



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:58 PM UTC

PDB ID : 3VW5 / pdb_00003vw5
Title : Crystal structure of sugar epimerase from ruminal bacterium
Authors : Fujiwara, T.; Saburi, W.; Tanaka, I.; Yao, M.
Deposited on : 2012-08-02
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

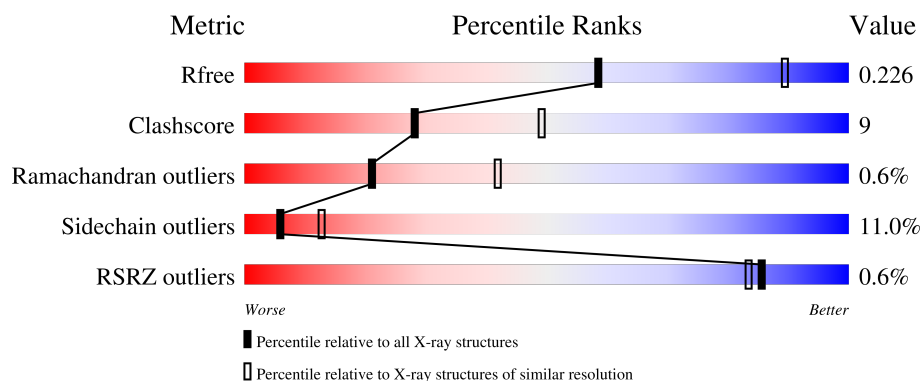
1 Overall quality at a glance

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	389	
1	B	389	
1	C	389	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9744 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 Cellobiose 2-epimerase.

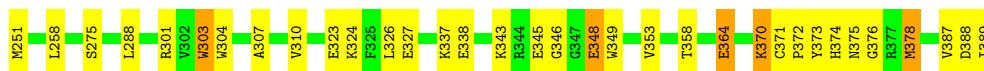
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			3196	2039	527	614	16			
1	B	389	Total	C	N	O	S	0	0	0
			3196	2039	527	614	16			
1	C	389	Total	C	N	O	S	0	0	0
			3196	2039	527	614	16			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	38	GLU	GLY	engineered mutation	UNP A8CF79
A	138	LEU	PRO	engineered mutation	UNP A8CF79
B	38	GLU	GLY	engineered mutation	UNP A8CF79
B	138	LEU	PRO	engineered mutation	UNP A8CF79
C	38	GLU	GLY	engineered mutation	UNP A8CF79
C	138	LEU	PRO	engineered mutation	UNP A8CF79

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	57	Total	O	0	0
			57	57		
2	B	59	Total	O	0	0
			59	59		
2	C	40	Total	O	0	0
			40	40		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.93Å 100.88Å 186.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.13 – 2.60 39.13 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.5 (39.13-2.60) 98.6 (39.13-2.60)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.173 , 0.222 0.177 , 0.226	Depositor DCC
R_{free} test set	2895 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	45.5	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.024 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9744	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.85% 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	0.52	0/3275	0.80	1/4427 (0.0%)
1	B	0.53	0/3275	0.82	3/4427 (0.1%)
1	C	0.53	0/3275	0.80	0/4427
All	All	0.52	0/9825	0.81	4/13281 (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
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	SER	N-CA-C	6.00	120.73	113.41
1	B	171	ASN	CB-CA-C	-5.72	109.45	117.23
1	B	126	SER	N-CA-C	5.39	119.94	112.45
1	B	179	MET	N-CA-C	-5.12	105.88	111.82

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	126	SER	Peptide
1	B	126	SER	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	3196	0	3032	49	0
1	B	3196	0	3032	48	0
1	C	3196	0	3032	72	0
2	A	57	0	0	1	0
2	B	59	0	0	2	0
2	C	40	0	0	0	0
All	All	9744	0	9096	167	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 9.

All (167) 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:C:307:ALA:HA	1:C:378:MET:HE3	1.40	1.01
1:C:126:SER:HB3	1:C:128:ASP:H	1.26	0.99
1:A:166:GLU:HG3	1:A:174:LYS:HD3	1.53	0.91
1:B:307:ALA:HA	1:B:378:MET:HE3	1.58	0.86
1:C:173:LEU:O	1:C:175:ALA:N	2.10	0.84
1:A:169:SER:HB2	1:A:174:LYS:H	1.42	0.84
1:C:212:MET:HA	1:C:216:VAL:HG13	1.62	0.82
1:C:163:VAL:HG12	1:C:164:ASP:H	1.50	0.75
1:C:126:SER:CB	1:C:128:ASP:H	1.99	0.75
1:A:32:TYR:CE1	1:A:364:GLU:HG2	2.22	0.74
1:A:163:VAL:HG12	1:A:164:ASP:H	1.55	0.72
1:C:45:LYS:HB3	1:C:50:HIS:CE1	2.24	0.72
1:A:166:GLU:O	1:A:167:ALA:HB3	1.90	0.71
1:C:45:LYS:HB3	1:C:50:HIS:HE1	1.55	0.71
1:C:110:TYR:HE1	1:C:184:HIS:CD2	2.10	0.70
1:C:173:LEU:C	1:C:175:ALA:H	2.00	0.70
1:A:32:TYR:HE1	1:A:364:GLU:HG2	1.54	0.70
1:A:212:MET:HA	1:A:216:VAL:HG13	1.73	0.70
1:C:166:GLU:O	1:C:168:LEU:N	2.25	0.69
1:B:32:TYR:CE1	1:B:364:GLU:HG2	2.26	0.69
1:B:32:TYR:HE1	1:B:364:GLU:HG2	1.60	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:348:GLU:HG3	1:C:349:TRP:H	1.59	0.67
1:B:10:LEU:HD13	1:B:64:LEU:HD11	1.77	0.66
1:C:170:GLU:HG2	1:C:171:ASN:OD1	1.97	0.64
1:A:246:GLU:HG2	1:A:250:LEU:HD22	1.79	0.64
1:C:348:GLU:CG	1:C:349:TRP:H	2.11	0.64
1:C:205:LEU:HD12	1:C:258:LEU:HD21	1.78	0.64
1:C:169:SER:OG	1:C:173:LEU:HA	1.98	0.63
1:C:192:LEU:HD13	1:C:201:VAL:HG21	1.81	0.63
1:C:310:VAL:HG21	1:C:378:MET:HE2	1.81	0.63
1:C:126:SER:HB3	1:C:128:ASP:N	2.06	0.62
1:A:166:GLU:HG3	1:A:174:LYS:CD	2.29	0.62
1:A:175:ALA:HB1	1:A:228:PHE:O	2.00	0.61
1:B:110:TYR:HE1	1:B:184:HIS:CD2	2.18	0.61
1:C:310:VAL:CG2	1:C:378:MET:HE2	2.31	0.61
1:B:348:GLU:HG3	1:B:349:TRP:H	1.66	0.61
1:C:32:TYR:CE1	1:C:364:GLU:HG2	2.35	0.61
1:A:180:ASN:HB2	2:A:447:HOH:O	2.01	0.59
1:A:183:LEU:HD11	1:A:250:LEU:HD23	1.84	0.59
1:B:192:LEU:HD13	1:B:201:VAL:HG21	1.83	0.59
1:A:52:ARG:HD3	1:A:371:CYS:SG	2.42	0.59
1:B:310:VAL:HG21	1:B:378:MET:HE2	1.83	0.59
1:A:166:GLU:O	1:A:167:ALA:CB	2.51	0.59
1:C:32:TYR:HE1	1:C:364:GLU:HG2	1.65	0.59
1:C:173:LEU:C	1:C:175:ALA:N	2.58	0.59
1:B:246:GLU:HG2	1:B:250:LEU:HD22	1.86	0.57
1:A:10:LEU:HD13	1:A:64:LEU:HD11	1.87	0.57
1:B:205:LEU:HD22	1:B:209:LEU:HD22	1.86	0.56
1:A:45:LYS:HB3	1:A:50:HIS:HE1	1.72	0.55
1:B:317:TYR:CG	1:B:326:LEU:HD13	2.43	0.54
1:B:165:ASN:HB3	1:B:175:ALA:O	2.07	0.54
1:C:348:GLU:HG3	1:C:349:TRP:N	2.23	0.54
1:A:45:LYS:HB3	1:A:50:HIS:CE1	2.44	0.53
1:C:205:LEU:HD13	1:C:258:LEU:HD11	1.91	0.53
1:B:45:LYS:HB3	1:B:50:HIS:HE1	1.74	0.53
1:B:154:GLU:HB2	1:B:165:ASN:ND2	2.24	0.53
1:B:326:LEU:HD11	1:B:387:VAL:HG21	1.90	0.53
1:C:122:TYR:CZ	1:C:126:SER:OG	2.60	0.53
1:C:171:ASN:OD1	1:C:171:ASN:N	2.41	0.53
1:B:171:ASN:ND2	1:B:293:ASP:OD2	2.42	0.52
1:B:304:TRP:HB2	1:B:374:HIS:CG	2.45	0.52
1:A:154:GLU:HB2	1:A:165:ASN:ND2	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:310:VAL:CG2	1:B:378:MET:HE2	2.40	0.52
1:A:153:ARG:NH2	1:B:140:GLU:OE1	2.40	0.52
1:C:246:GLU:HG2	1:C:250:LEU:HD22	1.91	0.52
1:C:213:ARG:HG2	1:C:213:ARG:HH11	1.76	0.51
1:A:169:SER:HB2	1:A:174:LYS:N	2.18	0.51
1:C:304:TRP:HB2	1:C:374:HIS:CG	2.44	0.51
1:C:310:VAL:HB	1:C:378:MET:HE2	1.92	0.51
1:B:180:ASN:HB2	2:B:448:HOH:O	2.10	0.51
1:B:352:GLU:OE2	1:B:363:LYS:HE2	2.11	0.51
1:C:212:MET:HA	1:C:216:VAL:CG1	2.39	0.50
1:B:95:MET:HE1	1:B:368:PRO:HG3	1.93	0.50
1:A:47:VAL:HG22	1:A:111:ASN:HD21	1.76	0.50
1:A:10:LEU:CD1	1:A:64:LEU:HD11	2.42	0.50
1:C:143:GLU:OE1	1:C:204:ARG:NE	2.46	0.49
1:B:187:GLU:OE1	1:B:377:ARG:NH1	2.42	0.49
1:B:45:LYS:HB3	1:B:50:HIS:CE1	2.47	0.49
1:C:310:VAL:CB	1:C:378:MET:HE2	2.42	0.49
1:B:105:THR:HB	1:B:158:ARG:HB2	1.94	0.48
1:C:20:LYS:HB3	1:C:20:LYS:NZ	2.28	0.48
1:C:97:ASP:OD1	1:C:97:ASP:C	2.55	0.48
1:B:348:GLU:CG	1:B:349:TRP:H	2.27	0.48
1:A:304:TRP:HB2	1:A:374:HIS:CG	2.49	0.48
1:C:110:TYR:CE1	1:C:184:HIS:CD2	2.98	0.48
1:C:304:TRP:HB2	1:C:374:HIS:CD2	2.49	0.48
1:C:159:GLN:H	1:C:159:GLN:HG2	1.46	0.48
1:B:179:MET:CE	1:B:212:MET:HE1	2.44	0.47
1:B:310:VAL:CB	1:B:378:MET:HE2	2.44	0.47
1:B:178:THR:HB	2:B:448:HOH:O	2.15	0.47
1:A:163:VAL:HG12	1:A:164:ASP:N	2.25	0.47
1:C:245:ILE:O	1:C:248:THR:OG1	2.33	0.47
1:C:247:ALA:HB1	1:C:251:MET:HE2	1.97	0.47
1:C:371:CYS:SG	1:C:373:TYR:HB3	2.55	0.47
1:A:119:LEU:HD13	1:A:134:LEU:HB3	1.97	0.47
1:A:205:LEU:HD13	1:A:258:LEU:HD11	1.97	0.47
1:A:179:MET:HE2	1:A:179:MET:HB2	1.64	0.46
1:A:251:MET:HE3	1:A:271:ASP:OD1	2.14	0.46
1:B:1:MET:HB3	1:B:389:ILE:O	2.15	0.46
1:C:159:GLN:HG3	1:C:161:ARG:HH21	1.80	0.46
1:A:348:GLU:CG	1:A:349:TRP:H	2.27	0.46
1:A:303:TRP:CZ2	1:A:370:LYS:HG3	2.51	0.46
1:B:356:ASP:O	1:B:357:HIS:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:LEU:HA	1:B:275:SER:HB2	1.98	0.45
1:C:348:GLU:CG	1:C:349:TRP:N	2.79	0.45
1:A:6:ILE:CG2	1:A:379:CYS:HA	2.47	0.45
1:B:128:ASP:OD2	1:B:130:GLU:N	2.49	0.45
1:C:165:ASN:HB3	1:C:175:ALA:O	2.17	0.45
1:A:192:LEU:HD13	1:A:201:VAL:HG21	1.99	0.45
1:C:375:ASN:OD1	1:C:376:GLY:N	2.49	0.45
1:C:81:LYS:HG3	1:C:134:LEU:HD21	1.98	0.45
1:B:303:TRP:CG	1:B:304:TRP:N	2.85	0.44
1:C:179:MET:HB2	1:C:179:MET:HE2	1.61	0.44
1:A:159:GLN:OE1	1:A:161:ARG:NH2	2.50	0.44
1:B:310:VAL:HB	1:B:378:MET:HE2	1.99	0.44
1:C:303:TRP:CG	1:C:304:TRP:N	2.86	0.44
1:A:250:LEU:HD12	1:A:250:LEU:HA	1.76	0.44
1:B:110:TYR:CE1	1:B:184:HIS:CD2	3.02	0.44
1:C:33:LEU:HD13	1:C:39:LEU:HD13	1.99	0.44
1:C:388:ASP:O	1:C:389:ILE:HB	2.18	0.44
1:C:301:ARG:HB2	1:C:353:VAL:O	2.18	0.44
1:C:179:MET:CE	1:C:212:MET:HE1	2.48	0.43
1:C:36:GLY:O	1:C:37:LEU:HB2	2.19	0.43
1:C:217:TYR:O	1:C:219:PRO:HD3	2.17	0.43
1:A:358:THR:HA	1:A:359:PRO:HD3	1.75	0.43
1:B:123:TYR:O	1:B:127:GLY:HA2	2.19	0.43
1:C:126:SER:CB	1:C:127:GLY:HA2	2.49	0.43
1:C:174:LYS:HB2	1:C:235:VAL:HG11	2.00	0.43
1:C:179:MET:HE1	1:C:212:MET:HE1	2.01	0.43
1:C:93:TYR:CD1	1:C:105:THR:HG22	2.54	0.43
1:C:112:ILE:HG22	1:C:142:ILE:HD11	2.01	0.42
1:B:156:PHE:CD2	1:B:156:PHE:N	2.87	0.42
1:C:85:ILE:CD1	1:C:134:LEU:HG	2.49	0.42
1:A:110:TYR:HE1	1:A:184:HIS:CD2	2.37	0.42
1:C:337:LYS:O	1:C:343:LYS:NZ	2.53	0.42
1:A:221:THR:HG22	1:A:223:ALA:H	1.84	0.42
1:B:172:GLY:HA3	1:B:173:LEU:HA	1.90	0.42
1:A:123:TYR:O	1:A:127:GLY:HA2	2.19	0.42
1:A:110:TYR:CE1	1:A:184:HIS:CD2	3.08	0.42
1:B:263:LEU:HD23	1:B:263:LEU:HA	1.94	0.42
1:B:250:LEU:HD12	1:B:250:LEU:HA	1.89	0.42
1:B:278:ILE:HD13	1:B:278:ILE:HA	1.82	0.42
1:C:106:MET:SD	1:C:108:HIS:HB2	2.60	0.42
1:A:213:ARG:HG2	1:A:214:ASP:OD1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:371:CYS:HB2	1:A:372:PRO:HD2	2.01	0.41
1:A:345:GLU:HA	1:A:346:GLY:HA2	1.78	0.41
1:C:94:TRP:HZ3	1:C:111:ASN:HB2	1.85	0.41
1:A:348:GLU:OE1	1:A:349:TRP:HB2	2.20	0.41
1:C:136:TYR:OH	1:C:198:ASN:HB2	2.20	0.41
1:A:303:TRP:CG	1:A:304:TRP:N	2.87	0.41
1:C:83:ASN:HB3	1:C:96:MET:HE2	2.03	0.41
1:A:162:LEU:CD2	1:B:137:ARG:HH21	2.33	0.41
1:A:348:GLU:CD	1:A:349:TRP:H	2.28	0.41
1:B:166:GLU:O	1:B:167:ALA:HB3	2.20	0.41
1:B:345:GLU:HA	1:B:346:GLY:HA2	1.89	0.41
1:C:18:TRP:CH2	1:C:372:PRO:HD3	2.56	0.41
1:B:136:TYR:OH	1:B:198:ASN:HB2	2.20	0.40
1:C:303:TRP:CZ2	1:C:370:LYS:HD3	2.56	0.40
1:A:48:ILE:HG13	1:A:111:ASN:OD1	2.21	0.40
1:A:123:TYR:O	1:A:127:GLY:CA	2.70	0.40
1:B:336:ILE:HD13	1:B:378:MET:HE1	2.02	0.40
1:C:18:TRP:CD1	1:C:372:PRO:HB3	2.56	0.40
1:C:18:TRP:NE1	1:C:372:PRO:HB3	2.36	0.40
1:C:345:GLU:HA	1:C:346:GLY:HA2	1.74	0.40
1:B:171:ASN:O	1:B:172:GLY:C	2.63	0.40
1:C:349:TRP:N	1:C:349:TRP:CD1	2.88	0.40
1:A:2:MET:HE2	1:A:2:MET:HB3	1.87	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	387/389 (100%)	371 (96%)	15 (4%)	1 (0%)	36	58
1	B	387/389 (100%)	369 (95%)	17 (4%)	1 (0%)	36	58

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	387/389 (100%)	365 (94%)	17 (4%)	5 (1%)	9	21
All	All	1161/1167 (100%)	1105 (95%)	49 (4%)	7 (1%)	21	42

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	174	LYS
1	B	172	GLY
1	A	167	ALA
1	C	166	GLU
1	C	171	ASN
1	C	196	ASP
1	C	167	ALA

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	333/333 (100%)	292 (88%)	41 (12%)	4	9
1	B	333/333 (100%)	302 (91%)	31 (9%)	8	18
1	C	333/333 (100%)	295 (89%)	38 (11%)	5	11
All	All	999/999 (100%)	889 (89%)	110 (11%)	6	13

All (110) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	SER
1	A	8	GLN
1	A	10	LEU
1	A	25	GLU
1	A	38	GLU
1	A	42	LYS
1	A	47	VAL
1	A	48	ILE

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Mol	Chain	Res	Type
1	A	129	LYS
1	A	134	LEU
1	A	137	ARG
1	A	159	GLN
1	A	162	LEU
1	A	163	VAL
1	A	169	SER
1	A	170	GLU
1	A	173	LEU
1	A	179	MET
1	A	183	LEU
1	A	192	LEU
1	A	194	LYS
1	A	199	GLU
1	A	200	LYS
1	A	205	LEU
1	A	209	LEU
1	A	216	VAL
1	A	221	THR
1	A	250	LEU
1	A	265	LYS
1	A	281	ILE
1	A	288	LEU
1	A	324	LYS
1	A	326	LEU
1	A	327	GLU
1	A	348	GLU
1	A	358	THR
1	A	364	GLU
1	A	371	CYS
1	A	384	THR
1	A	387	VAL
1	A	389	ILE
1	B	10	LEU
1	B	47	VAL
1	B	114	PHE
1	B	134	LEU
1	B	159	GLN
1	B	168	LEU
1	B	170	GLU
1	B	174	LYS
1	B	179	MET

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Mol	Chain	Res	Type
1	B	183	LEU
1	B	192	LEU
1	B	200	LYS
1	B	205	LEU
1	B	209	LEU
1	B	213	ARG
1	B	216	VAL
1	B	230	THR
1	B	250	LEU
1	B	281	ILE
1	B	288	LEU
1	B	296	GLU
1	B	324	LYS
1	B	326	LEU
1	B	331	SER
1	B	348	GLU
1	B	358	THR
1	B	364	GLU
1	B	370	LYS
1	B	378	MET
1	B	384	THR
1	B	387	VAL
1	C	10	LEU
1	C	25	GLU
1	C	47	VAL
1	C	48	ILE
1	C	101	LYS
1	C	126	SER
1	C	134	LEU
1	C	150	TYR
1	C	159	GLN
1	C	163	VAL
1	C	169	SER
1	C	171	ASN
1	C	174	LYS
1	C	179	MET
1	C	183	LEU
1	C	192	LEU
1	C	199	GLU
1	C	205	LEU
1	C	209	LEU
1	C	213	ARG

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Mol	Chain	Res	Type
1	C	216	VAL
1	C	221	THR
1	C	230	THR
1	C	250	LEU
1	C	275	SER
1	C	288	LEU
1	C	303	TRP
1	C	323	GLU
1	C	324	LYS
1	C	326	LEU
1	C	327	GLU
1	C	338	GLU
1	C	348	GLU
1	C	358	THR
1	C	364	GLU
1	C	370	LYS
1	C	378	MET
1	C	387	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	ASN
1	A	276	HIS
1	B	75	HIS
1	B	159	GLN
1	B	222	ASN
1	B	360	HIS
1	C	279	GLN

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 [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 [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	389/389 (100%)	-0.48	1 (0%) 90 88	31, 43, 63, 85	0
1	B	389/389 (100%)	-0.40	3 (0%) 82 80	33, 44, 64, 89	0
1	C	389/389 (100%)	-0.22	3 (0%) 82 80	35, 50, 67, 82	0
All	All	1167/1167 (100%)	-0.37	7 (0%) 85 83	31, 46, 65, 89	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	172	GLY	5.0
1	B	174	LYS	4.5
1	B	167	ALA	3.8
1	C	173	LEU	3.5
1	C	174	LYS	2.7
1	A	174	LYS	2.3
1	C	167	ALA	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.