



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:03 PM UTC

PDB ID : 6BUI / pdb_00006bui
Title : Crystal structure of a membrane protein, crystal form III
Authors : Ma, D.; Wang, Z.; Xu, W.
Deposited on : 2017-12-10
Resolution : 3.27 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

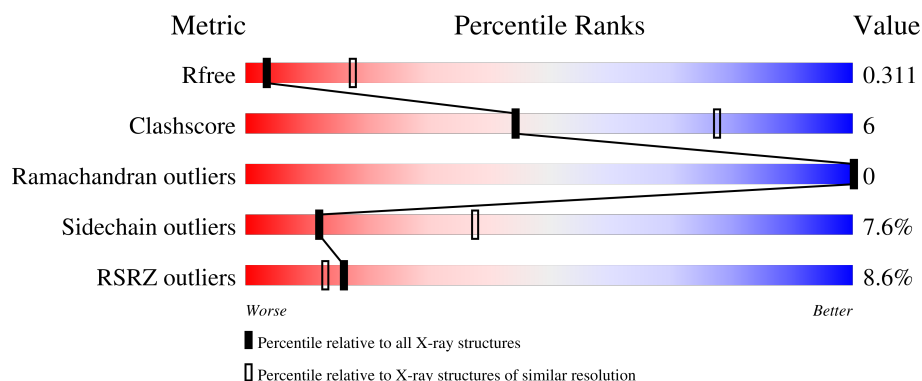
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.27 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	1303 (3.30-3.26)
Clashscore	190562	1354 (3.30-3.26)
Ramachandran outliers	187476	1334 (3.30-3.26)
Sidechain outliers	187428	1333 (3.30-3.26)
RSRZ outliers	180081	1303 (3.30-3.26)

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	425	6% 77% 18% ..
1	B	425	8% 75% 20% ..
1	C	425	5% 79% 16% ..
1	D	425	13% 77% 19% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 13792 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 Integral membrane protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	414	Total	C	N	O	S	0	0	0
			3448	2327	548	553	20			
1	A	414	Total	C	N	O	S	0	0	0
			3448	2327	548	553	20			
1	B	414	Total	C	N	O	S	0	0	0
			3448	2327	548	553	20			
1	D	414	Total	C	N	O	S	0	0	0
			3448	2327	548	553	20			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	416	LEU	-	expression tag	UNP Q5M4V4
C	417	GLU	-	expression tag	UNP Q5M4V4
C	418	HIS	-	expression tag	UNP Q5M4V4
C	419	HIS	-	expression tag	UNP Q5M4V4
C	420	HIS	-	expression tag	UNP Q5M4V4
C	421	HIS	-	expression tag	UNP Q5M4V4
C	422	HIS	-	expression tag	UNP Q5M4V4
C	423	HIS	-	expression tag	UNP Q5M4V4
C	424	HIS	-	expression tag	UNP Q5M4V4
C	425	HIS	-	expression tag	UNP Q5M4V4
A	416	LEU	-	expression tag	UNP Q5M4V4
A	417	GLU	-	expression tag	UNP Q5M4V4
A	418	HIS	-	expression tag	UNP Q5M4V4
A	419	HIS	-	expression tag	UNP Q5M4V4
A	420	HIS	-	expression tag	UNP Q5M4V4
A	421	HIS	-	expression tag	UNP Q5M4V4
A	422	HIS	-	expression tag	UNP Q5M4V4
A	423	HIS	-	expression tag	UNP Q5M4V4
A	424	HIS	-	expression tag	UNP Q5M4V4
A	425	HIS	-	expression tag	UNP Q5M4V4
B	416	LEU	-	expression tag	UNP Q5M4V4

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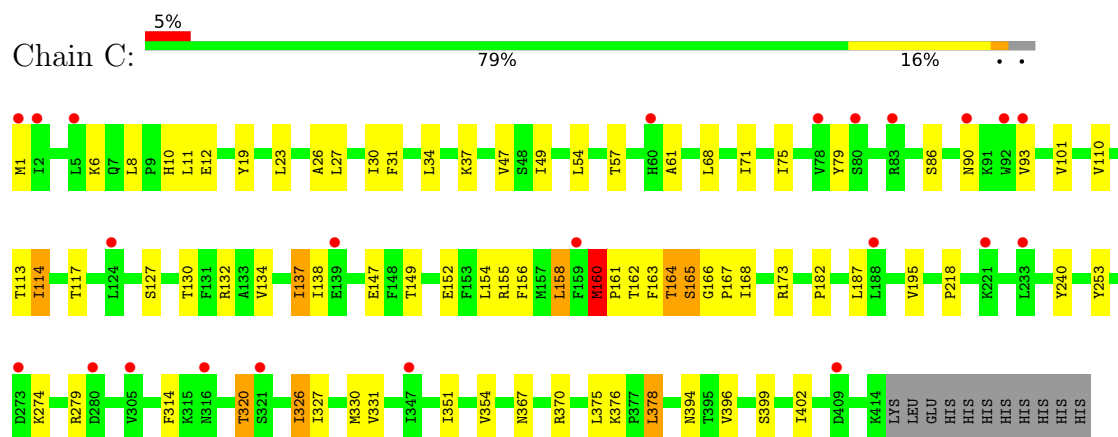
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Chain	Residue	Modelled	Actual	Comment	Reference
B	417	GLU	-	expression tag	UNP Q5M4V4
B	418	HIS	-	expression tag	UNP Q5M4V4
B	419	HIS	-	expression tag	UNP Q5M4V4
B	420	HIS	-	expression tag	UNP Q5M4V4
B	421	HIS	-	expression tag	UNP Q5M4V4
B	422	HIS	-	expression tag	UNP Q5M4V4
B	423	HIS	-	expression tag	UNP Q5M4V4
B	424	HIS	-	expression tag	UNP Q5M4V4
B	425	HIS	-	expression tag	UNP Q5M4V4
D	416	LEU	-	expression tag	UNP Q5M4V4
D	417	GLU	-	expression tag	UNP Q5M4V4
D	418	HIS	-	expression tag	UNP Q5M4V4
D	419	HIS	-	expression tag	UNP Q5M4V4
D	420	HIS	-	expression tag	UNP Q5M4V4
D	421	HIS	-	expression tag	UNP Q5M4V4
D	422	HIS	-	expression tag	UNP Q5M4V4
D	423	HIS	-	expression tag	UNP Q5M4V4
D	424	HIS	-	expression tag	UNP Q5M4V4
D	425	HIS	-	expression tag	UNP Q5M4V4

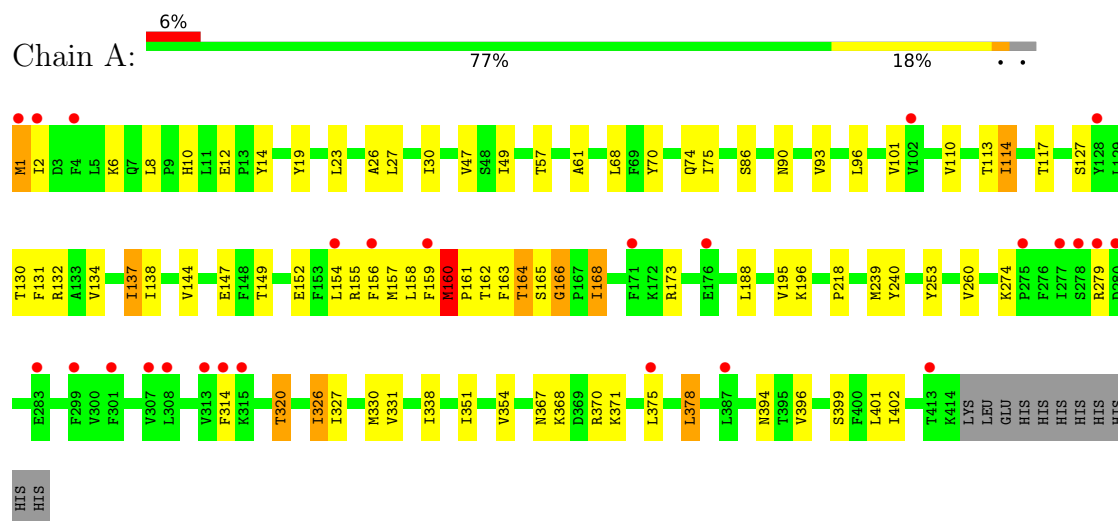
3 Residue-property plots [i](#)

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.

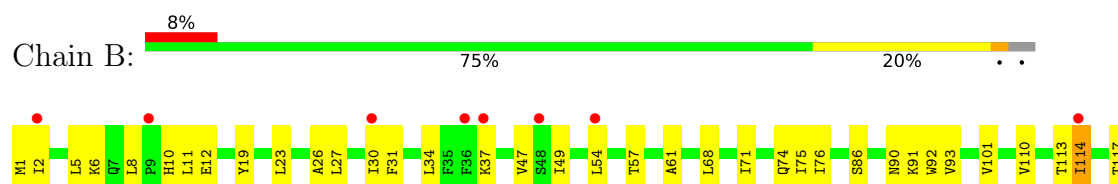
- Molecule 1: Integral membrane protein

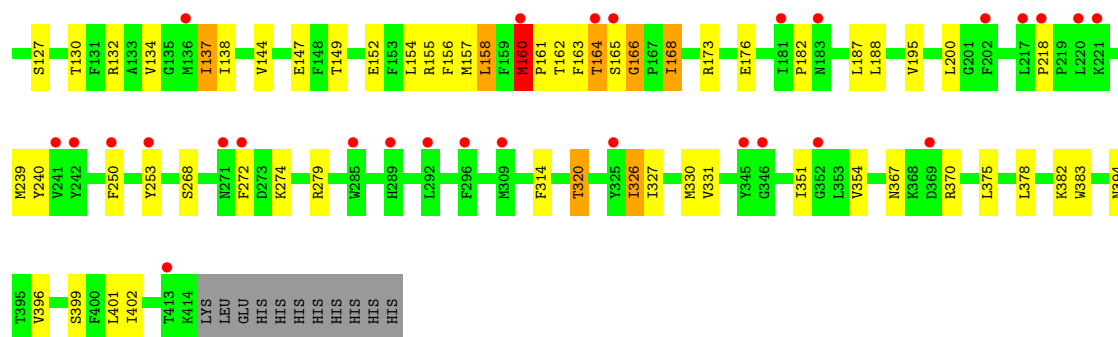


- Molecule 1: Integral membrane protein

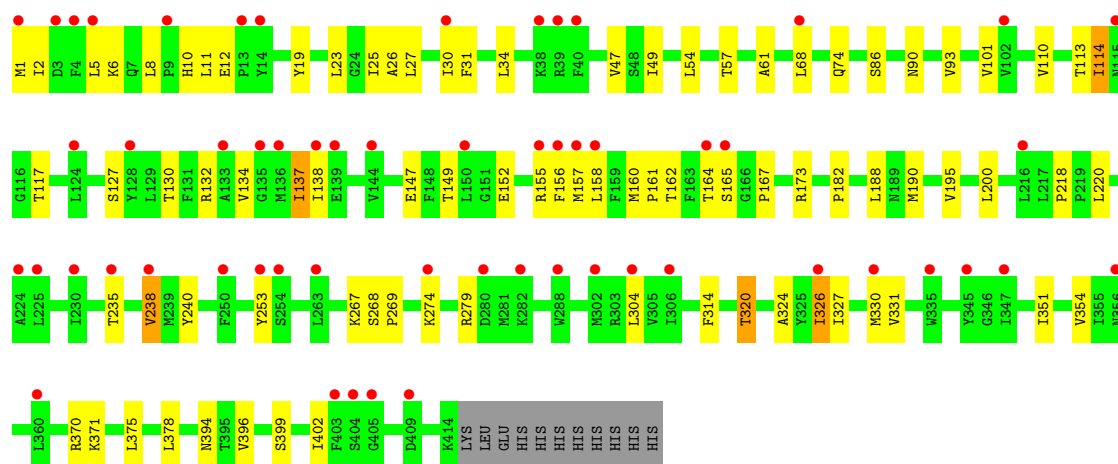
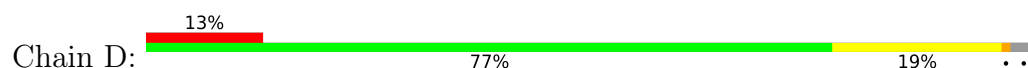


- Molecule 1: Integral membrane protein





● Molecule 1: Integral membrane protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	140.18Å 242.06Å 96.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	121.31 – 3.27 121.31 – 3.27	Depositor EDS
% Data completeness (in resolution range)	96.5 (121.31-3.27) 96.5 (121.31-3.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 3.25Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.280 , 0.300 0.285 , 0.311	Depositor DCC
R_{free} test set	2434 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	117.2	Xtriage
Anisotropy	0.374	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 75.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13792	wwPDB-VP
Average B, all atoms (Å ²)	151.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% 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	0.59	0/3556	1.09	4/4811 (0.1%)
1	B	0.58	0/3556	1.09	4/4811 (0.1%)
1	C	0.59	0/3556	1.10	2/4811 (0.0%)
1	D	0.56	0/3556	1.07	1/4811 (0.0%)
All	All	0.58	0/14224	1.09	11/19244 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	160	MET	N-CA-C	7.64	124.02	113.16
1	B	166	GLY	CA-C-N	6.65	128.15	119.84
1	B	166	GLY	C-N-CA	6.65	128.15	119.84
1	D	218	PRO	N-CA-C	6.43	118.54	110.70
1	A	218	PRO	N-CA-C	6.43	118.54	110.70
1	C	218	PRO	N-CA-C	6.41	118.52	110.70
1	B	218	PRO	N-CA-C	6.34	118.43	110.70
1	C	160	MET	N-CA-C	6.26	121.26	112.75
1	A	166	GLY	CA-C-N	6.13	127.51	119.84
1	A	166	GLY	C-N-CA	6.13	127.51	119.84
1	B	160	MET	N-CA-C	6.06	120.99	112.75

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	3448	0	3527	38	0
1	B	3448	0	3527	53	0
1	C	3448	0	3527	37	0
1	D	3448	0	3527	46	0
All	All	13792	0	14108	166	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 6.

All (166) 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:160:MET:HG3	1:B:161:PRO:CD	2.01	0.90
1:C:160:MET:HG3	1:C:161:PRO:CD	2.03	0.88
1:D:235:THR:HA	1:D:238:VAL:HG22	1.58	0.86
1:B:76:ILE:HD13	1:D:25:ILE:HG12	1.60	0.82
1:B:76:ILE:HG12	1:D:25:ILE:CD1	2.09	0.81
1:C:160:MET:HG3	1:C:161:PRO:N	1.96	0.80
1:B:160:MET:HG3	1:B:161:PRO:HD3	1.63	0.80
1:B:76:ILE:HG12	1:D:25:ILE:HD11	1.64	0.79
1:D:165:SER:O	1:D:253:TYR:CE2	2.35	0.79
1:B:76:ILE:CD1	1:D:25:ILE:HG12	2.13	0.78
1:C:160:MET:HG3	1:C:161:PRO:HD3	1.65	0.77
1:B:160:MET:HG3	1:B:161:PRO:N	1.95	0.77
1:B:162:THR:HA	1:B:166:GLY:HA3	1.72	0.71
1:B:168:ILE:N	1:B:168:ILE:HD12	2.06	0.71
1:A:131:PHE:CD2	1:A:168:ILE:HD11	2.26	0.70
1:B:76:ILE:CG2	1:D:25:ILE:HD13	2.26	0.66
1:B:162:THR:CA	1:B:166:GLY:HA3	2.26	0.65
1:C:90:ASN:HD22	1:C:93:VAL:HG23	1.62	0.65
1:D:90:ASN:HD22	1:D:93:VAL:HG23	1.63	0.64
1:A:90:ASN:HD22	1:A:93:VAL:HG23	1.62	0.64
1:B:90:ASN:HD22	1:B:93:VAL:HG23	1.63	0.64
1:D:130:THR:O	1:D:134:VAL:HG23	1.99	0.63
1:D:370:ARG:HD2	1:D:378:LEU:HD12	1.81	0.63
1:A:130:THR:O	1:A:134:VAL:HG23	1.99	0.63
1:D:165:SER:O	1:D:253:TYR:HE2	1.82	0.63
1:C:130:THR:O	1:C:134:VAL:HG23	1.99	0.62
1:B:76:ILE:HG21	1:D:25:ILE:HD13	1.80	0.62
1:B:130:THR:O	1:B:134:VAL:HG23	1.99	0.62
1:B:168:ILE:HD12	1:B:168:ILE:H	1.62	0.62
1:D:110:VAL:HG12	1:D:114:ILE:HD12	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:235:THR:HA	1:D:238:VAL:CG2	2.29	0.61
1:C:113:THR:HG22	1:C:113:THR:O	2.01	0.61
1:C:160:MET:CG	1:C:161:PRO:HD3	2.30	0.61
1:A:164:THR:HG23	1:A:253:TYR:OH	2.00	0.61
1:B:160:MET:CG	1:B:161:PRO:HD3	2.29	0.61
1:A:314:PHE:HB2	1:A:320:THR:HG23	1.82	0.60
1:B:314:PHE:HB2	1:B:320:THR:HG23	1.81	0.60
1:D:314:PHE:HB2	1:D:320:THR:HG23	1.82	0.60
1:C:314:PHE:HB2	1:C:320:THR:HG23	1.81	0.60
1:B:164:THR:HG23	1:B:253:TYR:OH	2.02	0.59
1:A:165:SER:O	1:A:165:SER:OG	2.21	0.59
1:D:113:THR:O	1:D:113:THR:HG22	2.02	0.58
1:C:110:VAL:HG12	1:C:114:ILE:HD12	1.86	0.57
1:A:162:THR:O	1:A:162:THR:OG1	2.20	0.56
1:B:113:THR:HG22	1:B:113:THR:O	2.06	0.55
1:B:101:VAL:HG21	1:B:137:ILE:HD13	1.89	0.55
1:C:327:ILE:O	1:C:331:VAL:HG23	2.07	0.55
1:D:26:ALA:HB1	1:D:47:VAL:HG21	1.89	0.55
1:A:131:PHE:CD2	1:A:168:ILE:CD1	2.89	0.55
1:A:113:THR:O	1:A:113:THR:HG22	2.08	0.54
1:A:110:VAL:HG12	1:A:114:ILE:HD12	1.89	0.54
1:B:149:THR:HB	1:B:152:GLU:H	1.72	0.54
1:B:168:ILE:H	1:B:168:ILE:CD1	2.20	0.54
1:C:26:ALA:HB1	1:C:47:VAL:HG21	1.90	0.54
1:B:327:ILE:O	1:B:331:VAL:HG23	2.08	0.53
1:A:149:THR:HB	1:A:152:GLU:H	1.72	0.53
1:B:110:VAL:HG12	1:B:114:ILE:HD12	1.90	0.53
1:D:149:THR:HB	1:D:152:GLU:H	1.72	0.53
1:C:101:VAL:HG21	1:C:137:ILE:HD13	1.90	0.53
1:C:149:THR:HB	1:C:152:GLU:H	1.72	0.53
1:D:101:VAL:HG21	1:D:137:ILE:HD13	1.89	0.53
1:D:351:ILE:HA	1:D:354:VAL:HG22	1.91	0.53
1:A:26:ALA:HB1	1:A:47:VAL:HG21	1.91	0.53
1:A:327:ILE:O	1:A:331:VAL:HG23	2.07	0.53
1:B:382:LYS:HE3	1:B:383:TRP:CZ3	2.43	0.53
1:C:351:ILE:HA	1:C:354:VAL:HG22	1.91	0.52
1:A:101:VAL:HG21	1:A:137:ILE:HD13	1.90	0.52
1:D:327:ILE:O	1:D:331:VAL:HG23	2.08	0.52
1:B:76:ILE:HG12	1:D:25:ILE:HD13	1.89	0.52
1:B:168:ILE:N	1:B:168:ILE:CD1	2.73	0.52
1:D:10:HIS:HE1	1:D:12:GLU:HG2	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:HIS:HE1	1:A:12:GLU:HG2	1.75	0.51
1:B:26:ALA:HB1	1:B:47:VAL:HG21	1.91	0.51
1:B:351:ILE:HA	1:B:354:VAL:HG22	1.91	0.51
1:C:164:THR:HG23	1:C:253:TYR:OH	2.11	0.51
1:A:351:ILE:HA	1:A:354:VAL:HG22	1.92	0.51
1:B:71:ILE:HG12	1:B:158:LEU:HD21	1.92	0.51
1:A:240:TYR:OH	1:A:394:ASN:ND2	2.44	0.50
1:B:10:HIS:HE1	1:B:12:GLU:HG2	1.76	0.50
1:B:163:PHE:C	1:B:163:PHE:CD1	2.90	0.50
1:A:6:LYS:HG3	1:A:61:ALA:HB3	1.93	0.50
1:C:71:ILE:HG12	1:C:158:LEU:HD21	1.94	0.50
1:C:6:LYS:HG3	1:C:61:ALA:HB3	1.93	0.50
1:C:10:HIS:HE1	1:C:12:GLU:HG2	1.75	0.49
1:B:165:SER:O	1:B:165:SER:OG	2.27	0.49
1:D:268:SER:OG	1:D:269:PRO:HD2	2.12	0.49
1:B:74:GLN:HG3	1:B:157:MET:HE1	1.94	0.49
1:A:75:ILE:HD11	1:A:154:LEU:HD21	1.94	0.49
1:C:240:TYR:OH	1:C:394:ASN:ND2	2.47	0.48
1:A:163:PHE:C	1:A:163:PHE:CD1	2.91	0.48
1:D:6:LYS:HG3	1:D:61:ALA:HB3	1.94	0.48
1:B:8:LEU:HD23	1:B:57:THR:HG21	1.95	0.48
1:C:163:PHE:CD1	1:C:163:PHE:C	2.91	0.48
1:C:75:ILE:HD11	1:C:154:LEU:HD21	1.96	0.48
1:A:27:LEU:HA	1:A:30:ILE:HD12	1.97	0.47
1:A:74:GLN:HG3	1:A:157:MET:HE1	1.96	0.47
1:D:27:LEU:HA	1:D:30:ILE:HD12	1.96	0.47
1:D:74:GLN:HG3	1:D:157:MET:HE1	1.95	0.47
1:B:27:LEU:HA	1:B:30:ILE:HD12	1.97	0.47
1:D:240:TYR:OH	1:D:394:ASN:ND2	2.45	0.47
1:D:326:ILE:O	1:D:330:MET:HB2	2.15	0.47
1:C:162:THR:HG22	1:C:167:PRO:HD2	1.97	0.47
1:B:75:ILE:HD11	1:B:154:LEU:HD21	1.96	0.47
1:D:162:THR:HG22	1:D:167:PRO:HD2	1.97	0.47
1:B:6:LYS:HG3	1:B:61:ALA:HB3	1.96	0.46
1:A:166:GLY:HA2	1:A:253:TYR:HE2	1.80	0.46
1:D:182:PRO:HG3	1:D:190:MET:HE1	1.96	0.46
1:C:27:LEU:HA	1:C:30:ILE:HD12	1.97	0.46
1:A:326:ILE:O	1:A:330:MET:HB2	2.15	0.46
1:B:326:ILE:O	1:B:330:MET:HB2	2.16	0.46
1:D:235:THR:CA	1:D:238:VAL:HG22	2.39	0.46
1:D:160:MET:HG3	1:D:161:PRO:HD3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:TYR:CZ	1:A:1:MET:HG3	2.51	0.46
1:C:326:ILE:O	1:C:330:MET:HB2	2.16	0.46
1:B:240:TYR:OH	1:B:394:ASN:ND2	2.46	0.45
1:A:399:SER:HA	1:A:402:ILE:HD12	1.98	0.45
1:B:162:THR:HB	1:B:166:GLY:HA3	1.99	0.45
1:B:399:SER:HA	1:B:402:ILE:HD12	1.98	0.45
1:D:8:LEU:HD23	1:D:57:THR:HG21	1.98	0.45
1:C:182:PRO:HG2	1:C:187:LEU:HD13	1.99	0.45
1:D:110:VAL:CG1	1:D:114:ILE:HD12	2.47	0.45
1:C:399:SER:HA	1:C:402:ILE:HD12	1.99	0.44
1:A:8:LEU:HD23	1:A:57:THR:HG21	1.99	0.44
1:D:399:SER:HA	1:D:402:ILE:HD12	1.99	0.44
1:C:165:SER:O	1:C:165:SER:OG	2.35	0.44
1:A:70:TYR:HE2	1:A:157:MET:HE2	1.83	0.44
1:A:159:PHE:CD2	1:A:161:PRO:HD2	2.53	0.44
1:A:131:PHE:HD2	1:A:168:ILE:HD11	1.79	0.43
1:B:182:PRO:HG2	1:B:187:LEU:HD13	2.00	0.43
1:C:367:ASN:ND2	1:C:378:LEU:H	2.17	0.43
1:C:8:LEU:HD23	1:C:57:THR:HG21	2.00	0.42
1:A:367:ASN:ND2	1:A:378:LEU:H	2.17	0.42
1:B:162:THR:CB	1:B:166:GLY:HA3	2.49	0.42
1:C:132:ARG:HG2	1:C:156:PHE:CZ	2.54	0.42
1:B:367:ASN:ND2	1:B:378:LEU:H	2.17	0.42
1:A:370:ARG:HG2	1:A:375:LEU:HD12	2.01	0.42
1:D:190:MET:HE2	1:D:267:LYS:HD3	2.01	0.42
1:B:91:LYS:HG3	1:B:92:TRP:CD1	2.55	0.42
1:A:160:MET:HE3	1:A:161:PRO:N	2.34	0.42
1:A:239:MET:HG3	1:A:401:LEU:HB2	2.02	0.42
1:D:160:MET:HG3	1:D:161:PRO:CD	2.50	0.42
1:D:110:VAL:HG12	1:D:114:ILE:CD1	2.48	0.42
1:D:132:ARG:HG2	1:D:156:PHE:CZ	2.55	0.42
1:C:19:TYR:HE2	1:C:23:LEU:HD22	1.85	0.41
1:A:159:PHE:CE2	1:A:161:PRO:HG2	2.55	0.41
1:A:132:ARG:HG2	1:A:156:PHE:CZ	2.54	0.41
1:B:370:ARG:HG2	1:B:375:LEU:HD12	2.03	0.41
1:B:11:LEU:HD12	1:B:54:LEU:HD13	2.02	0.41
1:D:19:TYR:HE2	1:D:23:LEU:HD22	1.86	0.41
1:C:110:VAL:CG1	1:C:114:ILE:HD12	2.50	0.41
1:B:31:PHE:HA	1:B:34:LEU:HD12	2.02	0.41
1:D:370:ARG:HG2	1:D:375:LEU:HD12	2.03	0.41
1:C:11:LEU:HD12	1:C:54:LEU:HD13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:370:ARG:HG2	1:C:375:LEU:HD12	2.02	0.41
1:B:19:TYR:HE1	1:B:23:LEU:HD22	1.85	0.41
1:D:31:PHE:HA	1:D:34:LEU:HD12	2.02	0.41
1:C:31:PHE:HA	1:C:34:LEU:HD12	2.02	0.41
1:A:14:TYR:HE1	1:A:260:VAL:HG21	1.86	0.41
1:D:220:LEU:HB3	1:D:238:VAL:HG13	2.03	0.41
1:D:304:LEU:HD23	1:D:324:ALA:HA	2.03	0.41
1:B:132:ARG:HG2	1:B:156:PHE:CZ	2.55	0.40
1:D:11:LEU:HD12	1:D:54:LEU:HD13	2.02	0.40
1:B:239:MET:HG3	1:B:401:LEU:HB2	2.03	0.40
1:C:166:GLY:HA2	1:C:253:TYR:HE2	1.86	0.40
1:A:19:TYR:HE1	1:A:23:LEU:HD22	1.87	0.40
1:B:250:PHE:HD2	1:B:272:PHE:CZ	2.40	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	412/425 (97%)	391 (95%)	21 (5%)	0	100	100
1	B	412/425 (97%)	392 (95%)	20 (5%)	0	100	100
1	C	412/425 (97%)	392 (95%)	20 (5%)	0	100	100
1	D	412/425 (97%)	391 (95%)	21 (5%)	0	100	100
All	All	1648/1700 (97%)	1566 (95%)	82 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	370/381 (97%)	339 (92%)	31 (8%)	10	34
1	B	370/381 (97%)	340 (92%)	30 (8%)	11	36
1	C	370/381 (97%)	344 (93%)	26 (7%)	14	41
1	D	370/381 (97%)	344 (93%)	26 (7%)	14	41
All	All	1480/1524 (97%)	1367 (92%)	113 (8%)	12	38

All (113) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	1	MET
1	C	37	LYS
1	C	49	ILE
1	C	68	LEU
1	C	86	SER
1	C	114	ILE
1	C	117	THR
1	C	127	SER
1	C	137	ILE
1	C	138	ILE
1	C	147	GLU
1	C	155	ARG
1	C	158	LEU
1	C	160	MET
1	C	164	THR
1	C	165	SER
1	C	168	ILE
1	C	173	ARG
1	C	195	VAL
1	C	274	LYS
1	C	279	ARG
1	C	320	THR
1	C	326	ILE
1	C	376	LYS

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Mol	Chain	Res	Type
1	C	378	LEU
1	C	396	VAL
1	A	1	MET
1	A	2	ILE
1	A	49	ILE
1	A	68	LEU
1	A	86	SER
1	A	96	LEU
1	A	114	ILE
1	A	117	THR
1	A	127	SER
1	A	137	ILE
1	A	138	ILE
1	A	144	VAL
1	A	147	GLU
1	A	155	ARG
1	A	158	LEU
1	A	160	MET
1	A	164	THR
1	A	168	ILE
1	A	173	ARG
1	A	188	LEU
1	A	195	VAL
1	A	196	LYS
1	A	274	LYS
1	A	279	ARG
1	A	320	THR
1	A	326	ILE
1	A	338	ILE
1	A	368	LYS
1	A	371	LYS
1	A	378	LEU
1	A	396	VAL
1	B	1	MET
1	B	2	ILE
1	B	5	LEU
1	B	37	LYS
1	B	49	ILE
1	B	68	LEU
1	B	86	SER
1	B	114	ILE
1	B	117	THR

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Mol	Chain	Res	Type
1	B	127	SER
1	B	137	ILE
1	B	138	ILE
1	B	144	VAL
1	B	147	GLU
1	B	155	ARG
1	B	158	LEU
1	B	160	MET
1	B	164	THR
1	B	168	ILE
1	B	173	ARG
1	B	176	GLU
1	B	188	LEU
1	B	195	VAL
1	B	200	LEU
1	B	268	SER
1	B	274	LYS
1	B	279	ARG
1	B	320	THR
1	B	326	ILE
1	B	396	VAL
1	D	1	MET
1	D	2	ILE
1	D	5	LEU
1	D	49	ILE
1	D	68	LEU
1	D	86	SER
1	D	114	ILE
1	D	117	THR
1	D	127	SER
1	D	137	ILE
1	D	138	ILE
1	D	147	GLU
1	D	155	ARG
1	D	158	LEU
1	D	164	THR
1	D	173	ARG
1	D	188	LEU
1	D	195	VAL
1	D	200	LEU
1	D	238	VAL
1	D	274	LYS

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Mol	Chain	Res	Type
1	D	279	ARG
1	D	320	THR
1	D	326	ILE
1	D	371	LYS
1	D	396	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	10	HIS
1	C	90	ASN
1	C	118	GLN
1	C	122	ASN
1	C	183	ASN
1	C	193	GLN
1	C	318	ASN
1	C	328	ASN
1	C	336	HIS
1	C	367	ASN
1	C	394	ASN
1	A	10	HIS
1	A	90	ASN
1	A	122	ASN
1	A	183	ASN
1	A	193	GLN
1	A	318	ASN
1	A	328	ASN
1	A	336	HIS
1	A	367	ASN
1	A	394	ASN
1	B	10	HIS
1	B	90	ASN
1	B	122	ASN
1	B	183	ASN
1	B	193	GLN
1	B	318	ASN
1	B	328	ASN
1	B	336	HIS
1	B	367	ASN
1	B	394	ASN
1	D	10	HIS
1	D	90	ASN

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Mol	Chain	Res	Type
1	D	115	ASN
1	D	118	GLN
1	D	122	ASN
1	D	183	ASN
1	D	193	GLN
1	D	318	ASN
1	D	328	ASN
1	D	336	HIS
1	D	356	ASN
1	D	367	ASN
1	D	394	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	414/425 (97%)	0.29	26 (6%) 26 18	86, 114, 163, 306	0
1	B	414/425 (97%)	0.47	36 (8%) 16 13	103, 138, 187, 236	0
1	C	414/425 (97%)	0.32	23 (5%) 30 20	86, 115, 171, 259	0
1	D	414/425 (97%)	0.86	57 (13%) 6 5	158, 217, 271, 289	0
All	All	1656/1700 (97%)	0.48	142 (8%) 16 13	86, 135, 249, 306	0

All (142) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	165	SER	9.0
1	D	40	PHE	8.6
1	D	253	TYR	7.7
1	A	277	ILE	7.6
1	A	278	SER	7.2
1	B	369	ASP	6.6
1	A	2	ILE	6.6
1	D	164	THR	6.5
1	B	36	PHE	5.5
1	B	250	PHE	5.1
1	A	314	PHE	4.6
1	C	273	ASP	4.4
1	D	306	ILE	4.4
1	A	313	VAL	4.4
1	D	144	VAL	4.2
1	D	39	ARG	4.2
1	C	60	HIS	4.1
1	D	136	MET	4.1
1	B	253	TYR	4.0
1	D	3	ASP	4.0
1	A	283	GLU	4.0

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Mol	Chain	Res	Type	RSRZ
1	C	92	TRP	3.9
1	B	9	PRO	3.9
1	D	250	PHE	3.9
1	C	80	SER	3.8
1	B	289	HIS	3.7
1	C	1	MET	3.7
1	A	387	LEU	3.7
1	B	272	PHE	3.7
1	B	2	ILE	3.6
1	A	156	PHE	3.6
1	D	38	LYS	3.5
1	D	274	LYS	3.5
1	D	14	TYR	3.5
1	C	221	LYS	3.4
1	B	220	LEU	3.4
1	B	114	ILE	3.4
1	D	4	PHE	3.4
1	D	326	ILE	3.3
1	D	356	ASN	3.3
1	C	2	ILE	3.3
1	D	135	GLY	3.3
1	A	280	ASP	3.3
1	B	217	LEU	3.3
1	C	139	GLU	3.3
1	B	37	LYS	3.1
1	A	176	GLU	3.1
1	A	275	PRO	3.1
1	C	83	ARG	3.0
1	D	230	ILE	3.0
1	A	171	PHE	3.0
1	D	302	MET	3.0
1	B	296	PHE	3.0
1	B	202	PHE	3.0
1	C	90	ASN	3.0
1	B	218	PRO	2.9
1	D	282	LYS	2.9
1	D	345	TYR	2.9
1	D	13	PRO	2.9
1	B	241	VAL	2.9
1	D	156	PHE	2.9
1	C	233	LEU	2.8
1	C	93	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	181	ILE	2.8
1	D	115	ASN	2.8
1	D	404	SER	2.8
1	C	305	VAL	2.8
1	D	238	VAL	2.8
1	B	345	TYR	2.7
1	D	133	ALA	2.7
1	C	5	LEU	2.7
1	C	188	LEU	2.7
1	A	299	PHE	2.7
1	D	139	GLU	2.7
1	A	128	TYR	2.7
1	B	309	MET	2.7
1	A	315	LYS	2.7
1	D	347	ILE	2.6
1	B	292	LEU	2.6
1	C	316	ASN	2.6
1	D	216	LEU	2.6
1	A	307	VAL	2.6
1	D	102	VAL	2.5
1	D	360	LEU	2.5
1	D	254	SER	2.5
1	D	409	ASP	2.5
1	B	136	MET	2.5
1	D	157	MET	2.5
1	C	280	ASP	2.5
1	B	271	ASN	2.5
1	A	413	THR	2.5
1	D	225	LEU	2.5
1	B	48	SER	2.4
1	A	301	PHE	2.4
1	A	279	ARG	2.4
1	D	124	LEU	2.4
1	D	138	ILE	2.4
1	C	124	LEU	2.4
1	D	68	LEU	2.4
1	B	164	THR	2.4
1	B	183	ASN	2.4
1	A	1	MET	2.4
1	A	159	PHE	2.4
1	D	150	LEU	2.4
1	D	9	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	242	TYR	2.3
1	D	128	TYR	2.3
1	C	78	VAL	2.3
1	B	221	LYS	2.3
1	B	160	MET	2.3
1	B	352	GLY	2.3
1	D	158	LEU	2.3
1	D	1	MET	2.3
1	D	224	ALA	2.3
1	C	321	SER	2.3
1	B	165	SER	2.3
1	A	4	PHE	2.2
1	C	347	ILE	2.2
1	B	285	TRP	2.2
1	B	346	GLY	2.2
1	D	304	LEU	2.2
1	B	30	ILE	2.2
1	D	155	ARG	2.2
1	C	409	ASP	2.2
1	D	263	LEU	2.1
1	D	330	MET	2.1
1	D	405	GLY	2.1
1	C	159	PHE	2.1
1	D	280	ASP	2.1
1	A	102	VAL	2.1
1	D	30	ILE	2.1
1	A	308	LEU	2.1
1	A	375	LEU	2.1
1	D	335	TRP	2.1
1	B	54	LEU	2.1
1	D	5	LEU	2.1
1	A	154	LEU	2.1
1	D	288	TRP	2.1
1	D	235	THR	2.1
1	B	325	TYR	2.0
1	B	413	THR	2.0
1	D	403	PHE	2.0

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

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.