



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:58 AM UTC

PDB ID : 7BOB / pdb_00007bob
Title : Exo-beta-1,4-mannosidase Op5Man5 from Opitutaceae bacterium strain TAV5
Authors : Kalyani, D.C.; Reichenbach, T.; Keskitalo, M.M.; Conrad, J.; Aspeborg, H.;
Divne, C.
Deposited on : 2021-01-24
Resolution : 2.20 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

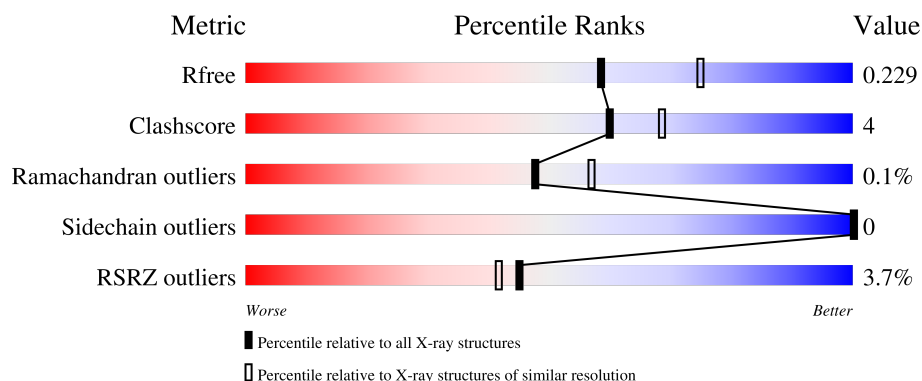
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

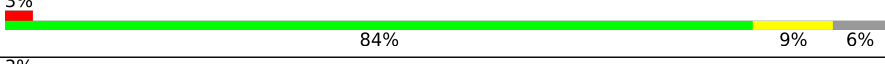
The reported resolution of this entry is 2.20 Å.

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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	601	
1	B	601	
1	C	601	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13498 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 Endo-beta-mannanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	564	Total	C	N	O	S	0	0	0
			4444	2833	785	803	23			
1	B	563	Total	C	N	O	S	0	0	0
			4440	2834	785	798	23			
1	C	560	Total	C	N	O	S	0	1	0
			4425	2824	782	796	23			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	595	ALA	-	expression tag	UNP W0J1H8
A	596	HIS	-	expression tag	UNP W0J1H8
A	597	HIS	-	expression tag	UNP W0J1H8
A	598	HIS	-	expression tag	UNP W0J1H8
A	599	HIS	-	expression tag	UNP W0J1H8
A	600	HIS	-	expression tag	UNP W0J1H8
A	601	HIS	-	expression tag	UNP W0J1H8
B	595	ALA	-	expression tag	UNP W0J1H8
B	596	HIS	-	expression tag	UNP W0J1H8
B	597	HIS	-	expression tag	UNP W0J1H8
B	598	HIS	-	expression tag	UNP W0J1H8
B	599	HIS	-	expression tag	UNP W0J1H8
B	600	HIS	-	expression tag	UNP W0J1H8
B	601	HIS	-	expression tag	UNP W0J1H8
C	595	ALA	-	expression tag	UNP W0J1H8
C	596	HIS	-	expression tag	UNP W0J1H8
C	597	HIS	-	expression tag	UNP W0J1H8
C	598	HIS	-	expression tag	UNP W0J1H8
C	599	HIS	-	expression tag	UNP W0J1H8
C	600	HIS	-	expression tag	UNP W0J1H8
C	601	HIS	-	expression tag	UNP W0J1H8

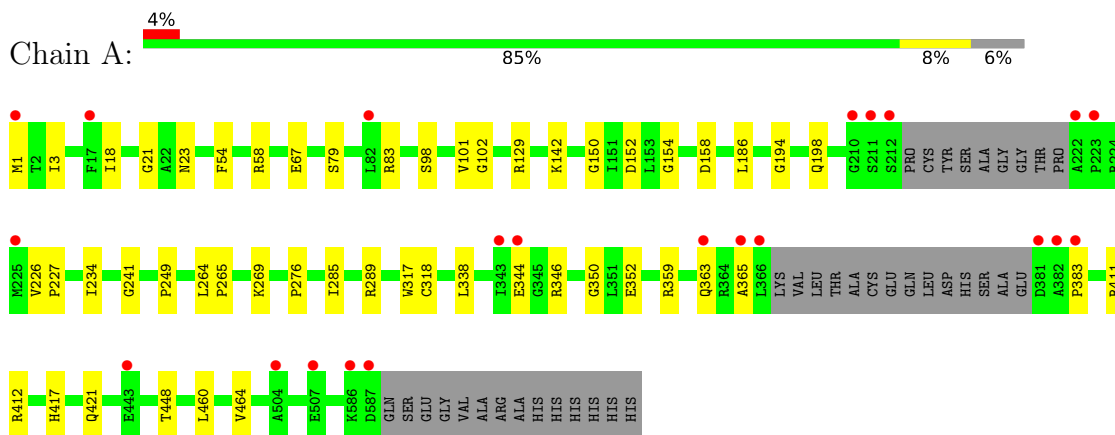
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	73	Total 73	O 73	0	0
2	B	56	Total 56	O 56	0	0
2	C	60	Total 60	O 60	0	0

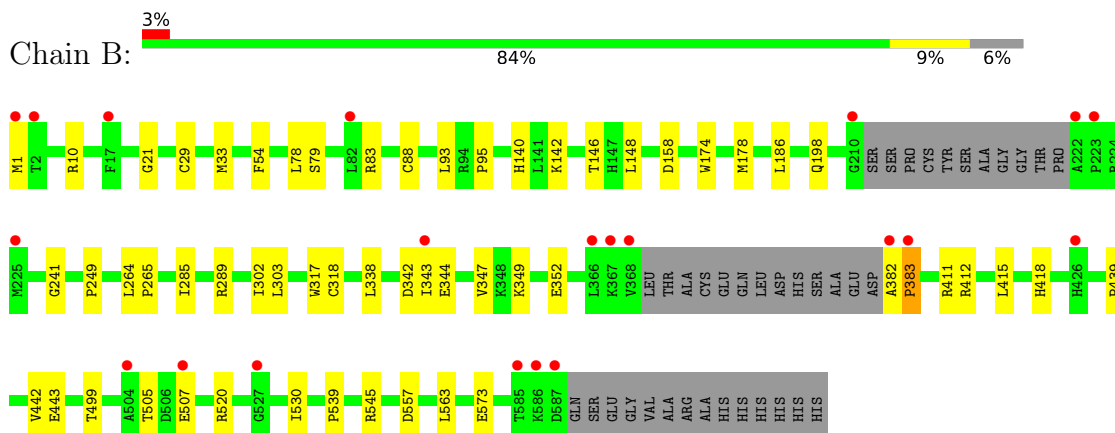
3 Residue-property plots

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.

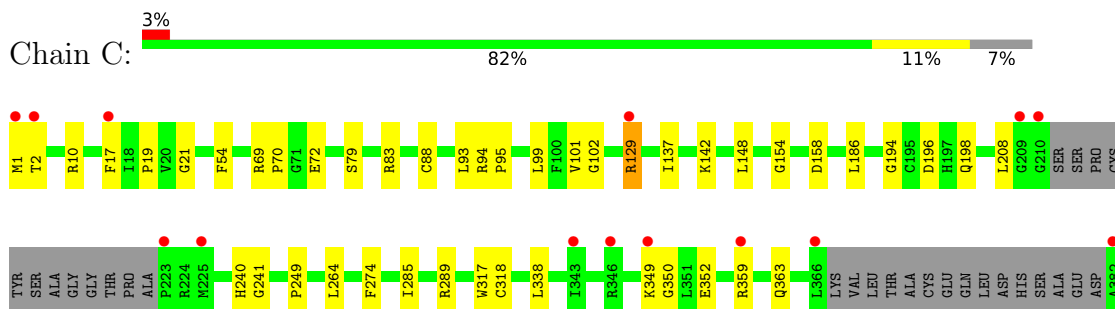
• Molecule 1: Endo-beta-mannanase

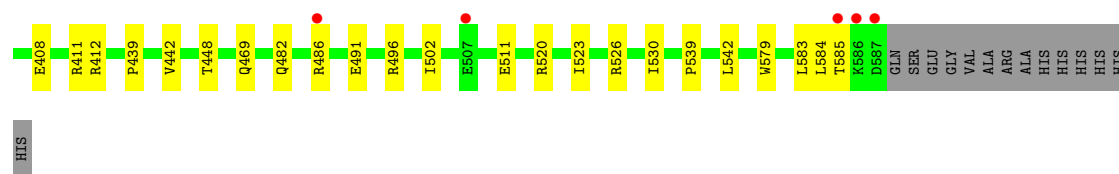


• Molecule 1: Endo-beta-mannanase



• Molecule 1: Endo-beta-mannanase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.14Å 162.45Å 168.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.87 – 2.20 48.87 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.87-2.20) 99.7 (48.87-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.15.2-3472	Depositor
R, R_{free}	0.201 , 0.227 0.204 , 0.229	Depositor DCC
R_{free} test set	2012 reflections (1.42%)	wwPDB-VP
Wilson B-factor (Å ²)	45.2	Xtriage
Anisotropy	0.702	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.005 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13498	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% 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	0.42	0/4579	0.58	0/6250
1	B	0.35	0/4575	0.52	0/6244
1	C	0.36	0/4563	0.55	0/6227
All	All	0.38	0/13717	0.55	0/18721

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	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	129	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	4444	0	4300	31	0
1	B	4440	0	4308	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	4425	0	4290	50	0
2	A	73	0	0	1	0
2	B	56	0	0	0	0
2	C	60	0	0	0	0
All	All	13498	0	12898	113	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 4.

All (113) 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:289:ARG:NH2	1:C:349:LYS:HZ1	1.38	1.20
1:C:289:ARG:HH21	1:C:349:LYS:NZ	1.46	1.11
1:C:289:ARG:NH2	1:C:349:LYS:NZ	2.02	1.00
1:B:505:THR:HG23	1:B:507:GLU:H	1.36	0.88
1:C:2:THR:HG23	1:C:10:ARG:HB2	1.54	0.88
1:A:241:GLY:HA3	1:A:264:LEU:HD13	1.63	0.78
1:A:344:GLU:OE1	1:A:346:ARG:HD3	1.87	0.74
1:B:412:ARG:NH2	1:B:557:ASP:OD2	2.22	0.72
1:A:129:ARG:NH1	2:A:701:HOH:O	2.22	0.72
1:C:289:ARG:CZ	1:C:349:LYS:NZ	2.57	0.68
1:B:343:ILE:H	1:B:343:ILE:HD12	1.57	0.68
1:C:289:ARG:CZ	1:C:349:LYS:HZ2	2.07	0.67
1:C:289:ARG:HE	1:C:349:LYS:HE3	1.60	0.67
1:B:249:PRO:HB2	1:B:411:ARG:HG3	1.77	0.66
1:A:344:GLU:OE1	1:A:346:ARG:CG	2.44	0.66
1:C:1:MET:HE1	1:C:17:PHE:CD2	2.31	0.66
1:C:17:PHE:CZ	1:C:94:ARG:HD2	2.30	0.66
1:C:241:GLY:HA3	1:C:264:LEU:HD13	1.80	0.63
1:C:249:PRO:HB2	1:C:411:ARG:HG3	1.82	0.62
1:A:344:GLU:OE1	1:A:346:ARG:CD	2.48	0.62
1:C:70:PRO:HB3	1:C:129:ARG:HE	1.65	0.61
1:B:241:GLY:HA3	1:B:264:LEU:HD13	1.84	0.59
1:C:289:ARG:NE	1:C:349:LYS:HZ2	2.02	0.57
1:B:499:THR:O	1:B:520:ARG:NH1	2.38	0.56
1:A:352:GLU:H	1:A:352:GLU:CD	2.14	0.56
1:C:289:ARG:HH21	1:C:349:LYS:HZ1	0.66	0.56
1:A:1:MET:HE1	1:A:150:GLY:H	1.72	0.55
1:A:344:GLU:OE1	1:A:346:ARG:HG3	2.07	0.54
1:B:79:SER:O	1:B:83:ARG:HG3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:PRO:HB2	1:A:411:ARG:HG3	1.89	0.54
1:A:23:ASN:OD1	1:A:58:ARG:HD3	2.08	0.54
1:B:289:ARG:HH11	1:B:349:LYS:HD2	1.73	0.54
1:C:289:ARG:NH2	1:C:349:LYS:HZ2	1.97	0.54
1:C:408:GLU:OE1	1:C:411:ARG:NH1	2.41	0.53
1:A:412:ARG:HD3	1:A:448:THR:O	2.09	0.53
1:C:158:ASP:OD2	1:C:198:GLN:HB2	2.09	0.52
1:C:289:ARG:HE	1:C:349:LYS:CE	2.21	0.52
1:B:78:LEU:HD22	1:B:140:HIS:CD2	2.44	0.52
1:A:67:GLU:OE2	1:A:129:ARG:NH1	2.43	0.52
1:A:142:LYS:HG3	1:A:186:LEU:HD21	1.92	0.51
1:C:469:GLN:HB3	1:C:585:THR:OG1	2.11	0.51
1:B:530:ILE:HD13	1:B:539:PRO:HG3	1.93	0.50
1:B:439:PRO:HA	1:B:442:VAL:O	2.13	0.49
1:B:382:ALA:N	1:B:383:PRO:HD3	2.28	0.49
1:C:496:ARG:HG3	1:C:520:ARG:HB3	1.95	0.49
1:A:350:GLY:N	1:A:352:GLU:OE2	2.43	0.49
1:B:21:GLY:HA2	1:B:54:PHE:CG	2.48	0.49
1:C:579:TRP:CZ2	1:C:583:LEU:HD11	2.48	0.49
1:C:289:ARG:NE	1:C:349:LYS:NZ	2.61	0.48
1:B:158:ASP:OD2	1:B:198:GLN:HB2	2.14	0.48
1:C:412:ARG:HD3	1:C:448:THR:O	2.13	0.48
1:B:418:HIS:NE2	1:B:573:GLU:OE2	2.35	0.48
1:B:174:TRP:O	1:B:178:MET:HG2	2.14	0.48
1:C:95:PRO:HG2	1:C:148:LEU:HD23	1.97	0.47
1:B:439:PRO:O	1:B:443:GLU:HG2	2.14	0.47
1:A:285:ILE:HA	1:A:338:LEU:HD22	1.96	0.47
1:C:285:ILE:HA	1:C:338:LEU:HD22	1.97	0.47
1:C:196:ASP:HB2	1:C:240:HIS:HB2	1.97	0.47
1:C:491:GLU:HG3	1:C:526:ARG:HG3	1.97	0.47
1:B:1:MET:HB3	1:B:10:ARG:O	2.15	0.47
1:A:21:GLY:HA2	1:A:54:PHE:CG	2.50	0.46
1:C:439:PRO:HA	1:C:442:VAL:O	2.15	0.46
1:C:530:ILE:HD13	1:C:539:PRO:HG3	1.98	0.46
1:B:1:MET:HE1	1:B:146:THR:HA	1.96	0.45
1:B:142:LYS:HB2	1:B:186:LEU:HD21	1.97	0.45
1:A:158:ASP:OD2	1:A:198:GLN:HB2	2.16	0.45
1:A:154:GLY:O	1:A:194:GLY:HA3	2.17	0.45
1:A:289:ARG:NH1	1:C:526:ARG:HH22	2.15	0.45
1:C:350:GLY:N	1:C:352:GLU:OE1	2.49	0.45
1:C:482:GLN:O	1:C:486:ARG:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:ARG:O	1:A:363:GLN:HG3	2.17	0.44
1:C:154:GLY:O	1:C:194:GLY:HA3	2.17	0.44
1:A:226:VAL:HG13	1:A:234:ILE:O	2.16	0.44
1:B:342:ASP:OD1	1:B:344:GLU:HG2	2.18	0.44
1:A:79:SER:O	1:A:83:ARG:HG3	2.18	0.44
1:C:10:ARG:HG2	1:C:10:ARG:HH11	1.83	0.43
1:B:285:ILE:HA	1:B:338:LEU:HD22	2.00	0.43
1:C:17:PHE:CE2	1:C:19:PRO:HG3	2.52	0.43
1:A:98:SER:HA	1:A:152:ASP:HB3	2.00	0.43
1:A:317:TRP:HA	1:A:318:CYS:HA	1.82	0.43
1:C:21:GLY:HA2	1:C:54:PHE:CG	2.54	0.43
1:C:99:LEU:HD22	1:C:137:ILE:HG21	2.00	0.43
1:C:142:LYS:HG3	1:C:186:LEU:HD21	2.01	0.42
1:C:359:ARG:NH1	1:C:363:GLN:OE1	2.52	0.42
1:A:265:PRO:O	1:A:269:LYS:HG3	2.20	0.42
1:B:264:LEU:HB3	1:B:265:PRO:HD3	2.01	0.42
1:B:29:CYS:SG	1:B:33:MET:HA	2.59	0.42
1:B:302:ILE:HG13	1:B:303:LEU:N	2.34	0.42
1:B:95:PRO:HG2	1:B:148:LEU:HD23	2.01	0.42
1:C:208:LEU:HB3	1:C:274:PHE:CD1	2.55	0.42
1:A:101:VAL:HA	1:A:102:GLY:HA3	1.73	0.42
1:C:584:LEU:HD23	1:C:584:LEU:HA	1.80	0.42
1:B:415:LEU:HD22	1:B:563:LEU:HG	2.01	0.41
1:C:289:ARG:HE	1:C:349:LYS:NZ	2.18	0.41
1:A:18:ILE:HG12	1:A:365:ALA:HB2	2.01	0.41
1:B:545:ARG:HE	1:B:545:ARG:HB3	1.62	0.41
1:C:88:CYS:HB3	1:C:93:LEU:O	2.20	0.41
1:C:317:TRP:HA	1:C:318:CYS:HA	1.82	0.41
1:B:186:LEU:HD23	1:B:186:LEU:HA	1.91	0.41
1:C:101:VAL:HA	1:C:102:GLY:HA3	1.76	0.41
1:B:88:CYS:HB3	1:B:93:LEU:O	2.20	0.41
1:B:289:ARG:NH1	1:B:349:LYS:HD2	2.36	0.41
1:C:69:ARG:NH2	1:C:72:GLU:OE1	2.46	0.41
1:B:317:TRP:HA	1:B:318:CYS:HA	1.66	0.41
1:A:417:HIS:O	1:A:421:GLN:HG3	2.21	0.41
1:A:460:LEU:O	1:A:464:VAL:HG23	2.21	0.41
1:B:347:VAL:HG11	1:B:352:GLU:HA	2.02	0.41
1:C:523:ILE:HD13	1:C:542:LEU:HD11	2.01	0.40
1:C:502:ILE:HG12	1:C:511:GLU:HG3	2.03	0.40
1:C:79:SER:O	1:C:83:ARG:HG3	2.22	0.40
1:A:3:ILE:HG21	1:A:276:PRO:HG3	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:VAL:HG13	1:A:227:PRO:HD2	2.03	0.40
1:C:317:TRP:CD1	1:C:317:TRP:C	3.00	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	558/601 (93%)	540 (97%)	17 (3%)	1 (0%)	43	51
1	B	557/601 (93%)	537 (96%)	19 (3%)	1 (0%)	43	51
1	C	555/601 (92%)	538 (97%)	17 (3%)	0	100	100
All	All	1670/1803 (93%)	1615 (97%)	53 (3%)	2 (0%)	48	57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	383	PRO
1	A	383	PRO

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	463/492 (94%)	463 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	462/492 (94%)	462 (100%)	0	100	100
1	C	461/492 (94%)	461 (100%)	0	100	100
All	All	1386/1476 (94%)	1386 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	140	HIS
1	A	565	GLN
1	B	140	HIS
1	B	199	GLN
1	B	231	GLN
1	B	544	HIS
1	C	251	GLN
1	C	421	GLN
1	C	529	GLN
1	C	570	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	564/601 (93%)	0.27	22 (3%) 43 40	42, 52, 73, 102	0
1	B	563/601 (93%)	0.45	21 (3%) 45 42	43, 58, 82, 110	0
1	C	560/601 (93%)	0.33	19 (3%) 48 45	31, 53, 75, 99	1 (0%)
All	All	1687/1803 (93%)	0.35	62 (3%) 45 42	31, 54, 78, 110	1 (0%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	7.0
1	B	382	ALA	6.7
1	C	382	ALA	6.6
1	A	343	ILE	6.5
1	B	222	ALA	6.3
1	C	366	LEU	5.6
1	C	343	ILE	5.6
1	B	1	MET	5.6
1	A	212	SER	4.8
1	C	1	MET	4.6
1	A	382	ALA	4.5
1	B	223	PRO	4.5
1	A	222	ALA	4.5
1	B	585	THR	4.4
1	C	587	ASP	4.4
1	A	225	MET	4.3
1	C	225	MET	4.3
1	B	343	ILE	4.2
1	B	210	GLY	4.2
1	B	225	MET	4.0
1	A	223	PRO	4.0
1	C	210	GLY	3.8
1	B	383	PRO	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	586	LYS	3.6
1	C	223	PRO	3.6
1	B	587	ASP	3.6
1	B	368	VAL	3.4
1	C	586	LYS	3.4
1	C	17	PHE	3.3
1	C	585	THR	3.3
1	A	381	ASP	3.2
1	A	366	LEU	3.2
1	A	211	SER	3.1
1	A	586	LYS	3.0
1	B	366	LEU	2.8
1	B	82	LEU	2.8
1	A	507	GLU	2.7
1	B	426	HIS	2.6
1	C	507	GLU	2.6
1	C	209	GLY	2.5
1	C	129	ARG	2.5
1	C	2	THR	2.5
1	C	359	ARG	2.5
1	A	344	GLU	2.4
1	A	210	GLY	2.4
1	A	82	LEU	2.3
1	A	587	ASP	2.3
1	B	527	GLY	2.3
1	A	363	GLN	2.2
1	A	443	GLU	2.2
1	A	17	PHE	2.2
1	B	507	GLU	2.2
1	C	349	LYS	2.2
1	A	383	PRO	2.2
1	A	365	ALA	2.2
1	B	504	ALA	2.2
1	C	486	ARG	2.1
1	B	17	PHE	2.1
1	C	346	ARG	2.1
1	A	504	ALA	2.1
1	B	2	THR	2.1
1	B	367	LYS	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.