



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2018 – 08:06 am GMT

PDB ID : 2E8G
Title : The structure of protein from *P. horikoshii* at 1.7 angstrom resolution
Authors : Gao, Y.G.; Yao, M.; Tanaka, I.
Deposited on : 2007-01-19
Resolution : 1.70 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 : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

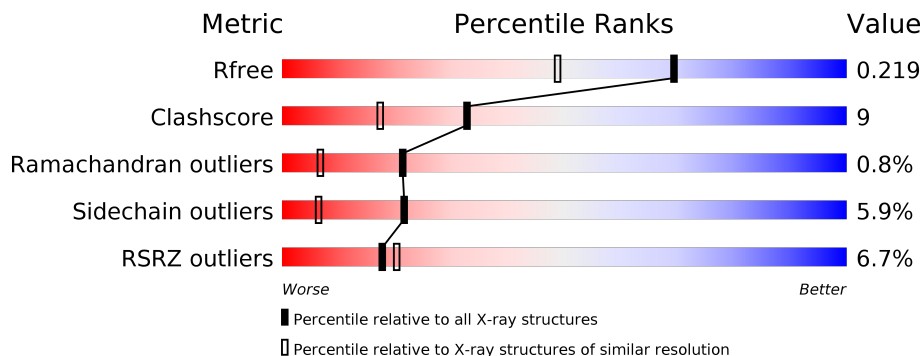
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	111664	3793 (1.70-1.70)
Clashscore	122126	4167 (1.70-1.70)
Ramachandran outliers	120053	4100 (1.70-1.70)
Sidechain outliers	120020	4100 (1.70-1.70)
RSRZ outliers	108989	3718 (1.70-1.70)

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. 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	241	<div> <div>10%</div> <div> <div></div> <div>83%</div> <div>15%</div> </div> <div></div> </div>
1	B	241	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>14%</div> </div> <div></div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4155 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 Hypothetical protein PH0536.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	0	0	0
			1891	1204	332	350	5			
1	B	240	Total	C	N	O	S	0	0	0
			1891	1204	332	350	5			

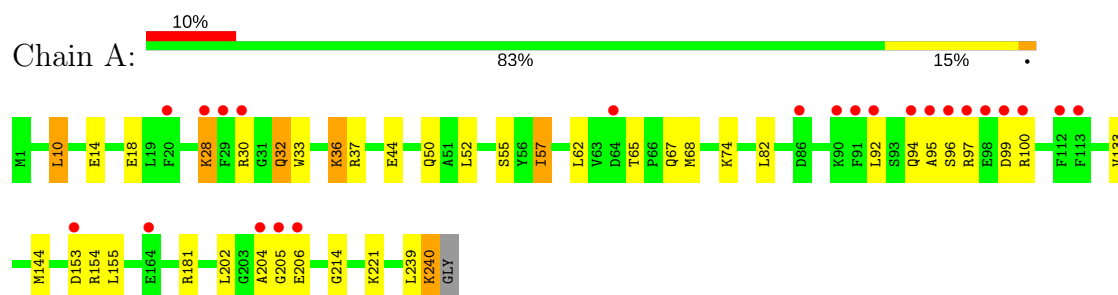
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	184	Total	O	0	0
			184	184		
2	B	189	Total	O	0	0
			189	189		

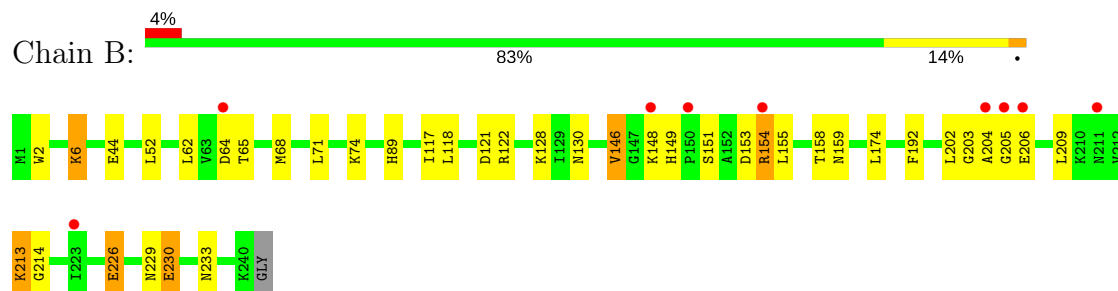
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Hypothetical protein PH0536



• Molecule 1: Hypothetical protein PH0536



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.19Å 71.69Å 69.19Å 90.00° 96.15° 90.00°	Depositor
Resolution (Å)	34.40 – 1.70 34.40 – 1.70	Depositor EDS
% Data completeness (in resolution range)	9.9 (34.40-1.70) 99.7 (34.40-1.70)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.85 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.199 , 0.237 0.219 , 0.219	Depositor DCC
R_{free} test set	5287 reflections (10.08%)	wwPDB-VP
Wilson B-factor (Å ²)	21.4	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4155	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.97 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.2732e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

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	0.65	0/1920	0.73	1/2587 (0.0%)
1	B	0.65	0/1920	0.72	1/2587 (0.0%)
All	All	0.65	0/3840	0.72	2/5174 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	121	ASP	CB-CG-OD1	5.84	123.56	118.30
1	A	10	LEU	CB-CG-CD1	5.55	120.43	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1891	0	1958	34	0
1	B	1891	0	1958	38	0
2	A	184	0	0	4	1
2	B	189	0	0	10	0
All	All	4155	0	3916	72	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 9.

All (72) 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:B:213:LYS:HD3	1:B:214:GLY:N	1.83	0.94
1:B:44:GLU:HB3	1:B:74:LYS:HE2	1.53	0.89
1:A:36:LYS:O	1:A:36:LYS:HE3	1.80	0.81
1:A:28:LYS:HE3	2:A:405:HOH:O	1.81	0.81
1:A:204:ALA:O	1:A:206:GLU:N	2.14	0.80
1:B:130:ASN:O	2:B:401:HOH:O	1.98	0.80
1:A:65:THR:HG23	1:A:68:MET:H	1.45	0.80
1:A:28:LYS:HD3	1:A:30:ARG:HG3	1.65	0.78
1:A:28:LYS:HG2	1:A:30:ARG:NH1	2.02	0.74
1:B:213:LYS:HD3	1:B:214:GLY:H	1.52	0.74
1:A:62:LEU:O	1:A:65:THR:HG22	1.88	0.73
1:B:149:HIS:HE1	1:B:192:PHE:O	1.73	0.72
1:B:65:THR:HG23	1:B:68:MET:H	1.54	0.72
1:A:95:ALA:HA	1:A:99:ASP:HB3	1.74	0.70
1:B:226:GLU:H	1:B:226:GLU:CD	1.96	0.67
1:B:213:LYS:HA	1:B:213:LYS:HE2	1.79	0.64
1:A:14:GLU:OE2	1:A:50:GLN:NE2	2.32	0.63
1:A:52:LEU:HD21	1:A:65:THR:HG21	1.79	0.63
1:B:203:GLY:HA2	1:B:209:LEU:CD1	2.29	0.63
1:B:130:ASN:HA	2:B:337:HOH:O	1.99	0.61
1:A:28:LYS:HE2	1:A:30:ARG:HH11	1.66	0.60
1:A:30:ARG:HG3	1:A:33:TRP:HE1	1.66	0.60
1:B:229:ASN:HB3	2:B:300:HOH:O	2.01	0.60
1:A:154:ARG:NH2	2:A:390:HOH:O	2.34	0.59
1:B:213:LYS:CD	1:B:214:GLY:H	2.16	0.59
1:A:28:LYS:HD3	1:A:30:ARG:CG	2.31	0.58
2:A:402:HOH:O	1:B:213:LYS:HD2	2.05	0.56
1:B:52:LEU:HD21	1:B:65:THR:HG21	1.87	0.56
1:A:65:THR:OG1	1:A:67:GLN:OE1	2.24	0.56
1:A:239:LEU:O	1:A:240:LYS:HB2	2.05	0.56
1:A:14:GLU:CD	1:A:50:GLN:NE2	2.59	0.55
1:B:149:HIS:CE1	1:B:192:PHE:O	2.58	0.55
1:A:28:LYS:HG2	1:A:30:ARG:HH11	1.70	0.55
1:B:203:GLY:HA2	1:B:209:LEU:HD13	1.88	0.55
1:A:92:LEU:O	1:A:95:ALA:HB3	2.05	0.54
1:B:149:HIS:HD2	1:B:151:SER:H	1.54	0.54
1:A:95:ALA:HB1	1:A:100:ARG:HA	1.89	0.54
1:B:213:LYS:CD	1:B:214:GLY:N	2.63	0.54
1:B:130:ASN:C	2:B:401:HOH:O	2.42	0.54
1:B:65:THR:CG2	1:B:68:MET:H	2.20	0.53
1:B:89:HIS:HE1	2:B:291:HOH:O	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:GLN:HE22	1:A:94:GLN:NE2	2.08	0.51
1:B:71:LEU:HG	1:B:117:ILE:HG23	1.93	0.51
1:A:30:ARG:NH1	1:A:94:GLN:HG2	2.26	0.51
1:B:213:LYS:CE	1:B:214:GLY:H	2.25	0.49
1:A:28:LYS:CG	1:A:30:ARG:NH1	2.73	0.49
1:B:52:LEU:CD2	1:B:65:THR:HG21	2.42	0.49
1:A:18:GLU:OE1	2:A:402:HOH:O	2.20	0.49
1:B:159:ASN:ND2	2:B:316:HOH:O	2.45	0.49
1:B:149:HIS:HD2	1:B:151:SER:N	2.11	0.48
1:B:153:ASP:O	1:B:154:ARG:HG2	2.13	0.47
1:B:204:ALA:O	1:B:206:GLU:N	2.47	0.47
1:A:28:LYS:HE2	1:A:30:ARG:HD2	1.96	0.47
1:A:28:LYS:CE	1:A:30:ARG:HD2	2.45	0.47
1:A:30:ARG:CZ	1:A:94:GLN:HG2	2.45	0.47
1:B:230:GLU:CG	2:B:428:HOH:O	2.63	0.46
1:B:230:GLU:HG3	2:B:428:HOH:O	2.16	0.46
1:A:214:GLY:HA2	1:A:221:LYS:CE	2.46	0.46
1:B:146:VAL:HG22	1:B:158:THR:HG22	1.99	0.45
1:B:62:LEU:O	1:B:65:THR:HG22	2.16	0.45
1:A:144:MET:HB2	1:A:144:MET:HE3	1.89	0.43
1:B:6:LYS:HB3	1:B:6:LYS:NZ	2.34	0.43
1:A:44:GLU:HB3	1:A:74:LYS:HE3	2.01	0.42
1:A:28:LYS:CG	1:A:30:ARG:HH11	2.31	0.42
1:A:181:ARG:HA	1:A:181:ARG:HD2	1.74	0.42
1:B:233:ASN:ND2	2:B:362:HOH:O	2.54	0.41
1:B:2:TRP:CH2	1:B:130:ASN:HB2	2.55	0.41
1:B:154:ARG:C	1:B:174:LEU:HD21	2.41	0.41
1:A:28:LYS:HG3	1:A:28:LYS:H	1.62	0.40
1:A:55:SER:O	1:A:57:ILE:HD12	2.20	0.40
1:B:2:TRP:CE3	1:B:128:LYS:HG2	2.56	0.40
1:B:130:ASN:HB3	2:B:274:HOH:O	2.21	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)
2:A:350:HOH:O	2:A:353:HOH:O[2_546]	2.10	0.10

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	238/241 (99%)	233 (98%)	2 (1%)	3 (1%)	13	2
1	B	238/241 (99%)	231 (97%)	6 (2%)	1 (0%)	36	19
All	All	476/482 (99%)	464 (98%)	8 (2%)	4 (1%)	21	7

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	205	GLY
1	B	205	GLY
1	A	97	ARG
1	A	96	SER

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	204/204 (100%)	192 (94%)	12 (6%)	21	6
1	B	204/204 (100%)	192 (94%)	12 (6%)	21	6
All	All	408/408 (100%)	384 (94%)	24 (6%)	21	6

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	28	LYS

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Mol	Chain	Res	Type
1	A	32	GLN
1	A	36	LYS
1	A	37	ARG
1	A	57	ILE
1	A	82	LEU
1	A	133	VAL
1	A	153	ASP
1	A	155	LEU
1	A	202	LEU
1	A	240	LYS
1	B	6	LYS
1	B	64	ASP
1	B	118	LEU
1	B	122	ARG
1	B	146	VAL
1	B	148	LYS
1	B	154	ARG
1	B	155	LEU
1	B	202	LEU
1	B	213	LYS
1	B	226	GLU
1	B	230	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	88	HIS
1	A	94	GLN
1	A	159	ASN
1	B	40	GLN
1	B	89	HIS
1	B	149	HIS
1	B	159	ASN
1	B	233	ASN

5.3.3 RNA ⓘ

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

6.1 Protein, DNA and RNA chains

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	240/241 (99%)	0.54	23 (9%) 8 9	14, 20, 32, 44	0
1	B	240/241 (99%)	0.41	9 (3%) 40 46	13, 20, 29, 36	0
All	All	480/482 (99%)	0.47	32 (6%) 18 20	13, 20, 31, 44	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	95	ALA	8.5
1	B	205	GLY	7.3
1	A	96	SER	7.1
1	B	204	ALA	5.9
1	B	64	ASP	4.9
1	A	204	ALA	4.4
1	B	154	ARG	4.4
1	A	98	GLU	4.1
1	A	153	ASP	4.1
1	A	97	ARG	3.9
1	A	29	PHE	3.9
1	A	94	GLN	3.9
1	A	205	GLY	3.8
1	A	91	PHE	3.7
1	A	206	GLU	3.2
1	B	148	LYS	3.0
1	B	206	GLU	3.0
1	A	30	ARG	2.8
1	A	92	LEU	2.7
1	A	86	ASP	2.7
1	A	90	LYS	2.6
1	A	99	ASP	2.3
1	A	100	ARG	2.3
1	A	164	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	150	PRO	2.2
1	A	113	PHE	2.2
1	B	223	ILE	2.1
1	A	64	ASP	2.1
1	A	28	LYS	2.1
1	B	211	ASN	2.1
1	A	20	PHE	2.0
1	A	112	PHE	2.0

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.