



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 08:33 AM BST

PDB ID : 6MD3  
Title : Structure of T. brucei RRP44 PIN domain  
Authors : Guimaraes, B.G.; Cesaro, G.  
Deposited on : 2018-09-03  
Resolution : 2.29 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	:	2.13.1
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)	:	2.13.1

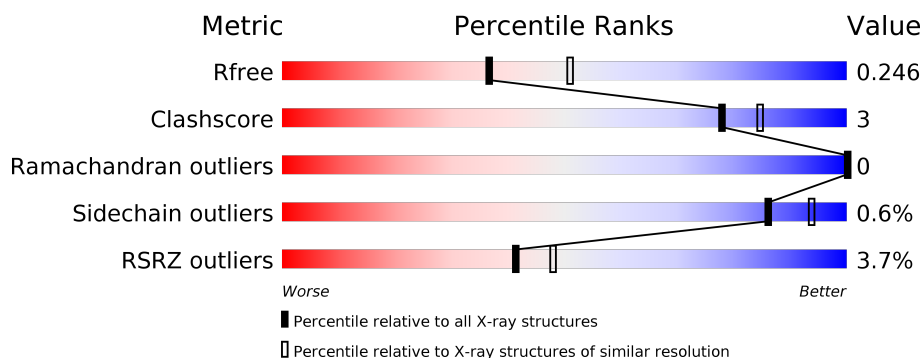
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.29 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	203	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>6%</div> <div>16%</div> </div> </div>
1	B	203	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>9%</div> <div>14%</div> </div> </div>
1	C	203	<div> <div>4%</div> <div> <div></div> <div>80%</div> <div>8%</div> <div>11%</div> </div> </div>
1	D	203	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>7%</div> <div>19%</div> </div> </div>
1	E	203	<div> <div>6%</div> <div> <div></div> <div>76%</div> <div>6%</div> <div>17%</div> </div> </div>
1	F	203	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>5%</div> <div>12%</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8635 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 Rrp44p homologue.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	18	0	0
			1303	813	233	243	14			
1	B	174	Total	C	N	O	S	11	0	0
			1332	828	239	251	14			
1	C	180	Total	C	N	O	S	7	0	0
			1379	856	248	261	14			
1	D	165	Total	C	N	O	S	37	0	0
			1265	793	226	232	14			
1	E	168	Total	C	N	O	S	39	0	0
			1290	806	233	237	14			
1	F	179	Total	C	N	O	S	22	0	0
			1369	849	246	260	14			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	219	LEU	-	expression tag	UNP Q95Z12
A	220	GLU	-	expression tag	UNP Q95Z12
A	221	HIS	-	expression tag	UNP Q95Z12
A	222	HIS	-	expression tag	UNP Q95Z12
A	223	HIS	-	expression tag	UNP Q95Z12
A	224	HIS	-	expression tag	UNP Q95Z12
A	225	HIS	-	expression tag	UNP Q95Z12
A	226	HIS	-	expression tag	UNP Q95Z12
B	219	LEU	-	expression tag	UNP Q95Z12
B	220	GLU	-	expression tag	UNP Q95Z12
B	221	HIS	-	expression tag	UNP Q95Z12
B	222	HIS	-	expression tag	UNP Q95Z12
B	223	HIS	-	expression tag	UNP Q95Z12
B	224	HIS	-	expression tag	UNP Q95Z12
B	225	HIS	-	expression tag	UNP Q95Z12
B	226	HIS	-	expression tag	UNP Q95Z12
C	219	LEU	-	expression tag	UNP Q95Z12

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Chain	Residue	Modelled	Actual	Comment	Reference
C	220	GLU	-	expression tag	UNP Q95Z12
C	221	HIS	-	expression tag	UNP Q95Z12
C	222	HIS	-	expression tag	UNP Q95Z12
C	223	HIS	-	expression tag	UNP Q95Z12
C	224	HIS	-	expression tag	UNP Q95Z12
C	225	HIS	-	expression tag	UNP Q95Z12
C	226	HIS	-	expression tag	UNP Q95Z12
D	219	LEU	-	expression tag	UNP Q95Z12
D	220	GLU	-	expression tag	UNP Q95Z12
D	221	HIS	-	expression tag	UNP Q95Z12
D	222	HIS	-	expression tag	UNP Q95Z12
D	223	HIS	-	expression tag	UNP Q95Z12
D	224	HIS	-	expression tag	UNP Q95Z12
D	225	HIS	-	expression tag	UNP Q95Z12
D	226	HIS	-	expression tag	UNP Q95Z12
E	219	LEU	-	expression tag	UNP Q95Z12
E	220	GLU	-	expression tag	UNP Q95Z12
E	221	HIS	-	expression tag	UNP Q95Z12
E	222	HIS	-	expression tag	UNP Q95Z12
E	223	HIS	-	expression tag	UNP Q95Z12
E	224	HIS	-	expression tag	UNP Q95Z12
E	225	HIS	-	expression tag	UNP Q95Z12
E	226	HIS	-	expression tag	UNP Q95Z12
F	219	LEU	-	expression tag	UNP Q95Z12
F	220	GLU	-	expression tag	UNP Q95Z12
F	221	HIS	-	expression tag	UNP Q95Z12
F	222	HIS	-	expression tag	UNP Q95Z12
F	223	HIS	-	expression tag	UNP Q95Z12
F	224	HIS	-	expression tag	UNP Q95Z12
F	225	HIS	-	expression tag	UNP Q95Z12
F	226	HIS	-	expression tag	UNP Q95Z12

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	2	Total Mn 2 2	0	0
2	E	2	Total Mn 2 2	0	0
2	B	2	Total Mn 2 2	0	0
2	C	2	Total Mn 2 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	Mn 2	0	0
2	F	2	Total 2	Mn 2	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total 1	Cl 1	0	0
4	E	1	Total 1	Cl 1	0	0
4	B	2	Total 2	Cl 2	0	0
4	C	1	Total 1	Cl 1	0	0
4	A	3	Total 3	Cl 3	0	0
4	F	3	Total 3	Cl 3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	113	Total 113	O 113	0	0

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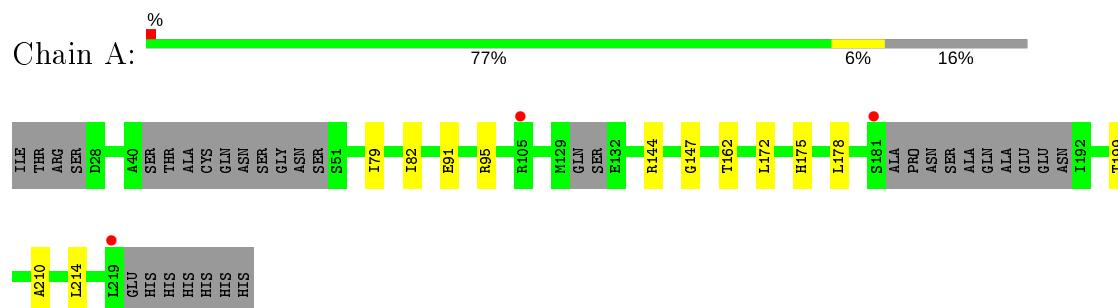
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	160	Total 160	O 160	0	0
5	C	130	Total 130	O 130	0	0
5	D	66	Total 66	O 66	0	0
5	E	64	Total 64	O 64	0	0
5	F	129	Total 129	O 129	0	0

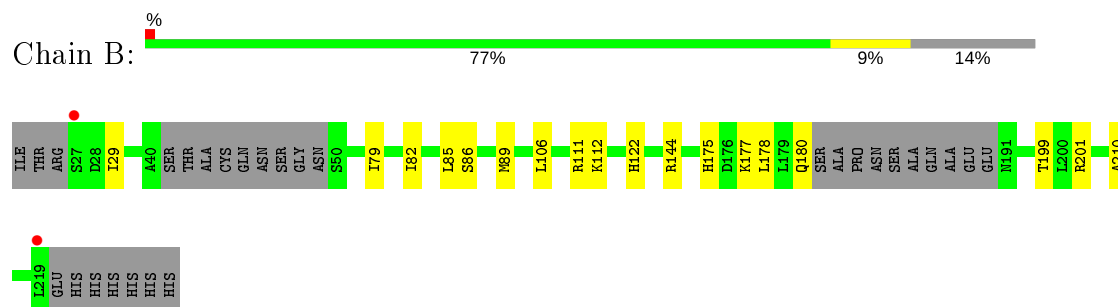
### 3 Residue-property plots [i](#)

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.

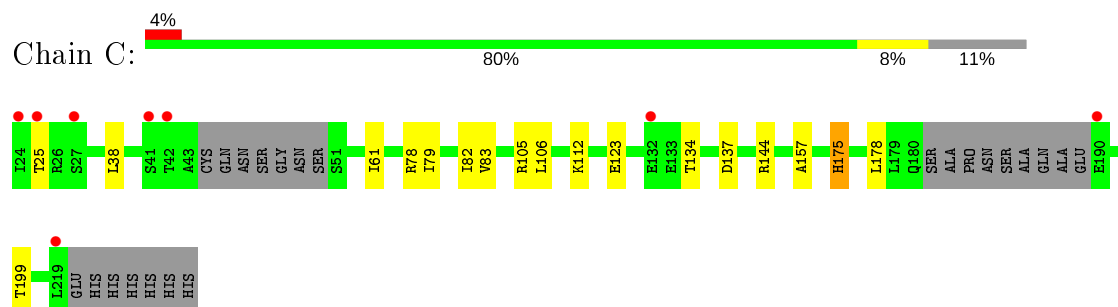
- Molecule 1: Rrp44p homologue



- Molecule 1: Rrp44p homologue

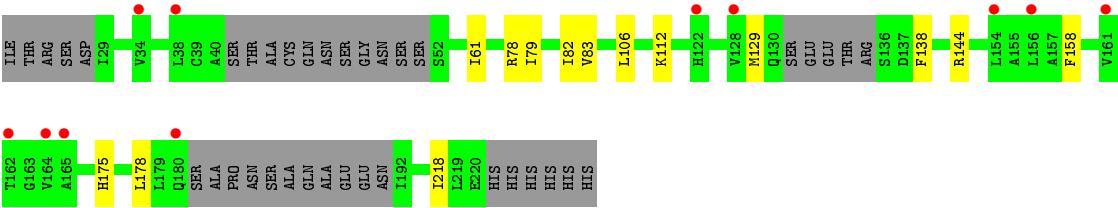


- Molecule 1: Rrp44p homologue

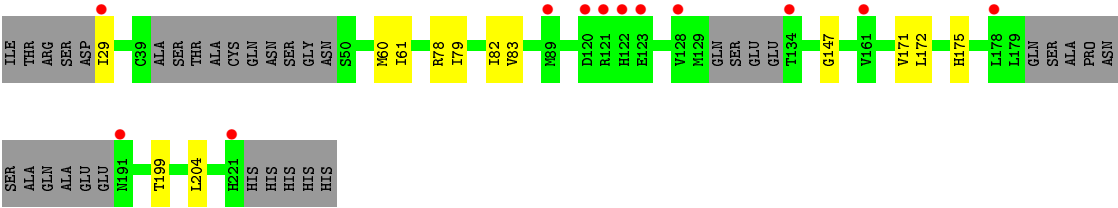
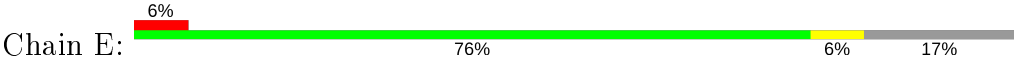


- Molecule 1: Rrp44p homologue

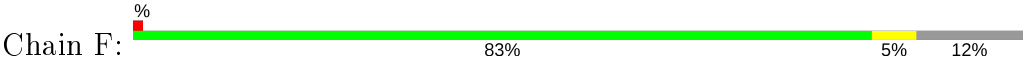




● Molecule 1: Rrp44p homologue



● Molecule 1: Rrp44p homologue





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.57Å 89.57Å 321.61Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.91 – 2.29 44.79 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.6 (45.91-2.29) 99.6 (44.79-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.38 (at 2.29Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, $R_{free}$	0.200 , 0.230 0.208 , 0.246	Depositor DCC
$R_{free}$ test set	3229 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.5	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 59.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.148 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8635	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.61% 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](#)

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

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	0.48	0/1318	0.67	0/1782
1	B	0.53	0/1348	0.70	0/1824
1	C	0.51	0/1395	0.69	0/1888
1	D	0.47	0/1280	0.67	0/1731
1	E	0.46	0/1306	0.65	0/1766
1	F	0.50	0/1386	0.66	0/1876
All	All	0.49	0/8033	0.68	0/10867

There are no bond length outliers.

There are no bond angle outliers.

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	1303	0	1316	7	0
1	B	1332	0	1341	10	0
1	C	1379	0	1390	12	0
1	D	1265	0	1284	8	0
1	E	1290	0	1306	9	0
1	F	1369	0	1371	6	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
4	A	3	0	0	0	0
4	B	2	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	3	0	0	0	0
5	A	113	0	0	0	0
5	B	160	0	0	1	0
5	C	130	0	0	0	0
5	D	66	0	0	0	0
5	E	64	0	0	2	0
5	F	129	0	0	0	0
All	All	8635	0	8008	46	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (46) 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:79:ILE:HG22	1:A:82:ILE:HD11	1.81	0.61
5:B:401:HOH:O	1:C:105:ARG:HG3	2.03	0.58
1:C:25:THR:HB	1:C:123:GLU:HG2	1.86	0.58
1:C:61:ILE:HG12	1:C:83:VAL:HB	1.85	0.57
1:B:177:LYS:HB3	1:B:180:GLN:HB2	1.88	0.56
1:B:79:ILE:HG22	1:B:82:ILE:HD11	1.87	0.56
1:D:61:ILE:HG12	1:D:83:VAL:HB	1.88	0.55
1:F:79:ILE:HG22	1:F:82:ILE:HD11	1.90	0.53
1:A:147:GLY:HA3	1:A:172:LEU:HD21	1.90	0.53
1:C:106:LEU:HD22	1:C:112:LYS:HD3	1.89	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:177:LYS:HB3	1:F:180:GLN:HB2	1.90	0.52
1:F:106:LEU:HD22	1:F:112:LYS:HD3	1.93	0.51
1:D:106:LEU:HD22	1:D:112:LYS:HD3	1.93	0.51
1:F:144:ARG:HD3	1:F:178:LEU:O	2.12	0.50
1:B:144:ARG:HD3	1:B:178:LEU:O	2.12	0.49
1:C:144:ARG:HD3	1:C:178:LEU:O	2.13	0.49
1:C:79:ILE:HG22	1:C:82:ILE:HD11	1.93	0.49
1:C:78:ARG:HD3	1:F:210:ALA:HB1	1.94	0.48
1:E:79:ILE:HG22	1:E:82:ILE:HD11	1.96	0.48
1:A:144:ARG:HD3	1:A:178:LEU:O	2.14	0.47
1:D:144:ARG:HD3	1:D:178:LEU:O	2.13	0.47
1:E:61:ILE:HG12	1:E:83:VAL:HB	1.96	0.47
1:B:85:LEU:O	1:B:89:MET:HG2	2.14	0.47
1:C:38:LEU:HD22	1:C:157:ALA:HB2	1.96	0.47
1:C:134:THR:HG22	1:C:137:ASP:CG	2.35	0.46
1:A:91:GLU:O	1:A:95:ARG:HG3	2.16	0.46
1:E:171:VAL:HB	5:E:453:HOH:O	2.17	0.45
1:A:214:LEU:HD21	1:E:79:ILE:HG13	1.99	0.44
1:C:175:HIS:HA	1:C:199:THR:HG22	1.99	0.44
1:F:175:HIS:HA	1:F:199:THR:HG22	1.99	0.44
1:B:29:ILE:HD13	1:B:122:HIS:HA	2.00	0.44
1:C:25:THR:CB	1:C:123:GLU:HG2	2.48	0.44
1:D:79:ILE:HG22	1:D:82:ILE:HD11	2.00	0.44
1:E:60:MET:HG3	5:E:453:HOH:O	2.18	0.43
1:E:147:GLY:HA3	1:E:172:LEU:HD21	2.01	0.43
1:A:175:HIS:HA	1:A:199:THR:HG22	2.01	0.42
1:B:111:ARG:HH22	1:D:158:PHE:HB3	1.85	0.42
1:D:129:MET:HG2	1:D:138:PHE:CD1	2.55	0.41
1:E:79:ILE:HD11	1:E:204:LEU:HD11	2.01	0.41
1:B:175:HIS:HA	1:B:199:THR:HG22	2.02	0.41
1:A:210:ALA:HB1	1:E:78:ARG:HD3	2.04	0.40
1:B:201:ARG:HG3	1:D:218:ILE:HD12	2.02	0.40
1:C:134:THR:HG22	1:C:137:ASP:OD2	2.22	0.40
1:B:210:ALA:HB1	1:D:78:ARG:HD3	2.03	0.40
1:B:106:LEU:HD11	1:B:112:LYS:HD3	2.02	0.40
1:E:175:HIS:HA	1:E:199:THR:HG22	2.02	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	162/203 (80%)	159 (98%)	3 (2%)	0	100	100
1	B	168/203 (83%)	166 (99%)	2 (1%)	0	100	100
1	C	174/203 (86%)	172 (99%)	2 (1%)	0	100	100
1	D	157/203 (77%)	153 (98%)	4 (2%)	0	100	100
1	E	160/203 (79%)	157 (98%)	3 (2%)	0	100	100
1	F	173/203 (85%)	170 (98%)	3 (2%)	0	100	100
All	All	994/1218 (82%)	977 (98%)	17 (2%)	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	146/174 (84%)	145 (99%)	1 (1%)	84	92
1	B	150/174 (86%)	149 (99%)	1 (1%)	84	92
1	C	155/174 (89%)	154 (99%)	1 (1%)	86	94
1	D	141/174 (81%)	140 (99%)	1 (1%)	84	92
1	E	145/174 (83%)	144 (99%)	1 (1%)	84	92
1	F	154/174 (88%)	154 (100%)	0	100	100
All	All	891/1044 (85%)	886 (99%)	5 (1%)	86	94

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	THR
1	B	86	SER
1	C	175	HIS
1	D	175	HIS
1	E	29	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	113	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 35 ligands modelled in this entry, 35 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	170/203 (83%)	0.15	3 (1%) 68 74	48, 64, 101, 123	4 (2%)
1	B	174/203 (85%)	0.06	2 (1%) 80 85	41, 53, 85, 119	3 (1%)
1	C	180/203 (88%)	0.20	8 (4%) 34 41	45, 55, 103, 115	2 (1%)
1	D	165/203 (81%)	0.43	11 (6%) 17 23	49, 76, 114, 181	9 (5%)
1	E	168/203 (82%)	0.42	12 (7%) 16 21	53, 80, 118, 148	10 (5%)
1	F	179/203 (88%)	0.05	2 (1%) 80 85	47, 58, 92, 121	6 (3%)
All	All	1036/1218 (85%)	0.21	38 (3%) 41 48	41, 64, 107, 181	34 (3%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	181	SER	7.9
1	D	128	VAL	5.5
1	D	156	LEU	5.2
1	E	123	GLU	4.5
1	C	190	GLU	4.2
1	C	42	THR	4.1
1	F	27	SER	3.4
1	D	34	VAL	3.3
1	A	105	ARG	3.3
1	A	219	LEU	3.1
1	D	164	VAL	3.1
1	D	122	HIS	3.1
1	E	134	THR	3.1
1	C	25	THR	3.0
1	D	180	GLN	3.0
1	C	41	SER	3.0
1	E	89	MET	3.0
1	B	219	LEU	2.9
1	E	221	HIS	2.8

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Mol	Chain	Res	Type	RSRZ
1	E	120	ASP	2.7
1	C	27	SER	2.7
1	E	121	ARG	2.6
1	D	161	VAL	2.6
1	E	122	HIS	2.6
1	C	219	LEU	2.6
1	E	178	LEU	2.6
1	E	161	VAL	2.4
1	E	29	ILE	2.3
1	E	191	ASN	2.3
1	D	38	LEU	2.3
1	B	27	SER	2.3
1	D	154	LEU	2.2
1	C	132	GLU	2.2
1	D	165	ALA	2.1
1	C	24	ILE	2.1
1	F	180	GLN	2.0
1	E	128	VAL	2.0
1	D	162	THR	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 monosaccharides 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(Å <sup>2</sup> )	Q<0.9
3	ZN	D	303	1/1	0.43	0.26	253,253,253,253	0
4	CL	B	306	1/1	0.60	0.13	96,96,96,96	0
2	MN	A	301	1/1	0.64	0.06	139,139,139,139	0
4	CL	A	307	1/1	0.70	0.13	97,97,97,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	CL	E	305	1/1	0.79	0.10	90,90,90,90	0
2	MN	E	301	1/1	0.80	0.06	167,167,167,167	0
4	CL	F	307	1/1	0.82	0.58	94,94,94,94	0
2	MN	F	301	1/1	0.82	0.10	131,131,131,131	0
4	CL	D	305	1/1	0.85	0.21	112,112,112,112	0
3	ZN	C	304	1/1	0.86	0.16	147,147,147,147	0
3	ZN	D	304	1/1	0.86	0.44	143,143,143,143	0
3	ZN	F	304	1/1	0.86	0.18	178,178,178,178	0
4	CL	A	306	1/1	0.88	0.14	100,100,100,100	0
2	MN	D	301	1/1	0.89	0.16	163,163,163,163	0
2	MN	D	302	1/1	0.89	0.42	78,78,78,78	0
4	CL	C	305	1/1	0.90	0.11	73,73,73,73	0
3	ZN	E	304	1/1	0.90	0.19	131,131,131,131	0
3	ZN	B	304	1/1	0.91	0.13	111,111,111,111	0
2	MN	E	302	1/1	0.92	0.40	132,132,132,132	0
4	CL	B	305	1/1	0.92	0.21	85,85,85,85	0
3	ZN	A	304	1/1	0.93	0.09	133,133,133,133	0
2	MN	B	301	1/1	0.95	0.15	96,96,96,96	0
4	CL	F	306	1/1	0.95	0.07	79,79,79,79	0
3	ZN	E	303	1/1	0.95	0.06	132,132,132,132	0
2	MN	C	301	1/1	0.96	0.09	135,135,135,135	0
4	CL	F	305	1/1	0.96	0.10	82,82,82,82	0
4	CL	A	305	1/1	0.97	0.07	76,76,76,76	0
2	MN	A	302	1/1	0.97	0.22	102,102,102,102	0
2	MN	B	302	1/1	0.98	0.09	68,68,68,68	0
2	MN	C	302	1/1	0.98	0.07	94,94,94,94	0
3	ZN	A	303	1/1	0.98	0.16	60,60,60,60	1
2	MN	F	302	1/1	0.98	0.16	89,89,89,89	0
3	ZN	C	303	1/1	0.99	0.16	60,60,60,60	1
3	ZN	F	303	1/1	1.00	0.12	75,75,75,75	0
3	ZN	B	303	1/1	1.00	0.11	70,70,70,70	0

## 6.5 Other polymers

There are no such residues in this entry.