



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:51 PM UTC

PDB ID : 3ZIL / pdb_00003zil
Title : Structure of the Wpl1 protein
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Deposited on : 2013-01-09
Resolution : 2.01 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

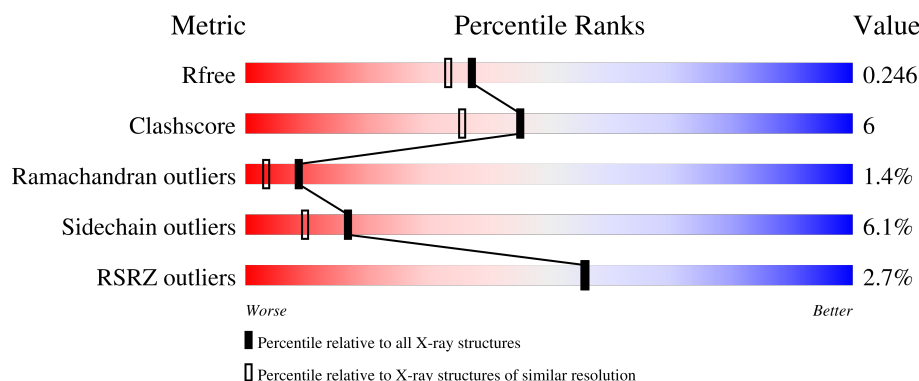
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

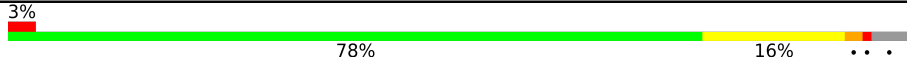
The reported resolution of this entry is 2.01 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	378	 3% 78% 16% ...
2	B	28	 18% 18% . 61%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3034 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called AAR187CP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2873	1843	480	536	14			

- Molecule 2 is a protein called AAL182WP.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	11	Total	C	N	O	0	0	0
			100	59	20	21			

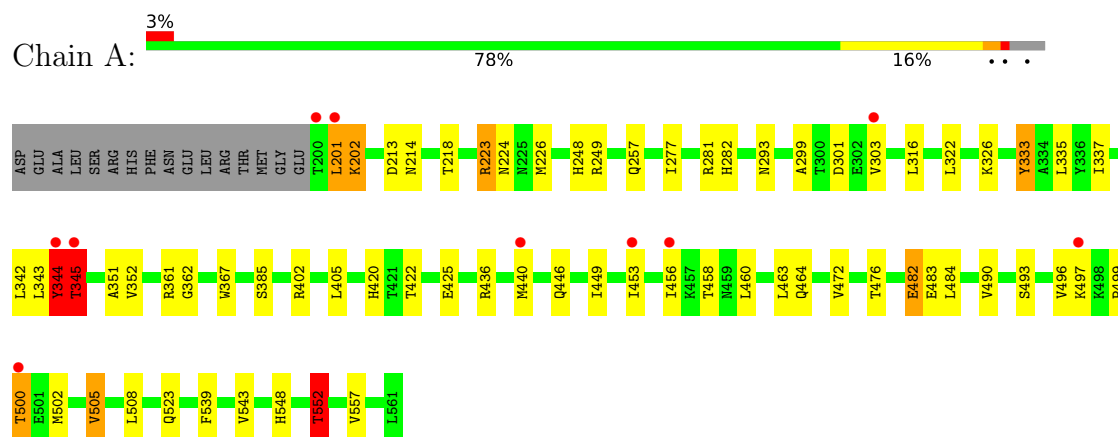
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	61	Total	O	0	0
			61	61		

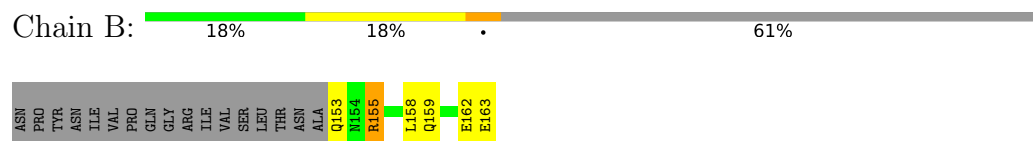
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: AAR187CP



• Molecule 2: AAL182WP



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.40Å 35.21Å 117.38Å 90.00° 101.89° 90.00°	Depositor
Resolution (Å)	29.12 – 2.01 29.12 – 2.01	Depositor EDS
% Data completeness (in resolution range)	98.6 (29.12-2.01) 98.7 (29.12-2.01)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.6	Depositor
R, R_{free}	0.200 , 0.246 0.199 , 0.246	Depositor DCC
R_{free} test set	1322 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.208	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3034	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.39	14/2927 (0.5%)	1.28	11/3976 (0.3%)
2	B	1.04	0/99	1.09	0/129
All	All	1.38	14/3026 (0.5%)	1.27	11/4105 (0.3%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	352	VAL	CA-C	6.55	1.59	1.52
1	A	425	GLU	C-O	6.45	1.31	1.24
1	A	248	HIS	CG-CD2	6.30	1.42	1.35
1	A	248	HIS	CE1-NE2	6.05	1.38	1.32
1	A	335	LEU	N-CA	5.89	1.53	1.46
1	A	282	HIS	CG-CD2	5.81	1.42	1.35
1	A	361	ARG	CZ-NH2	5.71	1.40	1.33
1	A	420	HIS	CG-CD2	5.59	1.42	1.35
1	A	463	LEU	CA-C	5.51	1.59	1.52
1	A	226	MET	N-CA	5.50	1.53	1.46
1	A	299	ALA	N-CA	5.22	1.52	1.46
1	A	351	ALA	N-CA	5.06	1.52	1.46
1	A	293	ASN	N-CA	5.05	1.52	1.46
1	A	385	SER	CA-C	5.03	1.59	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	343	LEU	N-CA-C	-7.61	102.88	113.37
1	A	345	THR	N-CA-C	-7.16	90.94	111.00
1	A	201	LEU	N-CA-C	-6.36	102.33	110.53
1	A	344	TYR	CA-C-N	6.09	132.66	121.70
1	A	344	TYR	C-N-CA	6.09	132.66	121.70
1	A	552	THR	N-CA-C	-6.08	104.73	111.36
1	A	500	THR	N-CA-C	-5.52	106.04	112.89
1	A	482	GLU	N-CA-C	5.11	117.51	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	436	ARG	CG-CD-NE	-5.10	100.77	112.00
1	A	436	ARG	N-CA-CB	5.06	117.35	110.01
1	A	218	THR	N-CA-C	-5.04	103.61	110.36

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2873	0	2917	34	0
2	B	100	0	98	3	0
3	A	61	0	0	1	0
All	All	3034	0	3015	35	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (35) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:TYR:HB3	1:A:345:THR:CA	1.75	1.16
1:A:344:TYR:HB3	1:A:345:THR:CB	1.80	1.09
1:A:446:GLN:HA	1:A:449:ILE:HD12	1.49	0.93
1:A:223:ARG:NH1	1:A:224:ASN:OD1	2.04	0.91
1:A:301:ASP:OD1	1:A:303:VAL:HG22	1.74	0.88
1:A:344:TYR:CB	1:A:345:THR:CB	2.52	0.87
1:A:223:ARG:HH11	1:A:223:ARG:HG2	1.40	0.85
1:A:344:TYR:HB3	1:A:345:THR:HA	1.55	0.85
1:A:201:LEU:O	1:A:202:LYS:HB3	1.84	0.75
1:A:548:HIS:O	1:A:552:THR:HB	1.89	0.73
1:A:342:LEU:C	1:A:344:TYR:N	2.57	0.62
1:A:223:ARG:HH11	1:A:223:ARG:CG	2.11	0.59
1:A:342:LEU:C	1:A:344:TYR:H	2.11	0.57
1:A:303:VAL:O	1:A:303:VAL:HG23	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:LEU:O	1:A:202:LYS:CB	2.55	0.55
1:A:214:ASN:HB2	1:A:257:GLN:OE1	2.09	0.53
1:A:496:VAL:HG23	1:A:497:LYS:HG3	1.90	0.53
2:B:153:GLN:HB3	2:B:155:ARG:NH1	2.25	0.52
1:A:422:THR:HG23	3:A:2054:HOH:O	2.10	0.50
1:A:508:LEU:HD21	2:B:158:LEU:HD11	1.94	0.50
1:A:277:ILE:HG13	1:A:277:ILE:O	2.13	0.49
1:A:453:ILE:HG22	1:A:453:ILE:O	2.13	0.48
1:A:202:LYS:HB3	1:A:202:LYS:HE2	1.76	0.47
1:A:362:GLY:HA2	1:A:367:TRP:HZ2	1.81	0.46
1:A:502:MET:HE3	1:A:502:MET:HB3	1.86	0.45
1:A:249:ARG:O	1:A:249:ARG:HG2	2.12	0.45
1:A:460:LEU:O	1:A:464:GLN:HG3	2.17	0.44
1:A:440:MET:HE3	1:A:483:GLU:HG2	1.99	0.44
1:A:458:THR:HG23	1:A:505:VAL:HG21	2.00	0.43
1:A:281:ARG:HH12	1:A:344:TYR:N	2.17	0.43
1:A:472:VAL:HG11	1:A:484:LEU:HD21	2.00	0.42
1:A:333:TYR:O	1:A:337:ILE:HG12	2.21	0.41
1:A:548:HIS:HB3	2:B:162:GLU:OE2	2.21	0.41
1:A:326:LYS:HE3	1:A:326:LYS:HB2	1.83	0.40
1:A:539:PHE:O	1:A:543:VAL:HG22	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	360/378 (95%)	350 (97%)	5 (1%)	5 (1%)	9	3
2	B	9/28 (32%)	8 (89%)	1 (11%)	0	100	100
All	All	369/406 (91%)	358 (97%)	6 (2%)	5 (1%)	9	3

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	202	LYS
1	A	344	TYR
1	A	499	PRO
1	A	213	ASP
1	A	345	THR

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	319/335 (95%)	302 (95%)	17 (5%)	20	13
2	B	11/26 (42%)	8 (73%)	3 (27%)	0	0
All	All	330/361 (91%)	310 (94%)	20 (6%)	17	10

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	223	ARG
1	A	316	LEU
1	A	322	LEU
1	A	333	TYR
1	A	344	TYR
1	A	402	ARG
1	A	405	LEU
1	A	456	ILE
1	A	476	THR
1	A	482	GLU
1	A	490	VAL
1	A	493	SER
1	A	500	THR
1	A	505	VAL
1	A	523	GLN
1	A	552	THR
1	A	557	VAL
2	B	155	ARG

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Mol	Chain	Res	Type
2	B	159	GLN
2	B	163	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	225	ASN
1	A	241	GLN
1	A	376	ASN
1	A	446	GLN
1	A	448	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	362/378 (95%)	-0.08	10 (2%) 55 54	12, 26, 50, 75	0
2	B	11/28 (39%)	0.19	0 100 100	21, 30, 70, 80	0
All	All	373/406 (91%)	-0.07	10 (2%) 56 56	12, 27, 51, 80	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	344	TYR	5.1
1	A	200	THR	3.6
1	A	201	LEU	3.1
1	A	345	THR	3.0
1	A	453	ILE	2.8
1	A	456	ILE	2.7
1	A	500	THR	2.4
1	A	303	VAL	2.3
1	A	497	LYS	2.1
1	A	440	MET	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.