



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 17, 2016 – 07:53 AM EST

PDB ID : 5GPD  
Title : Crystal structure of the binding domain of SREBP from fission yeast  
Authors : Gong, X.; Qian, H.W.; Wu, J.P.; Yan, N.  
Deposited on : 2016-08-01  
Resolution : 3.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028320  
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) : rb-20028320

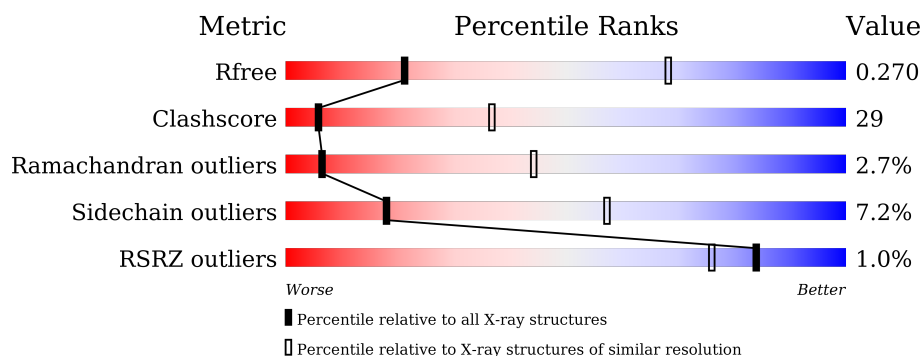
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	252	<div> <div></div> <div>50%</div> <div>27%</div> <div>18%</div> </div>
1	B	252	<div> <div>42%</div> <div>23%</div> <div>5%</div> <div>30%</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2938 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 Sterol regulatory element-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1573	1004	250	309	10			
1	B	177	Total	C	N	O	S	0	0	0
			1365	881	217	259	8			

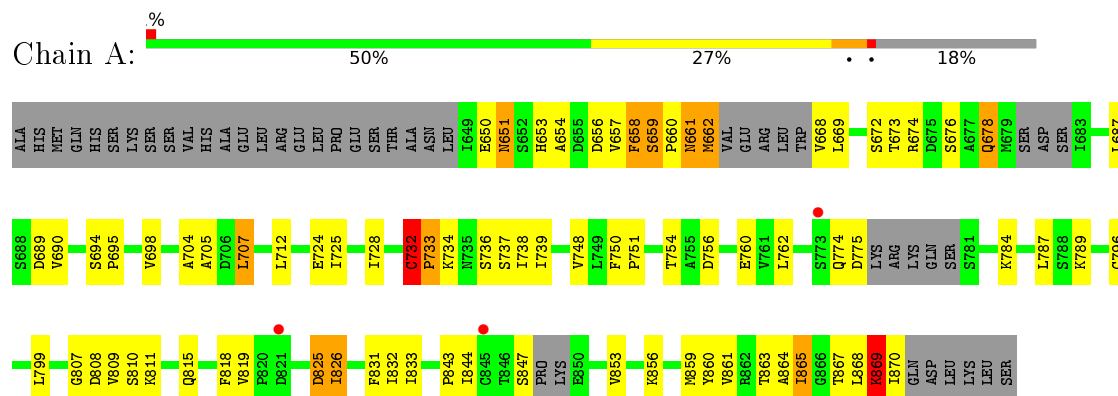
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	625	ALA	-	expression tag	UNP Q9UUD1
A	626	HIS	-	expression tag	UNP Q9UUD1
A	627	MET	-	expression tag	UNP Q9UUD1
A	644	SER	CYS	engineered mutation	UNP Q9UUD1
A	672	SER	CYS	engineered mutation	UNP Q9UUD1
B	625	ALA	-	expression tag	UNP Q9UUD1
B	626	HIS	-	expression tag	UNP Q9UUD1
B	627	MET	-	expression tag	UNP Q9UUD1
B	644	SER	CYS	engineered mutation	UNP Q9UUD1
B	672	SER	CYS	engineered mutation	UNP Q9UUD1

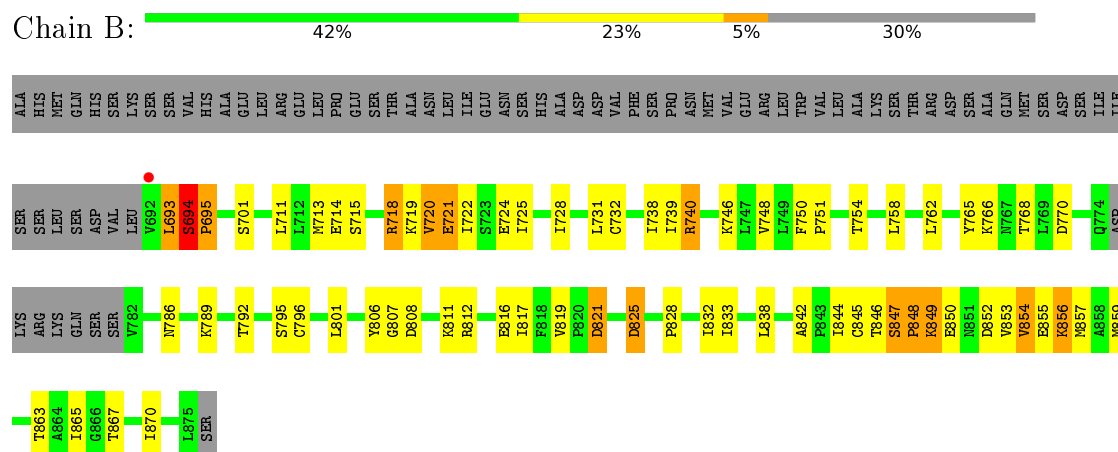
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Sterol regulatory element-binding protein 1



- Molecule 1: Sterol regulatory element-binding protein 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.90 Å 96.90 Å 151.59 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.97 – 3.50 34.97 – 3.50	Depositor EDS
% Data completeness (in resolution range)	98.2 (34.97-3.50) 92.0 (34.97-3.50)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.43 (at 3.47 Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, $R_{free}$	0.265 , 0.279 0.258 , 0.270	Depositor DCC
$R_{free}$ test set	477 reflections (4.80%)	DCC
Wilson B-factor (Å <sup>2</sup> )	138.5	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 117.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	2938	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	142.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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	0.36	1/1592 (0.1%)	0.54	2/2161 (0.1%)
1	B	0.61	6/1386 (0.4%)	0.57	3/1885 (0.2%)
All	All	0.49	7/2978 (0.2%)	0.55	5/4046 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	740	ARG	NE-CZ	-9.45	1.20	1.33
1	B	740	ARG	CZ-NH2	-9.02	1.21	1.33
1	B	740	ARG	CD-NE	-7.98	1.32	1.46
1	B	740	ARG	CZ-NH1	-7.26	1.23	1.33
1	B	848	PRO	N-CD	5.28	1.55	1.47
1	A	733	PRO	N-CD	5.03	1.54	1.47
1	B	695	PRO	N-CD	5.03	1.54	1.47

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	659	SER	C-N-CD	6.05	141.11	128.40
1	A	732	CYS	C-N-CD	6.02	141.05	128.40
1	B	694	SER	C-N-CD	6.02	141.04	128.40
1	B	847	SER	C-N-CD	5.58	140.11	128.40
1	B	740	ARG	NE-CZ-NH1	5.04	122.82	120.30

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	1573	0	1589	82	0
1	B	1365	0	1405	93	0
All	All	2938	0	2994	171	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 29.

All (171) 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:859:MET:O	1:A:863:THR:HG22	1.40	1.21
1:B:715:SER:CB	1:B:720:VAL:HG21	1.78	1.12
1:A:733:PRO:HG2	1:A:736:SER:HB3	1.36	1.08
1:B:842:ALA:O	1:B:846:THR:HG23	1.55	1.06
1:A:774:GLN:O	1:A:775:ASP:OD1	1.74	1.05
1:A:662:MET:HG2	1:A:695:PRO:HB3	1.38	1.05
1:B:842:ALA:O	1:B:846:THR:CG2	2.07	1.02
1:B:715:SER:HA	1:B:720:VAL:CG2	1.90	1.01
1:A:733:PRO:HD2	1:A:736:SER:OG	1.63	0.98
1:B:715:SER:HB3	1:B:720:VAL:HG21	1.48	0.95
1:B:819:VAL:HG11	1:B:856:LYS:HB3	1.48	0.94
1:B:758:LEU:HD23	1:B:762:LEU:CD1	1.98	0.92
1:B:758:LEU:HD23	1:B:762:LEU:HD13	1.50	0.91
1:B:722:ILE:HG13	1:B:750:PHE:CD1	2.06	0.91
1:B:714:GLU:OE1	1:B:718:ARG:NH1	2.03	0.90
1:A:748:VAL:HG21	1:A:833:ILE:HG12	1.55	0.88
1:B:711:LEU:HD11	1:B:720:VAL:CG1	2.05	0.86
1:A:809:VAL:HG22	1:B:806:TYR:CE1	2.11	0.85
1:B:715:SER:CA	1:B:720:VAL:HG21	2.05	0.85
1:A:662:MET:CG	1:A:695:PRO:HB3	2.08	0.84
1:A:825:ASP:O	1:A:826:ILE:HG23	1.78	0.84
1:A:733:PRO:HG2	1:A:736:SER:CB	2.08	0.81
1:A:733:PRO:CG	1:A:736:SER:HB3	2.09	0.81
1:B:722:ILE:HG13	1:B:750:PHE:CE1	2.15	0.81
1:B:724:GLU:O	1:B:728:ILE:HG13	1.80	0.80
1:B:811:LYS:HD3	1:B:844:ILE:HG22	1.63	0.80
1:A:869:LYS:H	1:A:869:LYS:HD2	1.47	0.79
1:B:715:SER:CA	1:B:720:VAL:CG2	2.60	0.79
1:A:733:PRO:CD	1:A:736:SER:OG	2.30	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:669:LEU:HD11	1:A:784:LYS:NZ	1.99	0.77
1:B:721:GLU:C	1:B:750:PHE:CE1	2.58	0.77
1:B:849:LYS:HA	1:B:849:LYS:HE2	1.66	0.77
1:B:694:SER:HB2	1:B:695:PRO:CD	2.15	0.77
1:A:662:MET:HG2	1:A:695:PRO:CB	2.14	0.77
1:A:859:MET:O	1:A:863:THR:CG2	2.30	0.76
1:B:758:LEU:CD2	1:B:762:LEU:CD1	2.64	0.75
1:B:849:LYS:HA	1:B:849:LYS:CE	2.16	0.74
1:B:849:LYS:HE2	1:B:849:LYS:CA	2.18	0.72
1:B:842:ALA:O	1:B:846:THR:HG21	1.88	0.72
1:A:650:GLU:O	1:A:654:ALA:HB3	1.90	0.71
1:A:661:ASN:HD22	1:A:661:ASN:H	1.37	0.71
1:B:715:SER:CB	1:B:720:VAL:CG2	2.63	0.71
1:B:714:GLU:CD	1:B:718:ARG:HH12	1.94	0.71
1:B:711:LEU:HD11	1:B:720:VAL:HG11	1.73	0.70
1:B:714:GLU:CD	1:B:718:ARG:NH1	2.44	0.70
1:A:811:LYS:NZ	1:A:847:SER:OG	2.24	0.70
1:B:758:LEU:CD2	1:B:762:LEU:HD11	2.22	0.70
1:A:658:PHE:CZ	1:A:687:LEU:HA	2.27	0.69
1:A:669:LEU:HD11	1:A:784:LYS:HZ2	1.55	0.69
1:A:787:LEU:HD21	1:A:818:PHE:HD2	1.58	0.69
1:B:721:GLU:C	1:B:750:PHE:HE1	1.94	0.68
1:A:825:ASP:O	1:A:826:ILE:CG2	2.40	0.67
1:A:658:PHE:HZ	1:A:687:LEU:HA	1.57	0.67
1:A:864:ALA:O	1:A:865:ILE:HG12	1.94	0.67
1:B:693:LEU:N	1:B:693:LEU:HD12	2.11	0.66
1:B:748:VAL:HG21	1:B:833:ILE:HG12	1.77	0.66
1:B:701:SER:HB2	1:B:738:ILE:HD13	1.79	0.65
1:B:847:SER:HB3	1:B:848:PRO:HD3	1.78	0.65
1:B:758:LEU:O	1:B:762:LEU:HD13	1.97	0.65
1:A:658:PHE:O	1:A:661:ASN:ND2	2.30	0.64
1:B:758:LEU:CD2	1:B:762:LEU:HD13	2.23	0.64
1:A:662:MET:CB	1:A:695:PRO:HB3	2.29	0.62
1:B:845:CYS:O	1:B:852:ASP:OD2	2.17	0.62
1:B:852:ASP:HB3	1:B:854:VAL:HG22	1.82	0.62
1:A:831:PHE:HE1	1:A:861:VAL:HG13	1.64	0.62
1:B:801:LEU:HD22	1:B:806:TYR:CE2	2.34	0.62
1:A:673:THR:O	1:A:676:SER:OG	2.17	0.61
1:A:661:ASN:HD21	1:A:698:VAL:HG21	1.65	0.61
1:B:693:LEU:HD12	1:B:693:LEU:H	1.66	0.60
1:A:657:VAL:HG13	1:A:678:GLN:OE1	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:853:VAL:HG12	1:B:853:VAL:O	2.02	0.59
1:A:657:VAL:CG1	1:A:678:GLN:OE1	2.51	0.58
1:B:758:LEU:HD23	1:B:762:LEU:HD11	1.80	0.58
1:A:819:VAL:HG21	1:A:856:LYS:HB3	1.85	0.58
1:A:865:ILE:O	1:A:865:ILE:HG22	2.03	0.57
1:A:864:ALA:C	1:A:865:ILE:HG12	2.25	0.57
1:B:812:ARG:NH1	1:B:816:GLU:OE1	2.34	0.57
1:B:738:ILE:H	1:B:738:ILE:HD12	1.70	0.57
1:A:733:PRO:CG	1:A:736:SER:CB	2.76	0.56
1:A:725:ILE:HG13	1:A:750:PHE:HE2	1.70	0.56
1:A:807:GLY:HA2	1:A:810:SER:HB3	1.88	0.56
1:B:715:SER:HA	1:B:720:VAL:HG22	1.86	0.55
1:B:751:PRO:HA	1:B:754:THR:HG23	1.87	0.55
1:A:844:ILE:HD12	1:A:844:ILE:H	1.70	0.55
1:A:707:LEU:HB3	1:A:728:ILE:HD12	1.89	0.55
1:B:795:SER:HB3	1:B:817:ILE:HD13	1.89	0.55
1:B:715:SER:HB2	1:B:720:VAL:HG21	1.82	0.54
1:B:722:ILE:N	1:B:750:PHE:CE1	2.75	0.54
1:A:668:VAL:O	1:A:668:VAL:HG12	2.08	0.54
1:B:713:MET:SD	1:B:828:PRO:HB3	2.47	0.54
1:A:661:ASN:HB2	1:A:690:VAL:HG13	1.89	0.53
1:A:756:ASP:O	1:A:760:GLU:HG3	2.08	0.53
1:B:711:LEU:CD1	1:B:720:VAL:HG11	2.37	0.53
1:B:806:TYR:O	1:B:808:ASP:N	2.41	0.53
1:A:868:LEU:HD23	1:A:870:ILE:HD12	1.91	0.53
1:A:825:ASP:C	1:A:826:ILE:HG23	2.28	0.53
1:B:721:GLU:O	1:B:750:PHE:CZ	2.61	0.53
1:B:722:ILE:HD12	1:B:722:ILE:H	1.73	0.53
1:B:722:ILE:HA	1:B:750:PHE:CZ	2.44	0.52
1:A:724:GLU:O	1:A:728:ILE:HG12	2.10	0.52
1:B:711:LEU:HD11	1:B:720:VAL:HG13	1.89	0.52
1:A:774:GLN:C	1:A:775:ASP:OD1	2.47	0.52
1:A:659:SER:N	1:A:660:PRO:CD	2.75	0.50
1:B:721:GLU:O	1:B:750:PHE:HZ	1.93	0.50
1:A:869:LYS:H	1:A:869:LYS:CD	2.14	0.49
1:A:669:LEU:HD11	1:A:784:LYS:HZ1	1.78	0.49
1:A:712:LEU:HB3	1:A:832:ILE:HG13	1.93	0.49
1:B:718:ARG:CG	1:B:719:LYS:N	2.76	0.49
1:A:662:MET:HB3	1:A:695:PRO:HG3	1.95	0.49
1:B:721:GLU:CA	1:B:750:PHE:HE1	2.25	0.49
1:A:656:ASP:C	1:A:658:PHE:H	2.16	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:792:THR:HG23	1:B:833:ILE:HD12	1.94	0.49
1:A:704:ALA:HB1	1:A:732:CYS:SG	2.54	0.48
1:B:711:LEU:HD23	1:B:725:ILE:HD13	1.95	0.48
1:A:704:ALA:CB	1:A:732:CYS:SG	3.01	0.48
1:A:811:LYS:O	1:A:815:GLN:HG3	2.14	0.48
1:A:809:VAL:HG22	1:B:806:TYR:HE1	1.75	0.47
1:A:661:ASN:H	1:A:661:ASN:ND2	2.09	0.47
1:B:725:ILE:HG13	1:B:750:PHE:CE2	2.50	0.47
1:B:863:THR:O	1:B:867:THR:HG23	2.14	0.47
1:A:662:MET:HG3	1:A:662:MET:H	1.48	0.46
1:B:711:LEU:CD1	1:B:720:VAL:CG1	2.87	0.46
1:B:718:ARG:HG2	1:B:719:LYS:N	2.31	0.46
1:A:818:PHE:CZ	1:B:770:ASP:HA	2.51	0.46
1:A:658:PHE:HE2	1:A:690:VAL:HG23	1.81	0.46
1:A:736:SER:O	1:A:739:ILE:HG22	2.16	0.45
1:B:720:VAL:C	1:B:721:GLU:HG3	2.35	0.45
1:B:721:GLU:C	1:B:750:PHE:CZ	2.89	0.45
1:A:672:SER:OG	1:A:687:LEU:HD11	2.17	0.45
1:B:853:VAL:HG13	1:B:856:LYS:HB2	1.99	0.45
1:A:815:GLN:HA	1:A:853:VAL:HG21	1.98	0.45
1:A:751:PRO:HA	1:A:754:THR:HG23	1.99	0.44
1:B:720:VAL:HG12	1:B:721:GLU:H	1.82	0.44
1:A:748:VAL:HG22	1:A:799:LEU:HD23	1.99	0.44
1:B:855:GLU:O	1:B:859:MET:HG3	2.18	0.43
1:A:687:LEU:O	1:A:687:LEU:HD23	2.18	0.43
1:B:865:ILE:H	1:B:865:ILE:HG13	1.65	0.43
1:B:832:ILE:HD13	1:B:832:ILE:HA	1.89	0.43
1:A:789:LYS:HA	1:A:789:LYS:HD3	1.85	0.43
1:B:765:TYR:HA	1:B:768:THR:HG22	2.00	0.43
1:B:693:LEU:N	1:B:693:LEU:CD1	2.79	0.43
1:B:694:SER:HB2	1:B:695:PRO:HD3	1.97	0.43
1:B:732:CYS:HB2	1:B:739:ILE:HD13	2.01	0.43
1:B:786:ASN:HB3	1:B:789:LYS:HG3	2.00	0.43
1:A:651:ASN:HA	1:A:651:ASN:HD22	1.52	0.42
1:A:657:VAL:HG12	1:A:657:VAL:O	2.18	0.42
1:A:860:TYR:O	1:A:863:THR:HG23	2.18	0.42
1:A:868:LEU:O	1:A:870:ILE:N	2.51	0.42
1:B:714:GLU:HB3	1:B:718:ARG:NH1	2.34	0.42
1:B:725:ILE:HG13	1:B:750:PHE:HE2	1.85	0.42
1:B:715:SER:HB2	1:B:720:VAL:CG2	2.45	0.42
1:A:733:PRO:CG	1:A:736:SER:OG	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:847:SER:HB3	1:B:848:PRO:CD	2.45	0.42
1:A:689:ASP:OD2	1:A:737:SER:HB2	2.20	0.42
1:A:656:ASP:C	1:A:658:PHE:N	2.73	0.41
1:B:766:LYS:HD3	1:B:766:LYS:HA	1.91	0.41
1:A:705:ALA:HB2	1:A:738:ILE:HG12	2.03	0.41
1:A:653:HIS:O	1:A:656:ASP:HB3	2.20	0.41
1:A:762:LEU:HD13	1:A:762:LEU:HA	1.80	0.41
1:B:722:ILE:O	1:B:725:ILE:N	2.53	0.41
1:A:809:VAL:CG2	1:B:806:TYR:CE1	2.94	0.41
1:A:659:SER:N	1:A:660:PRO:HD3	2.35	0.41
1:A:694:SER:O	1:A:698:VAL:HG23	2.21	0.40
1:A:748:VAL:HG23	1:A:796:CYS:SG	2.60	0.40
1:B:867:THR:HA	1:B:870:ILE:HG22	2.03	0.40
1:B:801:LEU:HD22	1:B:806:TYR:CD2	2.56	0.40
1:B:838:LEU:HD23	1:B:842:ALA:HB2	2.03	0.40
1:B:765:TYR:O	1:B:768:THR:HG22	2.21	0.40
1:B:792:THR:O	1:B:796:CYS:HB2	2.21	0.40
1:B:722:ILE:CG2	1:B:746:LYS:HD2	2.52	0.40
1:B:811:LYS:HD3	1:B:844:ILE:CG2	2.43	0.40

There are no symmetry-related clashes.

## 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	197/252 (78%)	170 (86%)	22 (11%)	5 (2%)	7	46
1	B	173/252 (69%)	157 (91%)	11 (6%)	5 (3%)	6	42
All	All	370/504 (73%)	327 (88%)	33 (9%)	10 (3%)	6	44

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	734	LYS
1	A	826	ILE
1	A	869	LYS
1	B	694	SER
1	B	821	ASP
1	B	850	GLU
1	A	865	ILE
1	B	807	GLY
1	A	843	PRO
1	B	825	ASP

### 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	180/228 (79%)	168 (93%)	12 (7%)	20	61
1	B	155/228 (68%)	143 (92%)	12 (8%)	16	54
All	All	335/456 (74%)	311 (93%)	24 (7%)	18	57

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	651	ASN
1	A	658	PHE
1	A	661	ASN
1	A	662	MET
1	A	674	ARG
1	A	678	GLN
1	A	707	LEU
1	A	732	CYS
1	A	808	ASP
1	A	825	ASP
1	A	867	THR
1	A	869	LYS
1	B	693	LEU
1	B	718	ARG

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Mol	Chain	Res	Type
1	B	720	VAL
1	B	721	GLU
1	B	731	LEU
1	B	740	ARG
1	B	821	ASP
1	B	825	ASP
1	B	849	LYS
1	B	854	VAL
1	B	856	LYS
1	B	857	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	651	ASN
1	A	661	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](#)

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 ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	207/252 (82%)	-0.18	3 (1%) 78 68	84, 137, 207, 280	0
1	B	177/252 (70%)	-0.29	1 (0%) 90 85	88, 136, 206, 269	0
All	All	384/504 (76%)	-0.23	4 (1%) 84 76	84, 137, 207, 280	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	692	VAL	3.0
1	A	821	ASP	2.3
1	A	845	CYS	2.1
1	A	773	SER	2.1

### 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.