-3817	0.6568
Au36	C102 Ge N9 O36 W10 C120 H216 Au36 S24	96-770-1506	0.6563
Na20 (Al20 Si76 O192)	C144 Cu6 O42 P3 Al20 Na20 O192 Si76	96-434-5367	0.6549
(C6 H12 (N H2)2)2 Zn Ga7 (H P O4)2 (P O4)6 (O H) F4 (H2 O)3	C12 H41 F4 Ga7 N4 O36 P8 Zn	96-710-8167	0.6538
	C64 H144 N4 O40 Si W12	96-151-4182	0.6533
	C96 N6 O36 Si W10	96-155-6348	0.6488
		96-152-7028	0.6479
		96-152-8075	0.6475
		96-431-9533	0.6469
		96-432-9779	0.6453



Pigment Green 7, Hostaperm Green GNX, Solid solution of CuPcC16 (main component), CuPcHCl15, CuPcH2Cl14, CuPcH3Cl13 and CuPcH4Cl12. Average composition Cu C32 N8 H1.2 Cl14.8 and 78 others...

C32 Cl16 Cu N8

96-210-8857

0.6451

Search-Match

Settings

Reference database used COD-inorg 2023.12.05
Automatic zeropoint adaptation Yes
Downgrade entries with low scaling factors Yes
Minimum figure-of-merit (FoM) 0.60
2theta window for peak corr. 0.30 deg.
Minimum rel. int. for peak corr. 0
Parameter/influence 2theta 0.50
Parameter/influence intensities 0.50
Parameter multiple/single phase(s) 0.50

Criteria for entries added by user

Reference:

Entry number:

96-901-5697;96-900-4985;96-900-8298;96-900-2780

Peak List

No.	2theta [°]	d [Å]	I/I0 (peak height)	Counts (peak area)	FWHM	Matched
1	5.59	15.7989	275.13	11.78	0.2800	
2	5.91	14.9423	310.76	32.32	0.6800	E
3	6.49	13.6082	156.68	6.71	0.2800	
4	6.85	12.8938	100.55	4.31	0.2800	
5	12.45	7.1039	279.97	18.84	0.4400	C
6	18.91	4.6891	215.04	11.84	0.3600	B
7	19.59	4.5279	194.03	5.93	0.2000	B
8	19.85	4.4692	268.85	27.96	0.6800	E
9	20.45	4.3394	202.04	59.32	1.9200	B,E
10	21.09	4.2091	364.39	57.95	1.0400	A,D
11	21.43	4.1431	201.83	24.69	0.8000	B
12	25.15	3.5381	345.05	19.00	0.3600	B,C
13	26.83	3.3202	1000.00	48.94	0.3200	A,B,E
14	28.49	3.1304	246.51	9.05	0.2400	B
15	28.79	3.0985	285.98	14.00	0.3200	B
16	31.29	2.8564	145.99	6.25	0.2800	E
17	33.29	2.6892	185.73	15.91	0.5600	D
18	34.79	2.5766	231.03	18.37	0.5200	B,D,E
19	35.25	2.5440	171.14	39.78	1.5200	D,E
20	35.95	2.4961	201.12	81.20	2.6400	B,C,D,E
21	36.83	2.4384	476.23	49.52	0.6800	A,B,D,E
22	39.77	2.2647	96.87	6.52	0.4400	A,B
23	40.07	2.2484	108.52	13.94	0.8400	D,E
24	40.41	2.2303	87.13	3.52	0.2639	A,B,E
25	41.19	2.1898	86.34	14.03	1.0627	B,C,D,E
26	42.75	2.1135	105.37	2.98	0.1849	A,B,C
27	50.39	1.8095	100.64	3.69	0.2400	A,B,D
28	53.29	1.7177	244.54	10.47	0.2800	B,D
29	53.67	1.7064	168.87	35.12	1.3600	B
30	58.97	1.5650	118.49	5.80	0.3200	B,C,D,E
31	59.21	1.5593	97.97	5.39	0.3600	B,D
32	60.21	1.5357	97.29	4.17	0.2800	A,B,C
33	61.43	1.5081	170.96	24.05	0.9200	B,D
34	63.63	1.4612	99.76	3.66	0.2400	B,E
35	64.97	1.4342	53.10	7.60	0.9365	B,C,E
36	68.57	1.3675	64.64	1.84	0.1859	A,B,C,D

Integrated Profile Areas

Based on calculated profile

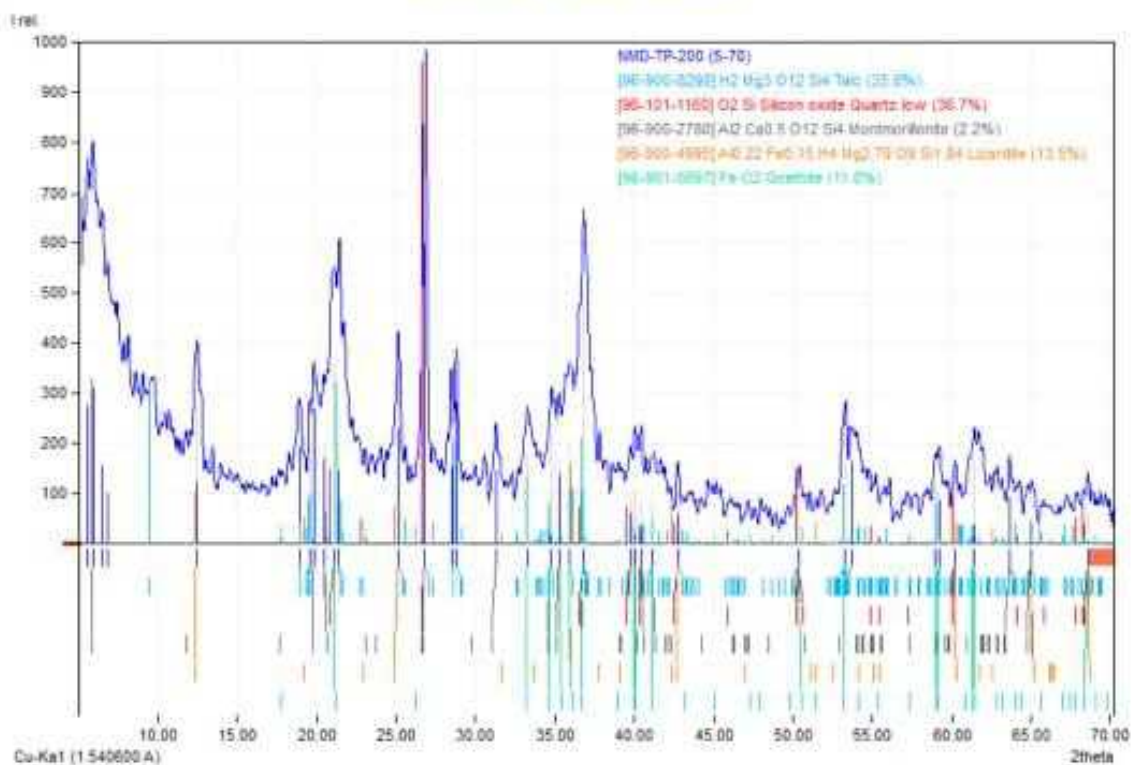
Profile area	Counts	Amount
Overall diffraction profile	81885	100.00%
Background radiation	49258	60.16%
Diffraction peaks	32627	39.84%
Peak area belonging to selected phases	20362	24.87%
Peak area of phase A (Taic)	6693	8.17%
Peak area of phase B (Silicon oxide Quartz low)	4352	5.31%
Peak area of phase C (Montmorillonite)	1680	2.05%
Peak area of phase D (Lizardite)	2273	2.78%
Phase E (Goethite)	5363	6.55%
Peak area	12265	14.98%



Peak Residuals

Peak data	Counts	Amount
Overall peak intensity	706	100.00%
Peak intensity belonging to selected phases	469	66.35%
Unidentified peak intensity	238	33.65%

Diffraction Pattern Graphics



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