



# Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 11:19 am BST

PDB ID : 5XQI  
Title : Crystal structure of full-length human Rogdi  
Authors : Lee, H.; Lee, C.  
Deposited on : 2017-06-07  
Resolution : 2.80 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](mailto: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.

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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)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

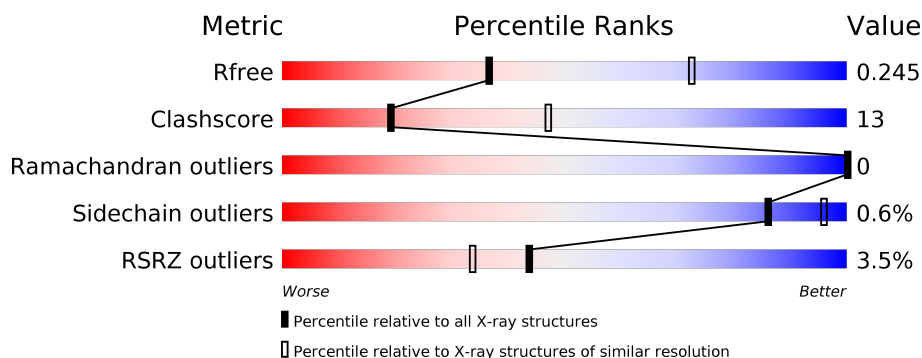
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	289	<div> <div> <div></div> <div>76%</div> <div>19%</div> <div>5%</div> </div> </div>
1	B	289	<div> <div> <div></div> <div>80%</div> <div>14%</div> <div>• •</div> </div> </div>
1	C	289	<div> <div> <div></div> <div>67%</div> <div>19%</div> <div>• 12%</div> </div> </div>
1	D	289	<div> <div> <div></div> <div>67%</div> <div>19%</div> <div>• 12%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8569 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 Protein rogdi homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2193	1399	385	398	11			
1	B	276	Total	C	N	O	S	0	0	0
			2197	1401	386	399	11			
1	C	254	Total	C	N	O	S	0	0	0
			2030	1304	349	367	10			
1	D	255	Total	C	N	O	S	0	0	0
			2043	1312	353	368	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP Q9GZN7
A	0	GLY	-	expression tag	UNP Q9GZN7
B	-1	SER	-	expression tag	UNP Q9GZN7
B	0	GLY	-	expression tag	UNP Q9GZN7
C	-1	SER	-	expression tag	UNP Q9GZN7
C	0	GLY	-	expression tag	UNP Q9GZN7
D	-1	SER	-	expression tag	UNP Q9GZN7
D	0	GLY	-	expression tag	UNP Q9GZN7

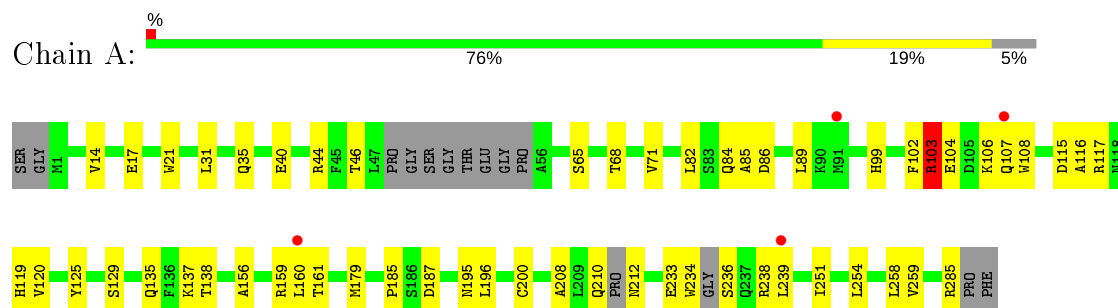
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	25	Total	O	0	0
			25	25		
2	B	36	Total	O	0	0
			36	36		
2	C	22	Total	O	0	0
			22	22		
2	D	23	Total	O	0	0
			23	23		

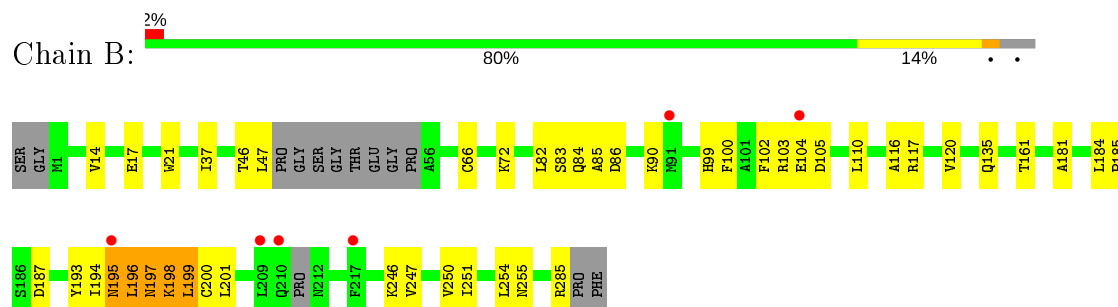
### 3 Residue-property plots [i](#)

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.

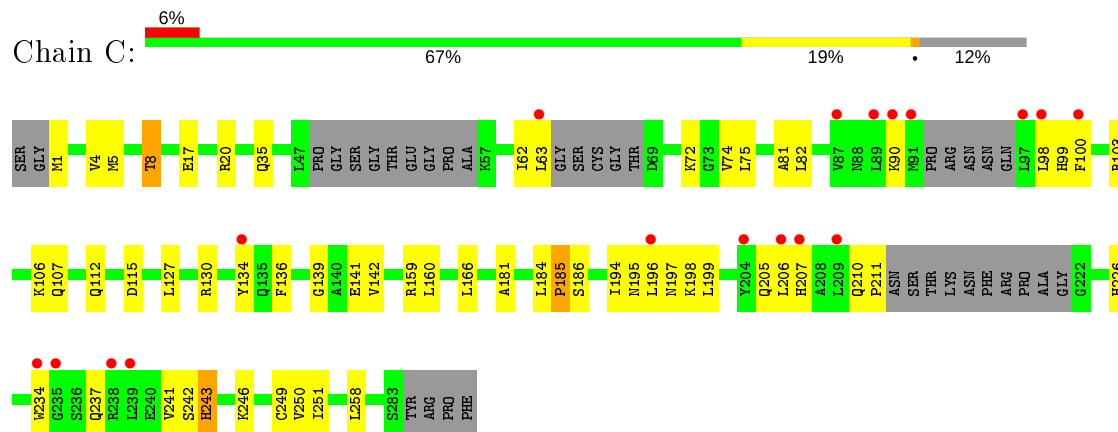
- Molecule 1: Protein rogdi homolog



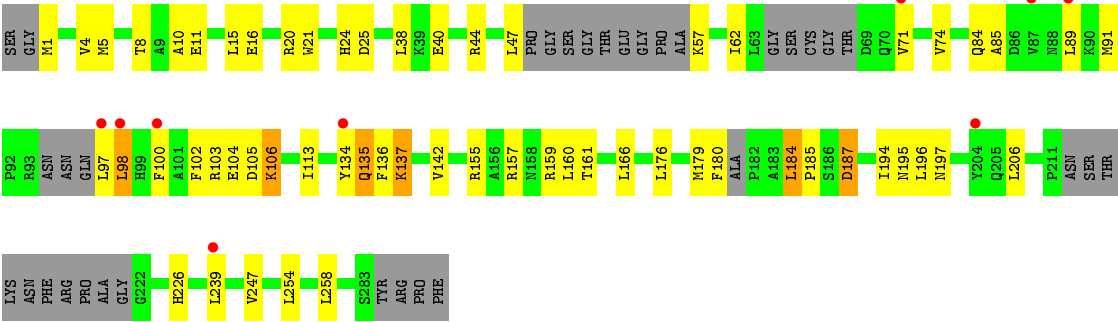
- Molecule 1: Protein rogdi homolog



- Molecule 1: Protein rogdi homolog



- Molecule 1: Protein rogdi homolog



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	169.37Å 169.37Å 220.32Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.17 – 2.80 36.17 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (36.17-2.80) 99.4 (36.17-2.80)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.79 (at 2.81Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.219 , 0.268 0.233 , 0.245	Depositor DCC
$R_{free}$ test set	1994 reflections (3.45%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	64.0	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 48.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.478 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8569	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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 [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	0.28	0/2233	0.60	4/3024 (0.1%)
1	B	0.28	0/2238	0.57	3/3032 (0.1%)
1	C	0.55	6/2066 (0.3%)	0.67	3/2798 (0.1%)
1	D	0.37	0/2079	0.71	8/2813 (0.3%)
All	All	0.38	6/8616 (0.1%)	0.64	18/11667 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	136	PHE	CA-CB	-8.94	1.34	1.53
1	C	136	PHE	CB-CG	-6.73	1.40	1.51
1	C	8	THR	C-O	-6.64	1.10	1.23
1	C	8	THR	CB-CG2	-5.66	1.33	1.52
1	C	136	PHE	CG-CD1	-5.28	1.30	1.38
1	C	211	PRO	N-CD	5.24	1.55	1.47

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	106	LYS	N-CA-CB	-9.27	93.91	110.60
1	B	196	LEU	N-CA-C	-8.71	87.47	111.00
1	D	135	GLN	N-CA-CB	-8.67	95.00	110.60
1	D	134	TYR	CB-CA-C	8.21	126.83	110.40
1	C	8	THR	OG1-CB-CG2	-7.16	93.52	110.00
1	A	103	ARG	N-CA-C	7.13	130.26	111.00
1	B	197	ASN	N-CA-C	7.01	129.94	111.00
1	D	197	ASN	N-CA-CB	6.80	122.85	110.60
1	A	103	ARG	CB-CA-C	-6.65	97.10	110.40
1	A	104	GLU	N-CA-CB	6.43	122.17	110.60
1	D	106	LYS	N-CA-C	6.33	128.08	111.00
1	B	195	ASN	N-CA-C	-6.04	94.71	111.00
1	D	187	ASP	CB-CG-OD1	5.82	123.54	118.30
1	C	210	GLN	C-N-CD	5.57	140.09	128.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	185	PRO	CB-CA-C	-5.41	98.48	112.00
1	D	98	LEU	CA-CB-CG	5.34	127.59	115.30
1	A	104	GLU	N-CA-C	-5.34	96.57	111.00
1	D	184	LEU	CB-CG-CD1	-5.26	102.06	111.00

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	2193	0	2223	41	0
1	B	2197	0	2227	58	0
1	C	2030	0	2068	49	3
1	D	2043	0	2083	84	3
2	A	25	0	0	0	0
2	B	36	0	0	1	0
2	C	22	0	0	0	0
2	D	23	0	0	2	0
All	All	8569	0	8601	220	3

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 (220) 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:D:159:ARG:NH1	2:D:301:HOH:O	1.61	1.30
1:D:196:LEU:O	1:D:258:LEU:HD11	1.19	1.27
1:D:4:VAL:O	1:D:8:THR:HG23	1.40	1.21
1:D:196:LEU:O	1:D:258:LEU:CD1	2.05	1.04
1:B:161:THR:CG2	1:B:196:LEU:HB3	1.89	1.02
1:D:103:ARG:HB2	1:D:106:LYS:HD2	1.39	1.02
1:B:161:THR:HG23	1:B:196:LEU:CB	1.91	1.00
1:D:84:GLN:HE22	1:D:102:PHE:H	1.10	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:THR:HG23	1:B:196:LEU:HB3	0.95	0.94
1:C:196:LEU:O	1:C:258:LEU:HD11	1.70	0.92
1:D:103:ARG:CB	1:D:106:LYS:HD2	2.00	0.90
1:B:197:ASN:C	1:B:198:LYS:HG2	1.90	0.90
1:B:103:ARG:CG	1:B:181:ALA:HB2	2.03	0.89
1:B:196:LEU:HD22	1:D:135:GLN:HG2	1.55	0.88
1:D:137:LYS:H	1:D:137:LYS:HD2	1.36	0.88
1:D:136:PHE:CE2	1:D:137:LYS:O	2.28	0.87
1:D:4:VAL:O	1:D:8:THR:CG2	2.23	0.86
1:A:161:THR:HG23	1:A:196:LEU:HD23	1.61	0.82
1:A:106:LYS:HD2	1:A:107:GLN:H	1.46	0.81
1:C:103:ARG:HG3	1:C:181:ALA:HB2	1.62	0.81
1:C:17:GLU:OE1	1:C:20:ARG:NH2	2.14	0.80
1:B:195:ASN:ND2	1:B:200:CYS:SG	2.54	0.79
1:B:197:ASN:O	1:B:198:LYS:HG2	1.83	0.79
1:A:84:GLN:OE1	1:A:102:PHE:HB2	1.83	0.79
1:C:197:ASN:HB3	1:C:258:LEU:HD12	1.62	0.79
1:B:103:ARG:HG3	1:B:181:ALA:HB2	1.62	0.78
1:C:4:VAL:O	1:C:8:THR:HG23	1.85	0.76
1:A:115:ASP:OD2	1:A:159:ARG:NH1	2.17	0.76
1:C:197:ASN:HB3	1:C:258:LEU:CD1	2.15	0.76
1:B:84:GLN:OE1	1:B:102:PHE:HB2	1.86	0.74
1:D:102:PHE:HE1	1:D:180:PHE:CZ	2.05	0.74
1:D:100:PHE:CD1	1:D:184:LEU:HD11	2.23	0.74
1:C:198:LYS:CG	1:C:250:VAL:HA	2.19	0.72
1:B:198:LYS:NZ	1:D:25:ASP:OD1	2.23	0.71
1:A:161:THR:O	1:A:195:ASN:O	2.07	0.70
1:D:137:LYS:N	1:D:137:LYS:HD2	2.06	0.70
1:C:207:HIS:CE1	1:C:242:SER:HB3	2.27	0.70
1:B:196:LEU:HD22	1:D:135:GLN:CG	2.21	0.70
1:A:82:LEU:HB3	1:A:108:TRP:HB3	1.73	0.70
1:A:82:LEU:O	1:A:102:PHE:HE2	1.76	0.69
1:D:137:LYS:NZ	1:D:142:VAL:HG22	2.07	0.69
1:B:66:CYS:O	1:B:90:LYS:NZ	2.25	0.69
1:C:205:GLN:O	1:C:205:GLN:HG3	1.92	0.68
1:D:184:LEU:HB3	1:D:185:PRO:HD2	1.76	0.67
1:A:65:SER:HB2	1:A:68:THR:HG22	1.74	0.67
1:B:82:LEU:O	1:B:102:PHE:HE2	1.75	0.67
1:B:197:ASN:OD1	1:B:255:ASN:OD1	2.11	0.67
1:D:137:LYS:HZ2	1:D:142:VAL:CG2	2.07	0.67
1:C:198:LYS:HG3	1:C:250:VAL:HA	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:130:ARG:NH1	1:C:141:GLU:OE1	2.29	0.66
1:A:160:LEU:HD21	1:A:254:LEU:HD22	1.77	0.65
1:D:137:LYS:HZ3	1:D:142:VAL:HG22	1.61	0.65
1:D:155:ARG:O	1:D:159:ARG:HG2	1.97	0.65
1:B:198:LYS:HD3	1:B:250:VAL:HG22	1.79	0.65
1:D:137:LYS:NZ	1:D:142:VAL:CG2	2.60	0.64
1:D:47:LEU:HB3	1:D:57:LYS:HE3	1.79	0.64
1:D:24:HIS:HB2	1:D:136:PHE:HB2	1.78	0.64
1:D:113:ILE:HD11	1:D:194:ILE:HD11	1.80	0.64
1:A:106:LYS:HD2	1:A:107:GLN:N	2.11	0.64
1:C:130:ARG:NH2	1:C:134:TYR:CE1	2.66	0.64
1:C:198:LYS:HG2	1:C:249:CYS:C	2.18	0.63
1:B:193:TYR:O	1:B:199:LEU:HD23	1.98	0.63
1:D:113:ILE:HG13	1:D:194:ILE:CD1	2.31	0.61
1:A:160:LEU:O	1:A:160:LEU:HD13	2.01	0.61
1:C:198:LYS:HG2	1:C:250:VAL:HA	1.82	0.60
1:C:198:LYS:HG2	1:C:250:VAL:N	2.16	0.60
1:B:196:LEU:HD12	1:B:196:LEU:O	2.01	0.60
1:B:100:PHE:CD2	1:B:184:LEU:HD23	2.37	0.59
1:C:166:LEU:O	1:C:226:HIS:NE2	2.30	0.59
1:D:103:ARG:CB	1:D:106:LYS:CD	2.78	0.59
1:B:200:CYS:HB3	1:B:246:LYS:NZ	2.18	0.59
1:A:103:ARG:HG3	1:A:179:MET:O	2.03	0.58
1:C:98:LEU:HD12	1:C:99:HIS:H	1.67	0.58
1:D:102:PHE:HZ	1:D:176:LEU:HD11	1.68	0.58
1:C:195:ASN:O	1:C:196:LEU:HB3	2.05	0.57
1:D:206:LEU:HD13	1:D:239:LEU:HD23	1.87	0.57
1:B:196:LEU:HD13	1:D:21:TRP:HH2	1.69	0.57
1:D:113:ILE:CG1	1:D:194:ILE:CD1	2.83	0.57
1:C:205:GLN:NE2	1:C:243:HIS:CE1	2.73	0.57
1:C:115:ASP:HB2	1:C:159:ARG:HG2	1.86	0.56
1:D:113:ILE:HG13	1:D:194:ILE:HD12	1.87	0.56
1:D:195:ASN:O	1:D:196:LEU:HB2	2.05	0.56
1:B:196:LEU:HD13	1:D:21:TRP:CH2	2.39	0.56
1:D:102:PHE:HE1	1:D:180:PHE:CE1	2.23	0.56
1:D:5:MET:O	1:D:8:THR:OG1	2.25	0.55
1:A:31:LEU:O	1:A:35:GLN:HG3	2.06	0.55
1:B:194:ILE:CD1	1:B:199:LEU:HG	2.37	0.55
1:B:196:LEU:CD1	1:D:21:TRP:CH2	2.90	0.54
1:D:84:GLN:NE2	1:D:102:PHE:H	1.93	0.54
1:B:103:ARG:HG2	1:B:181:ALA:HB2	1.84	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:199:LEU:HB2	1:C:251:ILE:HD11	1.89	0.54
1:B:82:LEU:O	1:B:102:PHE:CE2	2.59	0.53
1:B:200:CYS:HB3	1:B:246:LYS:HZ2	1.74	0.53
1:C:98:LEU:HG	1:C:100:PHE:CE1	2.43	0.53
1:D:102:PHE:HZ	1:D:176:LEU:CD1	2.21	0.53
1:B:84:GLN:HA	1:B:102:PHE:CD2	2.42	0.53
1:C:5:MET:HA	1:C:8:THR:OG1	2.09	0.53
1:C:198:LYS:HG2	1:C:250:VAL:CA	2.39	0.53
1:A:160:LEU:HD21	1:A:254:LEU:CD2	2.39	0.52
1:D:166:LEU:O	1:D:226:HIS:NE2	2.36	0.52
1:D:103:ARG:HG2	1:D:106:LYS:HE3	1.90	0.52
1:A:46:THR:HG22	1:A:117:ARG:HH21	1.75	0.52
1:D:100:PHE:CE1	1:D:184:LEU:HD11	2.44	0.52
1:B:83:SER:O	1:B:102:PHE:HD2	1.93	0.52
1:B:110:LEU:HD13	1:B:194:ILE:HD11	1.92	0.52
1:D:102:PHE:CE2	1:D:179:MET:HE2	2.45	0.51
1:D:102:PHE:CD2	1:D:179:MET:HE3	2.44	0.51
1:D:16:GLU:O	1:D:20:ARG:HG3	2.11	0.51
1:B:46:THR:HG22	1:B:117:ARG:HH22	1.75	0.51
1:D:137:LYS:H	1:D:137:LYS:CD	2.13	0.51
1:C:234:TRP:O	1:C:237:GLN:NE2	2.44	0.51
1:C:207:HIS:HE1	1:C:242:SER:HB3	1.74	0.51
1:B:83:SER:O	1:B:102:PHE:CD2	2.63	0.51
1:D:103:ARG:HB3	1:D:106:LYS:HD2	1.89	0.51
1:D:91:MET:HE3	1:D:98:LEU:HG	1.93	0.51
1:C:139:GLY:HA2	1:C:142:VAL:HG23	1.93	0.50
1:A:160:LEU:HD12	1:A:258:LEU:HD21	1.92	0.50
1:A:86:ASP:OD1	1:A:99:HIS:HE1	1.94	0.50
1:D:102:PHE:CE1	1:D:180:PHE:CE1	3.00	0.50
1:B:200:CYS:CB	1:B:246:LYS:HZ1	2.24	0.50
1:C:106:LYS:HD2	1:C:107:GLN:H	1.76	0.50
1:A:233:GLU:HB3	1:A:238:ARG:HG2	1.93	0.49
1:A:82:LEU:O	1:A:102:PHE:CE2	2.62	0.49
1:D:113:ILE:CD1	1:D:194:ILE:HD11	2.41	0.49
1:B:14:VAL:HG12	1:B:17:GLU:H	1.76	0.49
1:D:71:VAL:O	1:D:247:VAL:HG11	2.13	0.49
1:B:200:CYS:SG	1:B:246:LYS:NZ	2.74	0.48
1:A:208:ALA:HA	1:A:239:LEU:HD23	1.96	0.48
1:C:205:GLN:CG	1:C:243:HIS:H	2.25	0.48
1:A:185:PRO:HB2	1:A:187:ASP:OD1	2.13	0.48
1:D:161:THR:OG1	1:D:196:LEU:HD21	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:184:LEU:HD22	1:C:185:PRO:HD2	1.96	0.48
1:C:205:GLN:NE2	1:C:242:SER:OG	2.46	0.48
1:B:161:THR:HA	1:B:196:LEU:HA	1.95	0.48
1:D:89:LEU:HD23	1:D:91:MET:HE2	1.95	0.48
1:D:1:MET:O	1:D:5:MET:HG3	2.14	0.47
1:C:197:ASN:HB3	1:C:258:LEU:HD11	1.94	0.47
1:B:17:GLU:HG2	1:B:21:TRP:CZ2	2.50	0.47
1:D:136:PHE:CD2	1:D:137:LYS:O	2.67	0.47
1:A:14:VAL:HG12	1:A:17:GLU:H	1.79	0.47
1:D:103:ARG:HB3	1:D:106:LYS:CD	2.44	0.47
1:B:196:LEU:CD1	1:D:21:TRP:HH2	2.26	0.47
1:C:62:ILE:HG22	1:C:74:VAL:HG22	1.95	0.47
1:C:206:LEU:HD23	1:C:241:VAL:HA	1.96	0.47
1:B:194:ILE:HG22	1:B:195:ASN:O	2.15	0.47
1:B:196:LEU:HD12	1:D:21:TRP:CZ3	2.50	0.47
1:C:100:PHE:CD1	1:C:184:LEU:HD11	2.50	0.46
1:C:134:TYR:OH	1:C:141:GLU:HB3	2.15	0.46
1:B:72:LYS:NZ	2:B:304:HOH:O	2.48	0.46
1:D:102:PHE:CD2	1:D:179:MET:CE	2.97	0.46
1:D:102:PHE:C	1:D:104:GLU:H	2.18	0.46
1:D:157:ARG:NH1	2:D:302:HOH:O	2.30	0.46
1:A:137:LYS:HG3	1:A:138:THR:H	1.80	0.46
1:B:185:PRO:HB2	1:B:187:ASP:OD1	2.16	0.46
1:B:200:CYS:CB	1:B:246:LYS:NZ	2.78	0.46
1:D:187:ASP:O	1:D:187:ASP:OD1	2.33	0.46
1:B:103:ARG:HG3	1:B:181:ALA:CB	2.41	0.46
1:B:46:THR:HG22	1:B:117:ARG:NH2	2.32	0.45
1:D:137:LYS:N	1:D:137:LYS:CD	2.73	0.45
1:D:91:MET:CE	1:D:98:LEU:HG	2.45	0.45
1:A:161:THR:CG2	1:A:196:LEU:HD23	2.40	0.45
1:A:116:ALA:O	1:A:120:VAL:HG23	2.16	0.45
1:A:17:GLU:HG2	1:A:21:TRP:CZ2	2.52	0.45
1:C:1:MET:O	1:C:5:MET:HG3	2.16	0.45
1:C:90:LYS:O	1:C:90:LYS:HG3	2.14	0.45
1:B:196:LEU:CD1	1:D:21:TRP:CZ3	3.00	0.45
1:B:116:ALA:O	1:B:120:VAL:HG23	2.16	0.45
1:C:112:GLN:NE2	1:C:194:ILE:O	2.43	0.45
1:C:205:GLN:HG3	1:C:243:HIS:H	1.81	0.45
1:A:40:GLU:HG2	1:A:44:ARG:HH12	1.83	0.44
1:D:103:ARG:HB3	1:D:106:LYS:CG	2.47	0.44
1:A:259:VAL:HG13	1:C:17:GLU:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:VAL:HG22	1:A:89:LEU:HD22	2.00	0.44
1:C:81:ALA:HB1	1:C:107:GLN:HB2	2.00	0.44
1:C:160:LEU:HD23	1:C:160:LEU:HA	1.79	0.44
1:D:40:GLU:O	1:D:44:ARG:HG2	2.18	0.44
1:A:119:HIS:HB2	1:A:156:ALA:HB2	2.00	0.44
1:B:135:GLN:HB3	1:B:285:ARG:CZ	2.48	0.44
1:D:103:ARG:O	1:D:105:ASP:N	2.46	0.44
1:B:37:ILE:HD11	1:D:10:ALA:HB1	2.00	0.44
1:D:196:LEU:C	1:D:258:LEU:HD11	2.20	0.44
1:B:201:LEU:HB3	1:B:247:VAL:HG23	1.99	0.43
1:B:251:ILE:HB	1:B:254:LEU:HB2	2.00	0.43
1:C:5:MET:HE2	1:C:5:MET:HB2	1.77	0.43
1:A:106:LYS:NZ	1:A:179:MET:SD	2.88	0.43
1:D:62:ILE:HG22	1:D:74:VAL:HG22	1.99	0.43
1:A:125:TYR:O	1:A:129:SER:OG	2.37	0.43
1:A:210:GLN:HB2	1:A:212:ASN:OD1	2.18	0.43
1:B:84:GLN:CD	1:B:102:PHE:HB2	2.39	0.43
1:B:47:LEU:H	1:B:47:LEU:HG	1.73	0.43
1:B:86:ASP:OD1	1:B:99:HIS:CE1	2.72	0.43
1:A:160:LEU:HD12	1:A:258:LEU:CD2	2.49	0.42
1:C:246:LYS:HB2	1:C:246:LYS:HE2	1.74	0.42
1:D:254:LEU:HD23	1:D:254:LEU:HA	1.86	0.42
1:B:104:GLU:HB3	1:B:105:ASP:H	1.58	0.42
1:D:103:ARG:HB3	1:D:106:LYS:HG3	2.01	0.42
1:A:86:ASP:OD1	1:A:99:HIS:CE1	2.72	0.42
1:D:103:ARG:HA	1:D:103:ARG:HD3	1.82	0.42
1:D:103:ARG:C	1:D:105:ASP:H	2.23	0.42
1:B:85:ALA:HB3	1:B:102:PHE:CE1	2.55	0.42
1:A:195:ASN:HB2	1:A:200:CYS:SG	2.60	0.41
1:A:234:TRP:O	1:A:236:SER:N	2.53	0.41
1:C:75:LEU:HB3	1:C:82:LEU:HD11	2.02	0.41
1:A:251:ILE:HB	1:A:254:LEU:HB2	2.03	0.41
1:D:113:ILE:CG1	1:D:194:ILE:HD11	2.50	0.41
1:D:160:LEU:HD23	1:D:160:LEU:HA	1.79	0.41
1:A:135:GLN:HB3	1:A:285:ARG:NH1	2.35	0.41
1:D:38:LEU:HD23	1:D:38:LEU:HA	1.88	0.41
1:D:44:ARG:HD3	1:D:44:ARG:HA	1.86	0.41
1:D:5:MET:C	1:D:8:THR:HG1	2.24	0.41
1:A:85:ALA:HB3	1:A:102:PHE:CE1	2.56	0.41
1:C:63:LEU:O	1:C:72:LYS:HD3	2.20	0.41
1:C:35:GLN:HG2	1:C:127:LEU:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:ASN:O	1:B:198:LYS:CG	2.63	0.41
1:D:11:GLU:O	1:D:15:LEU:HD13	2.20	0.40
1:C:75:LEU:HD23	1:C:82:LEU:HD11	2.03	0.40
1:A:135:GLN:HB3	1:A:285:ARG:CZ	2.50	0.40
1:B:196:LEU:HD12	1:D:21:TRP:CH2	2.56	0.40
1:D:84:GLN:NE2	1:D:85:ALA:H	2.19	0.40
1:D:89:LEU:O	1:D:97:LEU:HA	2.22	0.40

All (3) 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:C:8:THR:CG2	1:D:8:THR:CG2[2_665]	1.70	0.50
1:C:8:THR:CG2	1:D:8:THR:CB[2_665]	1.71	0.49
1:C:8:THR:CB	1:D:8:THR:CB[2_665]	1.99	0.21

## 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	267/289 (92%)	261 (98%)	6 (2%)	0	100	100
1	B	270/289 (93%)	259 (96%)	11 (4%)	0	100	100
1	C	244/289 (84%)	231 (95%)	13 (5%)	0	100	100
1	D	243/289 (84%)	231 (95%)	12 (5%)	0	100	100
All	All	1024/1156 (89%)	982 (96%)	42 (4%)	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	239/248 (96%)	238 (100%)	1 (0%)	91	97
1	B	239/248 (96%)	237 (99%)	2 (1%)	81	94
1	C	222/248 (90%)	220 (99%)	2 (1%)	78	94
1	D	224/248 (90%)	223 (100%)	1 (0%)	91	97
All	All	924/992 (93%)	918 (99%)	6 (1%)	86	96

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	ARG
1	B	198	LYS
1	B	199	LEU
1	C	186	SER
1	C	243	HIS
1	D	137	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	35	GLN
1	C	197	ASN
1	C	205	GLN
1	C	207	HIS
1	C	237	GLN
1	C	243	HIS
1	C	255	ASN
1	D	84	GLN

### 5.3.3 RNA ⓘ

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 ⓘ

### 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, 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	275/289 (95%)	0.32	4 (1%) 73 68	37, 68, 122, 156	0
1	B	276/289 (95%)	0.35	6 (2%) 62 52	37, 70, 126, 156	0
1	C	254/289 (87%)	0.51	18 (7%) 16 9	36, 83, 132, 151	0
1	D	255/289 (88%)	0.47	9 (3%) 44 34	37, 82, 133, 159	0
All	All	1060/1156 (91%)	0.41	37 (3%) 44 34	36, 74, 130, 159	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	89	LEU	5.2
1	B	104	GLU	4.9
1	C	89	LEU	4.8
1	D	134	TYR	4.2
1	D	97	LEU	3.7
1	C	235	GLY	3.2
1	C	100	PHE	3.2
1	C	196	LEU	3.1
1	D	71	VAL	3.1
1	D	204	TYR	3.1
1	B	210	GLN	3.0
1	A	107	GLN	3.0
1	C	238	ARG	3.0
1	C	204	TYR	2.9
1	C	63	LEU	2.9
1	D	87	VAL	2.8
1	C	234	TRP	2.7
1	C	97	LEU	2.6
1	C	90	LYS	2.6
1	B	91	MET	2.5
1	B	217	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	195	ASN	2.4
1	C	87	VAL	2.4
1	A	91	MET	2.4
1	A	160	LEU	2.3
1	D	98	LEU	2.3
1	C	91	MET	2.3
1	C	98	LEU	2.3
1	C	207	HIS	2.2
1	D	100	PHE	2.2
1	B	209	LEU	2.2
1	C	206	LEU	2.1
1	D	239	LEU	2.1
1	A	239	LEU	2.1
1	C	134	TYR	2.1
1	C	209	LEU	2.0
1	C	239	LEU	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 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.