



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:10 AM UTC

PDB ID : 8X4G / pdb_00008x4g
Title : Crystal structure of the D117A mutant of DIMT1 from *Pyrococcus horikoshii*
Authors : Saha, S.; Kanaujia, S.P.
Deposited on : 2023-11-15
Resolution : 3.30 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

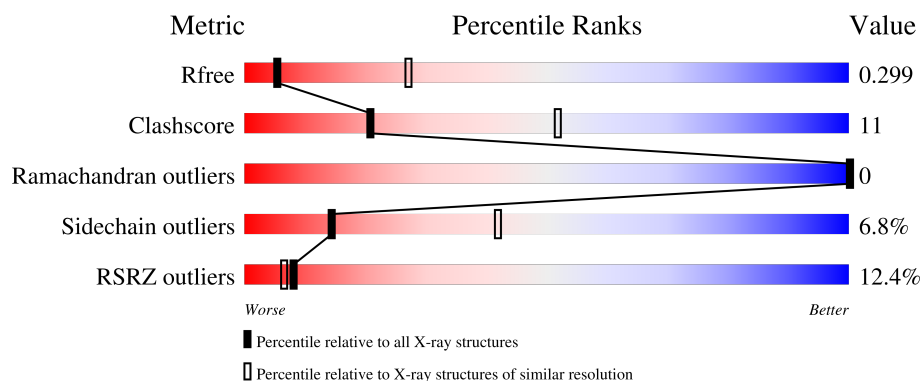
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	290	10% 66% 24% • 7%
1	B	290	13% 65% 26% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4397 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 Probable ribosomal RNA small subunit methyltransferase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			2192	1418	383	385	6			
1	B	270	Total	C	N	O	S	0	0	0
			2192	1418	383	385	6			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP O59487
A	2	SER	-	expression tag	UNP O59487
A	3	SER	-	expression tag	UNP O59487
A	4	ARG	-	expression tag	UNP O59487
A	5	ILE	-	expression tag	UNP O59487
A	6	ARG	-	expression tag	UNP O59487
A	7	ILE	-	expression tag	UNP O59487
A	8	ASN	-	expression tag	UNP O59487
A	9	THR	-	expression tag	UNP O59487
A	10	LYS	-	expression tag	UNP O59487
A	11	SER	-	expression tag	UNP O59487
A	12	LEU	-	expression tag	UNP O59487
A	13	LEU	-	expression tag	UNP O59487
A	14	VAL	-	expression tag	UNP O59487
A	15	GLN	-	expression tag	UNP O59487
A	16	PRO	-	expression tag	UNP O59487
A	17	GLY	-	expression tag	UNP O59487
A	18	TYR	-	expression tag	UNP O59487
A	19	SER	-	expression tag	UNP O59487
A	20	GLY	-	expression tag	UNP O59487
A	21	SER	-	expression tag	UNP O59487
A	22	LYS	-	expression tag	UNP O59487
A	117	ALA	ASP	engineered mutation	UNP O59487
B	1	MET	-	initiating methionine	UNP O59487
B	2	SER	-	expression tag	UNP O59487

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Chain	Residue	Modelled	Actual	Comment	Reference
B	3	SER	-	expression tag	UNP O59487
B	4	ARG	-	expression tag	UNP O59487
B	5	ILE	-	expression tag	UNP O59487
B	6	ARG	-	expression tag	UNP O59487
B	7	ILE	-	expression tag	UNP O59487
B	8	ASN	-	expression tag	UNP O59487
B	9	THR	-	expression tag	UNP O59487
B	10	LYS	-	expression tag	UNP O59487
B	11	SER	-	expression tag	UNP O59487
B	12	LEU	-	expression tag	UNP O59487
B	13	LEU	-	expression tag	UNP O59487
B	14	VAL	-	expression tag	UNP O59487
B	15	GLN	-	expression tag	UNP O59487
B	16	PRO	-	expression tag	UNP O59487
B	17	GLY	-	expression tag	UNP O59487
B	18	TYR	-	expression tag	UNP O59487
B	19	SER	-	expression tag	UNP O59487
B	20	GLY	-	expression tag	UNP O59487
B	21	SER	-	expression tag	UNP O59487
B	22	LYS	-	expression tag	UNP O59487
B	117	ALA	ASP	engineered mutation	UNP O59487

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	6	Total O 6 6	0	0
2	B	7	Total O 7 7	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.17Å 80.63Å 84.82Å 90.00° 114.86° 90.00°	Depositor
Resolution (Å)	64.52 – 3.30 64.52 – 3.30	Depositor EDS
% Data completeness (in resolution range)	93.6 (64.52-3.30) 93.6 (64.52-3.30)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.248 , 0.299 0.253 , 0.299	Depositor DCC
R_{free} test set	1074 reflections (9.74%)	wwPDB-VP
Wilson B-factor (Å ²)	27.9	Xtriage
Anisotropy	0.897	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	4397	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.94% 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 [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	0.54	0/2233	1.10	7/3005 (0.2%)
1	B	0.53	0/2233	1.08	6/3005 (0.2%)
All	All	0.54	0/4466	1.09	13/6010 (0.2%)

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

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	245	HIS	CA-CB-CG	9.73	123.53	113.80
1	B	245	HIS	CA-CB-CG	8.52	122.32	113.80
1	B	93	GLU	CB-CG-CD	5.99	122.78	112.60
1	A	93	GLU	CB-CG-CD	5.80	122.46	112.60
1	A	45	HIS	CA-CB-CG	5.62	119.42	113.80
1	A	135	TYR	CB-CA-C	5.61	121.82	110.38
1	B	135	TYR	CB-CA-C	5.44	121.48	110.38
1	A	25	ASP	CA-CB-CG	5.30	117.90	112.60
1	B	25	ASP	N-CA-C	-5.26	105.71	111.82
1	B	45	HIS	CA-CB-CG	5.17	118.97	113.80
1	B	95	ASP	CA-CB-CG	5.10	117.70	112.60
1	A	25	ASP	N-CA-C	-5.08	105.92	111.82
1	A	95	ASP	CA-CB-CG	5.07	117.67	112.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	GLU	Peptide
1	A	214	ARG	Sidechain

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	2192	0	2306	57	1
1	B	2192	0	2306	58	1
2	A	6	0	0	3	0
2	B	7	0	0	1	0
All	All	4397	0	4612	100	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 11.

All (100) 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:231:HIS:NE2	1:B:231:HIS:NE2	1.84	1.23
1:A:231:HIS:NE2	1:B:231:HIS:CE1	2.16	1.13
1:A:231:HIS:HE2	1:B:231:HIS:CE1	1.75	1.02
1:B:201:ARG:HH11	1:B:201:ARG:HB2	1.31	0.95
1:A:201:ARG:HH11	1:A:201:ARG:HB2	1.33	0.94
1:A:231:HIS:CE1	1:B:231:HIS:NE2	2.35	0.94
1:A:231:HIS:CE1	1:B:231:HIS:CE1	2.58	0.91
1:B:164:ARG:O	1:B:178:SER:OG	1.91	0.88
1:A:164:ARG:O	1:A:178:SER:OG	1.92	0.87
1:A:160:GLU:OE1	1:B:245:HIS:NE2	2.08	0.86
1:A:93:GLU:CD	2:A:301:HOH:O	2.19	0.84
1:A:245:HIS:NE2	1:B:160:GLU:OE1	2.12	0.82
1:A:93:GLU:OE1	2:A:301:HOH:O	2.01	0.77
1:A:160:GLU:OE1	1:B:245:HIS:CE1	2.40	0.74
1:A:231:HIS:HE2	1:B:231:HIS:HE2	0.72	0.71
1:A:233:ARG:O	1:A:267:ARG:NH1	2.24	0.70
1:B:233:ARG:O	1:B:267:ARG:NH1	2.25	0.68
1:A:27:LEU:HD12	1:A:75:LEU:HD12	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ILE:HG23	1:A:78:LEU:CD1	2.28	0.64
1:A:252:ASP:HB3	1:B:203:LYS:HG2	1.80	0.61
1:B:53:ILE:HG23	1:B:78:LEU:CD1	2.31	0.60
1:A:99:ILE:HG23	1:A:113:ILE:HG21	1.85	0.59
1:B:27:LEU:HD12	1:B:75:LEU:HD12	1.85	0.58
1:A:245:HIS:CE1	1:B:160:GLU:OE1	2.55	0.58
1:A:176:ARG:CZ	1:A:226:LYS:HE3	2.33	0.57
1:B:176:ARG:CZ	1:B:226:LYS:HE3	2.34	0.57
1:A:201:ARG:HH11	1:A:201:ARG:CB	2.13	0.56
1:B:201:ARG:HH11	1:B:201:ARG:CB	2.13	0.56
1:A:53:ILE:HG23	1:A:78:LEU:HD12	1.88	0.56
1:A:252:ASP:CB	1:B:203:LYS:HG2	2.36	0.55
1:A:27:LEU:HD12	1:A:75:LEU:CD1	2.37	0.54
1:A:167:ALA:HB2	1:A:173:ASN:HD21	1.73	0.53
1:B:93:GLU:HG2	1:B:98:ILE:HG21	1.90	0.53
1:A:93:GLU:HG2	1:A:98:ILE:HG21	1.91	0.53
1:A:69:LEU:O	1:A:129:VAL:HG13	2.09	0.52
1:B:53:ILE:HG23	1:B:78:LEU:HD12	1.90	0.52
1:A:39:ARG:NH2	1:A:45:HIS:CE1	2.78	0.51
1:B:39:ARG:NH2	1:B:45:HIS:CE1	2.78	0.51
1:B:167:ALA:HB2	1:B:173:ASN:HD21	1.74	0.51
1:B:244:ILE:CG2	1:B:249:VAL:HB	2.39	0.51
1:A:53:ILE:HG23	1:A:78:LEU:HD13	1.93	0.50
1:A:182:GLN:O	1:A:214:ARG:NH1	2.44	0.50
1:B:236:VAL:N	1:B:237:PRO:CD	2.75	0.50
1:B:69:LEU:O	1:B:129:VAL:HG13	2.12	0.49
1:B:244:ILE:HG23	1:B:249:VAL:HB	1.94	0.49
1:B:182:GLN:O	1:B:214:ARG:NH1	2.45	0.49
1:A:69:LEU:HD22	1:A:126:PHE:CE2	2.48	0.49
1:A:236:VAL:N	1:A:237:PRO:CD	2.76	0.49
1:B:27:LEU:HD12	1:B:75:LEU:CD1	2.43	0.49
1:A:244:ILE:CG2	1:A:249:VAL:HB	2.42	0.48
1:A:118:ALA:HA	1:A:121:VAL:HG22	1.96	0.48
1:B:68:ILE:HG21	1:B:82:LEU:HD13	1.95	0.48
1:A:226:LYS:O	1:A:230:GLN:N	2.48	0.47
1:A:74:GLY:HA2	2:A:302:HOH:O	2.14	0.47
1:B:23:MET:CE	1:B:50:GLU:HG2	2.45	0.47
1:A:195:LYS:HB2	1:A:202:PRO:HD2	1.96	0.47
1:A:199:TYR:HA	1:A:200:PRO:C	2.39	0.47
1:B:150:PHE:O	1:B:213:PRO:HG3	2.15	0.47
1:A:37:ARG:HB2	1:A:38:PRO:HD2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ILE:HG21	1:A:82:LEU:HD13	1.96	0.47
1:A:93:GLU:O	1:A:115:GLN:HA	2.15	0.46
1:A:221:ASN:O	1:A:225:VAL:HG23	2.15	0.46
1:A:39:ARG:HH21	1:A:45:HIS:CE1	2.33	0.46
1:B:199:TYR:HA	1:B:200:PRO:C	2.39	0.46
1:B:99:ILE:HG23	1:B:113:ILE:HG21	1.97	0.45
1:B:221:ASN:O	1:B:225:VAL:HG23	2.16	0.45
1:B:73:PRO:HD2	1:B:93:GLU:HB2	1.99	0.45
1:B:37:ARG:HB2	1:B:38:PRO:HD2	1.97	0.45
1:A:73:PRO:HD2	1:A:93:GLU:HB2	1.99	0.45
1:B:90:TYR:CG	1:B:124:PRO:HG3	2.51	0.45
1:B:195:LYS:HB2	1:B:202:PRO:HD2	1.98	0.45
1:B:226:LYS:O	1:B:230:GLN:N	2.50	0.45
1:A:258:ILE:O	1:A:261:VAL:HG12	2.17	0.44
1:B:69:LEU:HD22	1:B:126:PHE:CE2	2.52	0.44
1:B:53:ILE:HG23	1:B:78:LEU:HD13	1.98	0.44
1:A:251:LYS:HD3	1:A:255:ARG:NH1	2.33	0.44
1:B:93:GLU:O	1:B:115:GLN:HA	2.17	0.44
1:A:167:ALA:HB2	1:A:173:ASN:ND2	2.33	0.44
1:A:244:ILE:HG22	1:A:249:VAL:HB	2.00	0.43
1:B:176:ARG:NH1	1:B:226:LYS:HE3	2.34	0.43
1:A:245:HIS:CE1	1:B:160:GLU:CD	2.97	0.42
1:A:71:VAL:HG21	1:A:141:PHE:CE2	2.54	0.42
1:A:245:HIS:CE1	1:B:160:GLU:OE2	2.73	0.42
1:B:167:ALA:HB2	1:B:173:ASN:ND2	2.33	0.42
1:A:188:GLU:HB2	1:A:210:LEU:HB3	2.00	0.42
1:B:42:ILE:C	1:B:42:ILE:HD12	2.44	0.42
1:B:258:ILE:O	1:B:261:VAL:HG12	2.19	0.42
1:B:118:ALA:HA	1:B:121:VAL:HG22	2.02	0.41
1:A:90:TYR:CG	1:A:124:PRO:HG3	2.55	0.41
1:A:176:ARG:NH1	1:A:226:LYS:HE3	2.35	0.41
1:A:42:ILE:HD12	1:A:42:ILE:C	2.45	0.41
1:B:251:LYS:HD3	1:B:255:ARG:NH1	2.36	0.41
1:B:61:ASN:O	1:B:128:LYS:NZ	2.54	0.41
1:A:61:ASN:O	1:A:128:LYS:NZ	2.53	0.41
1:B:193:ILE:HD11	1:B:208:LEU:HB2	2.03	0.41
1:A:246:MET:O	1:B:164:ARG:NH1	2.54	0.41
1:B:39:ARG:HH21	1:B:45:HIS:CE1	2.39	0.40
1:B:64:GLU:O	1:B:87:LYS:HB2	2.21	0.40
1:B:93:GLU:CD	2:B:302:HOH:O	2.65	0.40
1:B:176:ARG:HG3	1:B:225:VAL:HG12	2.03	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:188:GLU:OE2	1:B:286:HIS:NE2[4_546]	1.89	0.31

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	268/290 (92%)	254 (95%)	14 (5%)	0	100	100
1	B	268/290 (92%)	255 (95%)	13 (5%)	0	100	100
All	All	536/580 (92%)	509 (95%)	27 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	242/260 (93%)	226 (93%)	16 (7%)	15	43
1	B	242/260 (93%)	225 (93%)	17 (7%)	14	41
All	All	484/520 (93%)	451 (93%)	33 (7%)	14	42

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	ILE
1	A	40	ASP
1	A	54	GLU
1	A	79	THR
1	A	172	ARG
1	A	179	LEU
1	A	186	ASN
1	A	201	ARG
1	A	214	ARG
1	A	219	VAL
1	A	222	GLU
1	A	236	VAL
1	A	238	ARG
1	A	240	LEU
1	A	254	ILE
1	A	281	GLU
1	B	36	ILE
1	B	40	ASP
1	B	50	GLU
1	B	54	GLU
1	B	79	THR
1	B	172	ARG
1	B	179	LEU
1	B	186	ASN
1	B	201	ARG
1	B	214	ARG
1	B	219	VAL
1	B	222	GLU
1	B	236	VAL
1	B	238	ARG
1	B	240	LEU
1	B	254	ILE
1	B	281	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	HIS
1	A	65	ASN
1	A	96	GLN
1	A	158	GLN
1	B	65	ASN
1	B	96	GLN

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Mol	Chain	Res	Type
1	B	158	GLN
1	B	265	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 ⓘ

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	270/290 (93%)	0.90	30 (11%) 10 9	20, 28, 55, 69	0
1	B	270/290 (93%)	1.04	37 (13%) 6 5	20, 28, 68, 80	0
All	All	540/580 (93%)	0.97	67 (12%) 8 6	20, 28, 60, 80	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	43	GLY	6.4
1	B	247	LEU	6.1
1	B	248	GLY	6.0
1	B	250	SER	5.8
1	A	245	HIS	5.6
1	B	44	GLN	5.5
1	B	245	HIS	5.4
1	A	41	SER	5.2
1	B	21	SER	5.1
1	A	43	GLY	5.0
1	B	41	SER	4.8
1	B	38	PRO	4.3
1	B	244	ILE	4.3
1	B	42	ILE	4.2
1	A	42	ILE	4.1
1	B	249	VAL	4.0
1	A	223	ASN	3.7
1	B	253	GLU	3.5
1	A	44	GLN	3.4
1	A	250	SER	3.3
1	B	243	SER	3.2
1	B	39	ARG	3.2
1	B	223	ASN	3.2
1	B	74	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	264	SER	3.1
1	A	253	GLU	3.1
1	B	40	ASP	3.0
1	A	135	TYR	3.0
1	B	254	ILE	3.0
1	B	51	ASP	2.9
1	B	259	ASN	2.8
1	A	21	SER	2.8
1	B	131	SER	2.7
1	B	258	ILE	2.6
1	A	227	ALA	2.6
1	A	39	ARG	2.5
1	A	38	PRO	2.5
1	B	290	SER	2.5
1	B	35	GLY	2.5
1	B	230	GLN	2.4
1	A	40	ASP	2.4
1	A	125	LYS	2.3
1	B	228	LEU	2.3
1	A	288	ILE	2.3
1	B	257	ILE	2.3
1	A	50	GLU	2.3
1	B	50	GLU	2.3
1	B	216	ASP	2.3
1	B	151	GLU	2.3
1	A	93	GLU	2.2
1	A	286	HIS	2.2
1	A	254	ILE	2.2
1	B	256	GLY	2.2
1	A	281	GLU	2.1
1	A	25	ASP	2.1
1	B	45	HIS	2.1
1	B	286	HIS	2.1
1	A	219	VAL	2.1
1	B	255	ARG	2.1
1	A	121	VAL	2.1
1	A	249	VAL	2.1
1	A	92	ILE	2.1
1	B	92	ILE	2.1
1	A	265	ASN	2.0
1	A	230	GLN	2.0
1	A	200	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	289	ILE	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.