



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 05:46 am BST

PDB ID : 4POF  
Title : PfMCM N-terminal domain without DNA  
Authors : Froelich, C.A.; Kang, S.; Epling, L.B.; Bell, S.P.; Enemark, E.J.  
Deposited on : 2014-02-25  
Resolution : 2.65 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

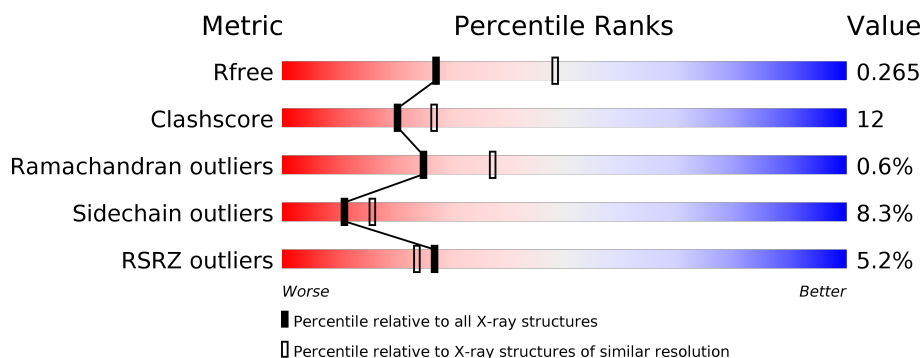
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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	1426 (2.66-2.62)
Clashscore	141614	1472 (2.66-2.62)
Ramachandran outliers	138981	1446 (2.66-2.62)
Sidechain outliers	138945	1446 (2.66-2.62)
RSRZ outliers	127900	1408 (2.66-2.62)

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	257	<div> <div>5%</div> <div> <div></div> <div>70%</div> <div>24%</div> <div>5%</div> </div> </div>
1	B	257	<div> <div>4%</div> <div> <div></div> <div>70%</div> <div>24%</div> <div>5%</div> </div> </div>
1	C	257	<div> <div>4%</div> <div> <div></div> <div>71%</div> <div>24%</div> <div>.</div> </div> </div>
1	D	257	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>21%</div> <div>5%</div> </div> </div>
1	E	257	<div> <div>7%</div> <div> <div></div> <div>72%</div> <div>23%</div> <div>.</div> </div> </div>
1	F	257	<div> <div>5%</div> <div> <div></div> <div>73%</div> <div>22%</div> <div>5%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12258 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 Cell division control protein 21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	0	0
			2042	1294	350	389	9			
1	B	255	Total	C	N	O	S	0	0	0
			2042	1294	350	389	9			
1	C	255	Total	C	N	O	S	0	0	0
			2042	1294	350	389	9			
1	D	255	Total	C	N	O	S	0	0	0
			2042	1294	350	389	9			
1	E	255	Total	C	N	O	S	0	0	0
			2042	1294	350	389	9			
1	F	255	Total	C	N	O	S	0	0	0
			2042	1294	350	389	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	EXPRESSION TAG	UNP Q8U3I4
A	1	VAL	-	EXPRESSION TAG	UNP Q8U3I4
B	0	SER	-	EXPRESSION TAG	UNP Q8U3I4
B	1	VAL	-	EXPRESSION TAG	UNP Q8U3I4
C	0	SER	-	EXPRESSION TAG	UNP Q8U3I4
C	1	VAL	-	EXPRESSION TAG	UNP Q8U3I4
D	0	SER	-	EXPRESSION TAG	UNP Q8U3I4
D	1	VAL	-	EXPRESSION TAG	UNP Q8U3I4
E	0	SER	-	EXPRESSION TAG	UNP Q8U3I4
E	1	VAL	-	EXPRESSION TAG	UNP Q8U3I4
F	0	SER	-	EXPRESSION TAG	UNP Q8U3I4
F	1	VAL	-	EXPRESSION TAG	UNP Q8U3I4

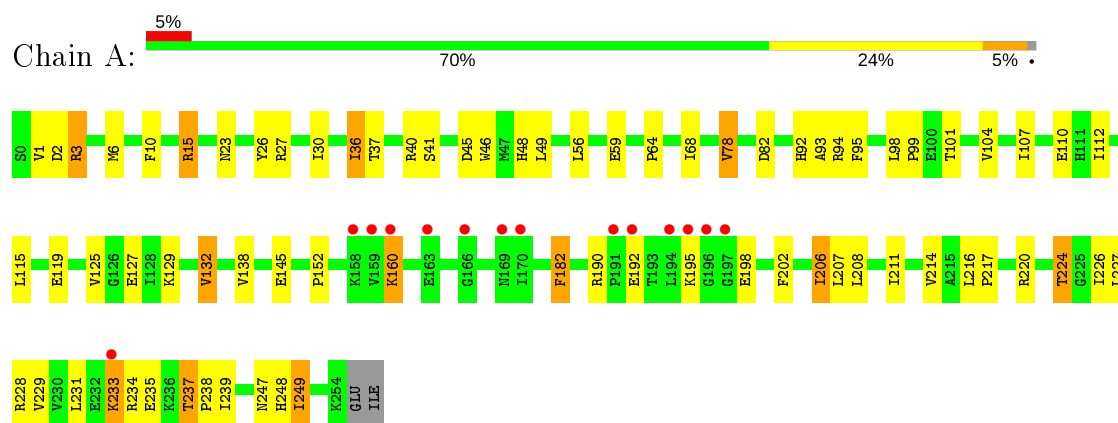
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total 1	Zn 1	0	0
2	E	1	Total 1	Zn 1	0	0
2	B	1	Total 1	Zn 1	0	0
2	C	1	Total 1	Zn 1	0	0
2	A	1	Total 1	Zn 1	0	0
2	F	1	Total 1	Zn 1	0	0

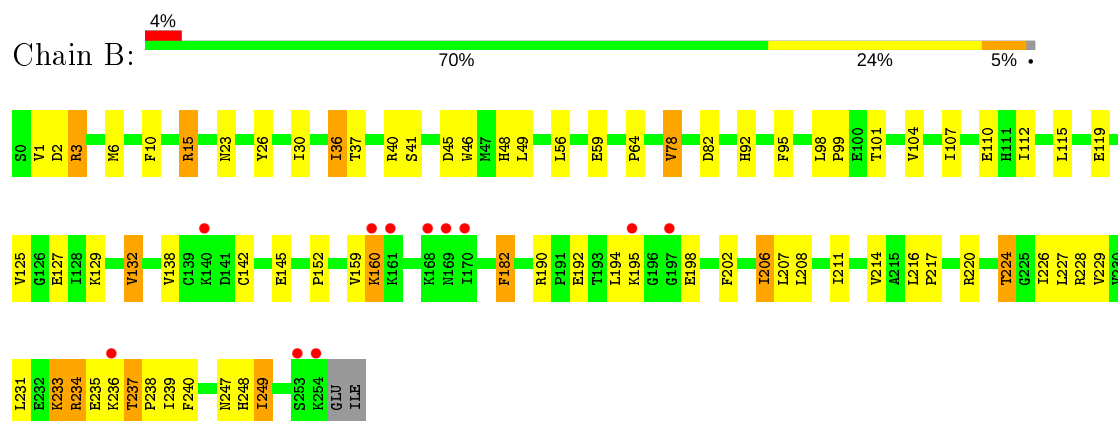
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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.

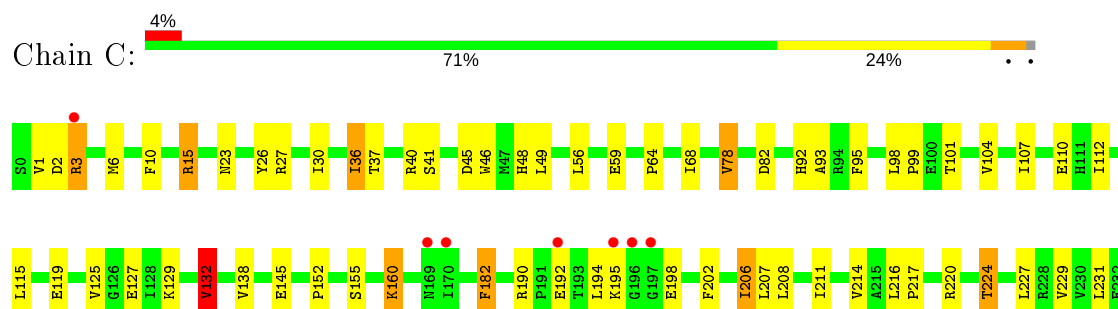
#### • Molecule 1: Cell division control protein 21

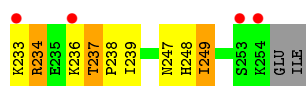


#### • Molecule 1: Cell division control protein 21

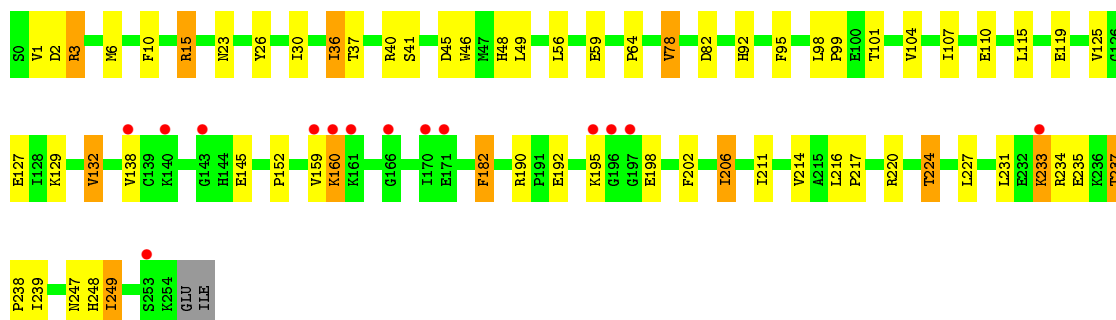
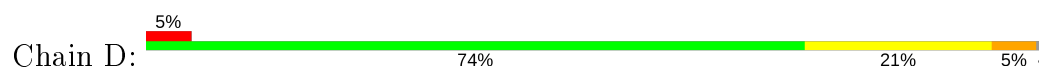


#### • Molecule 1: Cell division control protein 21

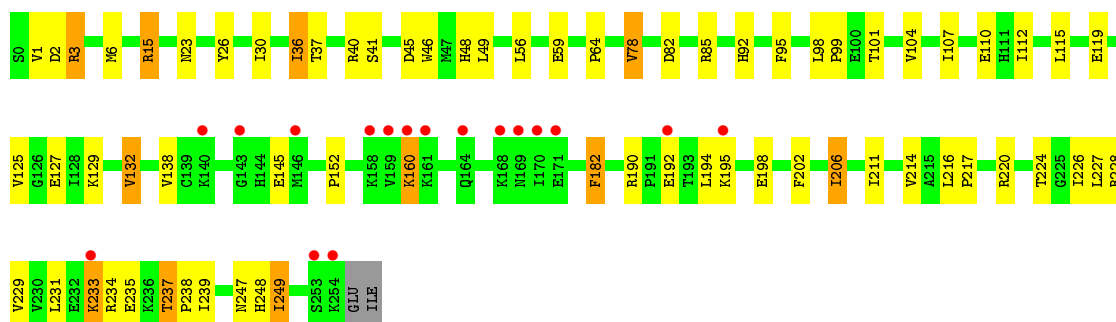




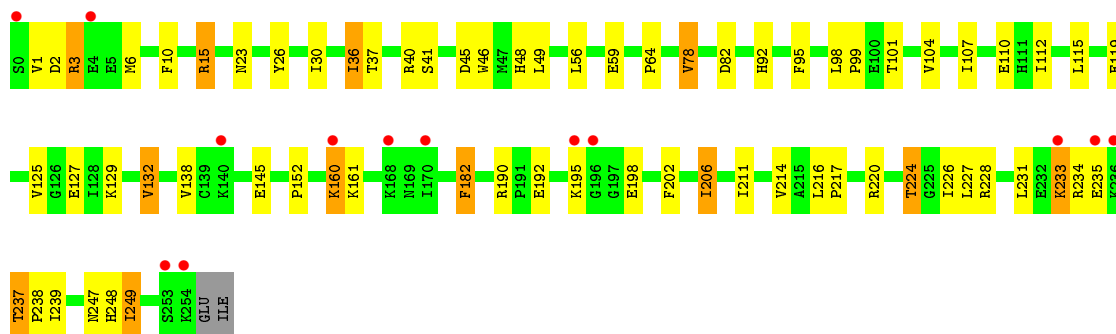
- Molecule 1: Cell division control protein 21



- Molecule 1: Cell division control protein 21



- Molecule 1: Cell division control protein 21



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.85Å 103.06Å 122.44Å 90.00° 119.85° 90.00°	Depositor
Resolution (Å)	47.33 – 2.65 47.33 – 2.65	Depositor EDS
% Data completeness (in resolution range)	98.6 (47.33-2.65) 98.6 (47.33-2.65)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.61 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, $R_{free}$	0.259 , 0.270 0.253 , 0.265	Depositor DCC
$R_{free}$ test set	3839 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.0	Xtriage
Anisotropy	1.153	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 34.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.022 for l,k,-h-l 0.022 for -h-l,k,h 0.000 for -h-l,-k,l 0.000 for h,-k,-h-l 0.000 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	12258	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.86 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2446e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.55	0/2077	0.69	1/2805 (0.0%)
1	B	0.52	0/2077	0.68	1/2805 (0.0%)
1	C	0.51	0/2077	0.68	1/2805 (0.0%)
1	D	0.49	0/2077	0.66	1/2805 (0.0%)
1	E	0.49	0/2077	0.66	1/2805 (0.0%)
1	F	0.57	0/2077	0.70	1/2805 (0.0%)
All	All	0.52	0/12462	0.68	6/16830 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	132	VAL	CB-CA-C	-6.20	99.62	111.40
1	E	132	VAL	CB-CA-C	-6.10	99.82	111.40
1	A	132	VAL	CB-CA-C	-5.90	100.20	111.40
1	C	132	VAL	CB-CA-C	-5.70	100.57	111.40
1	D	132	VAL	CB-CA-C	-5.66	100.64	111.40
1	F	132	VAL	CB-CA-C	-5.62	100.71	111.40

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	2042	0	2052	51	1
1	B	2042	0	2052	54	0
1	C	2042	0	2052	53	0
1	D	2042	0	2052	46	0
1	E	2042	0	2052	54	0
1	F	2042	0	2052	49	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
All	All	12258	0	12312	305	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 12.

All (305) 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:C:41:SER:HB3	1:C:92:HIS:HB2	1.46	0.97
1:B:211:ILE:HB	1:B:249:ILE:HD11	1.49	0.95
1:F:41:SER:HB3	1:F:92:HIS:HB2	1.52	0.92
1:B:41:SER:HB3	1:B:92:HIS:HB2	1.52	0.91
1:D:160:LYS:HD3	1:D:160:LYS:H	1.36	0.91
1:D:41:SER:HB3	1:D:92:HIS:HB2	1.51	0.91
1:E:160:LYS:H	1:E:160:LYS:HD3	1.37	0.90
1:C:160:LYS:HD3	1:C:160:LYS:H	1.36	0.90
1:E:211:ILE:HB	1:E:249:ILE:HD11	1.55	0.89
1:D:211:ILE:HB	1:D:249:ILE:HD11	1.54	0.88
1:E:41:SER:HB3	1:E:92:HIS:HB2	1.56	0.87
1:F:160:LYS:HD3	1:F:160:LYS:H	1.37	0.87
1:A:160:LYS:HD3	1:A:160:LYS:H	1.38	0.87
1:F:211:ILE:HB	1:F:249:ILE:HD11	1.56	0.86
1:C:211:ILE:HB	1:C:249:ILE:HD11	1.55	0.86
1:B:160:LYS:HD3	1:B:160:LYS:H	1.39	0.86
1:A:211:ILE:HB	1:A:249:ILE:HD11	1.58	0.85
1:A:41:SER:HB3	1:A:92:HIS:HB2	1.61	0.82
1:C:30:ILE:HD13	1:C:78:VAL:HG21	1.63	0.79
1:F:30:ILE:HD13	1:F:78:VAL:HG21	1.66	0.78
1:B:30:ILE:HD13	1:B:78:VAL:HG21	1.67	0.77
1:D:30:ILE:HD13	1:D:78:VAL:HG21	1.66	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:30:ILE:HD13	1:E:78:VAL:HG21	1.68	0.75
1:F:30:ILE:HD13	1:F:78:VAL:CG2	2.19	0.73
1:C:30:ILE:HD13	1:C:78:VAL:CG2	2.19	0.73
1:C:95:PHE:H	1:C:247:ASN:HD21	1.38	0.72
1:F:95:PHE:H	1:F:247:ASN:HD21	1.37	0.71
1:D:95:PHE:H	1:D:247:ASN:HD21	1.39	0.70
1:D:30:ILE:HD13	1:D:78:VAL:CG2	2.21	0.70
1:E:95:PHE:H	1:E:247:ASN:HD21	1.38	0.69
1:B:95:PHE:H	1:B:247:ASN:HD21	1.38	0.69
1:E:30:ILE:HD13	1:E:78:VAL:CG2	2.24	0.68
1:A:95:PHE:H	1:A:247:ASN:HD21	1.40	0.68
1:D:231:LEU:HD12	1:D:231:LEU:H	1.60	0.67
1:F:226:ILE:HG22	1:F:228:ARG:HD3	1.77	0.67
1:B:30:ILE:HD13	1:B:78:VAL:CG2	2.24	0.67
1:B:138:VAL:HG22	1:B:145:GLU:HG2	1.77	0.66
1:F:231:LEU:H	1:F:231:LEU:HD12	1.61	0.66
1:E:138:VAL:HG22	1:E:145:GLU:HG2	1.78	0.66
1:A:30:ILE:HD13	1:A:78:VAL:HG21	1.76	0.65
1:D:138:VAL:HG22	1:D:145:GLU:HG2	1.78	0.65
1:C:138:VAL:HG22	1:C:145:GLU:HG2	1.79	0.65
1:E:231:LEU:H	1:E:231:LEU:HD12	1.62	0.65
1:C:36:ILE:CD1	1:C:37:THR:H	2.10	0.64
1:A:138:VAL:HG22	1:A:145:GLU:HG2	1.78	0.64
1:C:231:LEU:HD12	1:C:231:LEU:H	1.62	0.64
1:D:40:ARG:HG3	1:D:40:ARG:HH21	1.63	0.64
1:F:36:ILE:CD1	1:F:37:THR:H	2.10	0.64
1:B:36:ILE:CD1	1:B:37:THR:H	2.10	0.63
1:F:138:VAL:HG22	1:F:145:GLU:HG2	1.80	0.63
1:A:30:ILE:HD13	1:A:78:VAL:CG2	2.28	0.63
1:E:36:ILE:CD1	1:E:37:THR:H	2.12	0.63
1:E:40:ARG:HH21	1:E:40:ARG:HG3	1.64	0.63
1:A:36:ILE:CD1	1:A:37:THR:H	2.12	0.62
1:E:190:ARG:HG2	1:E:192:GLU:HG2	1.81	0.62
1:B:190:ARG:HG2	1:B:192:GLU:HG2	1.81	0.62
1:F:190:ARG:HG2	1:F:192:GLU:HG2	1.81	0.62
1:C:190:ARG:HG2	1:C:192:GLU:HG2	1.82	0.62
1:B:49:LEU:HD11	1:B:56:LEU:HD23	1.82	0.62
1:A:195:LYS:HA	1:A:198:GLU:CB	2.30	0.61
1:D:36:ILE:CD1	1:D:37:THR:H	2.13	0.61
1:B:45:ASP:HB3	1:B:48:HIS:CD2	2.36	0.61
1:D:195:LYS:HA	1:D:198:GLU:CB	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:64:PRO:HB3	1:F:115:LEU:HB2	1.83	0.61
1:A:40:ARG:HG3	1:A:40:ARG:HH21	1.66	0.61
1:E:160:LYS:CD	1:E:160:LYS:H	2.11	0.61
1:C:195:LYS:HA	1:C:198:GLU:CB	2.31	0.61
1:D:160:LYS:H	1:D:160:LYS:CD	2.10	0.61
1:C:45:ASP:HB3	1:C:48:HIS:CD2	2.36	0.60
1:A:190:ARG:HG2	1:A:192:GLU:HG2	1.83	0.60
1:A:231:LEU:H	1:A:231:LEU:HD12	1.65	0.60
1:D:190:ARG:HG2	1:D:192:GLU:HG2	1.82	0.60
1:E:195:LYS:HA	1:E:198:GLU:CB	2.30	0.60
1:F:49:LEU:HD11	1:F:56:LEU:HD23	1.84	0.60
1:C:160:LYS:CD	1:C:160:LYS:H	2.10	0.60
1:A:226:ILE:HG22	1:A:228:ARG:HD3	1.84	0.60
1:A:45:ASP:HB3	1:A:48:HIS:CD2	2.37	0.60
1:B:95:PHE:H	1:B:247:ASN:ND2	1.99	0.59
1:F:195:LYS:HA	1:F:198:GLU:CB	2.32	0.59
1:B:195:LYS:HA	1:B:198:GLU:CB	2.33	0.59
1:B:231:LEU:H	1:B:231:LEU:HD12	1.67	0.59
1:C:49:LEU:HD11	1:C:56:LEU:HD23	1.83	0.59
1:C:40:ARG:HG3	1:C:40:ARG:HH21	1.68	0.59
1:B:226:ILE:HG22	1:B:228:ARG:HD3	1.85	0.59
1:F:40:ARG:HH21	1:F:40:ARG:HG3	1.67	0.59
1:F:41:SER:CB	1:F:92:HIS:HB2	2.31	0.58
1:D:95:PHE:H	1:D:247:ASN:ND2	2.01	0.58
1:E:45:ASP:HB3	1:E:48:HIS:CD2	2.38	0.58
1:F:95:PHE:H	1:F:247:ASN:ND2	2.02	0.58
1:E:95:PHE:H	1:E:247:ASN:ND2	2.01	0.58
1:C:41:SER:CB	1:C:92:HIS:HB2	2.30	0.58
1:D:49:LEU:HD11	1:D:56:LEU:HD23	1.85	0.58
1:D:64:PRO:HB3	1:D:115:LEU:HB2	1.86	0.57
1:A:49:LEU:HD11	1:A:56:LEU:HD23	1.87	0.57
1:C:26:TYR:OH	1:C:48:HIS:HD2	1.87	0.57
1:D:45:ASP:HB3	1:D:48:HIS:CD2	2.40	0.57
1:E:49:LEU:HD11	1:E:56:LEU:HD23	1.85	0.57
1:A:64:PRO:HB3	1:A:115:LEU:HB2	1.86	0.57
1:E:64:PRO:HB3	1:E:115:LEU:HB2	1.87	0.56
1:F:160:LYS:CD	1:F:160:LYS:H	2.11	0.56
1:B:64:PRO:HB3	1:B:115:LEU:HB2	1.87	0.56
1:C:95:PHE:H	1:C:247:ASN:ND2	2.03	0.56
1:B:95:PHE:O	1:B:224:THR:HG21	2.05	0.56
1:E:95:PHE:O	1:E:224:THR:HG21	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:LYS:CD	1:B:160:LYS:H	2.13	0.56
1:A:26:TYR:OH	1:A:48:HIS:HD2	1.88	0.56
1:D:41:SER:CB	1:D:92:HIS:HB2	2.29	0.56
1:A:95:PHE:H	1:A:247:ASN:ND2	2.04	0.56
1:C:64:PRO:HB3	1:C:115:LEU:HB2	1.88	0.55
1:B:40:ARG:HG3	1:B:40:ARG:HH21	1.71	0.55
1:B:41:SER:CB	1:B:92:HIS:HB2	2.31	0.55
1:F:214:VAL:HG11	1:F:249:ILE:HD13	1.88	0.55
1:D:26:TYR:OH	1:D:48:HIS:HD2	1.91	0.54
1:F:226:ILE:CG2	1:F:228:ARG:HD3	2.37	0.54
1:F:36:ILE:HD12	1:F:37:THR:H	1.72	0.54
1:A:98:LEU:HB3	1:A:99:PRO:HD2	1.89	0.53
1:E:224:THR:HG22	1:E:248:HIS:HB3	1.90	0.53
1:B:224:THR:HG22	1:B:248:HIS:HB3	1.90	0.53
1:B:98:LEU:HB3	1:B:99:PRO:HD2	1.91	0.53
1:E:41:SER:CB	1:E:92:HIS:HB2	2.34	0.53
1:D:98:LEU:HB3	1:D:99:PRO:HD2	1.89	0.53
1:E:36:ILE:HD12	1:E:37:THR:H	1.73	0.53
1:E:98:LEU:HB3	1:E:99:PRO:HD2	1.91	0.53
1:B:240:PHE:CE2	1:C:132:VAL:HG22	2.44	0.53
1:A:160:LYS:H	1:A:160:LYS:CD	2.13	0.52
1:F:3:ARG:HG2	1:F:3:ARG:HH21	1.74	0.52
1:F:26:TYR:OH	1:F:48:HIS:HD2	1.92	0.52
1:D:214:VAL:HG11	1:D:249:ILE:HD13	1.90	0.52
1:F:152:PRO:HB2	1:F:206:ILE:HD12	1.92	0.52
1:A:224:THR:HG22	1:A:248:HIS:HB3	1.92	0.52
1:A:237:THR:HG22	1:A:238:PRO:HD2	1.92	0.52
1:B:237:THR:HG22	1:B:238:PRO:HD2	1.90	0.51
1:C:237:THR:HG22	1:C:238:PRO:HD2	1.92	0.51
1:D:3:ARG:HH21	1:D:3:ARG:HG2	1.76	0.51
1:D:36:ILE:HD12	1:D:37:THR:H	1.74	0.51
1:B:214:VAL:HG11	1:B:249:ILE:HD13	1.91	0.51
1:C:98:LEU:HB3	1:C:99:PRO:HD2	1.92	0.51
1:B:3:ARG:HH21	1:B:3:ARG:HG2	1.76	0.51
1:C:30:ILE:CD1	1:C:78:VAL:HG21	2.35	0.51
1:D:30:ILE:CD1	1:D:78:VAL:HG21	2.40	0.51
1:E:226:ILE:HG22	1:E:228:ARG:HD3	1.92	0.51
1:C:95:PHE:O	1:C:224:THR:HG21	2.11	0.51
1:D:2:ASP:O	1:D:6:MET:HB2	2.11	0.51
1:F:195:LYS:HA	1:F:198:GLU:HB2	1.93	0.51
1:F:30:ILE:CD1	1:F:78:VAL:HG21	2.38	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:ILE:HD12	1:A:37:THR:H	1.75	0.50
1:A:3:ARG:HH21	1:A:3:ARG:HG2	1.76	0.50
1:E:195:LYS:HA	1:E:198:GLU:HB2	1.93	0.50
1:E:3:ARG:HG2	1:E:3:ARG:HH21	1.76	0.50
1:F:104:VAL:HA	1:F:107:ILE:CD1	2.41	0.50
1:F:125:VAL:HG22	1:F:217:PRO:HD3	1.93	0.50
1:A:195:LYS:HA	1:A:198:GLU:HB2	1.93	0.50
1:D:195:LYS:HA	1:D:198:GLU:HB2	1.92	0.50
1:E:214:VAL:HG11	1:E:249:ILE:HD13	1.94	0.50
1:C:214:VAL:HG11	1:C:249:ILE:HD13	1.93	0.50
1:D:224:THR:HG22	1:D:248:HIS:HB3	1.93	0.50
1:E:15:ARG:HH21	1:E:82:ASP:CG	2.15	0.50
1:A:1:VAL:HG23	1:A:59:GLU:HG3	1.93	0.50
1:C:104:VAL:HA	1:C:107:ILE:CD1	2.42	0.50
1:C:3:ARG:HH21	1:C:3:ARG:HG2	1.76	0.50
1:A:104:VAL:HA	1:A:107:ILE:CD1	2.41	0.49
1:C:195:LYS:HA	1:C:198:GLU:HB2	1.93	0.49
1:D:152:PRO:HB2	1:D:206:ILE:HD12	1.94	0.49
1:D:237:THR:HG22	1:D:238:PRO:HD2	1.94	0.49
1:A:127:GLU:HG3	1:A:129:LYS:HE3	1.93	0.49
1:B:26:TYR:OH	1:B:48:HIS:HD2	1.94	0.49
1:C:224:THR:HG22	1:C:248:HIS:HB3	1.94	0.49
1:A:95:PHE:O	1:A:224:THR:HG21	2.11	0.49
1:B:36:ILE:HD12	1:B:37:THR:H	1.75	0.49
1:E:182:PHE:CD2	1:E:182:PHE:C	2.86	0.49
1:F:46:TRP:CE3	1:F:99:PRO:HD3	2.47	0.49
1:A:152:PRO:HB2	1:A:206:ILE:HD12	1.95	0.49
1:E:237:THR:HG22	1:E:238:PRO:HD2	1.95	0.49
1:B:15:ARG:HH21	1:B:82:ASP:CG	2.16	0.49
1:D:95:PHE:O	1:D:224:THR:HG21	2.12	0.49
1:F:2:ASP:O	1:F:6:MET:HB2	2.12	0.49
1:E:127:GLU:HG3	1:E:129:LYS:HE3	1.95	0.49
1:C:36:ILE:HD12	1:C:37:THR:H	1.74	0.49
1:F:45:ASP:HB3	1:F:48:HIS:CD2	2.48	0.49
1:A:237:THR:HG22	1:A:239:ILE:HD13	1.94	0.49
1:B:182:PHE:CD2	1:B:182:PHE:C	2.85	0.49
1:C:2:ASP:O	1:C:6:MET:HB2	2.12	0.49
1:B:127:GLU:HG3	1:B:129:LYS:HE3	1.95	0.49
1:A:41:SER:CB	1:A:92:HIS:HB2	2.38	0.48
1:E:2:ASP:O	1:E:6:MET:HB2	2.13	0.48
1:B:195:LYS:HA	1:B:198:GLU:HB2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:95:PHE:O	1:F:224:THR:HG21	2.13	0.48
1:C:15:ARG:HH21	1:C:82:ASP:CG	2.17	0.48
1:F:127:GLU:HG3	1:F:129:LYS:HE3	1.95	0.48
1:A:15:ARG:HH21	1:A:82:ASP:CG	2.17	0.48
1:B:2:ASP:O	1:B:6:MET:HB2	2.14	0.48
1:D:1:VAL:HG23	1:D:59:GLU:HG3	1.95	0.48
1:E:26:TYR:OH	1:E:48:HIS:HD2	1.95	0.48
1:B:36:ILE:HD13	1:B:37:THR:H	1.79	0.48
1:C:182:PHE:CD2	1:C:182:PHE:C	2.86	0.48
1:F:237:THR:HG22	1:F:239:ILE:HD13	1.96	0.48
1:F:1:VAL:HG23	1:F:59:GLU:HG3	1.95	0.48
1:C:127:GLU:HG3	1:C:129:LYS:HE3	1.95	0.48
1:A:182:PHE:C	1:A:182:PHE:CD2	2.86	0.47
1:C:1:VAL:HG23	1:C:59:GLU:HG3	1.96	0.47
1:A:214:VAL:HG11	1:A:249:ILE:HD13	1.96	0.47
1:A:104:VAL:HA	1:A:107:ILE:HD11	1.96	0.47
1:D:125:VAL:HG22	1:D:217:PRO:HD3	1.97	0.47
1:E:1:VAL:HG23	1:E:59:GLU:HG3	1.97	0.47
1:E:237:THR:HG22	1:E:239:ILE:HD13	1.96	0.47
1:F:237:THR:HG22	1:F:238:PRO:HD2	1.96	0.47
1:C:125:VAL:HG22	1:C:217:PRO:HD3	1.96	0.47
1:D:15:ARG:HH21	1:D:82:ASP:CG	2.18	0.47
1:E:152:PRO:HB2	1:E:206:ILE:HD12	1.96	0.47
1:D:127:GLU:HG3	1:D:129:LYS:HE3	1.95	0.47
1:E:190:ARG:HD2	1:E:220:ARG:CZ	2.45	0.46
1:D:237:THR:HG22	1:D:239:ILE:HD13	1.97	0.46
1:F:224:THR:HG22	1:F:248:HIS:HB3	1.97	0.46
1:D:182:PHE:CD2	1:D:182:PHE:C	2.89	0.46
1:A:195:LYS:HA	1:A:198:GLU:HB3	1.98	0.46
1:F:104:VAL:HA	1:F:107:ILE:HD11	1.98	0.46
1:A:226:ILE:CG2	1:A:228:ARG:HD3	2.44	0.46
1:A:30:ILE:CD1	1:A:78:VAL:HG21	2.46	0.46
1:B:1:VAL:HG23	1:B:59:GLU:HG3	1.97	0.46
1:A:2:ASP:O	1:A:6:MET:HB2	2.15	0.46
1:D:160:LYS:HD3	1:D:160:LYS:N	2.18	0.46
1:E:30:ILE:CD1	1:E:78:VAL:HG21	2.42	0.46
1:F:182:PHE:C	1:F:182:PHE:CD2	2.90	0.46
1:E:112:ILE:HG12	1:E:229:VAL:HG13	1.97	0.46
1:C:36:ILE:HD13	1:C:37:THR:H	1.81	0.45
1:A:112:ILE:HG12	1:A:229:VAL:HG13	1.98	0.45
1:C:237:THR:HG22	1:C:239:ILE:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:98:LEU:HB3	1:F:99:PRO:HD2	1.99	0.45
1:D:224:THR:HG22	1:D:248:HIS:N	2.32	0.45
1:F:202:PHE:C	1:F:202:PHE:CD2	2.90	0.45
1:F:15:ARG:HH21	1:F:82:ASP:CG	2.19	0.45
1:A:190:ARG:HD2	1:A:220:ARG:CZ	2.47	0.45
1:B:226:ILE:CG2	1:B:228:ARG:HD3	2.46	0.45
1:A:94:ARG:HH11	1:A:94:ARG:HG2	1.81	0.45
1:B:190:ARG:HD2	1:B:220:ARG:CZ	2.47	0.45
1:C:104:VAL:HA	1:C:107:ILE:HD11	1.98	0.45
1:A:125:VAL:HG22	1:A:217:PRO:HD3	2.00	0.44
1:D:190:ARG:HD2	1:D:220:ARG:CZ	2.46	0.44
1:E:104:VAL:HA	1:E:107:ILE:CD1	2.47	0.44
1:A:46:TRP:CE3	1:A:99:PRO:HD3	2.52	0.44
1:B:125:VAL:HG22	1:B:217:PRO:HD3	1.99	0.44
1:E:224:THR:HG22	1:E:248:HIS:N	2.33	0.44
1:D:233:LYS:O	1:D:235:GLU:HG2	2.18	0.44
1:C:202:PHE:CD2	1:C:202:PHE:C	2.91	0.44
1:B:30:ILE:CD1	1:B:78:VAL:HG21	2.43	0.43
1:C:194:LEU:HB3	1:C:195:LYS:H	1.47	0.43
1:D:104:VAL:HA	1:D:107:ILE:CD1	2.48	0.43
1:B:104:VAL:HA	1:B:107:ILE:CD1	2.48	0.43
1:F:224:THR:HG22	1:F:248:HIS:N	2.33	0.43
1:B:224:THR:HG22	1:B:248:HIS:N	2.33	0.43
1:B:195:LYS:HA	1:B:198:GLU:HB3	2.01	0.43
1:C:46:TRP:CE3	1:C:99:PRO:HD3	2.54	0.43
1:C:224:THR:HG22	1:C:248:HIS:N	2.33	0.43
1:B:237:THR:HG22	1:B:239:ILE:HD13	1.99	0.43
1:E:125:VAL:HG22	1:E:217:PRO:HD3	1.99	0.43
1:A:224:THR:HG22	1:A:248:HIS:N	2.34	0.43
1:D:46:TRP:CE3	1:D:99:PRO:HD3	2.53	0.43
1:E:112:ILE:HG12	1:E:229:VAL:CG1	2.49	0.43
1:E:195:LYS:HA	1:E:198:GLU:HB3	1.99	0.43
1:A:207:LEU:O	1:A:208:LEU:HD23	2.19	0.42
1:D:202:PHE:C	1:D:202:PHE:CD2	2.92	0.42
1:F:195:LYS:HA	1:F:198:GLU:HB3	2.01	0.42
1:B:152:PRO:HB2	1:B:206:ILE:HD12	2.02	0.42
1:A:233:LYS:O	1:A:235:GLU:HG2	2.19	0.42
1:B:194:LEU:HB3	1:B:195:LYS:H	1.49	0.42
1:E:46:TRP:CE3	1:E:99:PRO:HD3	2.54	0.42
1:F:233:LYS:O	1:F:235:GLU:HG2	2.20	0.42
1:B:233:LYS:O	1:B:235:GLU:HG2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:194:LEU:HB3	1:E:195:LYS:H	1.47	0.42
1:C:112:ILE:HG12	1:C:229:VAL:HG13	2.02	0.42
1:B:236:LYS:O	1:C:155:SER:HB2	2.20	0.42
1:E:202:PHE:C	1:E:202:PHE:CD2	2.93	0.42
1:F:190:ARG:HD2	1:F:220:ARG:CZ	2.50	0.42
1:C:190:ARG:HD2	1:C:220:ARG:CZ	2.49	0.41
1:A:68:ILE:HG23	1:A:93:ALA:HB3	2.02	0.41
1:B:112:ILE:HG12	1:B:229:VAL:HG13	2.02	0.41
1:B:224:THR:CG2	1:B:248:HIS:HB3	2.50	0.41
1:C:207:LEU:O	1:C:208:LEU:HD23	2.20	0.41
1:D:195:LYS:HA	1:D:198:GLU:HB3	1.98	0.41
1:E:104:VAL:HA	1:E:107:ILE:HD11	2.03	0.41
1:E:233:LYS:O	1:E:235:GLU:HG2	2.20	0.41
1:F:125:VAL:CG2	1:F:217:PRO:HD3	2.50	0.41
1:E:160:LYS:N	1:E:160:LYS:HD3	2.19	0.41
1:C:234:ARG:HD2	1:C:234:ARG:H	1.85	0.41
1:C:234:ARG:HG2	1:C:236:LYS:H	1.86	0.41
1:F:160:LYS:HD3	1:F:160:LYS:N	2.19	0.41
1:C:125:VAL:CG2	1:C:217:PRO:HD3	2.51	0.41
1:E:237:THR:HG22	1:E:239:ILE:CD1	2.51	0.41
1:B:202:PHE:C	1:B:202:PHE:CD2	2.94	0.41
1:C:195:LYS:HA	1:C:198:GLU:HB3	2.00	0.41
1:D:36:ILE:HD13	1:D:37:THR:H	1.85	0.41
1:A:36:ILE:HD13	1:A:37:THR:H	1.83	0.41
1:C:152:PRO:HB2	1:C:206:ILE:HD12	2.02	0.41
1:A:202:PHE:CD2	1:A:202:PHE:C	2.94	0.40
1:B:46:TRP:CE3	1:B:99:PRO:HD3	2.56	0.40
1:E:224:THR:CG2	1:E:248:HIS:HB3	2.50	0.40
1:F:36:ILE:HD13	1:F:37:THR:H	1.83	0.40
1:B:104:VAL:HA	1:B:107:ILE:HD11	2.03	0.40
1:E:36:ILE:HD13	1:E:37:THR:H	1.85	0.40
1:E:226:ILE:CG2	1:E:228:ARG:HD3	2.51	0.40
1:B:207:LEU:O	1:B:208:LEU:HD23	2.22	0.40
1:D:224:THR:CG2	1:D:248:HIS:HB3	2.51	0.40
1:E:85:ARG:HA	1:E:85:ARG:HD3	1.97	0.40
1:B:234:ARG:HG2	1:B:236:LYS:H	1.85	0.40
1:C:68:ILE:HG23	1:C:93:ALA:HB3	2.04	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:192:GLU:O	1:F:161:LYS:NZ[2_545]	1.98	0.22

## 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	253/257 (98%)	237 (94%)	15 (6%)	1 (0%)	34	48
1	B	253/257 (98%)	236 (93%)	15 (6%)	2 (1%)	19	28
1	C	253/257 (98%)	235 (93%)	17 (7%)	1 (0%)	34	48
1	D	253/257 (98%)	235 (93%)	16 (6%)	2 (1%)	19	28
1	E	253/257 (98%)	237 (94%)	15 (6%)	1 (0%)	34	48
1	F	253/257 (98%)	233 (92%)	18 (7%)	2 (1%)	19	28
All	All	1518/1542 (98%)	1413 (93%)	96 (6%)	9 (1%)	25	37

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	LYS
1	B	233	LYS
1	C	233	LYS
1	D	233	LYS
1	E	233	LYS
1	F	233	LYS
1	F	112	ILE
1	B	159	VAL
1	D	159	VAL

### 5.3.2 Protein sidechains [i](#)

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	230/233 (99%)	210 (91%)	20 (9%)	10	14
1	B	230/233 (99%)	210 (91%)	20 (9%)	10	14
1	C	230/233 (99%)	210 (91%)	20 (9%)	10	14
1	D	230/233 (99%)	211 (92%)	19 (8%)	11	16
1	E	230/233 (99%)	213 (93%)	17 (7%)	13	21
1	F	230/233 (99%)	211 (92%)	19 (8%)	11	16
All	All	1380/1398 (99%)	1265 (92%)	115 (8%)	11	16

All (115) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ARG
1	A	10	PHE
1	A	15	ARG
1	A	23	ASN
1	A	27	ARG
1	A	36	ILE
1	A	78	VAL
1	A	101	THR
1	A	110	GLU
1	A	119	GLU
1	A	132	VAL
1	A	160	LYS
1	A	182	PHE
1	A	206	ILE
1	A	216	LEU
1	A	224	THR
1	A	227	LEU
1	A	234	ARG
1	A	237	THR
1	A	249	ILE
1	B	3	ARG
1	B	10	PHE
1	B	15	ARG
1	B	23	ASN
1	B	36	ILE
1	B	78	VAL
1	B	101	THR

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Mol	Chain	Res	Type
1	B	110	GLU
1	B	119	GLU
1	B	132	VAL
1	B	142	CYS
1	B	160	LYS
1	B	182	PHE
1	B	206	ILE
1	B	216	LEU
1	B	224	THR
1	B	227	LEU
1	B	234	ARG
1	B	237	THR
1	B	249	ILE
1	C	3	ARG
1	C	10	PHE
1	C	15	ARG
1	C	23	ASN
1	C	27	ARG
1	C	36	ILE
1	C	78	VAL
1	C	101	THR
1	C	110	GLU
1	C	119	GLU
1	C	132	VAL
1	C	160	LYS
1	C	182	PHE
1	C	206	ILE
1	C	216	LEU
1	C	224	THR
1	C	227	LEU
1	C	234	ARG
1	C	237	THR
1	C	249	ILE
1	D	3	ARG
1	D	10	PHE
1	D	15	ARG
1	D	23	ASN
1	D	36	ILE
1	D	78	VAL
1	D	101	THR
1	D	110	GLU
1	D	119	GLU

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Mol	Chain	Res	Type
1	D	132	VAL
1	D	160	LYS
1	D	182	PHE
1	D	206	ILE
1	D	216	LEU
1	D	224	THR
1	D	227	LEU
1	D	234	ARG
1	D	237	THR
1	D	249	ILE
1	E	3	ARG
1	E	15	ARG
1	E	23	ASN
1	E	36	ILE
1	E	78	VAL
1	E	101	THR
1	E	110	GLU
1	E	119	GLU
1	E	132	VAL
1	E	160	LYS
1	E	182	PHE
1	E	206	ILE
1	E	216	LEU
1	E	227	LEU
1	E	234	ARG
1	E	237	THR
1	E	249	ILE
1	F	3	ARG
1	F	10	PHE
1	F	15	ARG
1	F	23	ASN
1	F	36	ILE
1	F	78	VAL
1	F	101	THR
1	F	110	GLU
1	F	119	GLU
1	F	132	VAL
1	F	160	LYS
1	F	182	PHE
1	F	206	ILE
1	F	216	LEU
1	F	224	THR

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Mol	Chain	Res	Type
1	F	227	LEU
1	F	234	ARG
1	F	237	THR
1	F	249	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	A	48	HIS
1	A	63	ASN
1	A	117	GLN
1	A	188	GLN
1	A	247	ASN
1	B	12	ASN
1	B	48	HIS
1	B	63	ASN
1	B	117	GLN
1	B	188	GLN
1	B	247	ASN
1	C	12	ASN
1	C	48	HIS
1	C	63	ASN
1	C	117	GLN
1	C	188	GLN
1	C	247	ASN
1	D	12	ASN
1	D	48	HIS
1	D	63	ASN
1	D	117	GLN
1	D	188	GLN
1	D	247	ASN
1	E	12	ASN
1	E	48	HIS
1	E	63	ASN
1	E	117	GLN
1	E	188	GLN
1	E	247	ASN
1	F	12	ASN
1	F	48	HIS
1	F	63	ASN
1	F	117	GLN

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Mol	Chain	Res	Type
1	F	188	GLN
1	F	247	ASN

### 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](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	255/257 (99%)	0.28	14 (5%) 25 22	35, 65, 144, 162	0
1	B	255/257 (99%)	0.13	11 (4%) 35 31	39, 68, 135, 169	0
1	C	255/257 (99%)	0.22	11 (4%) 35 31	37, 69, 153, 170	0
1	D	255/257 (99%)	0.18	14 (5%) 25 22	40, 69, 159, 184	0
1	E	255/257 (99%)	0.25	17 (6%) 17 15	44, 73, 158, 182	0
1	F	255/257 (99%)	0.13	13 (5%) 28 24	35, 65, 142, 161	0
All	All	1530/1542 (99%)	0.20	80 (5%) 27 24	35, 68, 150, 184	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	254	LYS	10.0
1	D	233	LYS	7.0
1	A	197	GLY	5.8
1	C	253	SER	5.6
1	E	170	ILE	5.4
1	C	254	LYS	5.4
1	B	160	LYS	5.0
1	E	160	LYS	4.9
1	A	196	GLY	4.9
1	C	197	GLY	4.9
1	D	197	GLY	4.8
1	F	233	LYS	4.8
1	E	169	ASN	4.6
1	F	196	GLY	4.6
1	E	254	LYS	4.6
1	A	195	LYS	4.5
1	C	236	LYS	4.3
1	F	236	LYS	4.2
1	D	170	ILE	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	233	LYS	4.1
1	C	195	LYS	4.0
1	E	140	LYS	3.8
1	A	160	LYS	3.8
1	E	233	LYS	3.7
1	F	170	ILE	3.7
1	F	160	LYS	3.7
1	E	253	SER	3.7
1	D	253	SER	3.5
1	D	159	VAL	3.5
1	D	140	LYS	3.4
1	A	159	VAL	3.4
1	E	161	LYS	3.3
1	A	170	ILE	3.2
1	F	195	LYS	3.2
1	C	192	GLU	3.2
1	E	192	GLU	3.1
1	F	254	LYS	3.1
1	C	233	LYS	3.1
1	A	192	GLU	3.1
1	B	236	LYS	3.1
1	C	170	ILE	3.1
1	B	168	LYS	3.0
1	D	160	LYS	3.0
1	E	195	LYS	3.0
1	C	196	GLY	3.0
1	D	196	GLY	3.0
1	E	168	LYS	2.9
1	D	195	LYS	2.9
1	B	140	LYS	2.9
1	F	253	SER	2.8
1	E	164	GLN	2.8
1	B	169	ASN	2.7
1	E	171	GLU	2.6
1	D	166	GLY	2.6
1	F	168	LYS	2.6
1	B	195	LYS	2.5
1	B	161	LYS	2.5
1	F	140	LYS	2.5
1	A	163	GLU	2.4
1	F	235	GLU	2.4
1	B	197	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	143	GLY	2.4
1	B	170	ILE	2.3
1	A	194	LEU	2.3
1	E	159	VAL	2.3
1	F	0	SER	2.3
1	D	138	VAL	2.3
1	E	146	MET	2.2
1	D	171	GLU	2.2
1	C	169	ASN	2.1
1	A	166	GLY	2.1
1	D	161	LYS	2.1
1	B	253	SER	2.1
1	A	191	PRO	2.0
1	A	169	ASN	2.0
1	E	143	GLY	2.0
1	E	158	LYS	2.0
1	C	3	ARG	2.0
1	A	158	LYS	2.0
1	F	4	GLU	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	C	301	1/1	0.68	0.11	141,141,141,141	0
2	ZN	E	301	1/1	0.74	0.04	163,163,163,163	0
2	ZN	F	301	1/1	0.85	0.05	141,141,141,141	0
2	ZN	B	301	1/1	0.86	0.09	116,116,116,116	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	A	301	1/1	0.90	0.04	141,141,141,141	0
2	ZN	D	301	1/1	0.95	0.10	165,165,165,165	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.