



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:09 PM GMT

PDB ID : 4BP8  
Title : Oligopeptidase B from Trypanosoma brucei - open form  
Authors : Canning, P.; Rea, D.; Morty, R.; Fulop, V.  
Deposited on : 2013-05-23  
Resolution : 2.40 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

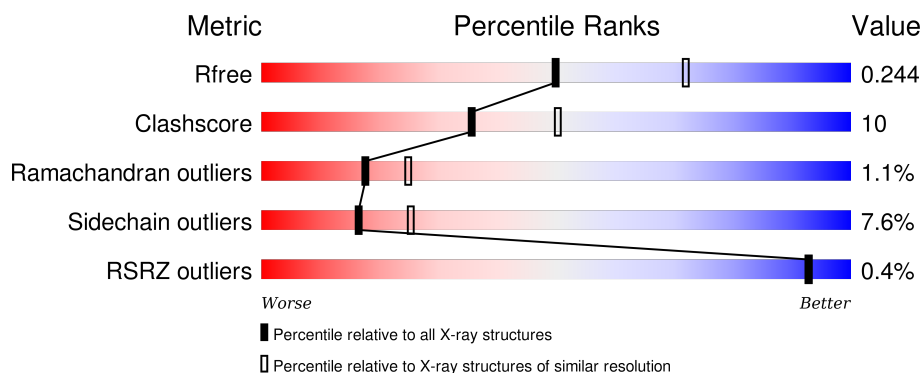
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	715	<div> <div></div> <div>80%15% . .</div> </div>
1	B	715	<div> <div></div> <div>79%17% . .</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12080 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 OLIGOPEPTIDASE B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	713	Total	C	N	O	S	Se	0	0	1
			5651	3577	979	1064	14	17			
1	B	713	Total	C	N	O	S	Se	0	0	1
			5651	3577	979	1064	14	17			

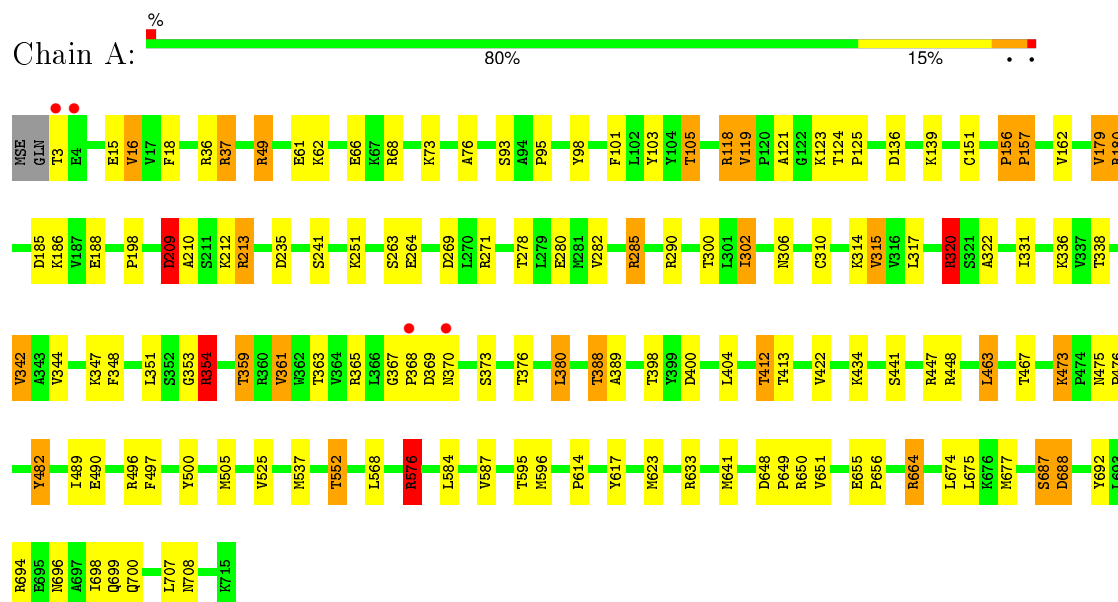
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	397	Total	O	0	0
			397	397		
2	B	381	Total	O	0	0
			381	381		

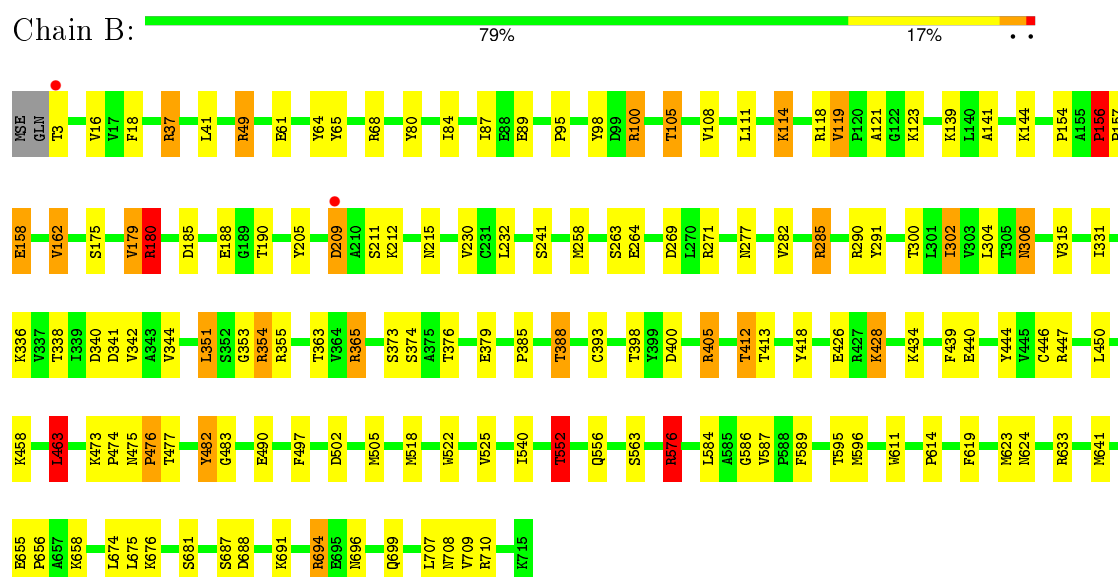
### 3 Residue-property plots

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: OLIGOPEPTIDASE B



#### • Molecule 1: OLIGOPEPTIDASE B



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	124.14Å 124.14Å 249.14Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	65.73 – 2.40 65.72 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.3 (65.73-2.40) 99.3 (65.72-2.40)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.80 (at 2.40Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.184 , 0.239 0.193 , 0.244	Depositor DCC
$R_{free}$ test set	3521 reflections (4.22%)	DCC
Wilson B-factor (Å <sup>2</sup> )	46.0	Xtriage
Anisotropy	0.463	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 35.2	EDS
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 86984 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	12080	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% 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, 0.375 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	0.76	2/5772 (0.0%)	0.90	16/7805 (0.2%)
1	B	0.75	1/5772 (0.0%)	0.86	15/7805 (0.2%)
All	All	0.76	3/11544 (0.0%)	0.88	31/15610 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	209	ASP	CB-CG	7.69	1.67	1.51
1	A	151	CYS	CB-SG	-6.77	1.70	1.82
1	A	209	ASP	CB-CG	5.54	1.63	1.51

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	576	ARG	NE-CZ-NH1	-12.12	114.24	120.30
1	A	354	ARG	NE-CZ-NH2	-10.82	114.89	120.30
1	A	156	PRO	C-N-CD	-10.74	96.96	120.60
1	B	49	ARG	NE-CZ-NH2	-9.89	115.36	120.30
1	B	463	LEU	CA-CB-CG	9.67	137.54	115.30
1	A	213	ARG	NE-CZ-NH2	-9.05	115.78	120.30
1	B	49	ARG	NE-CZ-NH1	8.92	124.76	120.30
1	A	354	ARG	NE-CZ-NH1	8.50	124.55	120.30
1	A	463	LEU	CA-CB-CG	6.84	131.03	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	576	ARG	NE-CZ-NH2	6.79	123.69	120.30
1	A	49	ARG	NE-CZ-NH2	-6.49	117.06	120.30
1	A	213	ARG	NE-CZ-NH1	6.23	123.41	120.30
1	B	710	ARG	NE-CZ-NH2	-6.22	117.19	120.30
1	B	482	TYR	CA-C-N	6.07	128.34	116.20
1	A	489	ILE	CG1-CB-CG2	-5.99	98.23	111.40
1	B	365	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	B	710	ARG	NE-CZ-NH1	5.90	123.25	120.30
1	A	320	ARG	NE-CZ-NH1	5.86	123.23	120.30
1	B	156	PRO	C-N-CD	-5.80	107.83	120.60
1	B	482	TYR	CB-CA-C	5.61	121.63	110.40
1	B	354	ARG	NE-CZ-NH1	5.57	123.08	120.30
1	B	354	ARG	NE-CZ-NH2	-5.37	117.62	120.30
1	B	576	ARG	NE-CZ-NH1	-5.25	117.68	120.30
1	B	209	ASP	CB-CA-C	5.24	120.88	110.40
1	A	320	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	B	552	THR	CB-CA-C	-5.19	97.58	111.60
1	A	482	TYR	CB-CA-C	5.19	120.77	110.40
1	A	482	TYR	CA-C-N	5.18	126.55	116.20
1	A	342	VAL	CB-CA-C	-5.17	101.58	111.40
1	B	482	TYR	C-N-CA	5.14	133.09	122.30
1	A	664	ARG	NE-CZ-NH2	-5.00	117.80	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	156	PRO	Peptide
1	A	482	TYR	Peptide
1	B	156	PRO	Peptide
1	B	179	VAL	Peptide
1	B	482	TYR	Peptide

## 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	5651	0	5536	99	0
1	B	5651	0	5536	118	0
2	A	397	0	0	19	0
2	B	381	0	0	21	0
All	All	12080	0	11072	216	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 10.

All (216) 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:576:ARG:HH11	1:A:576:ARG:CG	1.47	1.26
1:B:576:ARG:HH11	1:B:576:ARG:HG2	1.03	1.19
1:A:180:ARG:HD2	1:A:180:ARG:N	1.48	1.12
1:A:614:PRO:HB3	1:A:623:MSE:CE	1.78	1.12
1:A:576:ARG:NH1	1:A:576:ARG:HG2	1.46	1.11
1:A:614:PRO:CB	1:A:623:MSE:HE1	1.83	1.09
1:A:365:ARG:H	1:A:376:THR:HG21	1.17	1.08
1:A:614:PRO:HB3	1:A:623:MSE:HE1	1.28	1.05
1:A:180:ARG:HD2	1:A:180:ARG:H	0.99	1.02
1:A:365:ARG:H	1:A:376:THR:CG2	1.81	0.94
1:B:61:GLU:HG2	1:B:675:LEU:HD23	1.50	0.93
1:B:641:MSE:CE	1:B:699:GLN:HG3	2.02	0.90
1:B:576:ARG:NH1	1:B:576:ARG:HG2	1.87	0.89
1:B:576:ARG:HH11	1:B:576:ARG:CG	1.83	0.89
1:B:595:THR:HG22	1:B:596:MSE:HE2	1.56	0.88
1:B:641:MSE:HE2	1:B:699:GLN:HG3	1.56	0.88
1:A:180:ARG:N	1:A:180:ARG:CD	2.36	0.87
1:B:365:ARG:HB2	1:B:376:THR:HG21	1.58	0.86
1:B:68:ARG:HG2	1:B:674:LEU:HD21	1.57	0.86
1:B:388:THR:HG21	2:B:2247:HOH:O	1.73	0.86
1:B:302:ILE:HD11	1:B:342:VAL:HG11	1.58	0.86
1:B:483:GLY:HA2	2:B:2300:HOH:O	1.78	0.83
1:B:614:PRO:CA	1:B:623:MSE:HE1	2.08	0.83
1:B:641:MSE:HE2	1:B:699:GLN:CG	2.09	0.82
1:A:361:VAL:HG22	1:A:380:LEU:HD22	1.61	0.82
1:A:595:THR:HG22	1:A:596:MSE:HE2	1.62	0.81
1:A:388:THR:HG21	2:A:2255:HOH:O	1.82	0.80
1:A:365:ARG:N	1:A:376:THR:HG21	1.95	0.80
1:B:331:ILE:HD13	1:B:351:LEU:HD21	1.62	0.79
1:B:412:THR:HG22	1:B:413:THR:OG1	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:THR:HG21	1:A:389:ALA:H	1.48	0.78
1:A:587:VAL:O	1:A:651:VAL:HG21	1.82	0.77
1:B:100:ARG:HG3	1:B:100:ARG:HH11	1.51	0.75
1:A:210:ALA:HA	2:A:2154:HOH:O	1.87	0.75
1:A:447:ARG:HB2	2:A:2279:HOH:O	1.87	0.74
1:A:354:ARG:NH2	1:A:490:GLU:OE2	2.19	0.74
2:A:2358:HOH:O	1:B:576:ARG:HD2	1.87	0.74
1:A:576:ARG:HH11	1:A:576:ARG:HG2	0.64	0.74
1:B:304:LEU:CD2	1:B:341:ASP:HA	2.19	0.72
1:A:331:ILE:CD1	1:A:351:LEU:HD21	2.19	0.71
1:B:162:VAL:HG13	1:B:179:VAL:HG21	1.72	0.71
1:B:619:PHE:HB3	1:B:623:MSE:HE2	1.71	0.71
1:B:614:PRO:HB3	1:B:623:MSE:HE3	1.72	0.71
1:A:641:MSE:HE2	1:A:699:GLN:HG3	1.71	0.70
1:B:440:GLU:HB3	2:B:2272:HOH:O	1.91	0.70
1:A:61:GLU:HG2	1:A:675:LEU:HD23	1.73	0.69
1:B:388:THR:CG2	2:B:2246:HOH:O	2.39	0.69
1:A:359:THR:CG2	1:A:389:ALA:H	2.06	0.69
1:B:331:ILE:CD1	1:B:351:LEU:HD21	2.21	0.68
1:A:322:ALA:HB1	2:A:2100:HOH:O	1.93	0.68
1:A:595:THR:HG22	1:A:596:MSE:CE	2.23	0.68
1:B:388:THR:HG22	2:B:2246:HOH:O	1.95	0.67
1:B:290:ARG:H	1:B:306:ASN:HD21	1.42	0.67
1:B:641:MSE:HE1	1:B:699:GLN:HG3	1.76	0.67
1:B:269:ASP:OD1	1:B:271:ARG:HD3	1.95	0.66
1:A:213:ARG:NH2	1:A:235:ASP:O	2.26	0.66
1:A:331:ILE:HD13	1:A:351:LEU:HD21	1.78	0.66
1:B:290:ARG:H	1:B:306:ASN:ND2	1.93	0.66
1:A:373:SER:O	1:A:376:THR:HB	1.96	0.66
1:A:310:CYS:SG	1:A:314:LYS:HD3	2.37	0.65
1:B:304:LEU:HD22	1:B:341:ASP:HA	1.77	0.65
1:B:16:VAL:HG22	1:B:18:PHE:CE1	2.32	0.65
1:B:614:PRO:N	1:B:623:MSE:HE1	2.11	0.64
1:A:614:PRO:CA	1:A:623:MSE:HE1	2.27	0.64
1:A:269:ASP:OD1	1:A:271:ARG:HD3	1.96	0.64
1:A:15:GLU:OE2	1:A:36:ARG:HD3	1.97	0.64
1:B:614:PRO:HB3	1:B:623:MSE:CE	2.27	0.64
1:A:388:THR:CG2	2:A:2254:HOH:O	2.46	0.64
1:B:175:SER:HB3	1:B:188:GLU:OE2	1.98	0.63
1:A:617:TYR:HB3	1:B:576:ARG:HH12	1.63	0.63
1:A:412:THR:HG22	1:A:413:THR:OG1	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:576:ARG:HD2	2:A:2346:HOH:O	1.99	0.61
1:B:475:ASN:O	1:B:552:THR:HG23	2.00	0.61
1:B:584:LEU:HD22	1:B:699:GLN:HG2	1.81	0.61
1:A:348:PHE:CE2	1:A:422:VAL:HG21	2.35	0.61
1:B:209:ASP:HB2	1:B:215:ASN:ND2	2.15	0.61
1:A:576:ARG:NH1	1:A:576:ARG:CG	2.23	0.60
1:A:16:VAL:HG22	1:A:18:PHE:CE1	2.36	0.60
1:B:162:VAL:HG13	1:B:179:VAL:CG2	2.32	0.60
1:B:475:ASN:O	1:B:552:THR:CG2	2.49	0.60
1:B:302:ILE:HG12	1:B:344:VAL:HG21	1.83	0.60
1:B:614:PRO:HA	1:B:623:MSE:HE1	1.84	0.60
1:A:73:LYS:HE2	2:A:2069:HOH:O	2.01	0.59
1:A:614:PRO:CG	1:A:623:MSE:HE1	2.34	0.58
1:A:475:ASN:O	1:A:552:THR:CG2	2.51	0.58
1:B:476:PRO:HB2	1:B:505:MSE:HG2	1.85	0.57
1:B:105:THR:HG22	2:B:2097:HOH:O	2.05	0.57
1:B:95:PRO:O	1:B:405:ARG:NH2	2.35	0.57
1:A:475:ASN:O	1:A:552:THR:HG23	2.04	0.57
1:B:354:ARG:NH2	1:B:490:GLU:OE2	2.38	0.57
1:B:37:ARG:NH2	2:B:2038:HOH:O	2.38	0.56
1:B:400:ASP:HB2	2:B:2255:HOH:O	2.06	0.56
1:A:537:MSE:HE3	2:A:2333:HOH:O	2.06	0.56
1:A:434:LYS:HG3	2:A:2269:HOH:O	2.06	0.56
1:B:100:ARG:NH1	1:B:100:ARG:HG3	2.21	0.55
1:B:16:VAL:HG22	1:B:18:PHE:HE1	1.71	0.55
1:B:576:ARG:NH1	1:B:576:ARG:CG	2.54	0.55
1:A:280:GLU:OE1	1:A:320:ARG:HD3	2.06	0.55
1:A:119:VAL:HG13	1:A:123:LYS:HB2	1.88	0.55
1:B:64:TYR:O	1:B:68:ARG:HB2	2.06	0.55
1:A:614:PRO:CD	1:A:623:MSE:HE1	2.36	0.55
1:A:62:LYS:O	1:A:66:GLU:HG2	2.06	0.55
1:B:162:VAL:O	1:B:179:VAL:HG22	2.06	0.54
1:B:154:PRO:HA	1:B:162:VAL:HB	1.88	0.54
1:B:589:PHE:CZ	1:B:596:MSE:HE3	2.43	0.54
1:B:655:GLU:HB2	1:B:656:PRO:CD	2.36	0.54
1:A:264:GLU:OE2	1:A:285:ARG:HB2	2.07	0.54
1:B:655:GLU:HB2	1:B:656:PRO:HD3	1.90	0.54
1:A:315:VAL:HG13	1:A:331:ILE:HB	1.89	0.54
1:A:198:PRO:HD2	2:A:2146:HOH:O	2.08	0.54
1:B:474:PRO:HB2	1:B:556:GLN:NE2	2.23	0.53
1:B:68:ARG:HG2	1:B:674:LEU:CD2	2.33	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:TYR:CZ	1:A:700:GLN:HG2	2.43	0.53
1:B:614:PRO:CB	1:B:623:MSE:HE1	2.38	0.52
1:A:347:LYS:HE3	1:A:400:ASP:OD1	2.09	0.52
1:A:354:ARG:HH22	1:A:490:GLU:CD	2.12	0.52
1:B:426:GLU:HG3	2:B:2266:HOH:O	2.10	0.52
1:B:264:GLU:OE1	1:B:285:ARG:HB2	2.09	0.52
1:B:641:MSE:HE2	1:B:699:GLN:HG2	1.87	0.52
1:A:641:MSE:HE1	1:A:698:ILE:HG23	1.92	0.51
1:B:355:ARG:NH1	1:B:379:GLU:OE2	2.43	0.51
1:A:302:ILE:HG12	1:A:344:VAL:HG21	1.91	0.51
1:B:65:TYR:CZ	1:B:676:LYS:HD2	2.46	0.51
1:A:687:SER:HB2	1:A:692:TYR:CE1	2.46	0.51
1:B:258:MSE:HG2	2:B:2181:HOH:O	2.11	0.50
1:A:655:GLU:HB2	1:A:656:PRO:CD	2.42	0.50
1:B:331:ILE:HD13	1:B:351:LEU:CD2	2.38	0.50
1:A:388:THR:HG22	2:A:2254:HOH:O	2.11	0.50
1:B:365:ARG:H	1:B:376:THR:CG2	2.25	0.50
1:A:587:VAL:O	1:A:651:VAL:CG2	2.56	0.49
1:A:95:PRO:HA	1:A:103:TYR:O	2.12	0.49
1:B:614:PRO:CB	1:B:623:MSE:CE	2.91	0.49
1:A:476:PRO:HB2	1:A:505:MSE:HG2	1.94	0.49
1:B:190:THR:HG22	1:B:205:TYR:CD2	2.48	0.48
1:A:18:PHE:CE1	1:A:37:ARG:HD2	2.49	0.48
1:A:302:ILE:HG12	1:A:344:VAL:CG2	2.44	0.48
1:A:338:THR:HG21	2:A:2227:HOH:O	2.13	0.47
1:A:338:THR:O	1:A:353:GLY:HA3	2.13	0.47
1:B:3:THR:HG22	2:B:2368:HOH:O	2.13	0.47
1:B:80:TYR:CD2	1:B:694:ARG:NH2	2.81	0.47
1:A:157:PRO:HD2	2:A:2147:HOH:O	2.15	0.47
1:A:136:ASP:HB3	1:A:139:LYS:HG3	1.97	0.47
1:B:477:THR:OG1	1:B:552:THR:HG23	2.14	0.46
1:A:576:ARG:CD	2:A:2346:HOH:O	2.61	0.46
1:A:359:THR:HG21	1:A:388:THR:HA	1.96	0.46
1:B:385:PRO:HG2	1:B:447:ARG:HA	1.98	0.46
1:B:340:ASP:O	1:B:341:ASP:HB3	2.16	0.46
1:B:450:LEU:HD22	1:B:458:LYS:HB3	1.97	0.46
1:B:473:LYS:HB2	1:B:474:PRO:CD	2.45	0.46
1:B:681:SER:OG	1:B:691:LYS:NZ	2.48	0.46
1:A:68:ARG:HG2	1:A:674:LEU:HD21	1.97	0.46
1:B:446:CYS:HA	1:B:463:LEU:O	2.16	0.46
1:A:186:LYS:HE2	1:A:188:GLU:OE2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:VAL:CG2	1:B:18:PHE:CE1	2.97	0.45
1:A:677:MSE:HB3	2:A:2059:HOH:O	2.16	0.45
1:B:614:PRO:CA	1:B:623:MSE:CE	2.90	0.45
1:A:587:VAL:HB	1:A:651:VAL:HG23	1.98	0.45
1:B:477:THR:HB	1:B:552:THR:HG21	1.99	0.45
1:B:619:PHE:CB	1:B:623:MSE:HE2	2.45	0.45
1:A:655:GLU:HB2	1:A:656:PRO:HD3	1.99	0.45
1:A:361:VAL:HG22	1:A:380:LEU:CD2	2.39	0.45
1:A:269:ASP:OD1	1:A:271:ARG:CD	2.62	0.45
1:B:89:GLU:HB2	2:B:2086:HOH:O	2.17	0.44
1:B:483:GLY:CA	2:B:2300:HOH:O	2.51	0.44
1:B:209:ASP:HB2	1:B:215:ASN:HD21	1.81	0.44
1:A:93:SER:HB3	1:A:105:THR:HG22	1.98	0.44
1:B:418:TYR:HB2	1:B:428:LYS:O	2.18	0.44
1:B:619:PHE:HB3	1:B:623:MSE:CE	2.43	0.44
1:B:291:TYR:HA	1:B:304:LEU:O	2.18	0.44
1:B:61:GLU:HG3	1:B:675:LEU:HB3	1.98	0.44
1:A:650:ARG:O	1:A:651:VAL:HG22	2.18	0.44
1:A:209:ASP:HB3	2:A:2157:HOH:O	2.18	0.44
1:A:614:PRO:HB3	1:A:623:MSE:HE3	1.84	0.44
1:A:367:GLY:C	1:A:369:ASP:H	2.21	0.44
1:B:180:ARG:HB2	2:B:2129:HOH:O	2.17	0.44
1:B:306:ASN:C	1:B:306:ASN:HD22	2.22	0.43
1:B:540:ILE:HG21	1:B:576:ARG:HD3	2.00	0.43
1:B:100:ARG:HB3	2:B:2095:HOH:O	2.17	0.43
1:A:124:THR:HA	1:A:125:PRO:HD3	1.93	0.43
1:A:370:ASN:O	1:A:370:ASN:ND2	2.50	0.43
1:B:84:ILE:HA	1:B:87:ILE:HD12	2.00	0.43
1:A:388:THR:HG23	2:A:2254:HOH:O	2.12	0.43
1:B:563:SER:HA	1:B:587:VAL:O	2.18	0.43
1:B:230:VAL:HG11	1:B:277:ASN:ND2	2.33	0.43
1:B:540:ILE:CG2	1:B:576:ARG:HD3	2.48	0.42
1:B:212:LYS:HG2	2:B:2127:HOH:O	2.19	0.42
1:B:108:VAL:HB	1:B:111:LEU:HD12	2.01	0.42
1:A:584:LEU:HD22	1:A:699:GLN:HG2	2.00	0.42
1:B:434:LYS:HE3	2:B:2087:HOH:O	2.19	0.42
1:B:41:LEU:O	1:B:658:LYS:HD3	2.18	0.42
1:B:119:VAL:HG13	1:B:123:LYS:HB2	2.00	0.42
1:A:101:PHE:CD1	1:A:118:ARG:HG3	2.55	0.42
1:B:373:SER:O	1:B:376:THR:HB	2.19	0.42
1:B:518:MSE:HB2	1:B:522:TRP:CD2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:439:PHE:CE1	1:B:444:TYR:CE2	3.07	0.42
1:B:338:THR:O	1:B:353:GLY:HA3	2.20	0.42
1:A:179:VAL:HG13	1:A:180:ARG:HD3	2.01	0.42
1:B:114:LYS:HB3	1:B:114:LYS:HE3	1.79	0.42
1:B:141:ALA:O	1:B:144:LYS:HB2	2.20	0.42
1:B:158:GLU:HB2	2:B:2118:HOH:O	2.20	0.42
1:A:61:GLU:OE2	1:A:664:ARG:NH2	2.52	0.41
1:A:568:LEU:HD23	1:A:568:LEU:C	2.40	0.41
1:B:179:VAL:HG23	2:B:2120:HOH:O	2.20	0.41
1:B:475:ASN:O	1:B:552:THR:HG22	2.20	0.41
1:A:348:PHE:CE1	1:A:404:LEU:HD21	2.55	0.41
1:B:447:ARG:HB2	2:B:2279:HOH:O	2.20	0.41
1:A:448:ARG:HD3	2:A:2280:HOH:O	2.19	0.41
1:A:473:LYS:HE2	1:A:473:LYS:HB3	1.49	0.41
1:A:76:ALA:HB2	1:A:698:ILE:HD12	2.03	0.41
1:B:586:GLY:O	1:B:587:VAL:C	2.59	0.41
1:A:641:MSE:CE	1:A:699:GLN:HG3	2.46	0.41
1:B:180:ARG:HH11	1:B:180:ARG:HB3	1.86	0.41
1:A:648:ASP:OD1	1:A:649:PRO:HD2	2.21	0.40
1:B:624:ASN:HB3	2:B:2352:HOH:O	2.21	0.40
1:A:614:PRO:N	1:A:623:MSE:HE1	2.37	0.40
1:B:365:ARG:H	1:B:376:THR:HG21	1.85	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	711/715 (99%)	674 (95%)	29 (4%)	8 (1%)	17	25
1	B	711/715 (99%)	673 (95%)	30 (4%)	8 (1%)	17	25
All	All	1422/1430 (99%)	1347 (95%)	59 (4%)	16 (1%)	17	25

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	157	PRO
1	A	687	SER
1	B	157	PRO
1	A	121	ALA
1	A	285	ARG
1	A	688	ASP
1	B	180	ARG
1	B	285	ARG
1	B	121	ALA
1	B	687	SER
1	B	282	VAL
1	B	502	ASP
1	A	282	VAL
1	A	179	VAL
1	B	156	PRO
1	A	368	PRO

### 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	616/601 (102%)	567 (92%)	49 (8%)	15	23
1	B	616/601 (102%)	571 (93%)	45 (7%)	17	27
All	All	1232/1202 (102%)	1138 (92%)	94 (8%)	16	25

All (94) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	16	VAL
1	A	37	ARG
1	A	49	ARG
1	A	98	TYR
1	A	105	THR
1	A	118	ARG

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Mol	Chain	Res	Type
1	A	119	VAL
1	A	162	VAL
1	A	180	ARG
1	A	185	ASP
1	A	209	ASP
1	A	212	LYS
1	A	241	SER
1	A	251	LYS
1	A	263	SER
1	A	278	THR
1	A	290	ARG
1	A	300	THR
1	A	302	ILE
1	A	306	ASN
1	A	315	VAL
1	A	317	LEU
1	A	320	ARG
1	A	336	LYS
1	A	342	VAL
1	A	354	ARG
1	A	359	THR
1	A	361	VAL
1	A	363	THR
1	A	380	LEU
1	A	388	THR
1	A	398	THR
1	A	412	THR
1	A	441	SER
1	A	463	LEU
1	A	467	THR
1	A	473	LYS
1	A	496	ARG
1	A	497	PHE
1	A	525	VAL
1	A	552	THR
1	A	576	ARG
1	A	633	ARG
1	A	688	ASP
1	A	694	ARG
1	A	696	ASN
1	A	707	LEU
1	A	708	ASN

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Mol	Chain	Res	Type
1	B	37	ARG
1	B	49	ARG
1	B	98	TYR
1	B	100	ARG
1	B	105	THR
1	B	114	LYS
1	B	118	ARG
1	B	119	VAL
1	B	139	LYS
1	B	158	GLU
1	B	162	VAL
1	B	180	ARG
1	B	185	ASP
1	B	211	SER
1	B	232	LEU
1	B	241	SER
1	B	263	SER
1	B	300	THR
1	B	302	ILE
1	B	306	ASN
1	B	315	VAL
1	B	336	LYS
1	B	351	LEU
1	B	363	THR
1	B	374	SER
1	B	388	THR
1	B	393	CYS
1	B	398	THR
1	B	405	ARG
1	B	412	THR
1	B	428	LYS
1	B	463	LEU
1	B	476	PRO
1	B	497	PHE
1	B	525	VAL
1	B	552	THR
1	B	576	ARG
1	B	611	TRP
1	B	633	ARG
1	B	688	ASP
1	B	694	ARG
1	B	696	ASN

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Mol	Chain	Res	Type
1	B	707	LEU
1	B	708	ASN
1	B	709	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	224	GLN
1	A	306	ASN
1	A	370	ASN
1	A	643	GLN
1	A	699	GLN
1	B	224	GLN
1	B	306	ASN
1	B	370	ASN
1	B	556	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 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	696/715 (97%)	-0.31	4 (0%)	90 90	30, 45, 68, 90	0
1	B	696/715 (97%)	-0.32	2 (0%)	94 94	29, 46, 67, 92	0
All	All	1392/1430 (97%)	-0.31	6 (0%)	93 93	29, 45, 68, 92	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	THR	4.0
1	B	3	THR	3.0
1	A	368	PRO	2.8
1	A	4	GLU	2.2
1	B	209	ASP	2.1
1	A	370	ASN	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

There are no such residues in this entry.