



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:34 PM UTC

PDB ID : 3LCY / pdb_00003lcy
Title : Titin Ig tandem domains A164-A165
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Deposited on : 2010-01-12
Resolution : 2.50 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

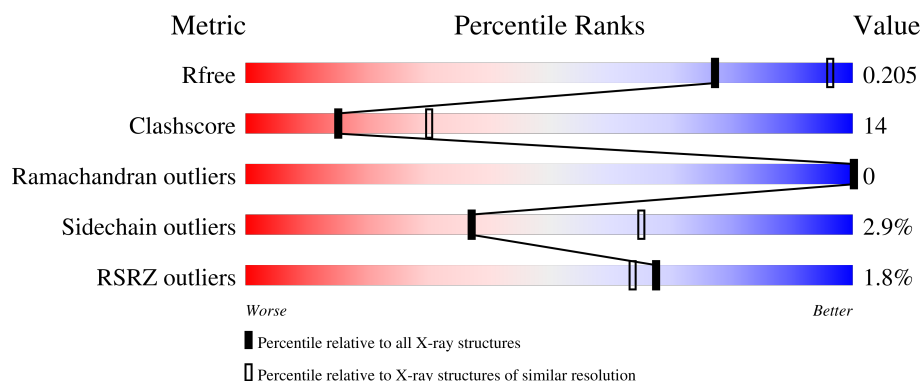
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	197	4% 64% 33% ...
1	B	197	2% 66% 29% ..
1	C	197	% 71% 25% ..
1	D	197	% 67% 31% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6542 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 Titin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	0	0
			1555	982	269	295	9			
1	B	196	Total	C	N	O	S	0	0	0
			1555	982	269	295	9			
1	C	195	Total	C	N	O	S	0	0	0
			1550	979	268	294	9			
1	D	195	Total	C	N	O	S	0	0	0
			1550	979	268	294	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q8WZ42
A	2	ALA	-	expression tag	UNP Q8WZ42
A	3	MET	-	expression tag	UNP Q8WZ42
B	1	GLY	-	expression tag	UNP Q8WZ42
B	2	ALA	-	expression tag	UNP Q8WZ42
B	3	MET	-	expression tag	UNP Q8WZ42
C	1	GLY	-	expression tag	UNP Q8WZ42
C	2	ALA	-	expression tag	UNP Q8WZ42
C	3	MET	-	expression tag	UNP Q8WZ42
D	1	GLY	-	expression tag	UNP Q8WZ42
D	2	ALA	-	expression tag	UNP Q8WZ42
D	3	MET	-	expression tag	UNP Q8WZ42

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	78	Total	O	0	0
			78	78		
2	B	94	Total	O	0	0
			94	94		

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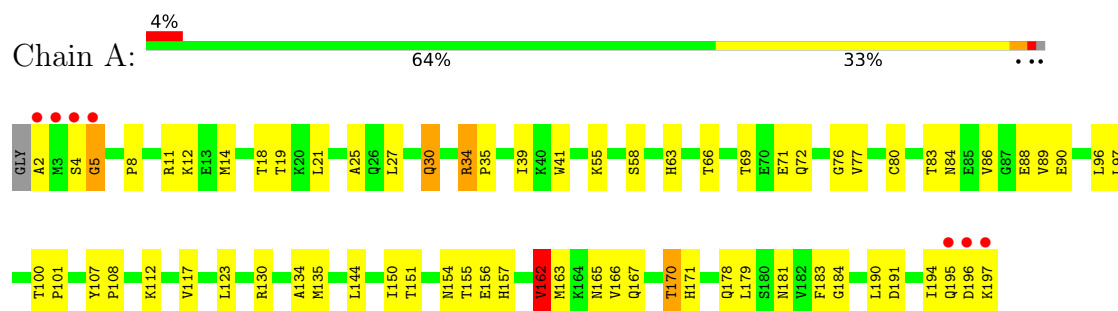
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	81	Total	O	0	0
			81	81		
2	D	79	Total	O	0	0
			79	79		

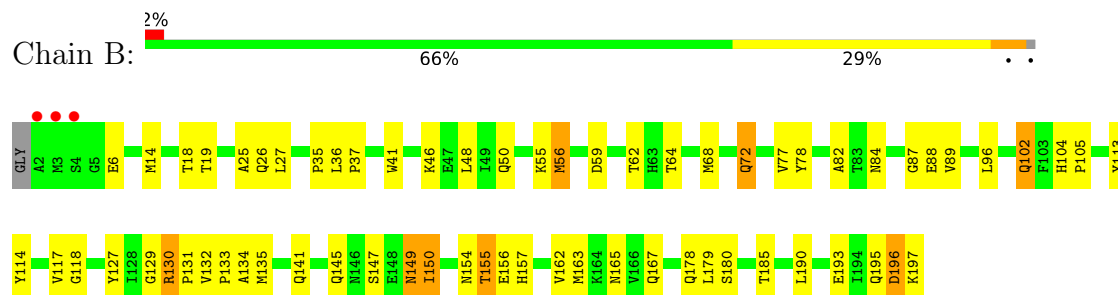
3 Residue-property plots [i](#)

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: Titin



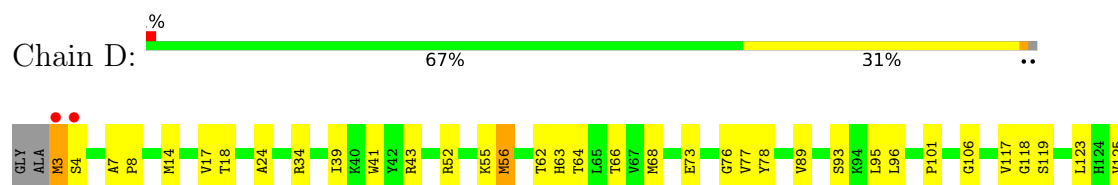
• Molecule 1: Titin



• Molecule 1: Titin



• Molecule 1: Titin



I128	G129	R130	P131
A134	M135	N146	M149
I152	E153	N154	T155
E156	H157	L161	K164
N165	V166	T170	K174
Q178	L179	S180	M181
T185	V186	I189	L190
D196	K197		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.09Å 103.82Å 116.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.59 – 2.50 36.59 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.4 (36.59-2.50) 99.6 (36.59-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 2.51Å)	Xtriage
Refinement program	PHENIX dev_276, REFMAC 5.5.0102	Depositor
R, R_{free}	0.203 , 0.240 (Not available) , 0.205	Depositor DCC
R_{free} test set	1666 reflections (3.45%)	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6542	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 6.34% 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	1.03	9/1586 (0.6%)	0.95	8/2142 (0.4%)
1	B	1.24	11/1586 (0.7%)	0.99	5/2142 (0.2%)
1	C	0.88	4/1581 (0.3%)	0.92	4/2135 (0.2%)
1	D	0.98	7/1581 (0.4%)	0.90	6/2135 (0.3%)
All	All	1.04	31/6334 (0.5%)	0.94	23/8554 (0.3%)

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	7	ALA	CA-CB	-9.01	1.39	1.53
1	B	133	PRO	C-O	-6.83	1.15	1.23
1	D	128	ILE	C-O	-6.79	1.16	1.24
1	B	162	VAL	C-O	-6.76	1.16	1.24
1	B	102	GLN	C-O	-6.61	1.15	1.23
1	D	64	THR	C-O	-6.49	1.16	1.24
1	B	130	ARG	C-O	-6.28	1.16	1.24
1	C	171	HIS	C-O	-6.18	1.16	1.24
1	A	69	THR	C-O	-5.91	1.16	1.23
1	B	134	ALA	CA-CB	-5.86	1.44	1.53
1	A	72	GLN	C-O	-5.83	1.16	1.24
1	C	128	ILE	C-O	-5.81	1.17	1.24
1	B	147	SER	C-O	-5.79	1.18	1.24
1	B	62	THR	C-O	-5.79	1.17	1.24
1	A	34	ARG	CA-C	-5.78	1.47	1.52
1	B	167	GLN	C-O	-5.71	1.16	1.23
1	A	76	GLY	C-O	-5.70	1.16	1.23
1	C	127	TYR	C-O	-5.69	1.16	1.23
1	B	113	TYR	C-O	-5.67	1.16	1.23
1	B	145	GLN	C-O	-5.65	1.16	1.23
1	A	71	GLU	C-O	-5.62	1.16	1.23
1	B	55	LYS	C-O	-5.51	1.17	1.24
1	C	173	GLY	C-O	-5.48	1.18	1.24
1	D	129	GLY	C-O	-5.46	1.16	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	162	VAL	C-O	-5.45	1.18	1.24
1	D	56	MET	C-O	-5.37	1.17	1.24
1	D	174	LYS	C-O	-5.37	1.17	1.24
1	A	30	GLN	C-O	-5.22	1.18	1.24
1	D	62	THR	C-O	-5.12	1.17	1.24
1	A	39	ILE	C-O	-5.05	1.18	1.24
1	A	34	ARG	C-O	-5.04	1.19	1.25

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	139	HIS	N-CA-C	-8.28	95.03	108.52
1	A	77	VAL	N-CA-C	7.91	119.19	109.30
1	C	183	PHE	N-CA-C	6.97	118.96	111.36
1	D	170	THR	N-CA-C	6.86	118.84	111.36
1	A	76	GLY	N-CA-C	6.29	121.14	112.52
1	B	77	VAL	N-CA-C	6.26	117.16	108.89
1	A	5	GLY	N-CA-C	-6.25	101.64	110.96
1	D	77	VAL	N-CA-C	6.25	117.64	109.58
1	B	196	ASP	N-CA-C	6.01	118.95	110.23
1	A	144	LEU	N-CA-C	5.96	118.90	109.96
1	C	134	ALA	N-CA-C	-5.94	100.32	109.76
1	D	76	GLY	N-CA-C	5.90	120.19	111.76
1	A	170	THR	N-CA-C	5.85	120.56	112.90
1	D	134	ALA	N-CA-C	-5.81	100.53	109.76
1	A	134	ALA	N-CA-C	-5.71	101.39	109.96
1	A	39	ILE	N-CA-C	5.50	116.23	108.36
1	C	5	GLY	N-CA-C	5.49	117.78	111.36
1	A	4	SER	N-CA-C	5.48	118.16	109.50
1	B	132	VAL	CA-C-N	5.37	125.57	120.31
1	B	132	VAL	C-N-CA	5.37	125.57	120.31
1	D	125	VAL	N-CA-C	5.08	115.07	107.51
1	D	106	GLY	N-CA-C	-5.07	108.45	114.48
1	B	150	ILE	N-CA-C	5.03	115.36	108.12

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	1555	0	1559	53	1
1	B	1555	0	1559	44	1
1	C	1550	0	1554	37	0
1	D	1550	0	1554	36	0
2	A	78	0	0	2	0
2	B	94	0	0	1	0
2	C	81	0	0	0	0
2	D	79	0	0	2	0
All	All	6542	0	6226	169	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 14.

All (169) 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:196:ASP:O	1:A:197:LYS:HG2	1.64	0.97
1:C:19:THR:HG21	1:C:25:ALA:HB2	1.49	0.94
1:C:155:THR:HG22	1:C:157:HIS:H	1.32	0.91
1:D:8:PRO:HB2	1:D:89:VAL:HG22	1.56	0.85
1:B:155:THR:HG22	1:B:157:HIS:H	1.41	0.85
1:A:117:VAL:HG23	1:A:195:GLN:O	1.78	0.84
1:A:196:ASP:O	1:A:197:LYS:CG	2.30	0.79
1:A:170:THR:HG23	1:A:171:HIS:CD2	2.19	0.78
1:B:117:VAL:HG23	1:B:195:GLN:O	1.85	0.75
1:A:150:ILE:CD1	1:A:163:MET:HG2	2.17	0.74
1:B:118:GLY:HA2	1:B:165:ASN:HD22	1.55	0.72
1:C:135:MET:HE3	1:C:154:ASN:HB3	1.70	0.72
1:B:150:ILE:CD1	1:B:163:MET:HG2	2.19	0.72
1:D:56:MET:HE1	2:D:281:HOH:O	1.90	0.71
1:A:8:PRO:O	1:A:89:VAL:HG11	1.90	0.71
1:C:20:LYS:HB3	1:C:23:GLU:HG3	1.75	0.69
1:A:167:GLN:HB2	1:A:170:THR:HG22	1.75	0.68
1:A:170:THR:HG23	1:A:171:HIS:HD2	1.57	0.68
1:B:150:ILE:HD13	1:B:163:MET:HG2	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:VAL:CG2	1:A:196:ASP:HA	2.23	0.67
1:A:117:VAL:HG23	1:A:195:GLN:C	2.19	0.67
1:A:8:PRO:O	1:A:89:VAL:CG1	2.42	0.67
1:A:12:LYS:HB3	1:A:30:GLN:HG3	1.78	0.66
1:B:41:TRP:CD1	1:B:56:MET:HG2	2.30	0.66
1:A:117:VAL:HG22	1:A:194:ILE:HG22	1.78	0.66
1:A:150:ILE:HD13	1:A:163:MET:HG2	1.77	0.66
1:A:14:MET:CE	1:A:27:LEU:HB3	2.25	0.65
1:A:100:THR:HG22	1:A:183:PHE:HB3	1.79	0.65
1:B:82:ALA:HB3	1:B:89:VAL:HG12	1.77	0.65
1:D:146:ASN:ND2	1:D:152:ILE:H	1.95	0.65
1:C:112:LYS:HG3	1:C:191:ASP:HB3	1.79	0.64
1:A:155:THR:HG22	1:A:157:HIS:H	1.63	0.64
1:B:149:ASN:O	1:B:149:ASN:ND2	2.30	0.64
1:D:41:TRP:CD1	1:D:56:MET:HG2	2.32	0.64
1:D:146:ASN:HD21	1:D:152:ILE:H	1.44	0.64
1:D:52:ARG:NH1	2:D:325:HOH:O	2.25	0.62
1:B:48:LEU:O	1:B:56:MET:HE1	1.98	0.62
1:A:41:TRP:CZ3	1:A:80:CYS:HB2	2.35	0.62
1:D:117:VAL:HG21	1:D:196:ASP:HB3	1.81	0.62
1:A:135:MET:HE3	1:A:154:ASN:HB3	1.82	0.61
1:D:135:MET:HE3	1:D:154:ASN:HB3	1.82	0.61
1:A:117:VAL:HG21	1:A:196:ASP:HA	1.82	0.60
1:B:50:GLN:HE22	1:B:56:MET:H	1.48	0.59
1:B:155:THR:HG22	1:B:157:HIS:N	2.15	0.59
1:C:112:LYS:HE3	1:C:193:GLU:OE1	2.03	0.59
1:A:196:ASP:C	1:A:197:LYS:CG	2.76	0.59
1:D:3:MET:N	1:D:34:ARG:HH22	2.01	0.58
1:B:118:GLY:HA2	1:B:165:ASN:ND2	2.19	0.58
1:D:155:THR:HG22	1:D:156:GLU:N	2.18	0.58
1:A:151:THR:HB	1:A:162:VAL:HG13	1.86	0.57
1:C:21:LEU:HD13	1:C:97:LEU:HD13	1.85	0.57
1:C:49:ILE:HD12	1:C:50:GLN:H	1.67	0.57
1:A:19:THR:HG21	1:A:25:ALA:HB2	1.86	0.57
1:C:155:THR:CG2	1:C:156:GLU:N	2.67	0.57
1:C:82:ALA:HB3	1:C:89:VAL:HG12	1.87	0.56
1:A:5:GLY:C	1:A:86:VAL:HG21	2.30	0.56
1:A:2:ALA:CB	2:A:242:HOH:O	2.53	0.55
1:C:155:THR:HG22	1:C:157:HIS:N	2.13	0.55
1:D:55:LYS:HB2	1:D:66:THR:HB	1.87	0.55
1:D:156:GLU:HG2	1:D:157:HIS:CD2	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:GLN:C	1:A:179:LEU:HD12	2.32	0.55
1:B:196:ASP:O	1:B:197:LYS:HB2	2.05	0.55
1:B:84:ASN:OD1	1:B:87:GLY:N	2.30	0.55
1:C:18:THR:HA	1:C:96:LEU:O	2.06	0.55
1:A:112:LYS:HG2	1:A:191:ASP:HB3	1.90	0.54
1:C:19:THR:HG21	1:C:25:ALA:CB	2.33	0.53
1:B:127:TYR:C	1:B:127:TYR:CD1	2.87	0.53
1:B:14:MET:HE3	1:B:27:LEU:HB3	1.90	0.53
1:A:135:MET:CE	1:A:154:ASN:HB3	2.39	0.52
1:B:72:GLN:OE1	1:B:129:GLY:N	2.28	0.52
1:C:130:ARG:HA	1:C:131:PRO:C	2.35	0.52
1:B:135:MET:HE3	1:B:154:ASN:HB3	1.92	0.52
1:A:58:SER:OG	1:A:63:HIS:ND1	2.38	0.52
1:B:155:THR:CG2	1:B:156:GLU:N	2.72	0.52
1:D:123:LEU:HD12	1:D:161:LEU:HD23	1.93	0.51
1:C:168:ARG:NH1	1:C:194:ILE:O	2.43	0.51
1:D:101:PRO:HB2	1:D:186:VAL:HG12	1.92	0.51
1:B:35:PRO:HG2	2:B:256:HOH:O	2.11	0.50
1:B:180:SER:OG	1:B:185:THR:HG23	2.11	0.50
1:B:19:THR:HG21	1:B:25:ALA:HB2	1.92	0.50
1:C:19:THR:HG22	1:C:95:LEU:HD11	1.94	0.50
1:D:180:SER:OG	1:D:185:THR:HG23	2.12	0.50
1:B:27:LEU:HD22	1:B:78:TYR:HB2	1.93	0.50
1:D:119:SER:O	1:D:166:VAL:HG22	2.11	0.50
1:C:178:GLN:C	1:C:179:LEU:HD12	2.37	0.49
1:D:18:THR:HA	1:D:96:LEU:O	2.12	0.49
1:A:167:GLN:HB2	1:A:170:THR:CG2	2.42	0.49
1:D:8:PRO:O	1:D:89:VAL:CG1	2.60	0.49
1:A:21:LEU:HD13	1:A:97:LEU:HD13	1.94	0.49
1:A:165:ASN:O	1:A:165:ASN:OD1	2.30	0.49
1:D:130:ARG:HA	1:D:131:PRO:C	2.37	0.49
1:A:117:VAL:HG23	1:A:196:ASP:HA	1.94	0.49
1:C:19:THR:CG2	1:C:95:LEU:HD11	2.43	0.49
1:A:155:THR:CG2	1:A:156:GLU:N	2.75	0.48
1:B:130:ARG:HA	1:B:131:PRO:C	2.37	0.48
1:C:101:PRO:HG2	1:C:184:GLY:O	2.13	0.48
1:A:83:THR:HG22	1:A:84:ASN:N	2.27	0.48
1:A:155:THR:HG22	1:A:156:GLU:N	2.28	0.48
1:B:190:LEU:HD12	1:B:190:LEU:N	2.29	0.48
1:C:135:MET:HB2	1:C:154:ASN:ND2	2.29	0.47
1:D:24:ALA:HB2	1:D:68:MET:CE	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:151:THR:HB	1:C:162:VAL:HG13	1.97	0.47
1:A:34:ARG:HA	1:A:35:PRO:C	2.38	0.47
1:A:107:TYR:N	1:A:108:PRO:HD3	2.30	0.47
1:C:148:GLU:HA	1:C:148:GLU:OE1	2.16	0.46
1:A:190:LEU:HD12	1:A:190:LEU:N	2.30	0.46
1:C:11:ARG:HD3	1:C:11:ARG:HA	1.69	0.46
1:D:14:MET:HE2	1:D:93:SER:OG	2.16	0.46
1:A:117:VAL:HG21	1:A:196:ASP:OD1	2.16	0.46
1:A:2:ALA:HB3	2:A:242:HOH:O	2.14	0.46
1:A:55:LYS:HB2	1:A:66:THR:HB	1.98	0.46
1:B:50:GLN:NE2	1:B:56:MET:H	2.13	0.46
1:C:84:ASN:OD1	1:C:84:ASN:C	2.59	0.46
1:C:107:TYR:N	1:C:108:PRO:HD3	2.31	0.46
1:D:178:GLN:C	1:D:179:LEU:HD12	2.41	0.46
1:A:97:LEU:HB2	1:A:130:ARG:HG3	1.97	0.46
1:A:178:GLN:O	1:A:179:LEU:HD12	2.16	0.45
1:C:164:LYS:CB	1:C:164:LYS:HZ3	2.28	0.45
1:A:155:THR:HG22	1:A:157:HIS:N	2.28	0.45
1:B:84:ASN:OD1	1:B:84:ASN:C	2.58	0.45
1:B:18:THR:HA	1:B:96:LEU:O	2.17	0.45
1:B:114:TYR:CD1	1:B:193:GLU:HB3	2.52	0.45
1:D:149:ASN:OD1	1:D:164:LYS:HE2	2.17	0.45
1:A:83:THR:HG23	1:A:88:GLU:HB3	1.99	0.45
1:D:155:THR:CG2	1:D:156:GLU:N	2.80	0.44
1:B:14:MET:CE	1:B:41:TRP:HZ3	2.30	0.44
1:B:135:MET:CE	1:B:154:ASN:HB3	2.48	0.44
1:C:151:THR:HB	1:C:162:VAL:CG1	2.48	0.44
1:B:46:LYS:HE3	1:B:46:LYS:HB2	1.89	0.44
1:C:36:LEU:HA	1:C:37:PRO:HD3	1.75	0.44
1:D:118:GLY:HA2	1:D:165:ASN:HD22	1.83	0.44
1:B:178:GLN:C	1:B:179:LEU:HD12	2.43	0.44
1:D:39:ILE:HG13	1:D:63:HIS:ND1	2.33	0.44
1:C:135:MET:CE	1:C:154:ASN:HB3	2.46	0.43
1:D:17:VAL:O	1:D:95:LEU:HD12	2.18	0.43
1:D:8:PRO:O	1:D:89:VAL:HG11	2.18	0.43
1:D:101:PRO:HD3	1:D:181:ASN:CG	2.43	0.43
1:B:50:GLN:HE21	1:B:50:GLN:HB2	1.62	0.43
1:B:104:HIS:HA	1:B:105:PRO:HD3	1.81	0.43
1:D:189:ILE:HD12	1:D:189:ILE:N	2.33	0.43
1:B:82:ALA:O	1:B:88:GLU:HA	2.19	0.42
1:D:24:ALA:HB2	1:D:68:MET:HE1	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:155:THR:HG23	1:C:156:GLU:N	2.35	0.42
1:A:166:VAL:HA	1:A:171:HIS:CD2	2.55	0.42
1:B:59:ASP:C	1:B:59:ASP:OD2	2.62	0.42
1:C:155:THR:HG23	1:C:156:GLU:H	1.85	0.42
1:D:3:MET:HB3	1:D:4:SER:H	1.60	0.42
1:C:179:LEU:CD1	1:C:179:LEU:N	2.83	0.42
1:B:102:GLN:NE2	1:B:105:PRO:HG3	2.34	0.42
1:A:90:GLU:O	1:B:141:GLN:HB2	2.20	0.41
1:C:26:GLN:HG3	1:C:64:THR:HG23	2.02	0.41
1:B:14:MET:HE3	1:B:41:TRP:HZ3	1.86	0.41
1:D:181:ASN:OD1	1:D:181:ASN:C	2.63	0.41
1:C:118:GLY:HA2	1:C:165:ASN:ND2	2.36	0.41
1:D:8:PRO:O	1:D:89:VAL:HG13	2.21	0.41
1:A:84:ASN:OD1	1:A:84:ASN:C	2.64	0.41
1:B:155:THR:HG22	1:B:156:GLU:N	2.36	0.41
1:D:43:ARG:HD3	1:D:78:TYR:CE1	2.56	0.41
1:A:101:PRO:HG2	1:A:184:GLY:O	2.20	0.41
1:A:165:ASN:O	1:A:165:ASN:CG	2.64	0.41
1:B:196:ASP:O	1:B:197:LYS:CB	2.67	0.41
1:B:36:LEU:HA	1:B:37:PRO:HD2	1.89	0.41
1:A:18:THR:HA	1:A:96:LEU:O	2.21	0.40
1:C:110:LYS:HB2	1:C:110:LYS:HE3	1.87	0.40
1:A:101:PRO:HD3	1:A:181:ASN:CG	2.47	0.40
1:C:96:LEU:HD12	1:C:96:LEU:N	2.35	0.40
1:D:190:LEU:HD12	1:D:190:LEU:N	2.36	0.40
1:B:68:MET:HA	1:B:68:MET:HE3	2.04	0.40
1:C:101:PRO:HD3	1:C:181:ASN:CG	2.46	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:A:11:ARG:NE	1:B:6:GLU:OE2[1_655]	2.15	0.05

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	194/197 (98%)	191 (98%)	3 (2%)	0	100	100
1	B	194/197 (98%)	192 (99%)	2 (1%)	0	100	100
1	C	193/197 (98%)	191 (99%)	2 (1%)	0	100	100
1	D	193/197 (98%)	191 (99%)	2 (1%)	0	100	100
All	All	774/788 (98%)	765 (99%)	9 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	171/171 (100%)	169 (99%)	2 (1%)	63	83
1	B	171/171 (100%)	165 (96%)	6 (4%)	32	58
1	C	171/171 (100%)	161 (94%)	10 (6%)	18	38
1	D	171/171 (100%)	169 (99%)	2 (1%)	63	83
All	All	684/684 (100%)	664 (97%)	20 (3%)	37	65

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	LEU
1	A	162	VAL
1	B	26	GLN
1	B	56	MET
1	B	64	THR
1	B	72	GLN
1	B	149	ASN
1	B	155	THR

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Mol	Chain	Res	Type
1	C	19	THR
1	C	23	GLU
1	C	49	ILE
1	C	86	VAL
1	C	89	VAL
1	C	136	THR
1	C	155	THR
1	C	156	GLU
1	C	169	LYS
1	C	179	LEU
1	D	3	MET
1	D	73	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	50	GLN
1	A	98	GLN
1	A	141	GLN
1	A	146	ASN
1	A	171	HIS
1	B	50	GLN
1	B	149	ASN
1	B	157	HIS
1	B	165	ASN
1	C	26	GLN
1	C	149	ASN
1	C	165	ASN
1	D	30	GLN
1	D	146	ASN
1	D	165	ASN

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	196/197 (99%)	-0.19	7 (3%) 46 41	19, 33, 65, 125	0
1	B	196/197 (99%)	-0.21	3 (1%) 72 68	19, 34, 60, 94	0
1	C	195/197 (98%)	-0.28	2 (1%) 79 76	20, 33, 51, 104	0
1	D	195/197 (98%)	-0.27	2 (1%) 79 76	19, 33, 54, 103	0
All	All	782/788 (99%)	-0.24	14 (1%) 67 64	19, 33, 57, 125	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	197	LYS	3.3
1	C	3	MET	3.3
1	B	4	SER	3.1
1	A	2	ALA	3.0
1	D	3	MET	2.9
1	C	4	SER	2.6
1	A	5	GLY	2.5
1	A	3	MET	2.5
1	B	3	MET	2.4
1	A	196	ASP	2.4
1	D	4	SER	2.3
1	A	4	SER	2.2
1	B	2	ALA	2.1
1	A	195	GLN	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.