



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 01:32 pm BST

PDB ID : 1Z81
Title : Crystal Structure of cyclophilin from Plasmodium yoelii.
Authors : Mulichak, A.; Alam, Z.; Amani, M.; Lew, J.; Wasney, G.; Sundstrom, M.; Arrowsmith, C.; Edwards, A.; Bochkarev, A.; Hui, R.; Vedadi, M.; Structural Genomics Consortium (SGC)
Deposited on : 2005-03-29
Resolution : 2.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

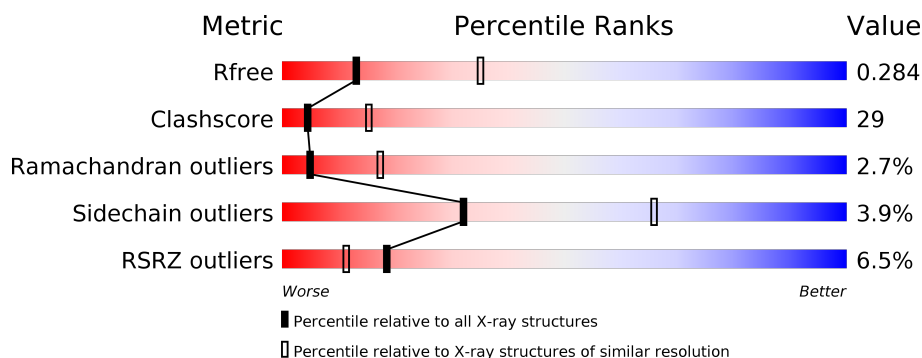
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 on 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	229	<div> <div>5%</div> <div> <div></div> <div>40%</div> <div>38%</div> <div>•</div> <div>19%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1455 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 cyclophilin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	186	Total	C	N	O	S	2	0	0
			1439	931	235	266	7			

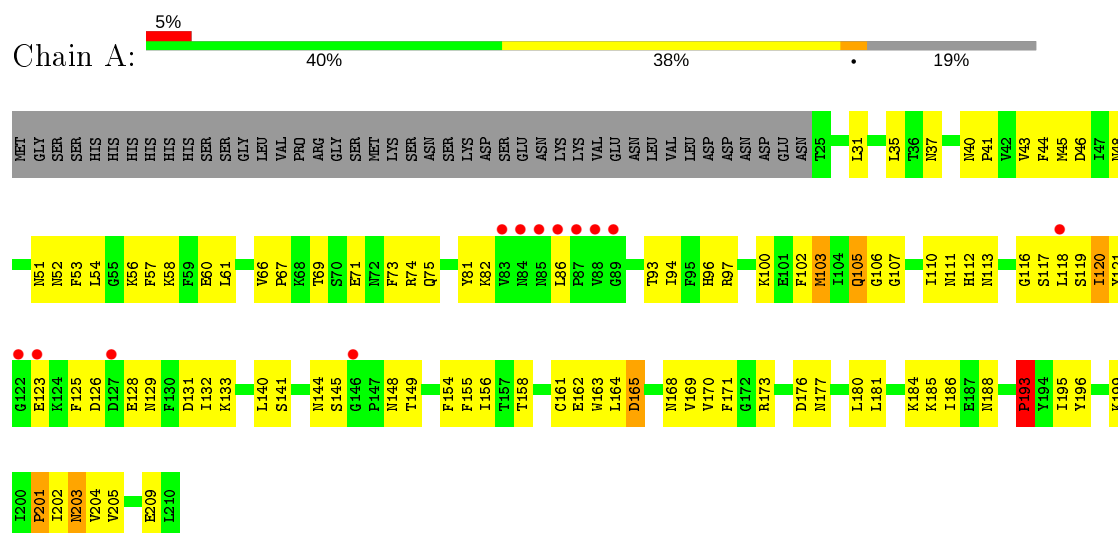
There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	CLONING ARTIFACT	UNP Q7RSH5
A	-17	GLY	-	CLONING ARTIFACT	UNP Q7RSH5
A	-16	SER	-	CLONING ARTIFACT	UNP Q7RSH5
A	-15	SER	-	CLONING ARTIFACT	UNP Q7RSH5
A	-14	HIS	-	CLONING ARTIFACT	UNP Q7RSH5
A	-13	HIS	-	CLONING ARTIFACT	UNP Q7RSH5
A	-12	HIS	-	CLONING ARTIFACT	UNP Q7RSH5
A	-11	HIS	-	CLONING ARTIFACT	UNP Q7RSH5
A	-10	HIS	-	CLONING ARTIFACT	UNP Q7RSH5
A	-9	HIS	-	CLONING ARTIFACT	UNP Q7RSH5
A	-8	SER	-	CLONING ARTIFACT	UNP Q7RSH5
A	-7	SER	-	CLONING ARTIFACT	UNP Q7RSH5
A	-6	GLY	-	CLONING ARTIFACT	UNP Q7RSH5
A	-5	LEU	-	CLONING ARTIFACT	UNP Q7RSH5
A	-4	VAL	-	CLONING ARTIFACT	UNP Q7RSH5
A	-3	PRO	-	CLONING ARTIFACT	UNP Q7RSH5
A	-2	ARG	-	CLONING ARTIFACT	UNP Q7RSH5
A	-1	GLY	-	CLONING ARTIFACT	UNP Q7RSH5
A	0	SER	-	CLONING ARTIFACT	UNP Q7RSH5

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	16	Total	O	0	0
			16	16		

- Molecule 1: cyclophilin



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	64.23 Å 64.23 Å 156.56 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 40.50 – 2.75	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.00-2.80) 96.4 (40.50-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.47 (at 2.77 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.257 , 0.285 0.256 , 0.284	Depositor DCC
R_{free} test set	493 reflections (5.48%)	wwPDB-VP
Wilson B-factor (Å ²)	73.4	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 68.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	1455	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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	0.37	0/1473	0.62	0/1999

There are no bond length outliers.

There are no bond angle outliers.

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	1439	0	1370	80	1
2	A	16	0	0	3	0
All	All	1455	0	1370	80	1

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 29.

All (80) 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:48:ASN:HD22	1:A:205:VAL:HG21	1.28	0.95
1:A:118:LEU:HD23	1:A:119:SER:N	1.89	0.88
1:A:81:TYR:HE2	1:A:120:ILE:HA	1.39	0.87
1:A:110:ILE:HG22	1:A:111:ASN:ND2	1.91	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:LYS:HA	1:A:86:LEU:O	1.76	0.85
1:A:66:VAL:HA	1:A:128:GLU:OE2	1.79	0.83
1:A:75:GLN:HG3	1:A:120:ILE:HG21	1.63	0.81
1:A:69:THR:HG22	1:A:125:PHE:HE2	1.48	0.77
1:A:48:ASN:HD22	1:A:205:VAL:CG2	1.97	0.76
1:A:110:ILE:HD12	1:A:117:SER:N	2.01	0.75
1:A:48:ASN:ND2	1:A:205:VAL:HG21	2.00	0.75
1:A:94:ILE:HG22	1:A:203:ASN:HB3	1.67	0.75
1:A:186:ILE:HG23	1:A:202:ILE:HD13	1.68	0.73
1:A:145:SER:OG	1:A:149:THR:HG21	1.89	0.73
1:A:81:TYR:CE2	1:A:120:ILE:HA	2.23	0.73
1:A:71:GLU:O	1:A:75:GLN:HG2	1.87	0.72
1:A:110:ILE:HB	1:A:116:GLY:HA3	1.73	0.71
1:A:132:ILE:HB	1:A:170:VAL:HB	1.71	0.70
1:A:161:CYS:SG	1:A:164:LEU:HD12	2.32	0.69
1:A:100:LYS:HE2	1:A:188:ASN:OD1	1.93	0.67
1:A:48:ASN:OD1	1:A:53:PHE:HA	1.95	0.67
1:A:203:ASN:H	1:A:203:ASN:ND2	1.95	0.65
1:A:199:LYS:O	1:A:201:PRO:HD3	1.97	0.64
1:A:61:LEU:HD21	1:A:73:PHE:CD1	2.34	0.62
1:A:144:ASN:HB2	1:A:149:THR:HB	1.83	0.61
1:A:111:ASN:O	1:A:112:HIS:HB2	2.00	0.60
1:A:93:THR:HG23	1:A:94:ILE:N	2.19	0.58
1:A:128:GLU:O	1:A:129:ASN:HB3	2.05	0.57
1:A:46:ASP:OD1	1:A:56:LYS:HE2	2.04	0.57
1:A:165:ASP:HA	2:A:214:HOH:O	2.04	0.56
1:A:177:ASN:O	1:A:181:LEU:HG	2.07	0.55
1:A:168:ASN:HA	2:A:213:HOH:O	2.07	0.54
1:A:37:ASN:HB3	1:A:40:ASN:ND2	2.23	0.54
1:A:141:SER:HB3	1:A:155:PHE:CZ	2.44	0.53
1:A:149:THR:O	1:A:149:THR:HG22	2.08	0.53
1:A:203:ASN:ND2	1:A:203:ASN:N	2.58	0.52
1:A:164:LEU:HD22	1:A:168:ASN:OD1	2.10	0.51
1:A:102:PHE:C	1:A:158:THR:HG22	2.30	0.51
1:A:105:GLN:HB2	1:A:155:PHE:HB3	1.94	0.50
1:A:180:LEU:O	1:A:184:LYS:HG2	2.12	0.50
1:A:126:ASP:HA	1:A:148:ASN:ND2	2.26	0.50
1:A:37:ASN:O	1:A:40:ASN:HB2	2.12	0.50
1:A:45:MET:HB3	1:A:204:VAL:HG13	1.93	0.50
1:A:126:ASP:HA	1:A:148:ASN:HD22	1.77	0.49
1:A:184:LYS:HA	1:A:184:LYS:HE2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LEU:HD21	1:A:73:PHE:HD1	1.77	0.49
1:A:107:GLY:HA3	1:A:154:PHE:HE1	1.78	0.49
1:A:195:ILE:O	1:A:196:TYR:HB2	2.12	0.49
1:A:51:ASN:C	1:A:52:ASN:HD22	2.16	0.48
1:A:81:TYR:CD2	1:A:120:ILE:HG22	2.48	0.48
1:A:81:TYR:CE2	1:A:120:ILE:HG22	2.48	0.48
1:A:169:VAL:N	2:A:213:HOH:O	2.34	0.48
1:A:133:LYS:HD2	1:A:165:ASP:OD2	2.13	0.48
1:A:96:HIS:CE1	1:A:106:GLY:HA2	2.49	0.47
1:A:110:ILE:HG22	1:A:111:ASN:HD21	1.75	0.47
1:A:203:ASN:HD22	1:A:203:ASN:H	1.62	0.47
1:A:35:LEU:HD22	1:A:60:GLU:HB2	1.98	0.46
1:A:140:LEU:HD23	1:A:156:ILE:HG12	1.96	0.46
1:A:48:ASN:OD1	1:A:53:PHE:CA	2.63	0.46
1:A:54:LEU:HD11	1:A:185:LYS:HD2	1.99	0.45
1:A:44:PHE:HA	1:A:57:PHE:O	2.17	0.45
1:A:110:ILE:C	1:A:111:ASN:HD22	2.20	0.45
1:A:67:PRO:HD2	1:A:128:GLU:OE2	2.17	0.45
1:A:111:ASN:HB3	1:A:113:ASN:OD1	2.18	0.44
1:A:162:GLU:O	1:A:164:LEU:N	2.51	0.44
1:A:43:VAL:HA	1:A:209:GLU:HA	2.00	0.44
1:A:107:GLY:HA3	1:A:154:PHE:CE1	2.52	0.44
1:A:110:ILE:HD12	1:A:117:SER:H	1.79	0.44
1:A:73:PHE:CG	1:A:171:PHE:HE2	2.36	0.44
1:A:119:SER:C	1:A:121:TYR:H	2.22	0.44
1:A:97:ARG:HG3	1:A:97:ARG:HH21	1.83	0.43
1:A:69:THR:HG22	1:A:125:PHE:CE2	2.39	0.43
1:A:102:PHE:CD2	1:A:103:MET:HG3	2.54	0.43
1:A:58:LYS:HD2	1:A:176:ASP:OD2	2.19	0.42
1:A:129:ASN:CG	1:A:131:ASP:OD2	2.58	0.42
1:A:118:LEU:HA	1:A:123:GLU:HA	2.02	0.41
1:A:110:ILE:HD11	1:A:118:LEU:HB2	2.00	0.41
1:A:186:ILE:HG23	1:A:202:ILE:CD1	2.44	0.41
1:A:40:ASN:HA	1:A:41:PRO:HD3	1.79	0.40
1:A:74:ARG:HH11	1:A:74:ARG:HG2	1.87	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:PRO:O	1:A:196:TYR:OH[8_665]	2.15	0.05

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	184/229 (80%)	157 (85%)	22 (12%)	5 (3%)	5 17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	163	TRP
1	A	165	ASP
1	A	193	PRO
1	A	201	PRO
1	A	120	ILE

5.3.2 Protein sidechains [i](#)

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	154/210 (73%)	148 (96%)	6 (4%)	32 66

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	LEU
1	A	103	MET
1	A	105	GLN
1	A	173	ARG
1	A	193	PRO
1	A	203	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	63	GLN
1	A	105	GLN
1	A	111	ASN
1	A	134	HIS
1	A	153	GLN
1	A	203	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 carbohydrates 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	186/229 (81%)	0.12	12 (6%) 18 11	42, 73, 101, 101	2 (1%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	88	VAL	5.9
1	A	85	ASN	4.1
1	A	122	GLY	3.7
1	A	84	ASN	3.6
1	A	118	LEU	2.9
1	A	87	PRO	2.8
1	A	146	GLY	2.7
1	A	123	GLU	2.5
1	A	83	VAL	2.4
1	A	86	LEU	2.3
1	A	127	ASP	2.2
1	A	89	GLY	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 carbohydrates 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.