



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:18 PM GMT

PDB ID : 3TJ1  
Title : Crystal Structure of RNA Polymerase I Transcription Initiation Factor Rrn3  
Authors : Blattner, C.; Jennebach, S.; Herzog, F.; Mayer, A.; Cheung, A.C.M.; Witte, G.; Lorenzen, K.; Hopfner, K-P.; Heck, A.J.R.; Aebersold, R.; Cramer, P.  
Deposited on : 2011-08-23  
Resolution : 2.85 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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

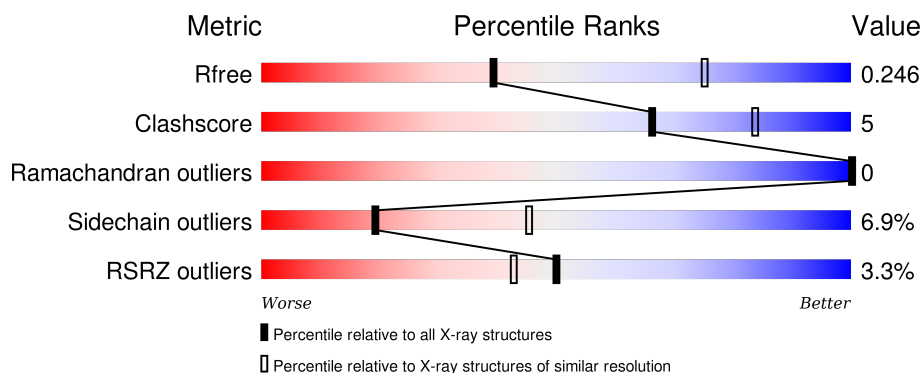
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.85 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))
$R_{free}$	91344	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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	649	<div> <div>3%</div> <div>61%</div> <div>13%</div> <div>•</div> <div>26%</div> </div>
1	B	649	<div> <div>2%</div> <div>61%</div> <div>12%</div> <div>•</div> <div>26%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8075 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 RNA polymerase I-specific transcription initiation factor RRN3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	483	Total	C	N	O	S	0	0	0
			3988	2583	654	730	21			
1	B	482	Total	C	N	O	S	0	0	0
			3984	2581	653	729	21			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	EXPRESSION TAG	UNP P36070
A	-20	GLY	-	EXPRESSION TAG	UNP P36070
A	-19	SER	-	EXPRESSION TAG	UNP P36070
A	-18	SER	-	EXPRESSION TAG	UNP P36070
A	-17	HIS	-	EXPRESSION TAG	UNP P36070
A	-16	HIS	-	EXPRESSION TAG	UNP P36070
A	-15	HIS	-	EXPRESSION TAG	UNP P36070
A	-14	HIS	-	EXPRESSION TAG	UNP P36070
A	-13	HIS	-	EXPRESSION TAG	UNP P36070
A	-12	HIS	-	EXPRESSION TAG	UNP P36070
A	-11	SER	-	EXPRESSION TAG	UNP P36070
A	-10	SER	-	EXPRESSION TAG	UNP P36070
A	-9	GLY	-	EXPRESSION TAG	UNP P36070
A	-8	LEU	-	EXPRESSION TAG	UNP P36070
A	-7	VAL	-	EXPRESSION TAG	UNP P36070
A	-6	PRO	-	EXPRESSION TAG	UNP P36070
A	-5	ARG	-	EXPRESSION TAG	UNP P36070
A	-4	GLY	-	EXPRESSION TAG	UNP P36070
A	-3	SER	-	EXPRESSION TAG	UNP P36070
A	-2	HIS	-	EXPRESSION TAG	UNP P36070
A	-1	MET	-	EXPRESSION TAG	UNP P36070
A	0	ALA	-	EXPRESSION TAG	UNP P36070
B	-21	MET	-	EXPRESSION TAG	UNP P36070
B	-20	GLY	-	EXPRESSION TAG	UNP P36070
B	-19	SER	-	EXPRESSION TAG	UNP P36070

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	SER	-	EXPRESSION TAG	UNP P36070
B	-17	HIS	-	EXPRESSION TAG	UNP P36070
B	-16	HIS	-	EXPRESSION TAG	UNP P36070
B	-15	HIS	-	EXPRESSION TAG	UNP P36070
B	-14	HIS	-	EXPRESSION TAG	UNP P36070
B	-13	HIS	-	EXPRESSION TAG	UNP P36070
B	-12	HIS	-	EXPRESSION TAG	UNP P36070
B	-11	SER	-	EXPRESSION TAG	UNP P36070
B	-10	SER	-	EXPRESSION TAG	UNP P36070
B	-9	GLY	-	EXPRESSION TAG	UNP P36070
B	-8	LEU	-	EXPRESSION TAG	UNP P36070
B	-7	VAL	-	EXPRESSION TAG	UNP P36070
B	-6	PRO	-	EXPRESSION TAG	UNP P36070
B	-5	ARG	-	EXPRESSION TAG	UNP P36070
B	-4	GLY	-	EXPRESSION TAG	UNP P36070
B	-3	SER	-	EXPRESSION TAG	UNP P36070
B	-2	HIS	-	EXPRESSION TAG	UNP P36070
B	-1	MET	-	EXPRESSION TAG	UNP P36070
B	0	ALA	-	EXPRESSION TAG	UNP P36070

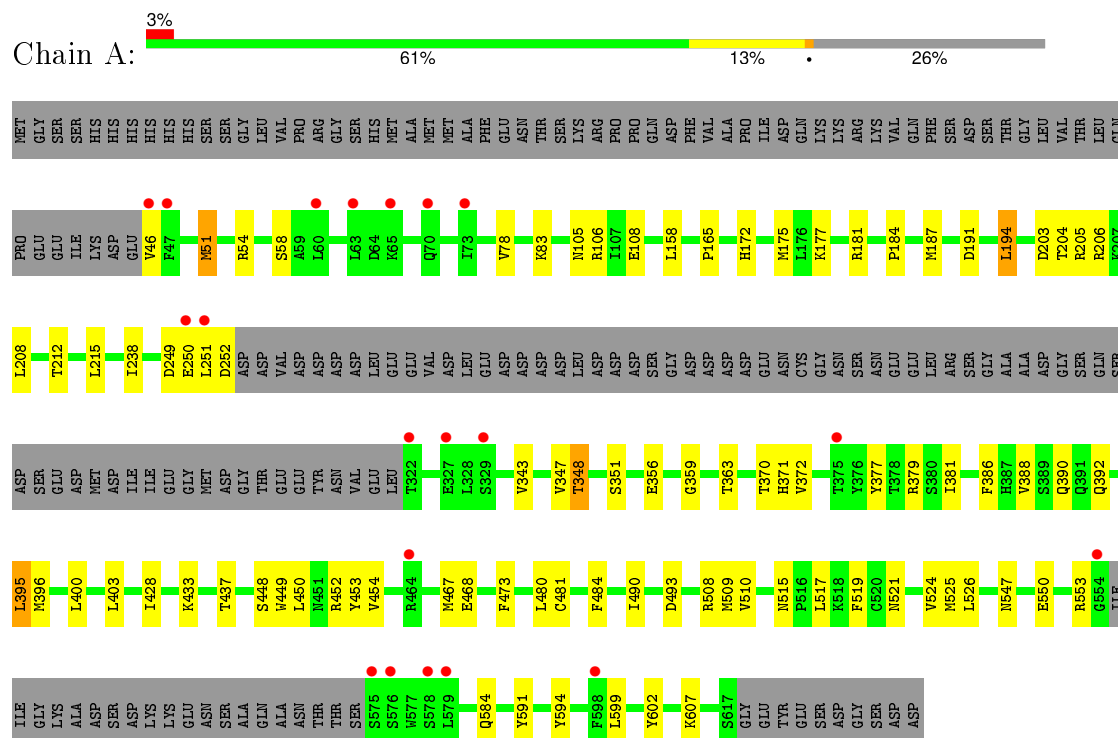
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	60	Total O 60 60	0	0
2	B	43	Total O 43 43	0	0

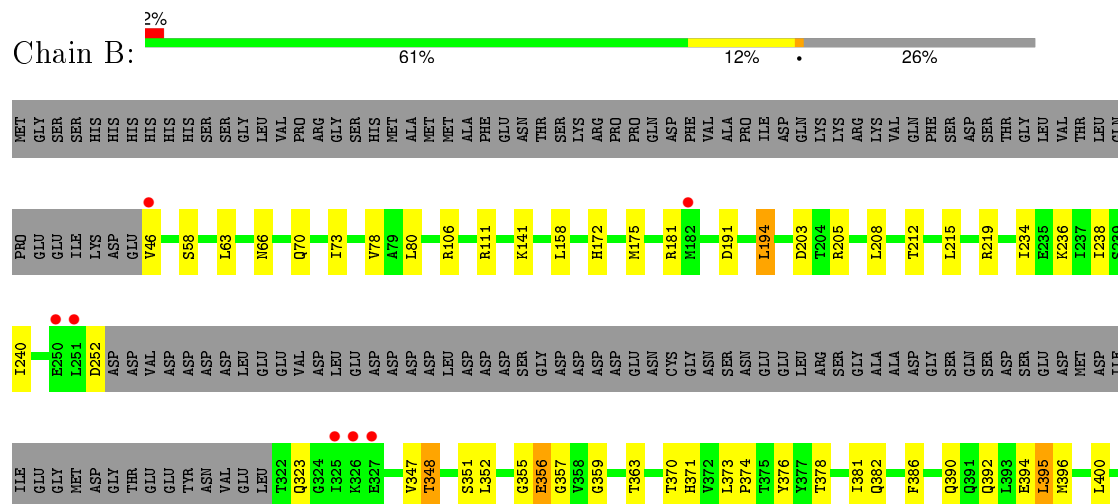
### 3 Residue-property plots

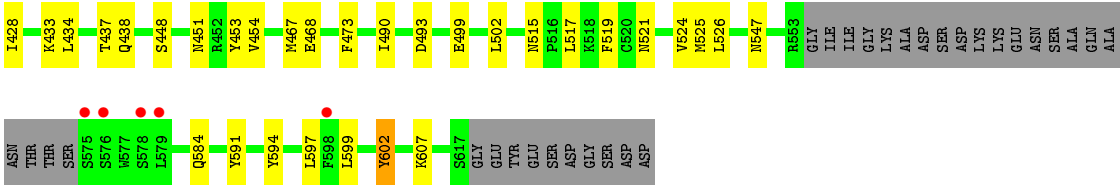
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: RNA polymerase I-specific transcription initiation factor RRN3



- Molecule 1: RNA polymerase I-specific transcription initiation factor RRN3





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.77Å 101.77Å 162.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.55 – 2.85 48.55 – 2.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.55-2.85) 100.0 (48.55-2.85)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.99 (at 2.86Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, $R_{free}$	0.208 , 0.242 0.218 , 0.246	Depositor DCC
$R_{free}$ test set	1900 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	47.5	Xtriage
Anisotropy	1.042	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 51.1	EDS
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.54$ , $\langle L^2 \rangle = 0.38$	Xtriage
Outliers	4 of 38054 reflections (0.011%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8075	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	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 53.50 % 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 4.1769e-05. 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.52	0/4078	0.71	0/5511
1	B	0.51	0/4074	0.69	0/5506
All	All	0.51	0/8152	0.70	0/11017

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

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	3988	0	3983	39	0
1	B	3984	0	3980	39	0
2	A	60	0	0	9	0
2	B	43	0	0	0	0
All	All	8075	0	7963	78	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 5.

All (78) 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:481:CYS:HA	2:A:670:HOH:O	1.54	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:480:LEU:HD22	2:A:669:HOH:O	1.60	1.00
1:A:347:VAL:HG11	1:A:388:VAL:HG23	1.64	0.80
1:A:450:LEU:HD12	2:A:669:HOH:O	1.81	0.79
1:A:400:LEU:HD21	1:A:428:ILE:HD11	1.72	0.70
1:B:400:LEU:HD21	1:B:428:ILE:HD11	1.73	0.69
1:A:208:LEU:O	1:A:212:THR:HG23	1.95	0.66
1:A:517:LEU:HA	1:A:525:MET:HE2	1.79	0.64
1:B:517:LEU:HA	1:B:525:MET:HE2	1.80	0.62
1:A:449:TRP:HA	2:A:677:HOH:O	1.99	0.62
1:A:343:VAL:O	1:A:347:VAL:HG12	2.00	0.62
1:A:400:LEU:CD2	1:A:428:ILE:HD11	2.29	0.62
1:B:400:LEU:CD2	1:B:428:ILE:HD11	2.30	0.62
1:B:378:THR:OG1	1:B:382:GLN:OE1	2.18	0.61
1:B:373:LEU:HB3	1:B:374:PRO:HD3	1.83	0.60
1:A:165:PRO:HD3	1:A:206:ARG:HH12	1.66	0.59
1:B:208:LEU:O	1:B:212:THR:HG23	2.03	0.58
1:A:348:THR:HG22	1:A:351:SER:H	1.68	0.57
1:B:376:TYR:CZ	1:B:591:TYR:HD2	2.25	0.55
1:B:390:GLN:NE2	1:B:433:LYS:H	2.05	0.55
1:B:396:MET:HE2	1:B:434:LEU:HG	1.88	0.55
1:B:348:THR:HG22	1:B:351:SER:H	1.72	0.55
1:A:396:MET:HE3	1:A:433:LYS:HB2	1.88	0.54
1:B:376:TYR:CZ	1:B:591:TYR:CD2	2.94	0.54
1:B:390:GLN:HG2	1:B:396:MET:HE3	1.91	0.53
1:B:390:GLN:HG2	1:B:396:MET:CE	2.40	0.52
1:A:521:ASN:HD22	1:A:524:VAL:H	1.56	0.52
1:B:521:ASN:HD22	1:B:524:VAL:H	1.56	0.52
1:A:484:PHE:HB3	2:A:670:HOH:O	2.10	0.52
1:A:377:TYR:HA	1:A:591:TYR:CE1	2.45	0.52
1:B:390:GLN:HE21	1:B:433:LYS:H	1.58	0.51
1:A:184:PRO:HA	2:A:683:HOH:O	2.10	0.51
1:B:451:ASN:O	1:B:454:VAL:HG12	2.11	0.50
1:B:370:THR:OG1	1:B:371:HIS:HD2	1.94	0.50
1:A:370:THR:OG1	1:A:371:HIS:HD2	1.95	0.50
1:A:386:PHE:HB2	1:A:594:TYR:CE2	2.48	0.49
1:B:515:ASN:HD21	1:B:547:ASN:HD21	1.61	0.48
1:A:515:ASN:HD21	1:A:547:ASN:HD21	1.62	0.48
1:A:359:GLY:O	1:A:363:THR:HG23	2.13	0.48
1:A:372:VAL:HG13	1:A:381:ILE:HG23	1.96	0.48
1:B:386:PHE:HB2	1:B:594:TYR:CE2	2.49	0.47
1:B:400:LEU:HD21	1:B:428:ILE:CD1	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:392:GLN:HB2	1:B:395:LEU:HD22	1.96	0.46
1:B:356:GLU:HG3	1:B:356:GLU:H	1.38	0.46
1:A:187:MET:HG3	2:A:683:HOH:O	2.15	0.45
1:A:348:THR:H	1:A:351:SER:HB3	1.81	0.45
1:A:51:MET:SD	1:A:54:ARG:NH2	2.89	0.45
1:B:467:MET:HG3	1:B:519:PHE:CZ	2.52	0.44
1:A:390:GLN:NE2	1:A:433:LYS:H	2.15	0.44
1:A:452:ARG:NE	2:A:677:HOH:O	2.50	0.44
1:B:597:LEU:HD21	1:B:602:TYR:CD1	2.53	0.43
1:A:191:ASP:HA	1:A:194:LEU:HD22	2.00	0.43
1:A:392:GLN:HB2	1:A:395:LEU:HD22	2.00	0.43
1:A:400:LEU:HD21	1:A:428:ILE:CD1	2.42	0.43
1:B:238:ILE:HG12	1:B:381:ILE:HG21	2.00	0.43
1:B:347:VAL:HG13	1:B:352:LEU:HD11	2.01	0.43
1:B:191:ASP:HA	1:B:194:LEU:HD22	2.01	0.43
1:B:70:GLN:HA	1:B:73:ILE:HD12	2.00	0.42
1:B:141:LYS:HG3	1:B:175:MET:HE1	2.01	0.42
1:B:236:LYS:O	1:B:240:ILE:HG12	2.19	0.42
1:B:355:GLY:C	1:B:357:GLY:H	2.22	0.42
1:B:396:MET:HB2	1:B:438:GLN:HE22	1.85	0.42
1:B:359:GLY:O	1:B:363:THR:HG23	2.20	0.42
1:B:234:ILE:HD12	1:B:381:ILE:HD11	2.02	0.42
1:B:158:LEU:HD11	1:B:175:MET:HG2	2.02	0.42
1:B:453:TYR:CE2	1:B:473:PHE:HB2	2.55	0.42
1:A:453:TYR:CE2	1:A:473:PHE:HB2	2.55	0.41
1:B:158:LEU:HD22	1:B:172:HIS:HA	2.02	0.41
1:B:499:GLU:O	1:B:502:LEU:HG	2.20	0.41
1:A:105:ASN:O	1:A:108:GLU:HB2	2.20	0.41
1:A:238:ILE:HG12	1:A:381:ILE:HG21	2.03	0.41
1:A:158:LEU:HD22	1:A:172:HIS:HA	2.03	0.41
1:A:177:LYS:HG2	1:A:181:ARG:HH22	1.86	0.41
1:B:66:ASN:OD1	1:B:111:ARG:NH1	2.53	0.41
1:A:403:LEU:HB3	2:A:672:HOH:O	2.20	0.41
1:A:473:PHE:CZ	1:A:509:MET:HE3	2.57	0.40
1:A:158:LEU:HD11	1:A:175:MET:HG2	2.03	0.40
1:A:467:MET:HG3	1:A:519:PHE:CZ	2.55	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	477/649 (74%)	453 (95%)	24 (5%)	0	100	100
1	B	476/649 (73%)	453 (95%)	23 (5%)	0	100	100
All	All	953/1298 (73%)	906 (95%)	47 (5%)	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	449/594 (76%)	415 (92%)	34 (8%)	16	40
1	B	449/594 (76%)	421 (94%)	28 (6%)	23	51
All	All	898/1188 (76%)	836 (93%)	62 (7%)	19	45

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	VAL
1	A	51	MET
1	A	58	SER
1	A	78	VAL
1	A	83	LYS
1	A	106	ARG
1	A	194	LEU
1	A	203	ASP

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Mol	Chain	Res	Type
1	A	204	THR
1	A	205	ARG
1	A	215	LEU
1	A	249	ASP
1	A	250	GLU
1	A	251	LEU
1	A	252	ASP
1	A	348	THR
1	A	356	GLU
1	A	379	ARG
1	A	395	LEU
1	A	437	THR
1	A	448	SER
1	A	454	VAL
1	A	468	GLU
1	A	490	ILE
1	A	493	ASP
1	A	508	ARG
1	A	510	VAL
1	A	526	LEU
1	A	550	GLU
1	A	553	ARG
1	A	584	GLN
1	A	599	LEU
1	A	602	TYR
1	A	607	LYS
1	B	46	VAL
1	B	58	SER
1	B	63	LEU
1	B	78	VAL
1	B	80	LEU
1	B	106	ARG
1	B	181	ARG
1	B	194	LEU
1	B	203	ASP
1	B	205	ARG
1	B	215	LEU
1	B	219	ARG
1	B	252	ASP
1	B	323	GLN
1	B	348	THR
1	B	356	GLU

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Mol	Chain	Res	Type
1	B	394	GLU
1	B	395	LEU
1	B	437	THR
1	B	448	SER
1	B	468	GLU
1	B	490	ILE
1	B	493	ASP
1	B	526	LEU
1	B	584	GLN
1	B	599	LEU
1	B	602	TYR
1	B	607	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	245	GLN
1	A	371	HIS
1	A	390	GLN
1	A	521	ASN
1	A	547	ASN
1	A	549	ASN
1	B	105	ASN
1	B	210	ASN
1	B	245	GLN
1	B	371	HIS
1	B	390	GLN
1	B	521	ASN
1	B	547	ASN
1	B	549	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no 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	483/649 (74%)	0.21	20 (4%) 41 34	33, 58, 103, 122	0
1	B	482/649 (74%)	0.13	12 (2%) 61 56	32, 60, 98, 118	0
All	All	965/1298 (74%)	0.17	32 (3%) 50 43	32, 59, 100, 122	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	575	SER	6.9
1	A	322	THR	4.2
1	A	375	THR	4.1
1	A	63	LEU	3.9
1	B	578	SER	3.9
1	B	576	SER	3.8
1	B	251	LEU	3.7
1	B	325	ILE	3.4
1	B	46	VAL	3.4
1	B	598	PHE	3.3
1	A	575	SER	3.1
1	A	578	SER	3.0
1	B	327	GLU	2.9
1	A	70	GLN	2.8
1	A	73	ILE	2.8
1	A	576	SER	2.8
1	A	250	GLU	2.7
1	A	251	LEU	2.7
1	A	554	GLY	2.6
1	A	579	LEU	2.6
1	B	182	MET	2.6
1	B	250	GLU	2.5
1	B	326	LYS	2.5
1	A	65	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	579	LEU	2.4
1	A	60	LEU	2.3
1	A	47	PHE	2.3
1	A	464	ARG	2.2
1	A	329	SER	2.0
1	A	327	GLU	2.0
1	A	598	PHE	2.0
1	A	46	VAL	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.