



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:02 AM UTC

PDB ID : 4QMB / pdb_00004qmb
Title : The structure of inorganic pyrophosphatase from Schistosoma japonicum
Authors : Wu, Q.F.
Deposited on : 2014-06-15
Resolution : 2.60 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

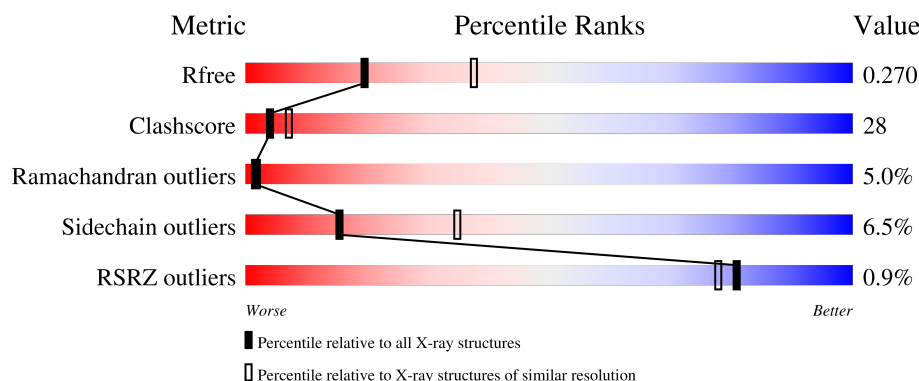
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	287	% 64% 26% 7% ..
1	B	287	% 67% 24% 5% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4618 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 SJCHGC07024 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2271	1458	377	424	12			
1	B	283	Total	C	N	O	S	0	0	0
			2259	1450	375	421	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	214	ALA	VAL	engineered mutation	UNP Q5DE13
B	214	ALA	VAL	engineered mutation	UNP Q5DE13

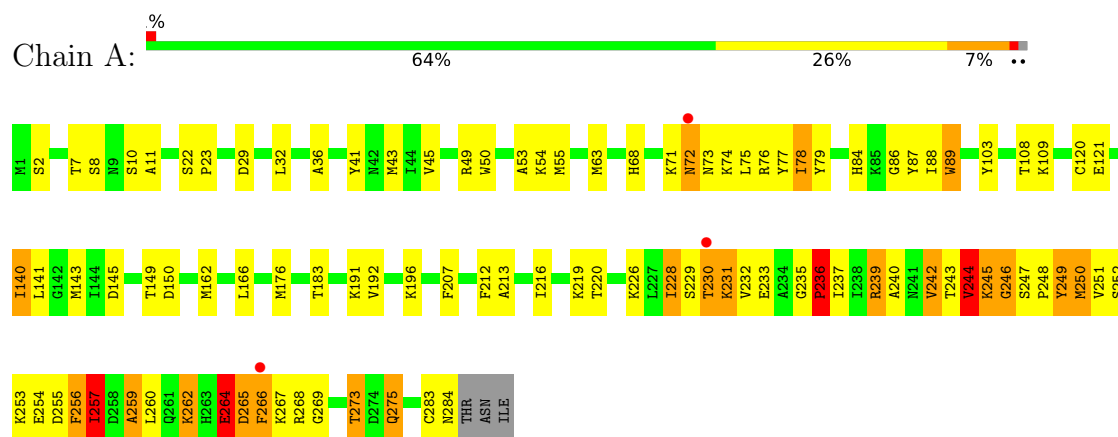
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	38	Total	O	0	0
			38	38		
2	B	50	Total	O	0	0
			50	50		

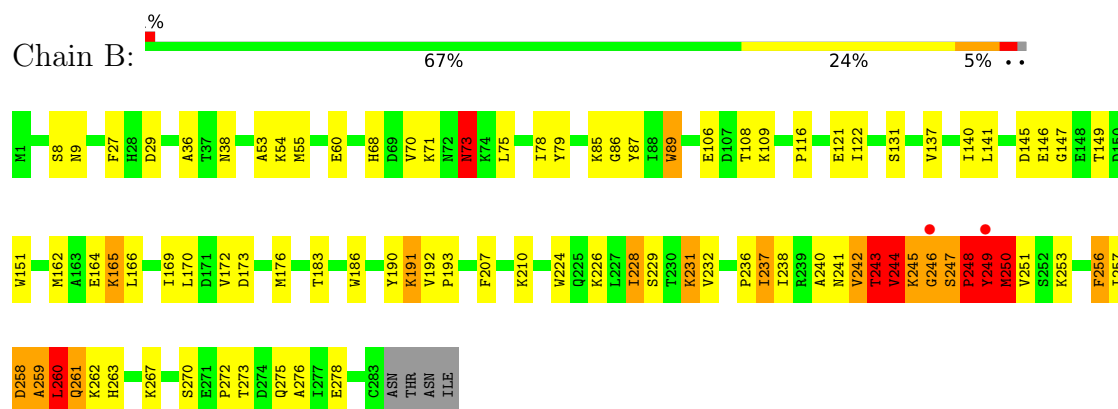
3 Residue-property plots [i](#)

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: SJCHGC07024 protein



• Molecule 1: SJCHGC07024 protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	75.75Å 75.75Å 122.96Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.88 – 2.60 37.88 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.6 (37.88-2.60) 98.1 (37.88-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	13.36 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.202 , 0.263 0.219 , 0.270	Depositor DCC
R_{free} test set	1985 reflections (8.21%)	wwPDB-VP
Wilson B-factor (Å ²)	47.8	Xtriage
Anisotropy	0.117	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l 0.468 for h,-h-k,-l 0.018 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4618	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.58% 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	0.67	0/2336	1.03	12/3168 (0.4%)
1	B	0.66	0/2323	0.96	9/3150 (0.3%)
All	All	0.66	0/4659	1.00	21/6318 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	4
All	All	0	6

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	243	THR	N-CA-C	-10.08	93.90	109.52
1	B	248	PRO	N-CA-C	-8.58	94.80	112.47
1	B	241	ASN	N-CA-C	8.01	122.03	110.10
1	A	72	ASN	N-CA-C	-6.87	104.85	113.23
1	A	249	TYR	CA-C-N	-6.40	109.32	121.54
1	A	249	TYR	C-N-CA	-6.40	109.32	121.54
1	A	192	VAL	CA-C-N	-6.14	113.30	119.56
1	A	192	VAL	C-N-CA	-6.14	113.30	119.56
1	A	257	ILE	N-CA-C	-6.09	96.66	109.34
1	A	262	LYS	CA-C-N	-5.93	114.89	122.77
1	A	262	LYS	C-N-CA	-5.93	114.89	122.77
1	B	173	ASP	N-CA-C	5.88	117.69	111.28
1	B	192	VAL	CA-C-N	-5.52	114.12	119.87
1	B	192	VAL	C-N-CA	-5.52	114.12	119.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	265	ASP	N-CA-C	5.50	122.52	110.80
1	A	259	ALA	N-CA-C	-5.34	104.71	111.11
1	A	255	ASP	N-CA-C	5.29	116.04	108.74
1	B	260	LEU	CA-CB-CG	5.20	134.51	116.30
1	A	264	GLU	N-CA-C	5.19	116.78	107.80
1	B	260	LEU	CA-C-N	-5.15	114.72	122.24
1	B	260	LEU	C-N-CA	-5.15	114.72	122.24

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	236	PRO	Peptide
1	A	244	VAL	Peptide
1	B	244	VAL	Peptide
1	B	249	TYR	Peptide
1	B	250	MET	Peptide
1	B	261	GLN	Peptide

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	2271	0	2198	132	0
1	B	2259	0	2195	119	0
2	A	38	0	0	8	0
2	B	50	0	0	2	0
All	All	4618	0	4393	248	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 28.

All (248) 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:236:PRO:O	1:A:237:ILE:HG22	1.16	1.32

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:LYS:N	1:B:248:PRO:HD2	1.40	1.31
1:B:245:LYS:HG3	1:B:249:TYR:CB	1.60	1.30
1:B:244:VAL:C	1:B:248:PRO:HD2	1.60	1.26
1:A:247:SER:HB2	1:A:249:TYR:CD2	1.71	1.24
1:A:248:PRO:CD	1:A:249:TYR:HA	1.67	1.21
1:A:249:TYR:O	1:A:250:MET:C	1.75	1.20
1:A:246:GLY:HA3	1:A:247:SER:OG	1.35	1.20
1:B:29:ASP:OD1	1:B:243:THR:O	1.61	1.17
1:B:247:SER:HA	1:B:249:TYR:HD1	1.06	1.17
1:B:245:LYS:HB2	1:B:249:TYR:HA	1.25	1.16
1:A:247:SER:CB	1:A:249:TYR:HD2	1.58	1.16
1:B:245:LYS:CG	1:B:249:TYR:CB	2.24	1.16
1:A:247:SER:CB	1:A:249:TYR:CD2	2.30	1.15
1:B:245:LYS:HG3	1:B:249:TYR:HB2	1.28	1.13
1:B:245:LYS:HA	1:B:246:GLY:C	1.71	1.12
1:B:243:THR:O	1:B:244:VAL:HB	1.51	1.10
1:A:248:PRO:HD2	1:A:249:TYR:CA	1.66	1.09
1:A:248:PRO:HB3	1:A:254:GLU:OE1	1.51	1.08
1:A:248:PRO:HD2	1:A:249:TYR:HA	1.16	1.07
1:A:247:SER:C	1:A:249:TYR:HA	1.82	1.04
1:A:236:PRO:O	1:A:237:ILE:CG2	2.06	1.02
1:B:245:LYS:H	1:B:248:PRO:CD	1.71	1.01
1:B:245:LYS:CG	1:B:249:TYR:HB3	1.89	1.00
1:B:247:SER:O	1:B:249:TYR:CD1	2.13	0.99
1:B:245:LYS:HG3	1:B:249:TYR:CA	1.93	0.99
1:B:247:SER:HA	1:B:249:TYR:CD1	1.98	0.98
1:A:249:TYR:O	1:A:251:VAL:N	1.96	0.97
1:A:248:PRO:N	1:A:249:TYR:HA	1.79	0.97
1:B:245:LYS:CB	1:B:249:TYR:HA	1.95	0.97
1:A:236:PRO:C	1:A:237:ILE:HG22	1.89	0.96
1:B:245:LYS:HG2	1:B:249:TYR:HB3	1.46	0.96
1:B:247:SER:CA	1:B:249:TYR:HD1	1.80	0.95
1:A:248:PRO:CD	1:A:249:TYR:CA	2.30	0.94
1:B:244:VAL:O	1:B:248:PRO:HD2	1.69	0.92
1:A:251:VAL:HG12	1:A:253:LYS:H	1.38	0.87
1:B:245:LYS:H	1:B:248:PRO:HD2	1.07	0.85
1:A:247:SER:OG	1:A:249:TYR:CD2	2.30	0.84
1:B:226:LYS:O	1:B:231:LYS:HG2	1.78	0.84
1:A:247:SER:OG	1:A:249:TYR:CE2	2.30	0.83
1:A:236:PRO:C	1:A:237:ILE:CG2	2.52	0.82
1:B:247:SER:O	1:B:249:TYR:CG	2.28	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:244:VAL:O	1:B:248:PRO:CG	2.30	0.80
1:A:29:ASP:OD1	1:A:244:VAL:HB	1.80	0.80
1:A:248:PRO:HB2	1:A:251:VAL:H	1.43	0.80
1:B:244:VAL:O	1:B:248:PRO:CD	2.30	0.80
1:B:244:VAL:O	1:B:248:PRO:HB2	1.80	0.79
1:B:244:VAL:C	1:B:248:PRO:CD	2.52	0.78
1:A:242:VAL:HG23	1:A:244:VAL:H	1.49	0.78
1:B:231:LYS:HD3	1:B:231:LYS:N	1.98	0.78
1:B:245:LYS:HB2	1:B:249:TYR:CA	2.13	0.78
1:B:231:LYS:HD3	1:B:231:LYS:H	1.48	0.77
1:B:245:LYS:CG	1:B:249:TYR:CA	2.60	0.77
1:A:240:ALA:O	1:A:242:VAL:HG12	1.85	0.77
1:A:244:VAL:HA	1:A:245:LYS:HB2	1.65	0.76
1:A:63:MET:HE1	1:A:262:LYS:HB3	1.68	0.75
1:A:244:VAL:HG22	1:A:245:LYS:HB2	1.67	0.74
1:B:245:LYS:N	1:B:248:PRO:CD	2.30	0.74
1:A:55:MET:HE1	1:A:68:HIS:ND1	2.02	0.73
1:A:226:LYS:NZ	2:A:308:HOH:O	2.20	0.73
1:B:164:GLU:HG3	1:B:165:LYS:HD2	1.70	0.73
1:A:275:GLN:N	2:A:303:HOH:O	2.20	0.73
1:A:247:SER:C	1:A:249:TYR:CA	2.60	0.72
1:B:55:MET:HE1	1:B:68:HIS:ND1	2.03	0.72
1:B:251:VAL:HG12	1:B:253:LYS:H	1.55	0.72
1:B:244:VAL:O	1:B:248:PRO:CB	2.37	0.71
1:B:245:LYS:CG	1:B:249:TYR:HA	2.18	0.71
1:B:106:GLU:O	1:B:109:LYS:NZ	2.23	0.70
1:B:240:ALA:O	1:B:242:VAL:HG12	1.91	0.70
1:A:259:ALA:O	1:A:262:LYS:HG2	1.91	0.70
1:B:245:LYS:HG3	1:B:249:TYR:HA	1.71	0.70
1:B:248:PRO:O	1:B:250:MET:N	2.25	0.70
1:B:242:VAL:HG21	1:B:248:PRO:HB2	1.74	0.69
1:B:8:SER:O	1:B:270:SER:OG	2.07	0.69
1:B:228:ILE:O	1:B:251:VAL:HG13	1.92	0.69
1:B:245:LYS:CG	1:B:249:TYR:HB2	2.06	0.69
1:A:140:ILE:HD11	1:A:212:PHE:HD2	1.58	0.68
1:A:231:LYS:HD3	1:A:233:GLU:OE1	1.94	0.68
1:B:242:VAL:HG21	1:B:248:PRO:CB	2.24	0.67
1:B:243:THR:O	1:B:244:VAL:CB	2.29	0.67
1:A:246:GLY:HA3	1:A:247:SER:CB	2.16	0.67
1:B:260:LEU:HG	1:B:261:GLN:OE1	1.94	0.67
1:B:244:VAL:HA	1:B:245:LYS:C	2.20	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:LYS:HA	1:B:247:SER:N	2.10	0.66
1:A:268:ARG:NH1	1:A:269:GLY:O	2.29	0.66
1:B:145:ASP:O	1:B:147:GLY:N	2.29	0.66
1:B:78:ILE:HD11	1:B:86:GLY:HA2	1.78	0.65
1:A:228:ILE:O	1:A:251:VAL:HG13	1.97	0.65
1:A:284:ASN:OD1	2:A:333:HOH:O	2.13	0.64
1:A:150:ASP:OD1	2:A:304:HOH:O	2.14	0.64
1:A:162:MET:HE3	1:A:166:LEU:HD11	1.78	0.64
1:A:244:VAL:HG22	1:A:245:LYS:CB	2.27	0.64
1:B:236:PRO:O	1:B:238:ILE:N	2.30	0.64
1:B:247:SER:C	1:B:249:TYR:CD1	2.75	0.64
1:B:78:ILE:HG12	1:B:87:TYR:CE1	2.33	0.64
1:B:245:LYS:CB	1:B:249:TYR:CA	2.70	0.64
1:A:257:ILE:O	2:A:324:HOH:O	2.15	0.63
1:A:140:ILE:HD11	1:A:212:PHE:CD2	2.33	0.63
1:B:258:ASP:O	1:B:260:LEU:N	2.32	0.63
1:A:235:GLY:N	1:A:236:PRO:HD2	2.14	0.62
1:B:245:LYS:CA	1:B:246:GLY:C	2.61	0.62
1:A:247:SER:CA	1:A:249:TYR:HA	2.29	0.62
1:A:244:VAL:CA	1:A:245:LYS:HB2	2.30	0.61
1:A:247:SER:OG	1:A:249:TYR:HE2	1.81	0.61
1:B:258:ASP:C	1:B:260:LEU:H	2.09	0.60
1:A:219:LYS:NZ	2:A:335:HOH:O	2.17	0.60
1:A:248:PRO:HB2	1:A:251:VAL:N	2.14	0.60
1:A:248:PRO:HB2	1:A:251:VAL:HG23	1.82	0.60
1:A:229:SER:C	1:A:231:LYS:H	2.10	0.59
1:A:246:GLY:N	1:A:247:SER:O	2.36	0.59
1:B:55:MET:HE1	1:B:68:HIS:CG	2.38	0.59
1:B:247:SER:CA	1:B:249:TYR:CD1	2.70	0.59
1:A:229:SER:C	1:A:231:LYS:N	2.59	0.58
1:A:78:ILE:HG12	1:A:87:TYR:CE1	2.38	0.58
1:A:78:ILE:HD11	1:A:86:GLY:HA2	1.86	0.58
1:B:8:SER:HB3	1:B:267:LYS:NZ	2.19	0.57
1:A:55:MET:HE1	1:A:68:HIS:CG	2.40	0.57
1:B:172:VAL:HG13	1:B:176:MET:HE3	1.87	0.57
1:A:71:LYS:O	1:A:72:ASN:HB2	2.05	0.56
1:A:247:SER:O	1:A:248:PRO:C	2.49	0.56
1:B:248:PRO:HB3	1:B:251:VAL:CG2	2.35	0.56
1:B:248:PRO:C	1:B:250:MET:N	2.63	0.56
1:A:246:GLY:CA	1:A:247:SER:OG	2.30	0.56
1:A:247:SER:CA	1:A:249:TYR:HD2	2.18	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:TRP:HB3	1:B:121:GLU:O	2.06	0.55
1:A:143:MET:HE2	1:A:145:ASP:HB2	1.87	0.55
1:A:253:LYS:HE2	2:A:331:HOH:O	2.07	0.55
1:A:256:PHE:CG	1:A:257:ILE:N	2.74	0.55
1:B:244:VAL:O	1:B:248:PRO:HG2	2.07	0.55
1:A:249:TYR:O	1:A:250:MET:O	2.23	0.55
1:A:84:HIS:HD2	2:B:347:HOH:O	1.91	0.54
1:A:54:LYS:O	1:A:55:MET:HE2	2.07	0.54
1:B:229:SER:O	1:B:251:VAL:HA	2.09	0.53
1:B:245:LYS:HG2	1:B:249:TYR:CB	2.15	0.53
1:B:247:SER:H	1:B:248:PRO:HD3	1.73	0.53
1:B:165:LYS:HD2	1:B:165:LYS:N	2.24	0.53
1:A:77:TYR:CZ	1:A:273:THR:HG21	2.44	0.53
1:A:53:ALA:HB1	1:A:55:MET:HE3	1.91	0.53
1:B:164:GLU:HG3	1:B:165:LYS:CD	2.38	0.53
1:B:258:ASP:C	1:B:260:LEU:N	2.65	0.53
1:A:246:GLY:N	1:A:247:SER:C	2.67	0.53
1:A:68:HIS:ND1	1:A:75:LEU:HD13	2.24	0.52
1:B:53:ALA:HB1	1:B:55:MET:HE3	1.91	0.52
1:B:210:LYS:NZ	2:B:332:HOH:O	2.41	0.52
1:A:229:SER:O	1:A:231:LYS:N	2.42	0.52
1:B:140:ILE:HD11	1:B:151:TRP:HB3	1.91	0.52
1:B:236:PRO:C	1:B:238:ILE:H	2.15	0.52
1:A:246:GLY:CA	1:A:247:SER:C	2.82	0.52
1:B:122:ILE:HD12	1:B:176:MET:HE1	1.90	0.52
1:B:29:ASP:OD1	1:B:244:VAL:HB	2.10	0.52
1:B:244:VAL:HG12	1:B:248:PRO:HG2	1.92	0.51
1:A:50:TRP:CE3	1:B:85:LYS:HD3	2.46	0.51
1:A:246:GLY:HA3	1:A:247:SER:HG	1.64	0.51
1:A:245:LYS:N	1:A:248:PRO:HA	2.25	0.50
1:B:162:MET:HE3	1:B:166:LEU:HD11	1.93	0.50
1:A:63:MET:SD	1:A:259:ALA:HB1	2.51	0.50
1:A:245:LYS:O	1:A:246:GLY:O	2.30	0.50
1:B:228:ILE:HG22	1:B:229:SER:N	2.27	0.50
1:A:247:SER:C	1:A:249:TYR:HD2	2.19	0.49
1:A:244:VAL:CG2	1:A:245:LYS:HB2	2.38	0.49
1:B:70:VAL:HG12	1:B:75:LEU:HD12	1.95	0.49
1:B:242:VAL:HG11	1:B:251:VAL:HG21	1.93	0.49
1:A:264:GLU:O	1:A:266:PHE:N	2.42	0.49
1:A:244:VAL:O	1:A:248:PRO:O	2.29	0.49
1:A:78:ILE:HD12	1:A:79:TYR:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:PRO:HG3	1:A:262:LYS:HD3	1.94	0.49
1:B:79:TYR:CD2	1:B:193:PRO:HG2	2.47	0.49
1:A:8:SER:HB3	1:A:267:LYS:HE2	1.95	0.49
1:A:140:ILE:CD1	1:A:213:ALA:HA	2.43	0.48
1:B:54:LYS:O	1:B:55:MET:HE2	2.13	0.48
1:A:88:ILE:HG13	1:A:89:TRP:CE2	2.48	0.48
1:A:226:LYS:O	1:A:231:LYS:HB2	2.13	0.48
1:B:169:ILE:HG23	1:B:170:LEU:HD12	1.95	0.48
1:B:261:GLN:HA	1:B:263:HIS:H	1.78	0.48
1:B:8:SER:HB3	1:B:267:LYS:HZ2	1.79	0.48
1:B:261:GLN:C	1:B:263:HIS:N	2.70	0.48
1:A:141:LEU:HB2	1:A:207:PHE:CE1	2.49	0.48
1:A:246:GLY:CA	1:A:247:SER:O	2.62	0.48
1:A:103:TYR:HB2	2:A:321:HOH:O	2.14	0.47
1:A:246:GLY:HA3	1:A:247:SER:O	2.14	0.47
1:B:245:LYS:H	1:B:247:SER:N	2.13	0.47
1:B:258:ASP:CG	1:B:259:ALA:N	2.71	0.47
1:A:10:SER:OG	1:A:11:ALA:N	2.48	0.47
1:A:63:MET:HE1	1:A:262:LYS:CB	2.43	0.47
1:B:89:TRP:CD1	1:B:122:ILE:HG22	2.50	0.47
1:B:191:LYS:HD3	1:B:191:LYS:HA	1.81	0.46
1:A:89:TRP:CZ2	1:A:183:THR:HG23	2.50	0.46
1:B:108:THR:HG21	1:B:149:THR:HG23	1.98	0.46
1:B:261:GLN:C	1:B:263:HIS:H	2.24	0.46
1:B:275:GLN:HG2	1:B:276:ALA:H	1.81	0.46
1:A:244:VAL:C	1:A:248:PRO:HA	2.41	0.46
1:A:162:MET:HE2	1:A:176:MET:HE3	1.97	0.45
1:B:247:SER:H	1:B:248:PRO:CD	2.29	0.45
1:B:228:ILE:HG23	1:B:251:VAL:HG22	1.97	0.45
1:A:249:TYR:C	1:A:251:VAL:N	2.67	0.45
1:A:216:ILE:O	1:A:220:THR:HG23	2.17	0.45
1:A:251:VAL:HB	1:A:254:GLU:N	2.32	0.45
1:A:260:LEU:C	1:A:262:LYS:H	2.24	0.45
1:B:141:LEU:HB2	1:B:207:PHE:CE1	2.51	0.45
1:A:50:TRP:CD2	1:B:85:LYS:HD3	2.51	0.45
1:A:230:THR:HA	1:A:252:SER:H	1.81	0.45
1:A:251:VAL:HB	1:A:254:GLU:H	1.80	0.45
1:A:245:LYS:O	1:A:246:GLY:C	2.60	0.44
1:B:73:ASN:OD1	1:B:73:ASN:N	2.51	0.44
1:A:247:SER:O	1:A:249:TYR:CD2	2.70	0.44
1:A:244:VAL:HG13	1:A:245:LYS:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:TRP:HB3	1:A:121:GLU:O	2.17	0.44
1:B:256:PHE:HA	1:B:257:ILE:HA	1.47	0.44
1:A:49:ARG:NH2	1:B:278:GLU:O	2.51	0.44
1:A:219:LYS:HD3	1:A:219:LYS:HA	1.82	0.44
1:A:247:SER:HA	1:A:249:TYR:HA	1.97	0.44
1:A:233:GLU:OE1	1:A:233:GLU:N	2.50	0.44
1:B:186:TRP:O	1:B:190:TYR:HB3	2.17	0.44
1:A:43:MET:HE2	1:A:45:VAL:HG22	1.99	0.44
1:B:242:VAL:HG21	1:B:248:PRO:HB3	1.99	0.43
1:B:55:MET:HE1	1:B:68:HIS:CE1	2.52	0.43
1:A:78:ILE:HD11	1:A:86:GLY:CA	2.47	0.43
1:B:242:VAL:CG2	1:B:248:PRO:HB2	2.46	0.43
1:A:247:SER:C	1:A:249:TYR:N	2.72	0.43
1:A:22:SER:HA	1:A:23:PRO:HD2	1.88	0.43
1:A:7:THR:O	1:A:10:SER:HB3	2.18	0.43
1:A:55:MET:HE1	1:A:68:HIS:CE1	2.53	0.43
1:A:247:SER:CB	1:A:249:TYR:CE2	2.90	0.42
1:A:140:ILE:HD13	1:A:213:ALA:HA	2.01	0.42
1:A:78:ILE:HD11	1:A:86:GLY:C	2.45	0.42
1:A:108:THR:HG21	1:A:149:THR:HG23	2.00	0.42
1:B:89:TRP:CZ2	1:B:183:THR:HG23	2.54	0.42
1:B:245:LYS:N	1:B:247:SER:N	2.68	0.42
1:A:191:LYS:HE3	1:A:196:LYS:HD3	2.02	0.42
1:A:242:VAL:HG23	1:A:243:THR:N	2.34	0.42
1:A:140:ILE:HD13	1:A:140:ILE:HG21	1.79	0.41
1:A:245:LYS:O	1:A:245:LYS:HG2	2.19	0.41
1:A:268:ARG:HD2	1:A:269:GLY:O	2.20	0.41
1:B:226:LYS:HA	1:B:231:LYS:HE2	2.03	0.41
1:A:89:TRP:CG	1:A:120:CYS:HB3	2.55	0.41
1:B:27:PHE:CE1	1:B:116:PRO:HG3	2.56	0.41
1:B:78:ILE:HD12	1:B:79:TYR:O	2.20	0.41
1:B:78:ILE:CD1	1:B:86:GLY:HA2	2.48	0.41
1:A:77:TYR:OH	1:A:273:THR:HG21	2.20	0.41
1:A:74:LYS:HB3	1:A:74:LYS:HE2	1.86	0.41
1:B:78:ILE:HD11	1:B:86:GLY:CA	2.46	0.41
1:A:235:GLY:N	1:A:236:PRO:CD	2.81	0.41
1:A:32:LEU:HD11	1:A:41:TYR:HD2	1.86	0.41
1:B:237:ILE:O	1:B:237:ILE:HG12	2.20	0.41
1:B:224:TRP:CZ2	1:B:228:ILE:HG13	2.57	0.40
1:B:9:ASN:ND2	1:B:272:PRO:HA	2.37	0.40
1:B:247:SER:N	1:B:248:PRO:CD	2.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:ARG:HD2	1:A:239:ARG:HA	1.55	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	282/287 (98%)	255 (90%)	16 (6%)	11 (4%)	2	3
1	B	281/287 (98%)	242 (86%)	22 (8%)	17 (6%)	1	1
All	All	563/574 (98%)	497 (88%)	38 (7%)	28 (5%)	1	2

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	PRO
1	A	245	LYS
1	A	250	MET
1	A	256	PHE
1	A	265	ASP
1	A	275	GLN
1	B	73	ASN
1	B	146	GLU
1	B	237	ILE
1	B	248	PRO
1	B	249	TYR
1	B	250	MET
1	B	258	ASP
1	B	259	ALA
1	A	246	GLY
1	A	266	PHE
1	B	231	LYS

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Mol	Chain	Res	Type
1	B	244	VAL
1	B	256	PHE
1	B	262	LYS
1	A	73	ASN
1	A	230	THR
1	B	36	ALA
1	B	273	THR
1	B	247	SER
1	B	260	LEU
1	A	36	ALA
1	B	246	GLY

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	245/251 (98%)	229 (94%)	16 (6%)	15	35
1	B	244/251 (97%)	228 (93%)	16 (7%)	15	34
All	All	489/502 (97%)	457 (94%)	32 (6%)	15	35

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	76	ARG
1	A	78	ILE
1	A	89	TRP
1	A	109	LYS
1	A	140	ILE
1	A	228	ILE
1	A	231	LYS
1	A	232	VAL
1	A	239	ARG
1	A	242	VAL
1	A	244	VAL

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Mol	Chain	Res	Type
1	A	257	ILE
1	A	264	GLU
1	A	273	THR
1	A	283	CYS
1	B	38	ASN
1	B	60	GLU
1	B	71	LYS
1	B	73	ASN
1	B	89	TRP
1	B	131	SER
1	B	137	VAL
1	B	165	LYS
1	B	191	LYS
1	B	228	ILE
1	B	232	VAL
1	B	242	VAL
1	B	243	THR
1	B	244	VAL
1	B	245	LYS
1	B	250	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	223	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	284/287 (98%)	-0.54	3 (1%) 78 74	24, 50, 104, 129	0
1	B	283/287 (98%)	-0.63	2 (0%) 84 82	24, 49, 100, 134	0
All	All	567/574 (98%)	-0.58	5 (0%) 81 78	24, 50, 103, 134	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	246	GLY	4.3
1	B	249	TYR	2.7
1	A	230	THR	2.3
1	A	72	ASN	2.3
1	A	266	PHE	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.