



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 01:56 AM EDT

PDB ID : 1SQJ / pdb_00001sqj
Title : Crystal Structure Analysis of Oligoxyloglucan reducing-end-specific cellobiohydrolase (OXG-RCBH)
Authors : Yaoi, K.; Kondo, H.; Noro, N.; Suzuki, M.; Tsuda, S.; Mitsuishi, Y.
Deposited on : 2004-03-19
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

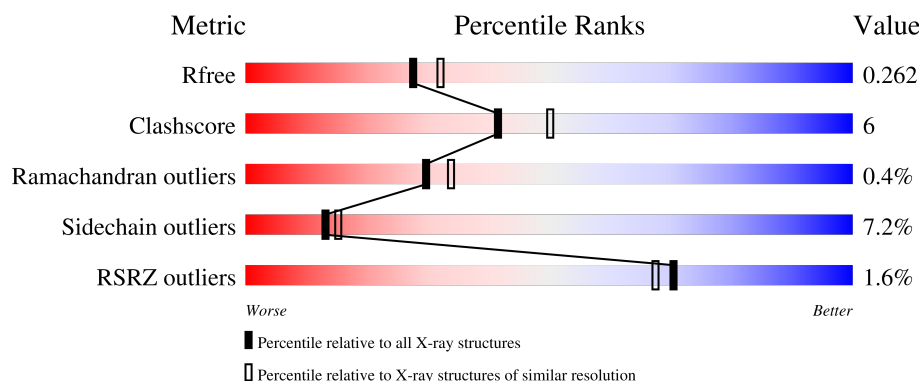
1 Overall quality at a glance

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	789	% 80% 16% ..
1	B	789	2% 82% 14% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11772 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 oligoxyloglucan reducing-end-specific cellobiohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	770	Total	C	N	O	S	0	0	0
			5853	3699	992	1143	19			
1	B	771	Total	C	N	O	S	0	0	0
			5856	3702	994	1141	19			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total	O	0	0
			19	19		
2	B	44	Total	O	0	0
			44	44		

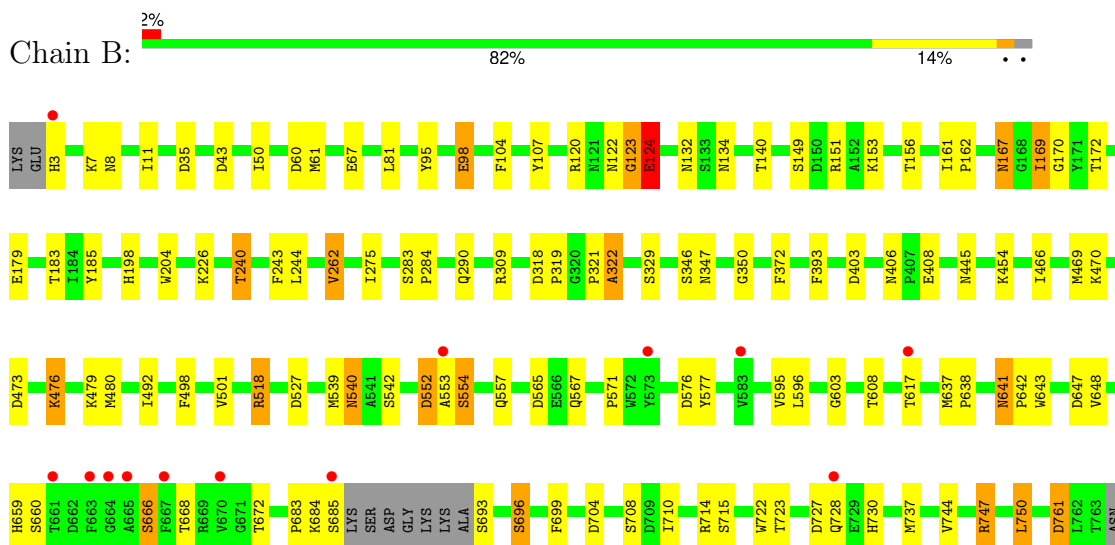
3 Residue-property plots

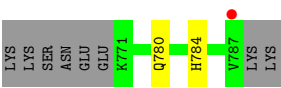
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: oligoxyloglucan reducing-end-specific cellobiohydrolase



- Molecule 1: oligoxyloglucan reducing-end-specific cellobiohydrolase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.08Å 146.05Å 212.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	87.2 (20.00-2.20) 87.1 (20.00-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.217 , 0.251 0.224 , 0.262	Depositor DCC
R_{free} test set	4300 reflections (4.39%)	wwPDB-VP
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.685	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 30.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11772	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% 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	0.75	2/6032 (0.0%)	0.96	3/8248 (0.0%)
1	B	0.75	1/6036 (0.0%)	0.98	4/8254 (0.0%)
All	All	0.75	3/12068 (0.0%)	0.97	7/16502 (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	2
1	B	0	3
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	466	ILE	CA-CB	6.12	1.62	1.54
1	A	411	MET	SD-CE	5.49	1.93	1.79
1	A	166	THR	CA-CB	5.28	1.59	1.53

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	165	PHE	N-CA-C	6.61	124.89	110.80
1	A	748	VAL	CB-CA-C	-6.16	101.45	110.62
1	B	124	GLU	CB-CA-C	5.80	121.96	110.42
1	B	617	THR	N-CA-C	-5.68	106.01	113.17
1	B	393	PHE	N-CA-C	-5.52	105.26	111.28
1	B	169	ILE	N-CA-C	-5.51	106.43	111.67
1	A	105	THR	CB-CA-C	5.02	118.42	109.38

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	122	ASN	Peptide
1	A	321	PRO	Peptide
1	B	122	ASN	Peptide
1	B	123	GLY	Peptide
1	B	321	PRO	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	5853	0	5490	72	0
1	B	5856	0	5498	76	0
2	A	19	0	0	0	0
2	B	44	0	0	0	0
All	All	11772	0	10988	146	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 (146) 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:290:GLN:HE22	1:A:322:ALA:HB3	1.06	1.08
1:B:647:ASP:OD2	1:B:659:HIS:HE1	1.48	0.96
1:A:205:GLU:OE1	1:B:198:HIS:HE1	1.49	0.93
1:A:290:GLN:NE2	1:A:322:ALA:HB3	1.85	0.91
1:A:385:HIS:HD2	1:A:387:THR:H	1.18	0.90
1:B:60:ASP:OD2	1:B:784:HIS:HD2	1.63	0.82
1:A:631:VAL:HG21	1:A:655:GLY:HA3	1.63	0.81
1:A:545:GLN:HE21	1:A:545:GLN:H	1.28	0.80
1:B:167:ASN:HD22	1:B:169:ILE:H	1.30	0.79
1:A:303:ASP:OD1	1:A:305:THR:HG22	1.82	0.79
1:A:543:HIS:HE1	1:A:546:GLY:O	1.66	0.78
1:A:23:HIS:HD2	1:A:25:LYS:H	1.30	0.78
1:A:205:GLU:OE1	1:B:198:HIS:CE1	2.38	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:672:THR:HG23	1:B:704:ASP:O	1.88	0.74
1:A:294:ALA:O	1:A:317:ARG:HD3	1.87	0.73
1:B:262:VAL:HG23	1:B:275:ILE:HG13	1.71	0.71
1:A:146:TRP:NE1	1:A:157:ASN:HD22	1.89	0.71
1:B:747:ARG:HD3	1:B:761:ASP:OD1	1.91	0.70
1:B:167:ASN:ND2	1:B:169:ILE:H	1.90	0.69
1:B:727:ASP:OD1	1:B:730:HIS:HD2	1.77	0.67
1:A:385:HIS:CD2	1:A:387:THR:H	2.07	0.67
1:B:240:THR:HG22	1:B:243:PHE:H	1.59	0.67
1:B:518:ARG:HD3	1:B:539:MET:O	1.96	0.66
1:B:647:ASP:OD2	1:B:659:HIS:CE1	2.40	0.66
1:A:266:ASN:HD22	1:A:268:THR:H	1.44	0.65
1:A:567:GLN:HE22	1:A:572:TRP:HE1	1.44	0.64
1:A:23:HIS:CD2	1:A:25:LYS:H	2.13	0.64
1:A:557:GLN:NE2	1:A:603:GLY:HA2	2.15	0.62
1:A:345:PRO:HG2	1:A:348:LEU:HD11	1.81	0.61
1:B:372:PHE:CD2	1:B:480:MET:HE1	2.35	0.61
1:B:683:PRO:HA	1:B:696:SER:HA	1.83	0.61
1:B:553:ALA:O	1:B:554:SER:HB3	2.01	0.61
1:A:454:LYS:HE3	1:A:642:PRO:O	2.01	0.60
1:B:167:ASN:HD21	1:B:170:GLY:H	1.49	0.60
1:B:132:ASN:C	1:B:132:ASN:HD22	2.11	0.59
1:B:641:ASN:HD22	1:B:643:TRP:H	1.51	0.58
1:A:545:GLN:H	1:A:545:GLN:NE2	1.97	0.58
1:B:553:ALA:O	1:B:554:SER:CB	2.52	0.57
1:A:185:TYR:CE2	1:A:244:LEU:HD21	2.39	0.57
1:B:660:SER:HB2	1:B:666:SER:O	2.05	0.56
1:B:98:GLU:OE1	1:B:151:ARG:NH2	2.38	0.56
1:B:60:ASP:OD2	1:B:784:HIS:CD2	2.51	0.56
1:B:262:VAL:CG2	1:B:275:ILE:HG13	2.36	0.55
1:B:322:ALA:HA	1:B:350:GLY:O	2.06	0.55
1:B:406:ASN:HD21	1:B:408:GLU:HB2	1.71	0.55
1:A:262:VAL:HG21	1:A:312:VAL:HG11	1.88	0.54
1:B:107:TYR:CE2	1:B:153:LYS:HG3	2.42	0.53
1:A:362:LYS:NZ	1:A:422:ASP:OD2	2.40	0.53
1:B:167:ASN:HD22	1:B:167:ASN:C	2.16	0.53
1:B:132:ASN:HD22	1:B:134:ASN:H	1.56	0.53
1:A:567:GLN:NE2	1:A:572:TRP:HE1	2.06	0.53
1:A:717:ASP:OD2	1:A:721:THR:OG1	2.23	0.52
1:A:453:PRO:HG2	1:A:457:ALA:O	2.09	0.52
1:A:543:HIS:CE1	1:A:546:GLY:O	2.55	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:552:ASP:OD1	1:B:553:ALA:O	2.29	0.51
1:B:699:PHE:CE1	1:B:714:ARG:HG3	2.45	0.51
1:A:107:TYR:CZ	1:A:153:LYS:HG2	2.46	0.51
1:B:683:PRO:CA	1:B:696:SER:HA	2.41	0.51
1:A:521:TYR:CE1	1:A:577:TYR:HB3	2.47	0.50
1:B:132:ASN:ND2	1:B:134:ASN:H	2.08	0.50
1:A:557:GLN:HE22	1:A:603:GLY:HA2	1.75	0.50
1:B:498:PHE:CD1	1:B:554:SER:HA	2.46	0.50
1:A:375:GLY:HA3	1:A:485:GLN:NE2	2.26	0.50
1:A:25:LYS:HD2	1:A:73:PRO:O	2.11	0.50
1:A:306:ASN:C	1:A:306:ASN:HD22	2.19	0.50
1:A:322:ALA:HA	1:A:350:GLY:O	2.12	0.50
1:B:557:GLN:OE1	1:B:603:GLY:HA2	2.12	0.50
1:A:631:VAL:CG2	1:A:655:GLY:HA3	2.38	0.49
1:B:183:THR:OG1	1:B:198:HIS:HD2	1.95	0.49
1:A:370:LEU:HD21	1:A:417:THR:HG21	1.95	0.49
1:A:146:TRP:CD1	1:A:157:ASN:HD22	2.30	0.49
1:B:185:TYR:CE2	1:B:244:LEU:HD11	2.47	0.49
1:A:287:TYR:HB3	1:A:322:ALA:HB2	1.95	0.48
1:A:123:GLY:HA3	1:A:236:LYS:HE2	1.96	0.48
1:B:167:ASN:ND2	1:B:170:GLY:H	2.10	0.48
1:B:518:ARG:CD	1:B:539:MET:O	2.62	0.47
1:A:54:ASP:OD2	1:A:755:ARG:NH1	2.47	0.47
1:B:95:TYR:HB3	1:B:104:PHE:CD2	2.49	0.47
1:A:137:TRP:CZ3	1:A:158:VAL:HG21	2.49	0.47
1:B:309:ARG:HA	1:B:329:SER:O	2.15	0.47
1:A:150:ASP:HB2	1:A:153:LYS:HE2	1.96	0.47
1:B:98:GLU:CD	1:B:151:ARG:HH22	2.23	0.47
1:B:283:SER:HA	1:B:284:PRO:C	2.39	0.47
1:B:750:LEU:N	1:B:750:LEU:HD12	2.30	0.46
1:B:124:GLU:OE2	1:B:124:GLU:HA	2.14	0.46
1:B:140:THR:O	1:B:170:GLY:HA3	2.15	0.46
1:A:439:ILE:HG13	1:A:439:ILE:O	2.15	0.46
1:A:699:PHE:CE2	1:A:748:VAL:HG22	2.51	0.46
1:A:287:TYR:CB	1:A:322:ALA:HB2	2.46	0.46
1:A:403:ASP:HB3	1:A:406:ASN:O	2.15	0.46
1:B:699:PHE:CZ	1:B:714:ARG:HG3	2.51	0.46
1:B:43:ASP:HB2	1:B:50:ILE:CD1	2.46	0.45
1:A:266:ASN:ND2	1:A:268:THR:H	2.13	0.45
1:B:715:SER:HB2	1:B:722:TRP:CE3	2.51	0.45
1:B:641:ASN:ND2	1:B:643:TRP:H	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:454:LYS:HE3	1:B:642:PRO:O	2.17	0.45
1:A:294:ALA:O	1:A:317:ARG:CD	2.61	0.44
1:B:498:PHE:HD1	1:B:554:SER:HA	1.82	0.44
1:B:132:ASN:HD21	1:B:134:ASN:HB2	1.82	0.44
1:B:750:LEU:N	1:B:750:LEU:CD1	2.81	0.44
1:A:95:TYR:HB3	1:A:104:PHE:CD1	2.53	0.44
1:B:608:THR:HG23	1:B:638:PRO:HG3	2.00	0.43
1:B:403:ASP:HB3	1:B:406:ASN:O	2.18	0.43
1:B:406:ASN:C	1:B:406:ASN:HD22	2.26	0.43
1:A:124:GLU:HA	1:A:124:GLU:OE1	2.18	0.43
1:B:162:PRO:HD3	1:B:204:TRP:HB2	2.01	0.43
1:A:631:VAL:HG13	1:A:658:PHE:CZ	2.54	0.43
1:A:746:GLY:HA3	1:A:762:LEU:HD22	2.01	0.43
1:A:785:CYS:O	1:A:786:TYR:CB	2.67	0.43
1:A:265:GLN:HG3	1:A:272:TRP:CE2	2.54	0.42
1:A:495:ALA:HB3	1:A:498:PHE:O	2.18	0.42
1:A:298:CYS:O	1:A:299:GLY:C	2.60	0.42
1:B:123:GLY:HA2	1:B:172:THR:O	2.19	0.42
1:B:346:SER:CB	1:B:347:ASN:HD22	2.33	0.42
1:B:518:ARG:HD3	1:B:539:MET:C	2.44	0.42
1:B:473:ASP:OD2	1:B:476:LYS:HG3	2.18	0.42
1:A:713:TYR:CE2	1:A:724:ARG:HD3	2.54	0.42
1:B:290:GLN:OE1	1:B:322:ALA:HB3	2.20	0.42
1:B:318:ASP:OD2	1:B:319:PRO:HA	2.20	0.42
1:A:249:ALA:HA	1:A:259:PHE:O	2.20	0.42
1:B:81:LEU:HB2	1:B:95:TYR:HB2	2.01	0.42
1:B:470:LYS:NZ	1:B:527:ASP:OD2	2.44	0.42
1:A:306:ASN:HD21	1:A:308:ASN:HB2	1.84	0.42
1:B:11:ILE:HG21	1:B:445:ASN:ND2	2.35	0.42
1:A:387:THR:O	1:A:388:PRO:C	2.62	0.42
1:B:43:ASP:HB2	1:B:50:ILE:HD12	2.02	0.42
1:A:98:GLU:OE1	1:A:151:ARG:NH2	2.50	0.41
1:A:371:ASP:OD1	1:A:373:ASN:N	2.45	0.41
1:A:743:LYS:HE2	1:A:743:LYS:HB2	1.91	0.41
1:B:540:ASN:HD22	1:B:542:SER:H	1.68	0.41
1:B:637:MET:HA	1:B:638:PRO:HD3	1.89	0.41
1:A:66:THR:HG23	1:A:83:GLN:NE2	2.36	0.41
1:A:601:GLN:HB2	1:A:645:ALA:HB1	2.02	0.41
1:A:483:ALA:HA	1:A:484:PRO:C	2.45	0.41
1:A:412:TYR:CZ	1:A:419:TRP:HB2	2.54	0.41
1:B:161:ILE:HA	1:B:162:PRO:HD3	1.97	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:571:PRO:HG2	1:B:595:VAL:HG11	2.02	0.41
1:A:191:PRO:HD3	1:A:233:GLN:OE1	2.20	0.41
1:A:771:LYS:HB3	1:A:771:LYS:HE3	1.86	0.41
1:A:146:TRP:NE1	1:A:157:ASN:ND2	2.64	0.41
1:B:123:GLY:CA	1:B:172:THR:O	2.69	0.41
1:B:576:ASP:O	1:B:577:TYR:C	2.63	0.41
1:A:40:TYR:CE1	1:A:442:ILE:HA	2.56	0.41
1:B:35:ASP:HB3	1:B:120:ARG:HD2	2.04	0.40
1:B:737:MET:HE3	1:B:737:MET:HB3	1.94	0.40
1:A:454:LYS:HD2	1:A:599:LYS:HG2	2.03	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	764/789 (97%)	720 (94%)	40 (5%)	4 (0%)	24	27
1	B	765/789 (97%)	727 (95%)	36 (5%)	2 (0%)	36	42
All	All	1529/1578 (97%)	1447 (95%)	76 (5%)	6 (0%)	30	34

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	PHE
1	A	322	ALA
1	B	322	ALA
1	A	124	GLU
1	A	289	ASN
1	B	124	GLU

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	608/625 (97%)	563 (93%)	45 (7%)	13	14
1	B	609/625 (97%)	566 (93%)	43 (7%)	13	16
All	All	1217/1250 (97%)	1129 (93%)	88 (7%)	13	15

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LEU
1	A	30	LEU
1	A	67	GLU
1	A	81	LEU
1	A	98	GLU
1	A	105	THR
1	A	124	GLU
1	A	161	ILE
1	A	163	ASP
1	A	167	ASN
1	A	226	LYS
1	A	244	LEU
1	A	265	GLN
1	A	266	ASN
1	A	312	VAL
1	A	315	LEU
1	A	317	ARG
1	A	328	LEU
1	A	344	SER
1	A	390	LEU
1	A	401	LEU
1	A	469	MET
1	A	507	SER
1	A	513	ASP
1	A	540	ASN
1	A	545	GLN
1	A	561	SER

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Mol	Chain	Res	Type
1	A	566	GLU
1	A	589	LYS
1	A	599	LYS
1	A	621	LYS
1	A	648	VAL
1	A	666	SER
1	A	668	THR
1	A	670	VAL
1	A	705	LYS
1	A	728	GLN
1	A	748	VAL
1	A	750	LEU
1	A	755	ARG
1	A	762	LEU
1	A	763	THR
1	A	764	ASN
1	A	770	GLU
1	A	781	LYS
1	B	3	HIS
1	B	7	LYS
1	B	8	ASN
1	B	61	MET
1	B	67	GLU
1	B	98	GLU
1	B	124	GLU
1	B	149	SER
1	B	156	THR
1	B	167	ASN
1	B	179	GLU
1	B	226	LYS
1	B	240	THR
1	B	262	VAL
1	B	469	MET
1	B	476	LYS
1	B	479	LYS
1	B	492	ILE
1	B	501	VAL
1	B	518	ARG
1	B	540	ASN
1	B	552	ASP
1	B	554	SER
1	B	565	ASP

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Mol	Chain	Res	Type
1	B	567	GLN
1	B	596	LEU
1	B	641	ASN
1	B	648	VAL
1	B	666	SER
1	B	668	THR
1	B	684	LYS
1	B	685	SER
1	B	693	SER
1	B	696	SER
1	B	708	SER
1	B	710	ILE
1	B	723	THR
1	B	728	GLN
1	B	744	VAL
1	B	747	ARG
1	B	750	LEU
1	B	761	ASP
1	B	780	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	HIS
1	A	83	GLN
1	A	157	ASN
1	A	242	ASN
1	A	265	GLN
1	A	266	ASN
1	A	290	GLN
1	A	306	ASN
1	A	308	ASN
1	A	373	ASN
1	A	385	HIS
1	A	485	GLN
1	A	510	HIS
1	A	540	ASN
1	A	543	HIS
1	A	545	GLN
1	A	557	GLN
1	A	567	GLN
1	A	601	GLN

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Mol	Chain	Res	Type
1	A	728	GLN
1	A	764	ASN
1	A	784	HIS
1	B	59	GLN
1	B	132	ASN
1	B	167	ASN
1	B	198	HIS
1	B	242	ASN
1	B	308	ASN
1	B	347	ASN
1	B	357	ASN
1	B	377	GLN
1	B	406	ASN
1	B	438	GLN
1	B	445	ASN
1	B	497	ASN
1	B	540	ASN
1	B	641	ASN
1	B	659	HIS
1	B	674	ASN
1	B	718	ASN
1	B	730	HIS
1	B	784	HIS

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	770/789 (97%)	0.05	11 (1%) 73 71	25, 36, 48, 63	0
1	B	771/789 (97%)	0.11	14 (1%) 67 64	23, 34, 55, 65	0
All	All	1541/1578 (97%)	0.08	25 (1%) 70 67	23, 35, 53, 65	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	786	TYR	4.9
1	A	165	PHE	3.6
1	B	665	ALA	3.6
1	A	164	ALA	3.5
1	A	764	ASN	3.4
1	B	583	VAL	3.3
1	B	787	VAL	3.3
1	B	685	SER	3.1
1	B	3	HIS	2.9
1	A	694	ALA	2.8
1	A	162	PRO	2.7
1	A	166	THR	2.7
1	B	573	TYR	2.6
1	B	664	GLY	2.5
1	B	667	PHE	2.5
1	A	565	ASP	2.4
1	B	663	PHE	2.3
1	B	670	VAL	2.3
1	A	781	LYS	2.3
1	B	728	GLN	2.3
1	A	123	GLY	2.2
1	A	513	ASP	2.1
1	B	617	THR	2.1
1	B	661	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	553	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.