



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 19, 2017 – 03:17 PM EDT

PDB ID : 3LCY  
Title : Titin Ig tandem domains A164-A165  
Authors : Chen, Q.; Groves, M.R.; Wilmanns, M.  
Deposited on : unknown  
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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345

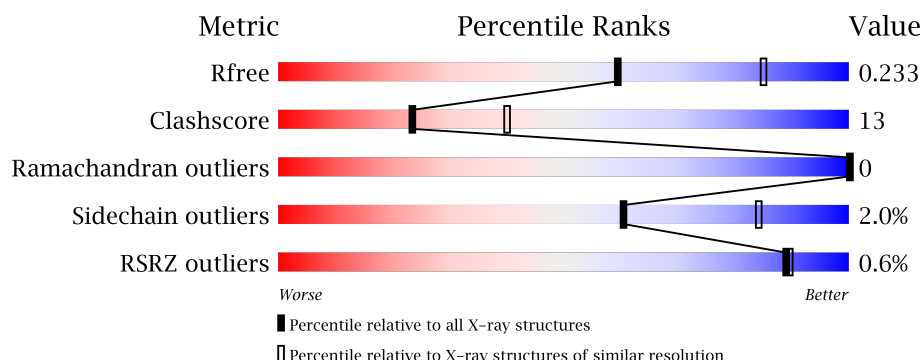
# 1 Overall quality at a glance

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))
$R_{free}$	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>2%</div> <div>69%</div> <div>30%</div> <div>..</div> </div>
1	B	197	<div> <div>74%</div> <div>25%</div> <div>..</div> </div>
1	C	197	<div> <div>%</div> <div>76%</div> <div>20%</div> <div>..</div> </div>
1	D	197	<div> <div>%</div> <div>74%</div> <div>24%</div> <div>..</div> </div>

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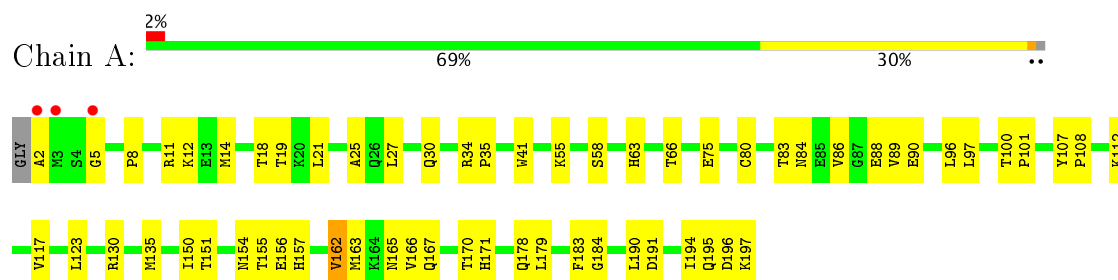
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	81	Total	O	0	0
			81	81		
2	D	79	Total	O	0	0
			79	79		

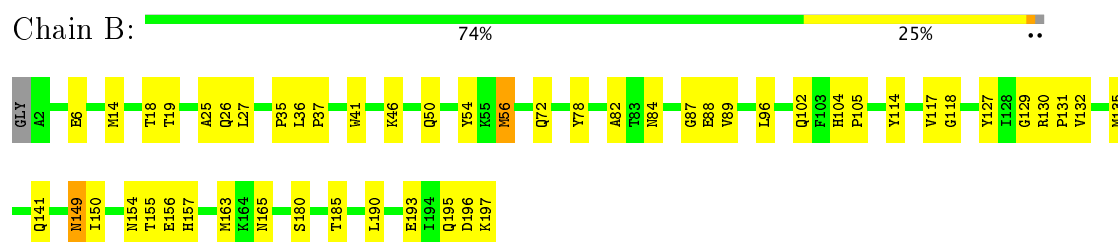
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

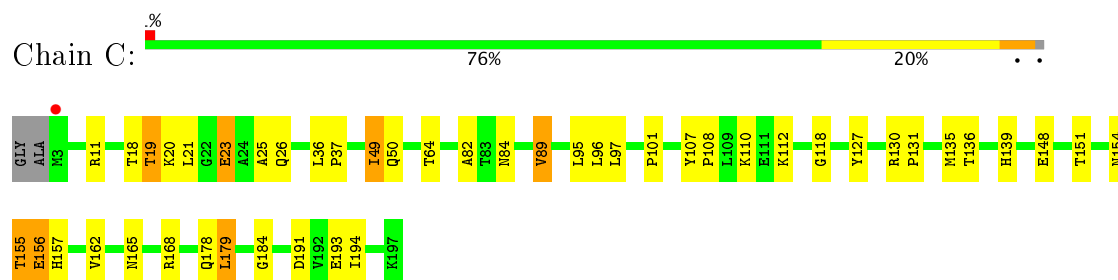
- Molecule 1: Titin



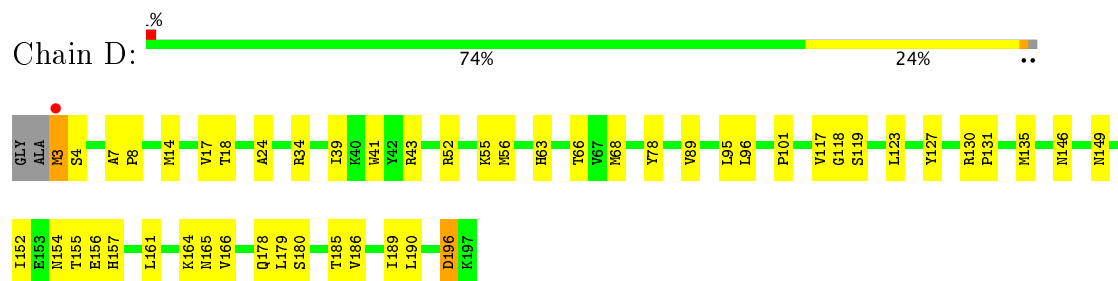
- Molecule 1: Titin



- Molecule 1: Titin



- Molecule 1: Titin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.09Å 103.82Å 116.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.59 – 2.50 43.06 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.4 (36.59-2.50) 99.4 (43.06-2.50)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.94 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0102, PHENIX dev_276	Depositor
R, $R_{free}$	0.203 , 0.240 0.203 , 0.233	Depositor DCC
$R_{free}$ test set	1665 reflections (5.33%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.8	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 39.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	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 <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.81	3/1586 (0.2%)	0.60	0/2142
1	B	1.00	2/1586 (0.1%)	0.64	0/2142
1	C	0.68	2/1581 (0.1%)	0.59	1/2135 (0.0%)
1	D	0.79	2/1581 (0.1%)	0.58	0/2135
All	All	0.83	9/6334 (0.1%)	0.60	1/8554 (0.0%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	7	ALA	CA-CB	-6.27	1.39	1.52
1	A	75	GLU	CD-OE2	-5.83	1.19	1.25
1	D	127	TYR	CD2-CE2	-5.57	1.30	1.39
1	B	54	TYR	CD2-CE2	-5.55	1.31	1.39
1	C	127	TYR	CD1-CE1	-5.42	1.31	1.39
1	A	162	VAL	CB-CG1	-5.41	1.41	1.52
1	B	132	VAL	CB-CG2	-5.39	1.41	1.52
1	A	162	VAL	CB-CG2	-5.16	1.42	1.52
1	C	127	TYR	CD2-CE2	-5.08	1.31	1.39

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	139	HIS	N-CA-C	-5.92	95.03	111.00

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	50	1
1	B	1555	0	1559	39	1
1	C	1550	0	1554	35	0
1	D	1550	0	1554	35	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	3	0
All	All	6542	0	6226	158	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 13.

All (158) 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:B:150:ILE:CD1	1:B:163:MET:HG2	2.19	0.72
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
1:A:117:VAL:CG2	1:A:196:ASP:HA	2.23	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:146:ASN:HD21	1:D:152:ILE:H	1.44	0.64
1:D:41:TRP:CD1	1:D:56:MET:HG2	2.32	0.64
1:A:117:VAL:HG23	1:A:195:GLN:C	2.19	0.62
1:D:52:ARG:NH1	2:D:325:HOH:O	2.25	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:117:VAL:HG21	1:A:196:ASP:HA	1.82	0.60
1:B:155:THR:HG22	1:B:157:HIS:N	2.15	0.59
1:B:50:GLN:HE22	1:B:56:MET:H	1.48	0.59
1:C:112:LYS:HE3	1:C:193:GLU:OE1	2.03	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:49:ILE:HD12	1:C:50:GLN:H	1.67	0.57
1:C:21:LEU:HD13	1:C:97:LEU:HD13	1.85	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: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:156:GLU:HG2	1:D:157:HIS:CD2	2.42	0.55
1:D:55:LYS:HB2	1:D:66:THR:HB	1.87	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:A:196:ASP:C	1:A:197:LYS:CG	2.76	0.54
1:C:19:THR:HG21	1:C:25:ALA:CB	2.33	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:A:58:SER:OG	1:A:63:HIS:ND1	2.38	0.52
1:A:5:GLY:C	1:A:86:VAL:HG21	2.30	0.52
1:B:155:THR:CG2	1:B:156:GLU:N	2.72	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:178:GLN:C	1:A:179:LEU:HD12	2.32	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:14:MET:HE3	1:B:27:LEU:HB3	1.93	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: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:165:ASN:O	1:A:165:ASN:OD1	2.30	0.49
1:A:21:LEU:HD13	1:A:97:LEU:HD13	1.94	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:D:56:MET:HE3	2:D:281:HOH:O	2.12	0.49
1:D:24:ALA:HB2	1:D:68:MET:HE3	1.95	0.49
1:A:155:THR:CG2	1:A:156:GLU:N	2.75	0.48
1:B:127:TYR:C	1:B:127:TYR:CD1	2.87	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:D:14:MET:HE3	1:D:41:TRP:HZ3	1.79	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
1:C:130:ARG:HA	1:C:131:PRO:C	2.35	0.47
1:C:151:THR:HB	1:C:162:VAL:HG13	1.97	0.47
1:A:107:TYR:N	1:A:108:PRO:HD3	2.30	0.47
1:A:135:MET:HE3	1:A:154:ASN:HB3	1.97	0.47
1:C:148:GLU:HA	1:C:148:GLU:OE1	2.16	0.46
1:A:190:LEU:N	1:A:190:LEU:HD12	2.30	0.46
1:C:11:ARG:HD3	1:C:11:ARG:HA	1.69	0.46
1:B:14:MET:HE3	1:B:41:TRP:HZ3	1.81	0.46
1:A:117:VAL:HG21	1:A:196:ASP:OD1	2.16	0.46
1:C:135:MET:CE	1:C:154:ASN:HB3	2.46	0.46
1:A:2:ALA:HB3	2:A:242:HOH:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:107:TYR:N	1:C:108:PRO:HD3	2.31	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:135:MET:HE1	1:C:154:ASN:HB3	1.98	0.45
1:D:56:MET:CE	2:D:281:HOH:O	2.64	0.45
1:A:155:THR:HG22	1:A:157:HIS:N	2.28	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:C:178:GLN:C	1:C:179:LEU:HD12	2.37	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:130:ARG:HA	1:D:131:PRO:C	2.37	0.44
1:D:155:THR:CG2	1:D:156:GLU:N	2.80	0.44
1:B:135:MET:CE	1:B:154:ASN:HB3	2.48	0.44
1:B:14:MET:CE	1:B:41:TRP:HZ3	2.30	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:130:ARG:HA	1:B:131:PRO:C	2.37	0.44
1:D:135:MET:CE	1:D:154:ASN:HB3	2.48	0.44
1:D:39:ILE:HG13	1:D:63:HIS:ND1	2.33	0.44
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: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:N	1:D:189:ILE:HD12	2.33	0.43
1:A:34:ARG:HA	1:A:35:PRO:C	2.38	0.42
1:B:82:ALA:O	1:B:88:GLU:HA	2.19	0.42
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: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:N	1:C:179:LEU:CD1	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:B:84:ASN:OD1	1:B:84:ASN:C	2.58	0.41
1:C:26:GLN:HG3	1:C:64:THR:HG23	2.02	0.41
1:C:118:GLY:HA2	1:C:165:ASN:ND2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:ASN:OD1	1:C:84:ASN:C	2.59	0.41
1:D:8:PRO:O	1:D:89:VAL:HG13	2.21	0.41
1:B:155:THR:HG22	1:B:156:GLU:N	2.36	0.41
1:D:178:GLN:C	1:D:179:LEU:HD12	2.41	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: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:C:96:LEU:N	1:C:96:LEU:HD12	2.35	0.40
1:D:190:LEU:HD12	1:D:190:LEU:N	2.36	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%)	170 (99%)	1 (1%)	89	97
1	B	171/171 (100%)	168 (98%)	3 (2%)	64	86
1	C	171/171 (100%)	163 (95%)	8 (5%)	30	54
1	D	171/171 (100%)	169 (99%)	2 (1%)	75	91
All	All	684/684 (100%)	670 (98%)	14 (2%)	60	84

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	LEU
1	B	26	GLN
1	B	56	MET
1	B	149	ASN
1	C	19	THR
1	C	23	GLU
1	C	49	ILE
1	C	89	VAL
1	C	136	THR
1	C	155	THR
1	C	156	GLU
1	C	179	LEU
1	D	3	MET
1	D	196	ASP

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

Mol	Chain	Res	Type
1	A	50	GLN
1	A	98	GLN
1	A	141	GLN
1	A	171	HIS
1	B	50	GLN

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Mol	Chain	Res	Type
1	B	149	ASN
1	B	165	ASN
1	B	178	GLN
1	C	26	GLN
1	C	98	GLN
1	C	149	ASN
1	C	157	HIS
1	C	165	ASN
1	D	30	GLN
1	D	146	ASN
1	D	165	ASN
1	D	178	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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	196/197 (99%)	-0.42	3 (1%) 74 75	19, 33, 65, 125	0
1	B	196/197 (99%)	-0.39	0 100 100	19, 34, 60, 94	0
1	C	195/197 (98%)	-0.42	1 (0%) 90 91	20, 33, 51, 104	0
1	D	195/197 (98%)	-0.48	1 (0%) 90 91	19, 33, 54, 103	0
All	All	782/788 (99%)	-0.43	5 (0%) 89 89	19, 33, 57, 125	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	3	MET	4.6
1	D	3	MET	3.4
1	A	2	ALA	2.7
1	A	3	MET	2.5
1	A	5	GLY	2.2

### 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 carbohydrates 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.