



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:34 AM UTC

PDB ID : 2Z8U / pdb_00002z8u
Title : Methanococcus jannaschii TBP
Authors : Adachi, N.; Senda, T.; Horikoshi, M.
Deposited on : 2007-09-10
Resolution : 1.90 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

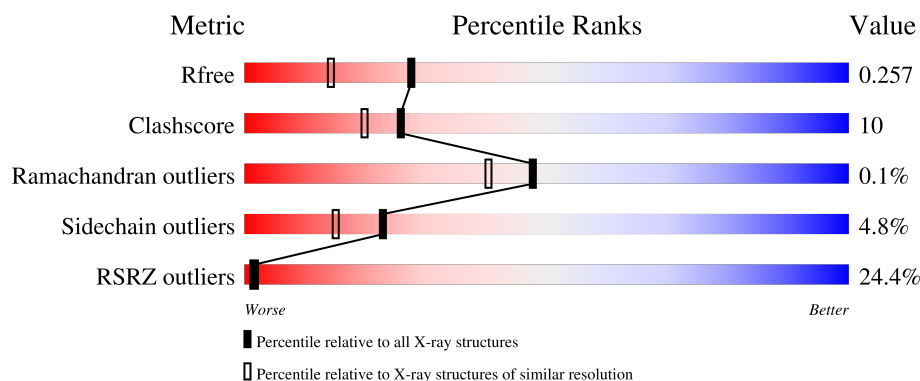
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	188	38% 63% 23% • 10%
1	B	188	11% 80% 12% • 7%
1	P	188	21% 77% 14% • 8%
1	Q	188	20% 74% 18% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5600 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 TATA-box-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	169	Total	C	N	O	S	0	0	0
			1293	826	209	253	5			
1	B	175	Total	C	N	O	S	0	0	0
			1348	864	215	264	5			
1	P	173	Total	C	N	O	S	0	0	0
			1334	854	213	262	5			
1	Q	175	Total	C	N	O	S	0	0	0
			1348	864	215	264	5			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP Q57930
A	-6	HIS	-	expression tag	UNP Q57930
A	-5	HIS	-	expression tag	UNP Q57930
A	-4	HIS	-	expression tag	UNP Q57930
A	-3	HIS	-	expression tag	UNP Q57930
A	-2	HIS	-	expression tag	UNP Q57930
A	-1	HIS	-	expression tag	UNP Q57930
A	0	PRO	-	expression tag	UNP Q57930
B	-7	MET	-	expression tag	UNP Q57930
B	-6	HIS	-	expression tag	UNP Q57930
B	-5	HIS	-	expression tag	UNP Q57930
B	-4	HIS	-	expression tag	UNP Q57930
B	-3	HIS	-	expression tag	UNP Q57930
B	-2	HIS	-	expression tag	UNP Q57930
B	-1	HIS	-	expression tag	UNP Q57930
B	0	PRO	-	expression tag	UNP Q57930
P	-7	MET	-	expression tag	UNP Q57930
P	-6	HIS	-	expression tag	UNP Q57930
P	-5	HIS	-	expression tag	UNP Q57930
P	-4	HIS	-	expression tag	UNP Q57930
P	-3	HIS	-	expression tag	UNP Q57930

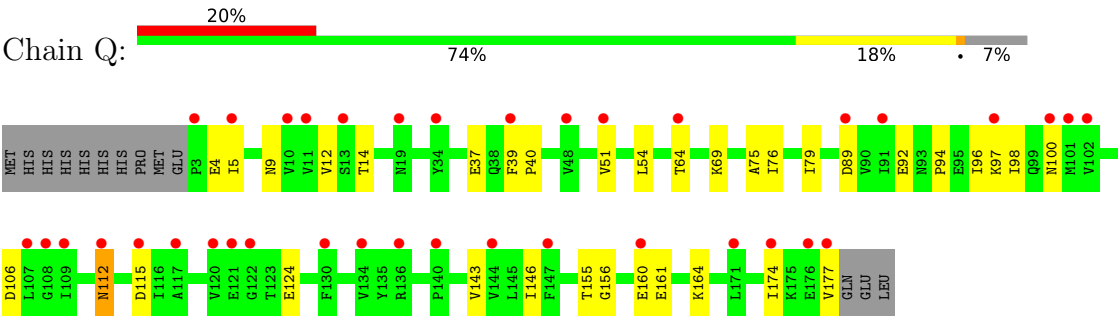
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Chain	Residue	Modelled	Actual	Comment	Reference
P	-2	HIS	-	expression tag	UNP Q57930
P	-1	HIS	-	expression tag	UNP Q57930
P	0	PRO	-	expression tag	UNP Q57930
Q	-7	MET	-	expression tag	UNP Q57930
Q	-6	HIS	-	expression tag	UNP Q57930
Q	-5	HIS	-	expression tag	UNP Q57930
Q	-4	HIS	-	expression tag	UNP Q57930
Q	-3	HIS	-	expression tag	UNP Q57930
Q	-2	HIS	-	expression tag	UNP Q57930
Q	-1	HIS	-	expression tag	UNP Q57930
Q	0	PRO	-	expression tag	UNP Q57930

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	58	Total O 58 58	0	0
2	B	94	Total O 94 94	0	0
2	P	53	Total O 53 53	0	0
2	Q	72	Total O 72 72	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.24Å 55.53Å 123.43Å 90.00° 91.05° 90.00°	Depositor
Resolution (Å)	50.00 – 1.90 50.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.00-1.90) 99.9 (50.00-1.90)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.226 , 0.259 0.225 , 0.257	Depositor DCC
R_{free} test set	2892 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.428	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for -k,-h,-l 0.014 for k,h,-l 0.025 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5600	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.70% 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	0.88	6/1303 (0.5%)	0.84	2/1756 (0.1%)
1	B	0.74	0/1363	0.83	0/1841
1	P	0.60	0/1348	0.82	0/1820
1	Q	0.64	0/1363	0.81	0/1841
All	All	0.72	6/5377 (0.1%)	0.83	2/7258 (0.0%)

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	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	86	ALA	C-N	11.62	1.51	1.33
1	A	41	GLY	C-O	7.87	1.39	1.23
1	A	41	GLY	C-N	7.35	1.44	1.33
1	A	25	VAL	C-O	-7.15	1.15	1.24
1	A	86	ALA	CA-C	6.99	1.62	1.52
1	A	86	ALA	C-O	5.81	1.31	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	130	PHE	CA-C-N	-5.44	114.37	120.14
1	A	130	PHE	C-N-CA	-5.44	114.37	120.14

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	86	ALA	Mainchain

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	1293	0	1372	44	0
1	B	1348	0	1417	29	0
1	P	1334	0	1400	20	0
1	Q	1348	0	1417	28	0
2	A	58	0	0	6	0
2	B	94	0	0	3	0
2	P	53	0	0	2	0
2	Q	72	0	0	3	0
All	All	5600	0	5606	104	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 10.

All (104) 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:100:ASN:ND2	1:B:124:GLU:HG2	1.57	1.18
1:A:100:ASN:HD21	1:B:124:GLU:CG	1.60	1.13
1:A:155:THR:HB	2:A:234:HOH:O	1.65	0.96
1:A:100:ASN:ND2	1:B:124:GLU:CG	2.23	0.94
1:B:38:GLN:HG2	2:B:270:HOH:O	1.68	0.93
1:P:126:GLU:HG3	1:P:129:GLN:HG2	1.52	0.92
2:A:234:HOH:O	1:B:124:GLU:HG3	1.74	0.88
1:P:126:GLU:CG	1:P:129:GLN:HG2	2.04	0.86
1:A:130:PHE:HE2	2:A:225:HOH:O	1.59	0.84
1:A:100:ASN:HD21	1:B:124:GLU:HG3	1.45	0.81
1:P:9:ASN:C	1:P:9:ASN:HD22	1.90	0.80
1:A:100:ASN:HD21	1:B:124:GLU:HG2	1.17	0.79
1:Q:39:PHE:CD1	1:Q:40:PRO:HD2	2.16	0.78
1:Q:100:ASN:HD21	1:Q:156:GLY:H	1.32	0.77
1:A:33:GLU:HG2	1:A:43:VAL:HB	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:69:LYS:HG3	1:Q:98:ILE:HD11	1.70	0.73
1:P:144:VAL:HG21	1:P:170:ILE:HD11	1.72	0.71
1:A:144:VAL:HG21	1:A:170:ILE:HD11	1.74	0.69
1:A:69:LYS:HE2	1:A:96:ILE:HG13	1.76	0.68
1:A:20:ILE:HD11	1:A:59:GLY:HA2	1.75	0.68
1:B:159:SER:OG	1:B:161:GLU:HG2	1.95	0.67
1:P:118:LEU:HD13	1:Q:64:THR:HG22	1.78	0.66
1:A:130:PHE:CD1	1:A:131:PRO:HD2	2.32	0.65
1:P:9:ASN:C	1:P:9:ASN:ND2	2.56	0.64
1:B:28:ILE:HD13	1:B:86:ALA:CB	2.28	0.63
1:Q:100:ASN:HD21	1:Q:156:GLY:N	1.96	0.63
1:A:125:TYR:CZ	1:A:127:PRO:HG3	2.34	0.63
1:A:32:ALA:HA	1:A:43:VAL:O	1.99	0.62
1:B:9:ASN:C	1:B:9:ASN:HD22	2.08	0.62
1:A:119:MET:HE2	1:B:43:VAL:HG21	1.80	0.61
1:A:100:ASN:HD22	1:B:124:GLU:HG2	1.62	0.61
1:P:126:GLU:HG2	1:P:129:GLN:HG2	1.82	0.60
1:B:112:ASN:C	1:B:112:ASN:HD22	2.11	0.59
1:A:146:ILE:HD13	1:A:174:ILE:HD11	1.85	0.58
1:A:130:PHE:CE2	1:A:131:PRO:O	2.57	0.57
2:P:187:HOH:O	1:Q:64:THR:HG21	2.04	0.57
1:B:97:LYS:HE2	1:B:99:GLN:NE2	2.22	0.55
1:P:69:LYS:HB2	1:P:98:ILE:HD11	1.89	0.55
1:A:20:ILE:CD1	1:A:59:GLY:HA2	2.37	0.54
1:Q:51:VAL:HG21	1:Q:75:ALA:HB1	1.90	0.54
1:A:119:MET:CE	1:B:43:VAL:HG21	2.38	0.53
1:Q:92:GLU:HB3	2:Q:240:HOH:O	2.09	0.53
1:B:9:ASN:C	1:B:9:ASN:ND2	2.66	0.53
1:Q:69:LYS:HE2	1:Q:96:ILE:O	2.09	0.52
1:P:136:ARG:HH11	1:Q:143:VAL:HG23	1.73	0.52
1:Q:39:PHE:CG	1:Q:40:PRO:HD2	2.44	0.52
1:A:98:ILE:HG21	1:A:101:MET:HE3	1.91	0.52
1:Q:12:VAL:HG12	1:Q:98:ILE:HD13	1.92	0.52
1:A:171:LEU:O	1:A:175:LYS:HG3	2.10	0.52
1:Q:39:PHE:HE2	1:Q:54:LEU:HD11	1.74	0.52
1:A:21:ASP:O	1:A:25:VAL:HG23	2.11	0.51
1:Q:5:ILE:HD13	1:Q:164:LYS:HG2	1.92	0.51
1:A:74:ILE:O	1:A:78:LYS:HG3	2.11	0.51
1:B:9:ASN:HD22	1:B:10:VAL:N	2.07	0.51
1:B:78:LYS:HG3	2:B:189:HOH:O	2.10	0.50
1:P:14:THR:OG1	1:P:94:PRO:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:VAL:HG23	1:B:124:GLU:CD	2.37	0.49
2:A:203:HOH:O	1:B:64:THR:HG21	2.11	0.49
1:A:130:PHE:CD2	1:A:131:PRO:O	2.65	0.49
1:P:73:GLU:O	1:P:77:LYS:HG3	2.13	0.49
1:Q:112:ASN:C	1:Q:112:ASN:HD22	2.21	0.48
1:Q:51:VAL:HG21	1:Q:75:ALA:CB	2.42	0.48
1:Q:76:ILE:O	1:Q:79:ILE:HG13	2.13	0.48
1:A:118:LEU:HD13	1:B:64:THR:HG22	1.96	0.48
1:A:10:VAL:HG12	1:A:101:MET:HG2	1.95	0.48
1:Q:97:LYS:NZ	2:Q:231:HOH:O	2.46	0.48
1:A:124:GLU:CD	1:B:102:VAL:HG23	2.39	0.47
1:Q:69:LYS:HG3	1:Q:98:ILE:CD1	2.44	0.47
1:A:145:LEU:HD22	2:A:225:HOH:O	2.15	0.46
1:B:28:ILE:HD13	1:B:86:ALA:HB2	1.97	0.46
1:Q:146:ILE:HD13	1:Q:174:ILE:HD11	1.97	0.46
1:Q:5:ILE:CD1	1:Q:164:LYS:HA	2.46	0.46
1:A:30:GLU:O	1:A:31:ASN:HB2	2.16	0.46
1:P:9:ASN:HD22	1:P:10:VAL:N	2.11	0.46
1:A:69:LYS:CG	1:A:96:ILE:HD11	2.46	0.46
1:Q:89:ASP:HB2	2:Q:251:HOH:O	2.16	0.45
1:B:30:GLU:O	1:B:31:ASN:HB2	2.16	0.45
1:B:97:LYS:HE2	1:B:99:GLN:HE22	1.81	0.45
1:P:112:ASN:HD22	1:P:115:ASP:H	1.65	0.45
1:A:51:VAL:HG21	1:A:75:ALA:CB	2.46	0.45
1:A:119:MET:HE2	1:B:43:VAL:CG2	2.44	0.45
1:A:69:LYS:HG2	1:A:96:ILE:HD11	1.99	0.44
1:A:112:ASN:ND2	1:A:115:ASP:H	2.15	0.44
1:P:88:ILE:HG22	1:P:90:VAL:HG13	1.99	0.44
1:P:33:GLU:HB3	1:P:43:VAL:HB	1.99	0.44
1:A:145:LEU:HB3	2:A:225:HOH:O	2.16	0.44
1:P:97:LYS:HD3	1:P:99:GLN:HE22	1.83	0.44
1:A:112:ASN:H	1:A:178:GLN:NE2	2.16	0.43
1:B:15:LYS:HE3	1:B:58:SER:O	2.18	0.43
1:P:112:ASN:ND2	1:P:115:ASP:H	2.16	0.43
1:Q:14:THR:OG1	1:Q:94:PRO:HB2	2.18	0.43
1:A:9:ASN:OD1	1:A:9:ASN:C	2.62	0.42
1:Q:100:ASN:ND2	1:Q:156:GLY:CA	2.82	0.42
1:A:112:ASN:H	1:A:178:GLN:HE22	1.68	0.42
1:P:124:GLU:HG3	1:Q:155:THR:HB	2.01	0.42
1:A:129:GLN:HE21	1:B:151:LYS:HE2	1.84	0.42
1:Q:112:ASN:ND2	1:Q:115:ASP:H	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:LEU:HG	1:A:88:ILE:HD12	2.02	0.41
1:Q:69:LYS:CE	1:Q:96:ILE:O	2.68	0.41
1:P:102:VAL:HG23	1:Q:124:GLU:CD	2.46	0.41
1:A:112:ASN:HD22	1:A:115:ASP:H	1.67	0.41
1:B:62:ASN:ND2	2:B:230:HOH:O	2.53	0.41
1:A:75:ALA:O	1:A:79:ILE:HG23	2.20	0.41
1:P:57:ARG:HB3	2:P:204:HOH:O	2.21	0.40

There are no symmetry-related clashes.

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	165/188 (88%)	162 (98%)	2 (1%)	1 (1%)	21	13
1	B	173/188 (92%)	169 (98%)	4 (2%)	0	100	100
1	P	171/188 (91%)	168 (98%)	3 (2%)	0	100	100
1	Q	173/188 (92%)	172 (99%)	1 (1%)	0	100	100
All	All	682/752 (91%)	671 (98%)	10 (2%)	1 (0%)	48	40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	32	ALA

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	146/165 (88%)	136 (93%)	10 (7%)	14	7
1	B	152/165 (92%)	146 (96%)	6 (4%)	28	21
1	P	150/165 (91%)	145 (97%)	5 (3%)	33	26
1	Q	152/165 (92%)	144 (95%)	8 (5%)	20	12
All	All	600/660 (91%)	571 (95%)	29 (5%)	23	15

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ILE
1	A	9	ASN
1	A	15	LYS
1	A	18	ASP
1	A	19	ASN
1	A	46	LEU
1	A	79	ILE
1	A	83	LEU
1	A	112	ASN
1	A	173	THR
1	B	9	ASN
1	B	20	ILE
1	B	35	GLU
1	B	109	ILE
1	B	112	ASN
1	B	174	ILE
1	P	5	ILE
1	P	9	ASN
1	P	35	GLU
1	P	109	ILE
1	P	112	ASN
1	Q	4	GLU
1	Q	9	ASN
1	Q	37	GLU
1	Q	106	ASP
1	Q	112	ASN
1	Q	160	GLU
1	Q	161	GLU
1	Q	177	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	112	ASN
1	A	129	GLN
1	B	9	ASN
1	B	31	ASN
1	B	38	GLN
1	B	112	ASN
1	P	9	ASN
1	P	99	GLN
1	P	112	ASN
1	P	129	GLN
1	Q	100	ASN
1	Q	112	ASN
1	Q	129	GLN

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 ⓘ

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	169/188 (89%)	2.17	72 (42%) 0 0	27, 38, 48, 60	0
1	B	175/188 (93%)	1.21	20 (11%) 10 10	24, 37, 49, 52	0
1	P	173/188 (92%)	1.45	40 (23%) 2 2	33, 39, 51, 61	0
1	Q	175/188 (93%)	1.40	37 (21%) 2 2	31, 38, 50, 56	0
All	All	692/752 (92%)	1.55	169 (24%) 2 1	24, 38, 50, 61	0

All (169) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	ILE	8.0
1	A	86	ALA	7.6
1	A	29	LEU	6.4
1	A	10	VAL	6.3
1	A	96	ILE	5.8
1	A	7	ILE	5.6
1	A	74	ILE	5.5
1	A	90	VAL	4.9
1	B	130	PHE	4.8
1	A	9	ASN	4.8
1	A	25	VAL	4.7
1	Q	177	VAL	4.6
1	P	39	PHE	4.5
1	A	8	VAL	4.5
1	A	6	LYS	4.4
1	P	124	GLU	4.4
1	A	3	PRO	4.4
1	A	109	ILE	4.4
1	A	65	GLY	4.3
1	A	61	VAL	4.2
1	A	28	ILE	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	66	ALA	4.1
1	Q	109	ILE	4.0
1	A	100	ASN	4.0
1	Q	3	PRO	4.0
1	A	68	SER	3.9
1	A	26	ALA	3.9
1	Q	39	PHE	3.9
1	P	40	PRO	3.9
1	Q	108	GLY	3.8
1	A	55	ILE	3.8
1	A	80	ILE	3.8
1	A	67	LYS	3.7
1	A	11	VAL	3.7
1	A	175	LYS	3.6
1	A	177	VAL	3.6
1	A	22	LEU	3.6
1	P	120	VAL	3.5
1	A	174	ILE	3.5
1	P	102	VAL	3.5
1	A	79	ILE	3.5
1	A	12	VAL	3.5
1	A	4	GLU	3.4
1	A	20	ILE	3.4
1	A	130	PHE	3.4
1	A	48	VAL	3.4
1	A	57	ARG	3.3
1	Q	117	ALA	3.3
1	B	18	ASP	3.3
1	P	130	PHE	3.3
1	P	109	ILE	3.2
1	Q	91	ILE	3.2
1	P	5	ILE	3.2
1	B	177	VAL	3.1
1	P	101	MET	3.1
1	Q	112	ASN	3.1
1	P	98	ILE	3.1
1	A	88	ILE	3.0
1	P	118	LEU	3.0
1	Q	134	VAL	3.0
1	A	108	GLY	3.0
1	A	87	GLY	3.0
1	Q	136	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	42	LEU	2.9
1	B	40	PRO	2.8
1	B	39	PHE	2.8
1	A	91	ILE	2.8
1	A	41	GLY	2.8
1	P	119	MET	2.8
1	A	178	GLN	2.8
1	A	78	LYS	2.8
1	B	49	PRO	2.8
1	Q	171	LEU	2.8
1	A	23	GLU	2.7
1	B	100	ASN	2.7
1	A	33	GLU	2.7
1	P	105	ALA	2.7
1	Q	19	ASN	2.7
1	B	90	VAL	2.7
1	B	122	GLY	2.6
1	A	176	GLU	2.6
1	A	82	GLU	2.6
1	Q	13	SER	2.6
1	P	91	ILE	2.6
1	P	174	ILE	2.6
1	B	38	GLN	2.6
1	P	134	VAL	2.6
1	Q	48	VAL	2.6
1	Q	5	ILE	2.6
1	P	150	GLY	2.6
1	B	48	VAL	2.6
1	A	70	GLU	2.5
1	A	89	ASP	2.5
1	Q	34	TYR	2.5
1	P	77	LYS	2.5
1	P	48	VAL	2.5
1	Q	51	VAL	2.5
1	A	135	TYR	2.5
1	Q	107	LEU	2.5
1	Q	120	VAL	2.5
1	Q	130	PHE	2.4
1	P	74	ILE	2.4
1	P	131	PRO	2.4
1	P	112	ASN	2.4
1	Q	174	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	101	MET	2.4
1	B	50	LYS	2.4
1	A	53	LEU	2.3
1	A	71	GLU	2.3
1	Q	176	GLU	2.3
1	A	72	ALA	2.3
1	Q	140	PRO	2.3
1	A	43	VAL	2.3
1	A	63	CYS	2.3
1	P	4	GLU	2.3
1	A	107	LEU	2.3
1	A	19	ASN	2.3
1	A	27	MET	2.3
1	P	170	ILE	2.3
1	P	97	LYS	2.3
1	A	13	SER	2.3
1	A	49	PRO	2.3
1	Q	122	GLY	2.3
1	P	110	GLU	2.3
1	Q	144	VAL	2.3
1	B	44	CYS	2.3
1	B	3	PRO	2.2
1	A	18	ASP	2.2
1	Q	115	ASP	2.2
1	A	51	VAL	2.2
1	A	46	LEU	2.2
1	A	73	GLU	2.2
1	P	66	ALA	2.2
1	A	93	ASN	2.2
1	P	169	LYS	2.2
1	A	134	VAL	2.2
1	P	53	LEU	2.2
1	P	106	ASP	2.2
1	P	108	GLY	2.2
1	B	5	ILE	2.2
1	Q	101	MET	2.2
1	B	143	VAL	2.2
1	Q	10	VAL	2.2
1	A	83	LEU	2.2
1	P	136	ARG	2.2
1	Q	100	ASN	2.1
1	P	7	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	52	ALA	2.1
1	P	80	ILE	2.1
1	Q	89	ASP	2.1
1	A	173	THR	2.1
1	Q	64	THR	2.1
1	P	70	GLU	2.1
1	P	36	PRO	2.1
1	B	35	GLU	2.1
1	Q	97	LYS	2.1
1	A	59	GLY	2.1
1	P	107	LEU	2.1
1	Q	121	GLU	2.0
1	Q	147	PHE	2.0
1	B	41	GLY	2.0
1	P	144	VAL	2.0
1	Q	11	VAL	2.0
1	P	157	LEU	2.0
1	B	32	ALA	2.0
1	B	138	ASP	2.0
1	Q	160	GLU	2.0
1	P	116	ILE	2.0
1	Q	102	VAL	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 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.