



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 09:41 AM EDT

PDB ID : 2Q74 / pdb_00002q74
Title : Mycobacterium tuberculosis SuhB
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Deposited on : 2007-06-06
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

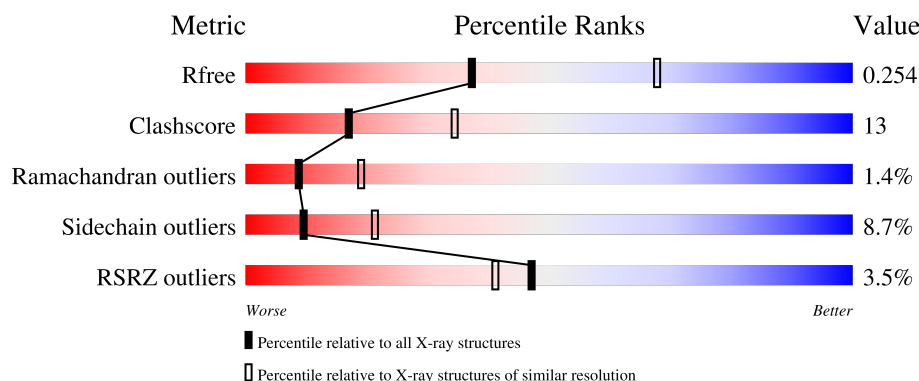
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

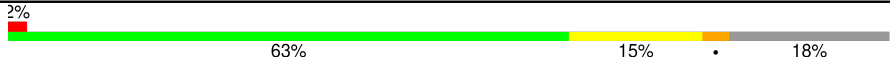
The reported resolution of this entry is 2.60 Å.

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	299	
1	B	299	
1	C	299	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5466 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 Inositol-1-monophosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1784	1115	325	339	5			
1	B	245	Total	C	N	O	S	0	0	0
			1784	1115	325	338	6			
1	C	243	Total	C	N	O	S	0	0	0
			1774	1109	323	336	6			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	291	ALA	-	expression tag	UNP P65165
A	292	LEU	-	expression tag	UNP P65165
A	293	GLU	-	expression tag	UNP P65165
A	294	HIS	-	expression tag	UNP P65165
A	295	HIS	-	expression tag	UNP P65165
A	296	HIS	-	expression tag	UNP P65165
A	297	HIS	-	expression tag	UNP P65165
A	298	HIS	-	expression tag	UNP P65165
A	299	HIS	-	expression tag	UNP P65165
B	291	ALA	-	expression tag	UNP P65165
B	292	LEU	-	expression tag	UNP P65165
B	293	GLU	-	expression tag	UNP P65165
B	294	HIS	-	expression tag	UNP P65165
B	295	HIS	-	expression tag	UNP P65165
B	296	HIS	-	expression tag	UNP P65165
B	297	HIS	-	expression tag	UNP P65165
B	298	HIS	-	expression tag	UNP P65165
B	299	HIS	-	expression tag	UNP P65165
C	291	ALA	-	expression tag	UNP P65165
C	292	LEU	-	expression tag	UNP P65165
C	293	GLU	-	expression tag	UNP P65165
C	294	HIS	-	expression tag	UNP P65165
C	295	HIS	-	expression tag	UNP P65165

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Chain	Residue	Modelled	Actual	Comment	Reference
C	296	HIS	-	expression tag	UNP P65165
C	297	HIS	-	expression tag	UNP P65165
C	298	HIS	-	expression tag	UNP P65165
C	299	HIS	-	expression tag	UNP P65165

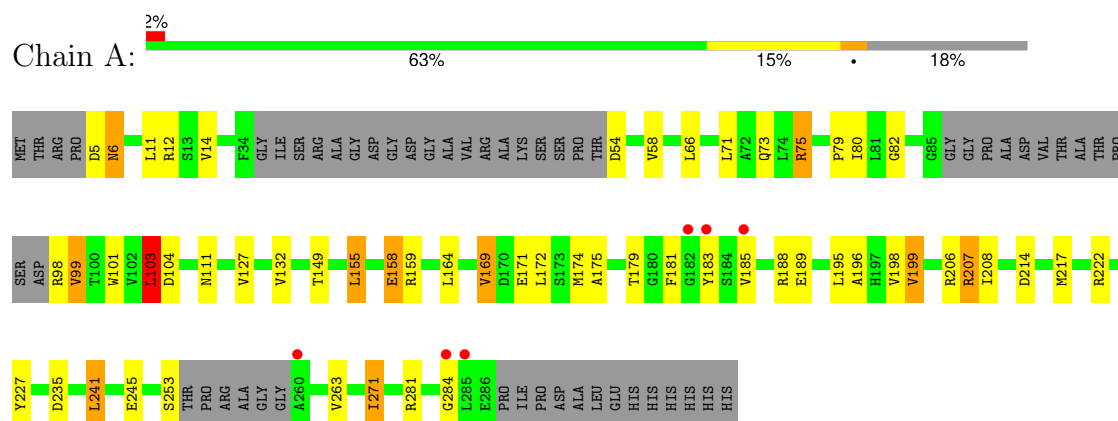
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	41	Total O 41 41	0	0
2	B	31	Total O 31 31	0	0
2	C	52	Total O 52 52	0	0

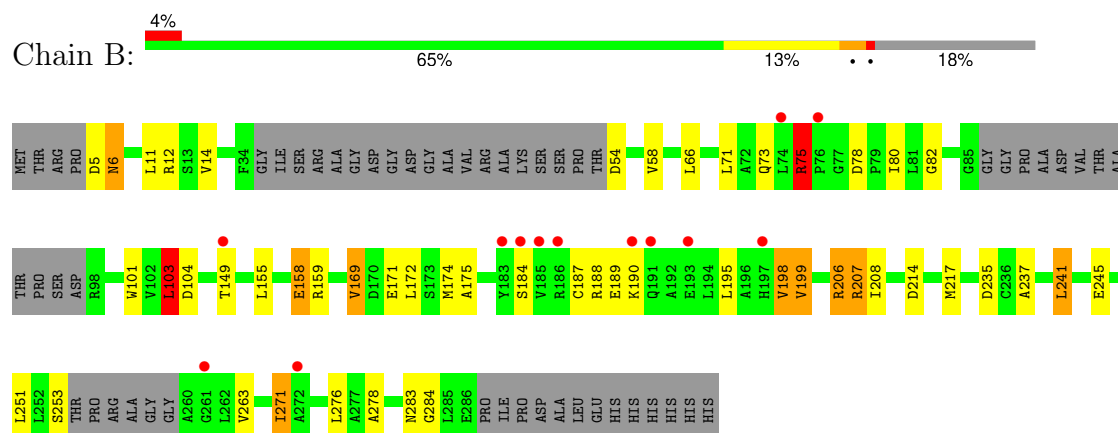
3 Residue-property plots [i](#)

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.

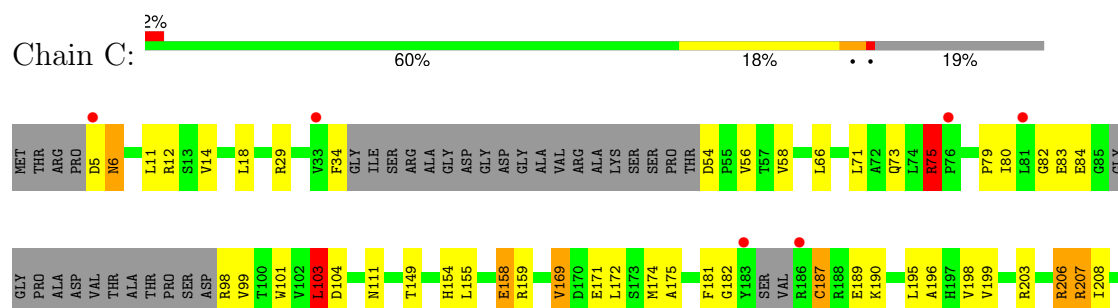
• Molecule 1: Inositol-1-monophosphatase



• Molecule 1: Inositol-1-monophosphatase



• Molecule 1: Inositol-1-monophosphatase





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	101.40Å 185.60Å 107.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.88 – 2.60 19.88 – 2.60	Depositor EDS
% Data completeness (in resolution range)	93.2 (19.88-2.60) 92.3 (19.88-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.28 (at 2.59Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.223 , 0.245 0.217 , 0.254	Depositor DCC
R_{free} test set	1469 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	68.9	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 65.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.015 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.021 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5466	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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	0.77	1/1806 (0.1%)	0.91	4/2457 (0.2%)
1	B	0.60	0/1806	0.85	4/2457 (0.2%)
1	C	0.74	0/1795	0.90	3/2440 (0.1%)
All	All	0.71	1/5407 (0.0%)	0.89	11/7354 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	199	VAL	CA-CB	7.84	1.58	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	103	LEU	N-CA-C	6.79	119.64	108.99
1	A	188	ARG	N-CA-C	6.69	118.65	111.36
1	A	103	LEU	N-CA-C	5.97	118.48	109.23
1	B	103	LEU	N-CA-C	5.85	118.17	108.99
1	C	75	ARG	CA-C-N	5.71	126.98	119.84
1	C	75	ARG	C-N-CA	5.71	126.98	119.84
1	B	75	ARG	CA-C-N	5.53	126.75	119.84
1	B	75	ARG	C-N-CA	5.53	126.75	119.84
1	A	75	ARG	CA-C-N	5.40	126.59	119.84
1	A	75	ARG	C-N-CA	5.40	126.59	119.84
1	B	199	VAL	CB-CA-C	-5.09	108.87	113.70

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	1784	0	1769	43	0
1	B	1784	0	1769	43	1
1	C	1774	0	1764	57	0
2	A	41	0	0	5	0
2	B	31	0	0	9	1
2	C	52	0	0	16	0
All	All	5466	0	5302	135	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 13.

All (135) 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:195:LEU:CD2	2:A:312:HOH:O	1.93	1.17
1:C:29:ARG:HD3	2:C:287:HOH:O	1.50	1.08
1:B:199:VAL:HG11	2:B:316:HOH:O	1.68	0.92
1:A:195:LEU:HD23	2:A:312:HOH:O	1.56	0.92
1:B:278:ALA:HB1	2:B:310:HOH:O	1.72	0.90
1:C:56:VAL:HG12	2:C:310:HOH:O	1.81	0.80
1:A:171:GLU:C	1:A:271:ILE:HD11	2.09	0.77
1:C:171:GLU:C	1:C:271:ILE:HD11	2.10	0.77
1:C:251:LEU:HD23	2:C:323:HOH:O	1.86	0.75
1:B:237:ALA:HB1	2:B:306:HOH:O	1.86	0.74
1:B:171:GLU:C	1:B:271:ILE:HD11	2.13	0.74
1:C:227:TYR:HB2	2:C:322:HOH:O	1.89	0.72
1:A:111:ASN:OD1	1:C:206:ARG:NH1	2.22	0.72
1:C:182:GLY:HA3	1:C:230:GLY:O	1.91	0.70
1:C:79:PRO:HD2	1:C:99:VAL:O	1.92	0.70
1:A:132:VAL:HG22	2:A:318:HOH:O	1.91	0.69
1:C:265:VAL:HG13	2:C:322:HOH:O	1.94	0.67
1:C:208:ILE:CD1	1:C:217:MET:HE1	2.26	0.66
1:A:281:ARG:HH22	1:B:283:ASN:CG	2.06	0.64
1:C:265:VAL:HG22	2:C:322:HOH:O	1.98	0.64
1:A:206:ARG:NH1	1:C:111:ASN:OD1	2.33	0.62
1:C:18:LEU:HD23	2:C:338:HOH:O	1.99	0.61
2:B:303:HOH:O	1:C:187:CYS:HB2	2.00	0.61
1:C:34:PHE:CD1	2:C:330:HOH:O	2.51	0.60
1:A:281:ARG:HH22	1:B:283:ASN:ND2	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:54:ASP:O	1:C:58:VAL:HG23	2.03	0.59
1:A:196:ALA:O	1:C:189:GLU:HG3	2.01	0.59
1:B:174:MET:HA	1:B:174:MET:HE2	1.83	0.59
1:B:54:ASP:O	1:B:58:VAL:HG23	2.01	0.59
1:C:187:CYS:O	1:C:187:CYS:SG	2.61	0.58
1:A:54:ASP:O	1:A:58:VAL:HG23	2.03	0.58
1:A:71:LEU:HD13	1:A:80:ILE:HD12	1.86	0.58
1:C:208:ILE:HD12	1:C:217:MET:HE1	1.86	0.58
1:B:208:ILE:CD1	1:B:217:MET:HE1	2.35	0.57
1:B:5:ASP:C	2:B:302:HOH:O	2.47	0.57
1:C:169:VAL:HB	1:C:271:ILE:HD12	1.87	0.57
1:B:14:VAL:HB	2:B:297:HOH:O	2.05	0.57
1:A:189:GLU:HG3	1:C:196:ALA:O	2.04	0.57
1:C:83:GLU:HB3	2:C:299:HOH:O	2.04	0.57
1:A:208:ILE:CD1	1:A:217:MET:HE1	2.34	0.57
1:C:71:LEU:HD13	1:C:80:ILE:HD12	1.87	0.56
1:C:174:MET:HA	1:C:174:MET:HE2	1.86	0.56
1:B:206:ARG:CD	2:B:311:HOH:O	2.53	0.56
1:B:206:ARG:NE	2:B:311:HOH:O	2.38	0.56
1:B:169:VAL:HB	1:B:271:ILE:HD12	1.87	0.56
1:A:208:ILE:HD12	1:A:217:MET:HE1	1.86	0.55
1:A:169:VAL:HB	1:A:271:ILE:HD12	1.87	0.55
1:A:214:ASP:HA	1:A:217:MET:HE3	1.87	0.55
1:B:187:CYS:O	1:B:189:GLU:N	2.40	0.55
1:A:195:LEU:HD21	2:A:312:HOH:O	1.82	0.55
1:A:207:ARG:HB3	1:C:207:ARG:HB3	1.90	0.54
1:C:171:GLU:O	1:C:271:ILE:HD11	2.07	0.53
1:C:6:ASN:HB2	2:C:306:HOH:O	2.08	0.53
1:B:171:GLU:O	1:B:271:ILE:HD11	2.08	0.53
1:A:14:VAL:CG2	1:A:101:TRP:CH2	2.92	0.52
1:B:71:LEU:HD13	1:B:80:ILE:HD12	1.91	0.52
1:C:260:ALA:HB2	2:C:325:HOH:O	2.09	0.52
1:C:195:LEU:O	1:C:199:VAL:HG23	2.09	0.52
1:B:208:ILE:HD12	1:B:217:MET:HE1	1.91	0.51
1:A:79:PRO:HD2	1:A:99:VAL:O	2.11	0.51
1:A:171:GLU:O	1:A:271:ILE:HD11	2.09	0.51
1:A:207:ARG:HD2	2:C:291:HOH:O	2.10	0.51
1:C:34:PHE:CG	2:C:330:HOH:O	2.62	0.51
1:C:208:ILE:HD13	1:C:217:MET:HE1	1.90	0.50
1:C:14:VAL:CG2	1:C:101:TRP:CH2	2.94	0.50
1:A:99:VAL:HG22	1:A:127:VAL:HG22	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:MET:HA	1:A:174:MET:HE2	1.94	0.49
1:B:206:ARG:HD3	2:B:311:HOH:O	2.11	0.49
1:B:66:LEU:C	1:B:66:LEU:HD23	2.38	0.49
1:B:14:VAL:CG2	1:B:101:TRP:CH2	2.96	0.48
1:B:104:ASP:OD2	1:B:235:ASP:OD1	2.32	0.48
1:C:5:ASP:O	1:C:6:ASN:HB2	2.14	0.48
1:B:214:ASP:HA	1:B:217:MET:HE3	1.94	0.48
1:A:195:LEU:O	1:A:199:VAL:HG23	2.14	0.47
1:B:175:ALA:HB2	1:B:271:ILE:HD13	1.95	0.47
1:C:174:MET:O	1:C:203:ARG:HD2	2.14	0.47
1:B:207:ARG:HH11	1:B:207:ARG:HB2	1.80	0.47
1:B:195:LEU:O	1:B:199:VAL:HG23	2.15	0.46
1:A:99:VAL:HG22	1:A:127:VAL:CG2	2.45	0.46
1:C:241:LEU:HD22	1:C:245:GLU:HG2	1.98	0.46
1:B:207:ARG:HH11	1:B:207:ARG:CB	2.28	0.46
1:A:104:ASP:OD2	1:A:235:ASP:OD1	2.33	0.45
1:C:82:GLY:HA2	1:C:103:LEU:HD23	1.97	0.45
1:A:181:PHE:CZ	1:A:207:ARG:NH2	2.85	0.45
1:B:251:LEU:HD21	1:B:276:LEU:HD11	1.99	0.45
1:C:182:GLY:CA	1:C:230:GLY:O	2.64	0.45
1:B:207:ARG:HH11	1:B:207:ARG:CG	2.30	0.45
1:A:82:GLY:HA2	1:A:103:LEU:HD23	1.98	0.45
1:B:253:SER:HB3	1:B:263:VAL:HB	1.99	0.44
1:A:179:THR:HG22	1:A:227:TYR:O	2.17	0.44
1:A:207:ARG:CG	1:A:207:ARG:HH11	2.30	0.44
1:C:75:ARG:HA	1:C:75:ARG:HD2	1.92	0.44
1:C:207:ARG:HH11	1:C:207:ARG:CB	2.31	0.44
1:C:207:ARG:HH11	1:C:207:ARG:CG	2.30	0.44
1:A:14:VAL:CG2	1:A:101:TRP:HH2	2.30	0.44
1:A:172:LEU:N	1:A:271:ILE:HD11	2.32	0.44
1:A:207:ARG:HH11	1:A:207:ARG:CB	2.31	0.43
1:B:208:ILE:HD13	1:B:217:MET:HE1	2.00	0.43
1:B:241:LEU:HD22	1:B:245:GLU:HG2	2.01	0.43
1:B:14:VAL:HG23	1:B:101:TRP:CH2	2.54	0.43
1:C:253:SER:HB3	1:C:263:VAL:HB	2.00	0.43
1:B:172:LEU:N	1:B:271:ILE:HD11	2.34	0.43
1:A:158:GLU:HG2	1:A:159:ARG:N	2.33	0.43
1:C:158:GLU:HG2	1:C:159:ARG:N	2.34	0.43
1:A:66:LEU:C	1:A:66:LEU:HD23	2.44	0.43
1:C:175:ALA:HB2	1:C:271:ILE:HD13	2.01	0.43
1:C:14:VAL:HG23	1:C:101:TRP:CH2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:VAL:HG23	1:A:101:TRP:CH2	2.54	0.42
1:C:214:ASP:HA	1:C:217:MET:HE3	2.01	0.42
1:B:158:GLU:HG2	1:B:159:ARG:N	2.34	0.42
1:B:195:LEU:HA	1:B:198:VAL:HG13	2.00	0.42
1:C:66:LEU:C	1:C:66:LEU:HD23	2.45	0.42
1:B:82:GLY:HA2	1:B:103:LEU:HD23	2.02	0.42
1:B:75:ARG:NH1	1:B:78:ASP:OD2	2.53	0.42
1:C:181:PHE:CZ	1:C:207:ARG:NH2	2.88	0.42
1:A:5:ASP:O	1:A:6:ASN:HB2	2.18	0.42
1:A:241:LEU:HD22	1:A:245:GLU:HG2	2.03	0.41
1:C:172:LEU:N	1:C:271:ILE:HD11	2.35	0.41
1:C:14:VAL:CG2	1:C:101:TRP:HH2	2.33	0.41
1:C:207:ARG:HH11	1:C:207:ARG:HB2	1.86	0.41
1:A:222:ARG:NE	2:A:308:HOH:O	2.54	0.41
1:C:104:ASP:OD2	1:C:235:ASP:OD1	2.38	0.41
1:C:229:HIS:ND1	1:C:263:VAL:HG22	2.35	0.41
1:C:187:CYS:HA	2:C:300:HOH:O	2.20	0.41
1:C:251:LEU:HD21	1:C:276:LEU:HD11	2.03	0.41
1:A:155:LEU:HB2	1:A:164:LEU:HD21	2.03	0.41
1:B:189:GLU:CB	1:C:190:LYS:HE2	2.50	0.41
1:B:190:LYS:HE3	2:C:300:HOH:O	2.20	0.41
1:B:75:ARG:HA	1:B:75:ARG:HD2	1.90	0.40
1:A:253:SER:HB3	1:A:263:VAL:HB	2.03	0.40
1:C:82:GLY:C	1:C:84:GLU:H	2.29	0.40
1:C:154:HIS:HD2	2:C:298:HOH:O	2.04	0.40
1:B:5:ASP:O	1:B:6:ASN:HB2	2.21	0.40
1:B:14:VAL:CG2	1:B:101:TRP:HH2	2.34	0.40
1:A:175:ALA:HB2	1:A:271:ILE:HD13	2.02	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:B:208:ILE:C	2:B:311:HOH:O[3_655]	2.17	0.03

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	237/299 (79%)	222 (94%)	11 (5%)	4 (2%)	7	15
1	B	237/299 (79%)	221 (93%)	12 (5%)	4 (2%)	7	15
1	C	233/299 (78%)	218 (94%)	13 (6%)	2 (1%)	14	30
All	All	707/897 (79%)	661 (94%)	36 (5%)	10 (1%)	9	19

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	B	6	ASN
1	C	6	ASN
1	A	185	VAL
1	B	188	ARG
1	A	183	TYR
1	B	184	SER
1	C	284	GLY
1	A	284	GLY
1	B	284	GLY

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	172/217 (79%)	157 (91%)	15 (9%)	9	21
1	B	172/217 (79%)	158 (92%)	14 (8%)	11	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	172/217 (79%)	156 (91%)	16 (9%)	8	18
All	All	516/651 (79%)	471 (91%)	45 (9%)	9	21

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	12	ARG
1	A	73	GLN
1	A	75	ARG
1	A	98	ARG
1	A	99	VAL
1	A	103	LEU
1	A	149	THR
1	A	155	LEU
1	A	158	GLU
1	A	169	VAL
1	A	198	VAL
1	A	207	ARG
1	A	241	LEU
1	A	271	ILE
1	B	11	LEU
1	B	12	ARG
1	B	73	GLN
1	B	75	ARG
1	B	103	LEU
1	B	149	THR
1	B	155	LEU
1	B	158	GLU
1	B	169	VAL
1	B	198	VAL
1	B	206	ARG
1	B	207	ARG
1	B	241	LEU
1	B	271	ILE
1	C	11	LEU
1	C	12	ARG
1	C	73	GLN
1	C	75	ARG
1	C	98	ARG
1	C	103	LEU
1	C	149	THR

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Mol	Chain	Res	Type
1	C	155	LEU
1	C	158	GLU
1	C	169	VAL
1	C	187	CYS
1	C	198	VAL
1	C	206	ARG
1	C	207	ARG
1	C	241	LEU
1	C	271	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	126	GLN
1	A	232	GLN
1	A	280	GLN
1	A	283	ASN
1	B	126	GLN
1	B	191	GLN
1	B	232	GLN
1	B	280	GLN
1	B	283	ASN
1	C	126	GLN
1	C	232	GLN
1	C	280	GLN
1	C	283	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

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

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	245/299 (81%)	0.13	6 (2%) 59 54	52, 69, 87, 94	0
1	B	245/299 (81%)	0.45	13 (5%) 32 26	52, 69, 87, 92	0
1	C	243/299 (81%)	0.12	7 (2%) 53 48	52, 69, 86, 93	0
All	All	733/897 (81%)	0.23	26 (3%) 47 41	52, 69, 87, 94	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	183	TYR	5.2
1	B	185	VAL	4.3
1	A	182	GLY	4.1
1	B	191	GLN	3.4
1	B	76	PRO	3.2
1	B	184	SER	3.2
1	A	285	LEU	2.9
1	A	284	GLY	2.8
1	B	272	ALA	2.6
1	B	261	GLY	2.6
1	C	76	PRO	2.6
1	B	197	HIS	2.5
1	B	190	LYS	2.5
1	C	285	LEU	2.5
1	B	193	GLU	2.5
1	A	185	VAL	2.5
1	B	149	THR	2.5
1	A	260	ALA	2.5
1	C	183	TYR	2.4
1	C	5	ASP	2.3
1	B	186	ARG	2.3
1	C	81	LEU	2.1
1	B	183	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	33	VAL	2.1
1	C	186	ARG	2.0
1	B	74	LEU	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.