



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:29 PM UTC

PDB ID : 5GLV / pdb_00005glv
Title : Tl-gal
Authors : Jang, S.B.; Hwang, E.Y.
Deposited on : 2016-07-12
Resolution : 1.80 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

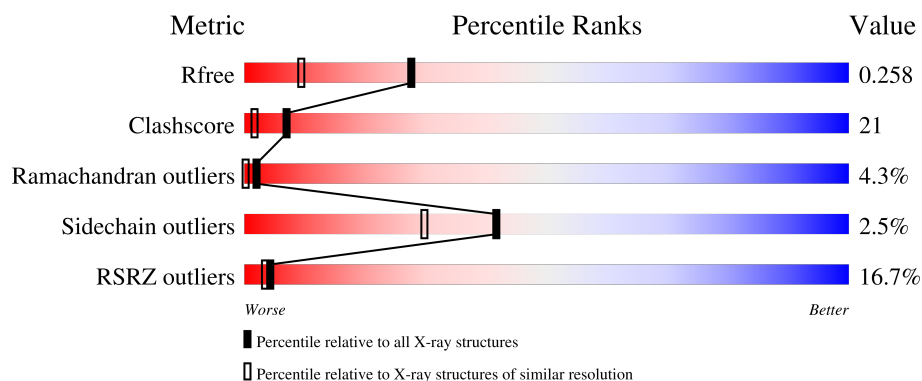
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	284	17% 65% 27% 6%
1	B	284	15% 64% 29% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4971 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.

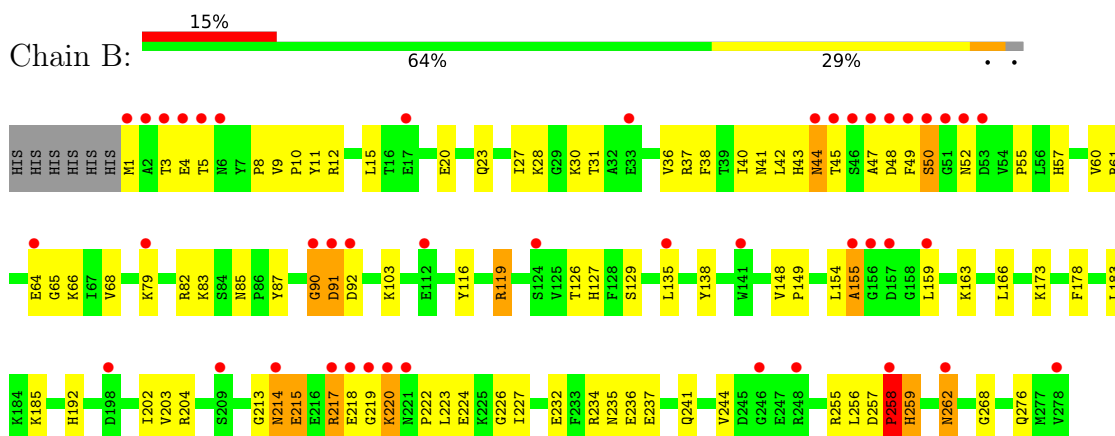
- Molecule 1 is a protein called galectin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	0	0
			2226	1414	384	425	3			
1	B	278	Total	C	N	O	S	0	0	0
			2226	1414	384	425	3			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	261	Total	O	0	0
			261	261		
2	B	258	Total	O	0	0
			258	258		

- Molecule 1: galectin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.42Å 84.00Å 78.36Å 90.00° 109.67° 90.00°	Depositor
Resolution (Å)	30.00 – 1.80 30.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	93.9 (30.00-1.80) 94.5 (30.00-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 1.80Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.258 0.223 , 0.258	Depositor DCC
R_{free} test set	3356 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	17.4	Xtriage
Anisotropy	0.567	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4971	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.99% 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	0.34	0/2281	0.88	6/3079 (0.2%)
1	B	0.34	0/2281	0.88	5/3079 (0.2%)
All	All	0.34	0/4562	0.88	11/6158 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	90	GLY	N-CA-C	-8.38	99.60	112.84
1	A	43	HIS	N-CA-C	6.35	118.60	109.14
1	B	237	GLU	N-CA-C	6.10	118.43	111.11
1	B	183	LEU	N-CA-C	6.02	119.52	109.95
1	A	221	ASN	CA-C-N	5.77	125.87	119.87
1	A	221	ASN	C-N-CA	5.77	125.87	119.87
1	A	42	LEU	N-CA-C	-5.46	99.62	108.52
1	B	43	HIS	N-CA-C	5.46	117.28	109.14
1	A	183	LEU	N-CA-C	5.39	118.06	109.81
1	B	258	PRO	N-CA-C	-5.37	101.41	112.47
1	A	255	ARG	N-CA-C	-5.32	107.16	113.97

There are no chirality outliers.

There are no planarity outliers.

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	2226	0	2167	93	0
1	B	2226	0	2167	92	0
2	A	261	0	0	9	0
2	B	258	0	0	11	0
All	All	4971	0	4334	185	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 21.

All (185) 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:217:ARG:HH11	1:A:217:ARG:HB2	1.02	1.13
1:A:217:ARG:HB2	1:A:217:ARG:NH1	1.71	1.05
1:A:45:THR:HG22	1:A:46:SER:H	1.25	0.99
1:A:38:PHE:CZ	1:A:60:VAL:HG11	2.15	0.82
1:A:45:THR:HG22	1:A:46:SER:N	1.95	0.80
1:B:38:PHE:CZ	1:B:60:VAL:HG11	2.17	0.80
1:A:224:GLU:O	1:A:227:ILE:HG22	1.80	0.80
1:A:38:PHE:CE1	1:A:60:VAL:HG11	2.17	0.79
1:B:224:GLU:O	1:B:227:ILE:HG22	1.81	0.79
1:B:173:LYS:HD3	1:B:226:GLY:HA3	1.67	0.77
1:A:217:ARG:HH11	1:A:217:ARG:CB	1.92	0.76
1:B:224:GLU:HB3	1:B:227:ILE:HG21	1.69	0.74
1:A:119:ARG:HD3	2:A:394:HOH:O	1.86	0.74
1:B:38:PHE:CE1	1:B:60:VAL:HG11	2.22	0.72
1:B:214:ASN:O	1:B:215:GLU:HB2	1.89	0.72
1:B:217:ARG:HH11	1:B:218:GLU:N	1.88	0.72
1:B:148:VAL:HG23	2:B:446:HOH:O	1.91	0.71
1:A:173:LYS:HD3	1:A:226:GLY:HA3	1.72	0.70
1:A:64:GLU:HG2	1:A:66:LYS:HD3	1.75	0.69
1:A:71:THR:HG23	2:A:494:HOH:O	1.92	0.69
1:B:224:GLU:HB3	1:B:227:ILE:CG2	2.23	0.68
1:B:217:ARG:HH11	1:B:218:GLU:H	1.40	0.67
1:A:224:GLU:HB3	1:A:227:ILE:HG21	1.76	0.67
1:A:45:THR:CG2	1:A:46:SER:H	2.02	0.66
1:B:119:ARG:HD3	2:B:395:HOH:O	1.96	0.66
1:A:44:ASN:HD22	1:A:52:ASN:HB3	1.60	0.64
1:B:220:LYS:HB3	2:B:515:HOH:O	1.96	0.64
1:A:79:LYS:HG3	2:A:447:HOH:O	1.96	0.64
1:B:60:VAL:O	1:B:60:VAL:HG13	1.97	0.64
1:A:79:LYS:HD2	1:A:79:LYS:O	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:VAL:HA	1:A:149:PRO:C	2.22	0.63
1:B:45:THR:HA	1:B:49:PHE:CE1	2.34	0.63
1:B:79:LYS:O	1:B:79:LYS:HD2	1.97	0.63
1:A:224:GLU:HB3	1:A:227:ILE:CG2	2.29	0.62
1:A:202:ILE:HD11	1:A:223:LEU:HD23	1.81	0.62
1:B:148:VAL:HA	1:B:149:PRO:C	2.24	0.61
1:A:41:ASN:ND2	1:A:57:HIS:ND1	2.46	0.61
1:A:235:ASN:C	1:A:235:ASN:HD22	2.07	0.61
1:B:218:GLU:HA	2:B:449:HOH:O	2.00	0.61
1:A:4:GLU:HG3	1:A:5:THR:H	1.66	0.60
1:B:217:ARG:O	1:B:218:GLU:HB2	2.01	0.60
1:B:64:GLU:OE2	1:B:82:ARG:NH1	2.35	0.59
1:B:217:ARG:HH11	1:B:217:ARG:HB2	1.66	0.59
1:B:68:VAL:HG22	1:B:82:ARG:NH1	2.18	0.59
1:B:103:LYS:HB2	1:B:116:TYR:O	2.02	0.58
1:B:45:THR:HA	1:B:49:PHE:CZ	2.38	0.58
1:A:223:LEU:HD22	1:A:223:LEU:N	2.19	0.58
1:B:41:ASN:ND2	1:B:57:HIS:ND1	2.48	0.58
1:A:184:LYS:HD2	1:A:256:LEU:HD23	1.85	0.58
1:B:37:ARG:HB3	1:B:61:ARG:HA	1.86	0.57
1:B:90:GLY:O	1:B:91:ASP:HB2	2.04	0.57
1:B:257:ASP:C	1:B:258:PRO:O	2.47	0.57
1:B:223:LEU:HD22	1:B:223:LEU:N	2.20	0.57
1:B:185:LYS:HE2	1:B:262:ASN:OD1	2.05	0.56
1:A:164:SER:OG	1:A:232:GLU:HG3	2.05	0.56
1:A:262:ASN:HD22	1:A:262:ASN:H	1.53	0.56
1:A:38:PHE:CE2	1:A:60:VAL:HG11	2.41	0.56
1:A:214:ASN:O	1:A:215:GLU:CB	2.53	0.56
1:B:38:PHE:CE2	1:B:60:VAL:HG11	2.40	0.56
1:B:27:ILE:HD12	1:B:40:ILE:HD13	1.88	0.55
1:B:262:ASN:H	1:B:262:ASN:HD22	1.54	0.55
1:B:222:PRO:HG2	1:B:244:VAL:HG11	1.89	0.55
1:A:15:LEU:HD12	1:A:15:LEU:N	2.22	0.55
1:A:232:GLU:CD	1:A:234:ARG:HE	2.15	0.55
1:A:44:ASN:HD22	1:A:52:ASN:CB	2.20	0.54
1:B:37:ARG:HD2	1:B:61:ARG:HG2	1.90	0.54
1:B:217:ARG:CZ	1:B:218:GLU:HG2	2.38	0.54
1:A:44:ASN:O	1:A:45:THR:CB	2.56	0.54
1:B:31:THR:OG1	1:B:90:GLY:O	2.26	0.53
1:B:85:ASN:ND2	1:B:87:TYR:H	2.06	0.53
1:A:64:GLU:O	1:A:66:LYS:HD2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:GLY:O	1:B:214:ASN:O	2.26	0.53
1:A:222:PRO:HG2	1:A:244:VAL:HG11	1.90	0.53
1:B:202:ILE:HD11	1:B:223:LEU:HD23	1.91	0.52
1:A:60:VAL:O	1:A:60:VAL:HG13	2.08	0.52
1:A:4:GLU:HG3	1:A:5:THR:N	2.25	0.52
1:A:234:ARG:HB2	1:A:241:GLN:HB2	1.90	0.52
1:A:38:PHE:CD1	1:A:60:VAL:HG11	2.45	0.52
1:A:65:GLY:C	1:A:66:LYS:HD2	2.35	0.51
1:B:37:ARG:HB2	1:B:60:VAL:O	2.10	0.51
1:B:15:LEU:HD13	1:B:126:THR:C	2.35	0.51
1:A:88:LYS:HE2	2:A:497:HOH:O	2.11	0.51
1:A:12:ARG:HD2	1:A:129:SER:HB3	1.93	0.50
1:A:126:THR:OG1	1:A:127:HIS:HD2	1.93	0.50
1:A:103:LYS:HB2	1:A:116:TYR:O	2.12	0.50
1:A:220:LYS:O	1:A:221:ASN:HB3	2.12	0.50
1:B:255:ARG:C	1:B:256:LEU:HD12	2.36	0.50
1:B:154:LEU:O	1:B:155:ALA:C	2.55	0.50
1:A:48:ASP:HA	2:A:479:HOH:O	2.12	0.50
1:B:276:GLN:HG3	2:B:478:HOH:O	2.10	0.49
1:A:201:ALA:CA	1:A:221:ASN:HD22	2.26	0.49
1:A:257:ASP:C	1:A:258:PRO:O	2.54	0.49
1:A:9:VAL:HA	1:A:10:PRO:C	2.38	0.49
1:B:15:LEU:N	1:B:15:LEU:HD12	2.28	0.49
1:B:49:PHE:HE1	1:B:52:ASN:HB2	1.76	0.49
1:B:217:ARG:HD3	1:B:218:GLU:HG2	1.95	0.49
1:B:258:PRO:O	1:B:259:HIS:O	2.31	0.49
1:A:7:TYR:HB2	1:A:136:ILE:HB	1.95	0.48
1:B:36:VAL:O	1:B:37:ARG:HB3	2.13	0.48
1:A:198:ASP:OD1	1:A:199:GLU:HG3	2.13	0.48
1:B:49:PHE:CD1	1:B:50:SER:N	2.81	0.48
1:A:136:ILE:HG21	1:A:139:ILE:HD11	1.95	0.48
1:B:30:LYS:HG2	1:B:92:ASP:OD2	2.14	0.48
1:B:257:ASP:OD2	1:B:258:PRO:O	2.33	0.47
1:B:217:ARG:HB2	1:B:217:ARG:NH1	2.29	0.47
1:A:71:THR:CG2	1:A:119:ARG:HE	2.27	0.47
1:A:19:PHE:HD1	1:A:23:GLN:HE21	1.63	0.47
1:A:66:LYS:HB3	1:A:66:LYS:HE3	1.69	0.47
1:B:12:ARG:HD2	1:B:129:SER:HB3	1.97	0.47
1:A:148:VAL:CG2	2:A:448:HOH:O	2.63	0.47
1:A:154:LEU:O	1:A:155:ALA:C	2.58	0.47
1:A:173:LYS:HD2	2:A:525:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:VAL:CG2	1:B:82:ARG:NH1	2.78	0.47
1:A:45:THR:HA	1:A:52:ASN:ND2	2.29	0.47
1:B:11:TYR:O	1:B:129:SER:HA	2.15	0.47
1:B:203:VAL:HG13	1:B:215:GLU:HG2	1.97	0.47
1:B:44:ASN:HB3	1:B:52:ASN:HB3	1.97	0.46
1:A:184:LYS:HE3	1:A:256:LEU:HB3	1.96	0.46
1:B:219:GLY:O	1:B:220:LYS:CB	2.63	0.46
1:A:30:LYS:CE	1:A:135:LEU:HD23	2.46	0.46
1:A:236:GLU:CD	1:A:241:GLN:HG3	2.40	0.46
1:A:148:VAL:HG23	2:A:448:HOH:O	2.15	0.46
1:A:202:ILE:HD11	1:A:223:LEU:CD2	2.45	0.46
1:A:82:ARG:O	1:A:83:LYS:HD2	2.15	0.46
1:B:126:THR:OG1	1:B:127:HIS:HD2	1.99	0.46
1:B:256:LEU:HD12	1:B:256:LEU:N	2.31	0.46
1:A:262:ASN:H	1:A:262:ASN:ND2	2.12	0.46
1:B:217:ARG:HD3	1:B:217:ARG:C	2.41	0.46
1:B:42:LEU:O	1:B:55:PRO:HD2	2.16	0.46
1:B:257:ASP:O	1:B:258:PRO:O	2.33	0.45
1:A:159:LEU:HD22	1:A:235:ASN:HB2	1.98	0.45
1:A:262:ASN:HD22	1:A:262:ASN:N	2.10	0.45
1:B:9:VAL:HA	1:B:10:PRO:C	2.42	0.45
1:A:68:VAL:CG2	1:A:82:ARG:HH11	2.30	0.45
1:B:227:ILE:HG23	2:B:430:HOH:O	2.17	0.45
1:B:38:PHE:CD1	1:B:60:VAL:HG11	2.51	0.44
1:A:160:ALA:O	1:A:163:LYS:HB2	2.18	0.44
1:B:1:MET:SD	1:B:1:MET:C	3.01	0.44
1:A:85:ASN:ND2	1:A:87:TYR:H	2.15	0.44
1:B:3:THR:O	1:B:5:THR:HG23	2.17	0.44
1:B:20:GLU:H	1:B:23:GLN:NE2	2.15	0.44
1:A:20:GLU:H	1:A:23:GLN:NE2	2.15	0.44
1:A:257:ASP:O	1:A:258:PRO:O	2.35	0.44
1:A:223:LEU:HD22	1:A:223:LEU:H	1.83	0.44
1:B:217:ARG:NH1	1:B:218:GLU:H	2.11	0.44
1:B:30:LYS:HE3	1:B:135:LEU:HD23	2.00	0.43
1:B:148:VAL:CG2	2:B:500:HOH:O	2.66	0.43
1:A:178:PHE:HA	1:A:268:GLY:HA3	2.00	0.43
1:A:214:ASN:O	1:A:215:GLU:HB2	2.17	0.43
1:B:192:HIS:O	1:B:204:ARG:HA	2.19	0.43
1:B:163:LYS:HE3	2:B:326:HOH:O	2.19	0.43
1:A:71:THR:HG21	1:A:119:ARG:HE	1.83	0.43
1:B:202:ILE:HD11	1:B:223:LEU:CD2	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:PRO:O	1:A:259:HIS:O	2.37	0.42
1:B:8:PRO:HG2	2:B:353:HOH:O	2.17	0.42
1:B:103:LYS:HB2	1:B:116:TYR:C	2.43	0.42
1:A:235:ASN:C	1:A:235:ASN:ND2	2.73	0.42
1:A:217:ARG:O	1:A:218:GLU:O	2.37	0.42
1:A:15:LEU:N	1:A:15:LEU:CD1	2.82	0.42
1:A:218:GLU:HA	2:A:380:HOH:O	2.18	0.42
1:B:15:LEU:HD13	1:B:126:THR:O	2.19	0.42
1:B:235:ASN:HD22	1:B:236:GLU:N	2.16	0.42
1:B:44:ASN:O	1:B:126:THR:OG1	2.37	0.42
1:B:217:ARG:NH1	1:B:218:GLU:N	2.63	0.42
1:A:217:ARG:NH1	1:A:217:ARG:CB	2.63	0.42
1:B:232:GLU:OE2	1:B:234:ARG:NE	2.50	0.42
1:A:220:LYS:O	1:A:221:ASN:CB	2.68	0.42
1:B:227:ILE:HD13	2:B:496:HOH:O	2.19	0.42
1:B:262:ASN:HD22	1:B:262:ASN:N	2.14	0.42
1:B:79:LYS:HG3	2:B:427:HOH:O	2.20	0.42
1:B:234:ARG:HB2	1:B:241:GLN:HB2	2.00	0.42
1:A:192:HIS:O	1:A:204:ARG:HA	2.20	0.41
1:A:256:LEU:HD12	1:A:256:LEU:N	2.35	0.41
1:B:65:GLY:O	1:B:66:LYS:HD2	2.20	0.41
1:A:185:LYS:HG2	1:A:262:ASN:OD1	2.20	0.41
1:B:178:PHE:HA	1:B:268:GLY:HA3	2.02	0.41
1:B:28:LYS:HB2	1:B:138:TYR:HB3	2.03	0.41
1:A:11:TYR:C	1:A:12:ARG:HG2	2.46	0.41
1:A:44:ASN:O	1:A:45:THR:OG1	2.38	0.41
1:A:38:PHE:CD1	1:A:60:VAL:CG1	3.04	0.41
1:A:45:THR:HA	1:A:52:ASN:HD22	1.85	0.41
1:B:38:PHE:O	1:B:60:VAL:HG12	2.21	0.41
1:A:4:GLU:O	1:A:6:ASN:N	2.54	0.40
1:A:121:PRO:O	1:A:124:SER:HB3	2.21	0.40
1:B:82:ARG:O	1:B:83:LYS:HD2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	276/284 (97%)	255 (92%)	9 (3%)	12 (4%)	2	0
1	B	276/284 (97%)	252 (91%)	12 (4%)	12 (4%)	2	0
All	All	552/568 (97%)	507 (92%)	21 (4%)	24 (4%)	2	0

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	45	THR
1	B	155	ALA
1	B	214	ASN
1	B	215	GLU
1	B	220	LYS
1	A	5	THR
1	A	155	ALA
1	A	215	GLU
1	A	218	GLU
1	A	259	HIS
1	B	44	ASN
1	B	259	HIS
1	A	2	ALA
1	A	258	PRO
1	B	48	ASP
1	B	50	SER
1	B	258	PRO
1	A	48	ASP
1	A	50	SER
1	B	4	GLU
1	B	47	ALA
1	B	91	ASP
1	A	220	LYS
1	A	51	GLY

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	240/246 (98%)	233 (97%)	7 (3%)	37	25
1	B	240/246 (98%)	235 (98%)	5 (2%)	47	36
All	All	480/492 (98%)	468 (98%)	12 (2%)	42	30

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	LYS
1	A	71	THR
1	A	159	LEU
1	A	166	LEU
1	A	217	ARG
1	A	235	ASN
1	A	262	ASN
1	B	119	ARG
1	B	159	LEU
1	B	166	LEU
1	B	217	ARG
1	B	262	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	41	ASN
1	A	44	ASN
1	A	52	ASN
1	A	85	ASN
1	A	127	HIS
1	A	214	ASN
1	A	221	ASN
1	A	235	ASN
1	A	276	GLN
1	B	23	GLN
1	B	41	ASN
1	B	43	HIS
1	B	85	ASN
1	B	127	HIS

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Mol	Chain	Res	Type
1	B	235	ASN
1	B	276	GLN

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	278/284 (97%)	0.91	49 (17%) 4 3	11, 21, 46, 76	0
1	B	278/284 (97%)	0.89	44 (15%) 5 4	10, 20, 45, 77	0
All	All	556/568 (97%)	0.90	93 (16%) 4 3	10, 20, 46, 77	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	THR	10.2
1	B	49	PHE	10.0
1	B	219	GLY	9.3
1	B	3	THR	8.7
1	A	2	ALA	8.4
1	B	44	ASN	7.7
1	B	218	GLU	7.5
1	A	47	ALA	7.5
1	A	49	PHE	7.4
1	B	2	ALA	7.4
1	B	45	THR	7.0
1	B	51	GLY	6.9
1	A	46	SER	6.9
1	A	155	ALA	6.8
1	B	155	ALA	6.8
1	B	47	ALA	6.7
1	B	6	ASN	6.6
1	B	220	LYS	6.3
1	B	5	THR	6.2
1	A	48	ASP	6.2
1	A	218	GLU	6.2
1	B	46	SER	6.0
1	A	51	GLY	5.8
1	A	52	ASN	5.4

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Mol	Chain	Res	Type	RSRZ
1	B	1	MET	5.3
1	A	45	THR	5.2
1	A	220	LYS	5.1
1	A	3	THR	5.1
1	A	50	SER	4.9
1	B	50	SER	4.9
1	B	4	GLU	4.8
1	B	90	GLY	4.7
1	A	4	GLU	4.7
1	B	214	ASN	4.7
1	A	221	ASN	4.6
1	B	48	ASP	4.3
1	A	219	GLY	4.3
1	A	278	VAL	4.2
1	A	44	ASN	4.0
1	A	214	ASN	4.0
1	B	258	PRO	4.0
1	A	1	MET	3.8
1	A	18	PRO	3.7
1	B	91	ASP	3.7
1	B	217	ARG	3.5
1	B	156	GLY	3.4
1	A	173	LYS	3.3
1	B	52	ASN	3.2
1	B	157	ASP	3.2
1	A	258	PRO	3.1
1	A	91	ASP	3.0
1	A	256	LEU	3.0
1	B	141	TRP	3.0
1	A	6	ASN	3.0
1	A	7	TYR	2.9
1	A	217	ARG	2.9
1	B	33	GLU	2.9
1	A	112	GLU	2.9
1	B	221	ASN	2.8
1	A	209	SER	2.8
1	A	12	ARG	2.8
1	B	248	ARG	2.7
1	A	247	GLU	2.7
1	B	79	LYS	2.7
1	B	112	GLU	2.7
1	B	53	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	209	SER	2.6
1	B	262	ASN	2.6
1	A	34	ASP	2.5
1	B	17	GLU	2.5
1	A	53	ASP	2.4
1	A	222	PRO	2.4
1	A	157	ASP	2.4
1	A	259	HIS	2.4
1	A	37	ARG	2.3
1	B	64	GLU	2.3
1	B	278	VAL	2.3
1	A	33	GLU	2.3
1	B	198	ASP	2.3
1	B	159	LEU	2.3
1	A	185	LYS	2.3
1	A	223	LEU	2.3
1	A	262	ASN	2.2
1	B	92	ASP	2.2
1	B	246	GLY	2.2
1	A	210	GLY	2.1
1	A	198	ASP	2.1
1	B	124	SER	2.1
1	A	154	LEU	2.1
1	B	135	LEU	2.1
1	A	19	PHE	2.1
1	A	227	ILE	2.1
1	A	156	GLY	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.