



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 09:09 am GMT

PDB ID : 3ND0  
Title : X-ray crystal structure of a slow cyanobacterial Cl<sup>-</sup>/H<sup>+</sup> antiporter  
Authors : Jayaram, H.; Robertson, J.; Wu, F.; Williams, C.; Miller, C.  
Deposited on : 2010-06-06  
Resolution : 3.20 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

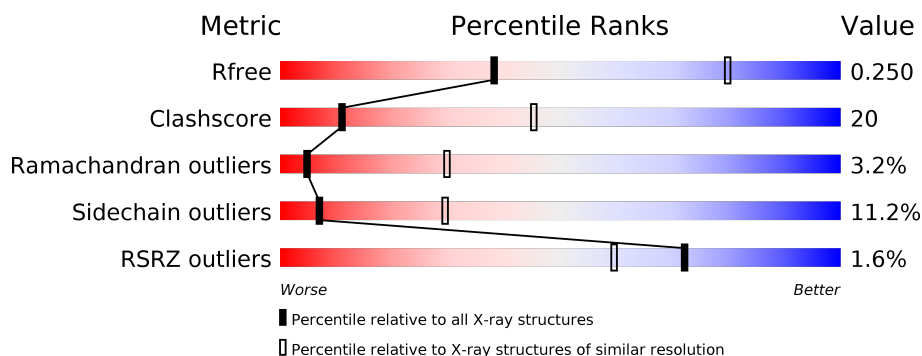
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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))
$R_{free}$	100719	1015 (3.22-3.18)
Clashscore	112137	1009 (3.20-3.20)
Ramachandran outliers	110173	1118 (3.22-3.18)
Sidechain outliers	110143	1117 (3.22-3.18)
RSRZ outliers	101464	1020 (3.22-3.18)

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	466	<div> <div>2%</div> <div> <div></div> <div>54%</div> <div>30%</div> <div>7%</div> <div>9%</div> </div> </div>
1	B	466	<div> <div>2%</div> <div> <div></div> <div>58%</div> <div>27%</div> <div>6%</div> <div>9%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6400 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 Sll0855 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	425	Total	C	N	O	S	0	0	0
			3199	2110	533	536	20			
1	B	425	Total	C	N	O	S	0	0	0
			3199	2110	533	536	20			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	452	SER	-	EXPRESSION TAG	UNP P73745
A	453	LEU	-	EXPRESSION TAG	UNP P73745
A	454	VAL	-	EXPRESSION TAG	UNP P73745
A	455	PRO	-	EXPRESSION TAG	UNP P73745
A	456	ARG	-	EXPRESSION TAG	UNP P73745
A	457	GLY	-	EXPRESSION TAG	UNP P73745
A	458	SER	-	EXPRESSION TAG	UNP P73745
A	459	GLY	-	EXPRESSION TAG	UNP P73745
A	460	GLY	-	EXPRESSION TAG	UNP P73745
A	461	HIS	-	EXPRESSION TAG	UNP P73745
A	462	HIS	-	EXPRESSION TAG	UNP P73745
A	463	HIS	-	EXPRESSION TAG	UNP P73745
A	464	HIS	-	EXPRESSION TAG	UNP P73745
A	465	HIS	-	EXPRESSION TAG	UNP P73745
A	466	HIS	-	EXPRESSION TAG	UNP P73745
B	452	SER	-	EXPRESSION TAG	UNP P73745
B	453	LEU	-	EXPRESSION TAG	UNP P73745
B	454	VAL	-	EXPRESSION TAG	UNP P73745
B	455	PRO	-	EXPRESSION TAG	UNP P73745
B	456	ARG	-	EXPRESSION TAG	UNP P73745
B	457	GLY	-	EXPRESSION TAG	UNP P73745
B	458	SER	-	EXPRESSION TAG	UNP P73745
B	459	GLY	-	EXPRESSION TAG	UNP P73745
B	460	GLY	-	EXPRESSION TAG	UNP P73745
B	461	HIS	-	EXPRESSION TAG	UNP P73745

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Chain	Residue	Modelled	Actual	Comment	Reference
B	462	HIS	-	EXPRESSION TAG	UNP P73745
B	463	HIS	-	EXPRESSION TAG	UNP P73745
B	464	HIS	-	EXPRESSION TAG	UNP P73745
B	465	HIS	-	EXPRESSION TAG	UNP P73745
B	466	HIS	-	EXPRESSION TAG	UNP P73745

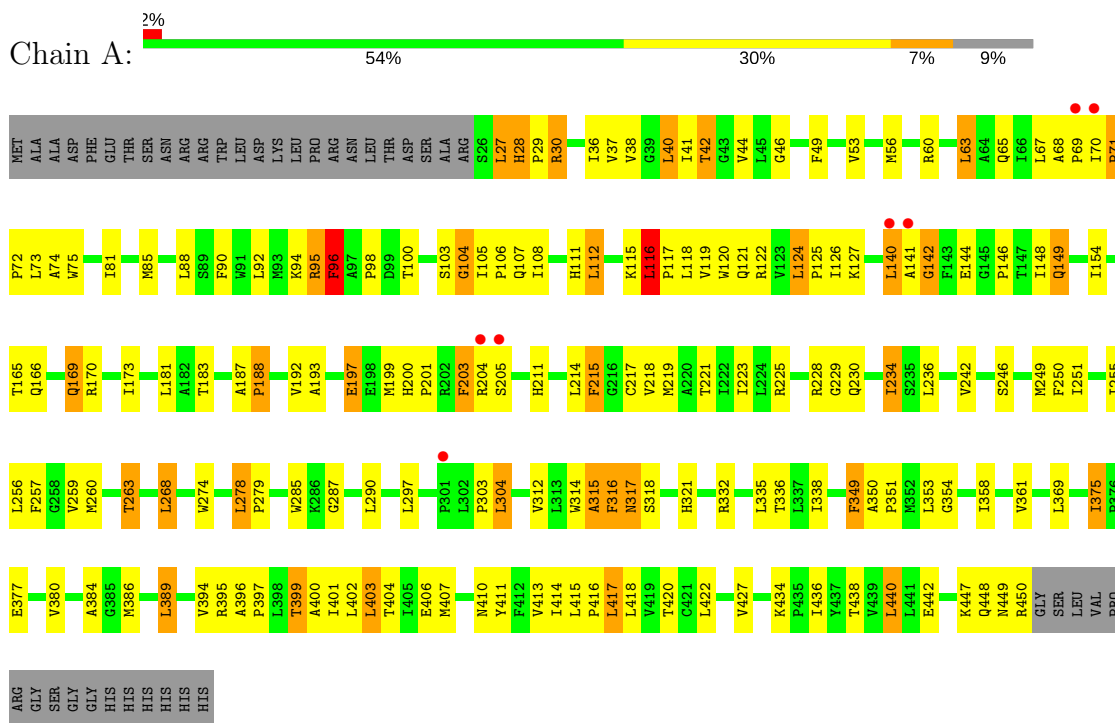
- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Cl 1 1	0	0
2	A	1	Total Cl 1 1	0	0

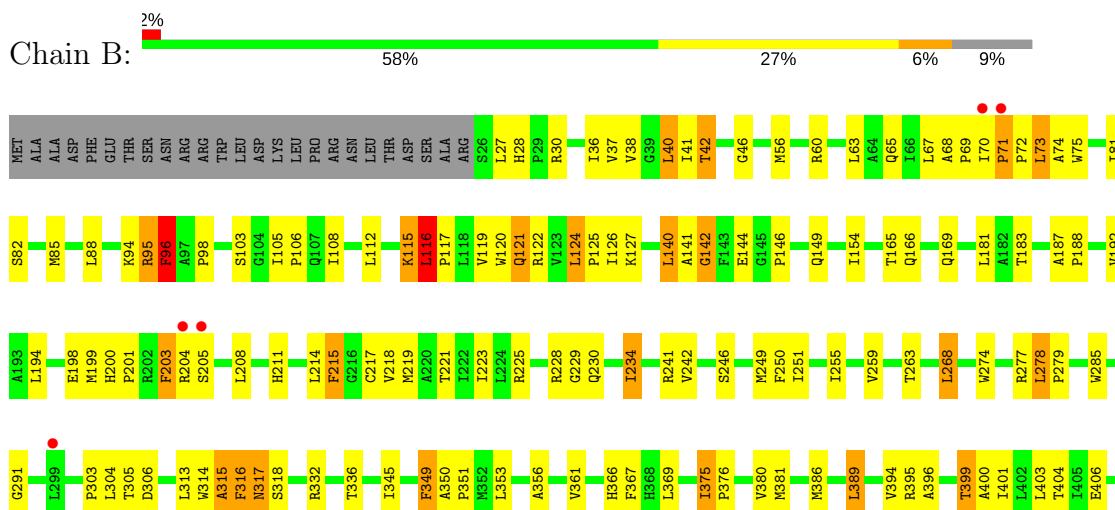
### 3 Residue-property plots [i](#)

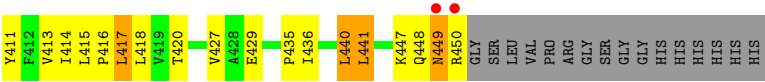
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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 ( $RSRZ > 2$ ). 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: Sll0855 protein



#### • Molecule 1: Sll0855 protein





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	203.79Å 203.79Å 96.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	65.22 – 3.20 66.71 – 3.20	Depositor EDS
% Data completeness (in resolution range)	91.4 (65.22-3.20) 91.4 (66.71-3.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.26	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 3.19Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, $R_{free}$	0.243 , 0.269 0.227 , 0.250	Depositor DCC
$R_{free}$ test set	1748 reflections (4.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	82.2	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 55.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.052 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.86	EDS
Total number of atoms	6400	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.25	0/3276	0.42	0/4457
1	B	0.24	0/3276	0.42	0/4457
All	All	0.25	0/6552	0.42	0/8914

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	3199	0	3360	133	0
1	B	3199	0	3360	135	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
All	All	6400	0	6720	263	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 20.

All (263) 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:B:225:ARG:HD3	1:B:230:GLN:HG3	1.33	1.08
1:A:225:ARG:HD3	1:A:230:GLN:HG3	1.36	1.04
1:A:71:PRO:HB2	1:A:72:PRO:CD	1.99	0.91
1:B:71:PRO:HB2	1:B:72:PRO:CD	2.03	0.88
1:B:95:ARG:HH22	1:B:450:ARG:HH22	1.18	0.88
1:A:218:VAL:HG22	1:B:418:LEU:HD13	1.57	0.86
1:A:71:PRO:HB2	1:A:72:PRO:HD2	1.58	0.84
1:A:278:LEU:HB3	1:A:279:PRO:HA	1.60	0.83
1:A:278:LEU:HB3	1:A:279:PRO:CA	2.09	0.82
1:A:95:ARG:HH22	1:A:450:ARG:HH22	1.23	0.82
1:A:418:LEU:HD13	1:B:218:VAL:HG22	1.64	0.79
1:A:144:GLU:HG3	1:A:349:PHE:HB2	1.64	0.79
1:B:71:PRO:HB2	1:B:72:PRO:HD2	1.66	0.76
1:B:278:LEU:HB3	1:B:279:PRO:CA	2.16	0.76
1:B:144:GLU:HG3	1:B:349:PHE:HB2	1.70	0.74
1:B:375:ILE:HG22	1:B:375:ILE:O	1.88	0.73
1:A:395:ARG:O	1:A:395:ARG:HG2	1.89	0.73
1:A:188:PRO:HG3	1:A:221:THR:HG21	1.72	0.72
1:B:188:PRO:HG3	1:B:221:THR:HG21	1.71	0.72
1:A:192:VAL:HG21	1:A:214:LEU:HD23	1.73	0.71
1:A:375:ILE:CG2	1:A:375:ILE:O	2.39	0.70
1:B:395:ARG:HG2	1:B:395:ARG:O	1.91	0.70
1:B:116:LEU:HD23	1:B:117:PRO:HD2	1.74	0.70
1:A:375:ILE:HG22	1:A:375:ILE:O	1.92	0.69
1:B:303:PRO:HG3	1:B:314:TRP:CG	2.27	0.69
1:B:278:LEU:HB3	1:B:279:PRO:HA	1.74	0.69
1:B:268:LEU:HD12	1:B:436:ILE:HD11	1.74	0.68
1:B:259:VAL:O	1:B:263:THR:HG23	1.94	0.68
1:A:141:ALA:HB1	1:A:142:GLY:HA3	1.76	0.66
1:A:68:ALA:N	1:A:69:PRO:HD2	2.10	0.66
1:A:303:PRO:HG3	1:A:314:TRP:CG	2.30	0.66
1:A:127:LYS:HD2	1:A:146:PRO:HA	1.79	0.65
1:B:95:ARG:NH2	1:B:450:ARG:HH12	1.95	0.65
1:B:68:ALA:N	1:B:69:PRO:HD2	2.10	0.65
1:A:278:LEU:CB	1:A:279:PRO:HA	2.27	0.65
1:B:115:LYS:HD2	1:B:115:LYS:H	1.62	0.65
1:B:95:ARG:HH22	1:B:450:ARG:NH2	1.93	0.65
1:A:401:ILE:HG23	1:A:417:LEU:HD13	1.79	0.64
1:A:111:HIS:ND1	1:A:118:LEU:HB3	2.11	0.64
1:A:259:VAL:O	1:A:263:THR:HG23	1.96	0.64
1:B:375:ILE:O	1:B:375:ILE:CG2	2.46	0.64
1:B:94:LYS:HB3	1:B:98:PRO:HG3	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:411:TYR:CE1	1:A:414:ILE:HD12	2.32	0.64
1:A:251:ILE:O	1:A:255:ILE:HG12	1.97	0.63
1:A:394:VAL:O	1:A:395:ARG:HB3	1.98	0.62
1:B:192:VAL:HG21	1:B:214:LEU:HD23	1.81	0.62
1:A:422:LEU:HD11	1:B:215:PHE:CE1	2.33	0.62
1:B:141:ALA:HB1	1:B:142:GLY:HA3	1.82	0.61
1:B:278:LEU:HB3	1:B:279:PRO:C	2.21	0.61
1:A:28:HIS:NE2	1:A:30:ARG:HB2	2.15	0.61
1:B:278:LEU:CB	1:B:279:PRO:HA	2.30	0.61
1:A:413:VAL:O	1:A:413:VAL:HG12	2.00	0.60
1:A:278:LEU:CB	1:A:279:PRO:CA	2.80	0.60
1:B:71:PRO:CB	1:B:72:PRO:CD	2.79	0.60
1:A:225:ARG:HA	1:A:229:GLY:HA3	1.83	0.60
1:B:401:ILE:HD13	1:B:418:LEU:HD23	1.83	0.60
1:A:71:PRO:CB	1:A:72:PRO:HD2	2.31	0.60
1:B:127:LYS:HD2	1:B:146:PRO:HA	1.84	0.60
1:B:416:PRO:O	1:B:420:THR:HG22	2.02	0.59
1:B:278:LEU:CB	1:B:279:PRO:CA	2.80	0.59
1:A:268:LEU:HD12	1:A:436:ILE:HD11	1.83	0.59
1:A:72:PRO:O	1:A:75:TRP:HB2	2.03	0.58
1:A:71:PRO:CB	1:A:72:PRO:CD	2.78	0.58
1:B:394:VAL:O	1:B:395:ARG:HB3	2.03	0.58
1:A:416:PRO:O	1:A:420:THR:HG22	2.02	0.58
1:B:396:ALA:HB1	1:B:399:THR:HG23	1.86	0.58
1:A:95:ARG:NH2	1:A:450:ARG:HH22	1.96	0.58
1:B:380:VAL:HG13	1:B:413:VAL:HG13	1.85	0.58
1:B:67:LEU:HB2	1:B:74:ALA:HB2	1.84	0.57
1:A:201:PRO:HG2	1:A:203:PHE:CE1	2.40	0.57
1:A:315:ALA:HB1	1:A:361:VAL:HG21	1.86	0.57
1:B:28:HIS:CE1	1:B:30:ARG:HB2	2.39	0.57
1:A:36:ILE:HD11	1:A:215:PHE:CD1	2.40	0.56
1:A:95:ARG:HH22	1:A:450:ARG:NH2	2.00	0.56
1:B:199:MET:HG2	1:B:199:MET:O	2.05	0.56
1:B:225:ARG:HA	1:B:229:GLY:HA3	1.85	0.56
1:B:71:PRO:CB	1:B:72:PRO:HD2	2.34	0.56
1:B:214:LEU:O	1:B:218:VAL:HG23	2.05	0.56
1:B:72:PRO:O	1:B:75:TRP:HB2	2.05	0.56
1:A:404:THR:OG1	1:A:417:LEU:HD11	2.05	0.56
1:A:40:LEU:HG	1:A:219:MET:HE2	1.85	0.56
1:B:187:ALA:HB1	1:B:406:GLU:OE1	2.05	0.56
1:A:37:VAL:O	1:A:41:ILE:HG12	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:396:ALA:HB1	1:A:399:THR:CG2	2.37	0.55
1:B:242:VAL:HG11	1:B:416:PRO:HG2	1.88	0.55
1:A:400:ALA:O	1:A:404:THR:HG23	2.06	0.55
1:A:108:ILE:O	1:A:112:LEU:HB2	2.07	0.55
1:A:67:LEU:HB2	1:A:74:ALA:HB2	1.88	0.54
1:B:350:ALA:HB3	1:B:351:PRO:HD3	1.89	0.54
1:B:217:CYS:O	1:B:221:THR:HG23	2.07	0.54
1:B:41:ILE:HB	1:B:154:ILE:HG21	1.89	0.54
1:B:96:PHE:O	1:B:122:ARG:HD2	2.08	0.54
1:A:214:LEU:O	1:A:218:VAL:HG23	2.08	0.53
1:B:413:VAL:O	1:B:413:VAL:HG12	2.08	0.53
1:B:251:ILE:O	1:B:255:ILE:HG12	2.08	0.53
1:B:255:ILE:HD13	1:B:427:VAL:HG21	1.88	0.53
1:B:201:PRO:HG2	1:B:203:PHE:CE1	2.43	0.53
1:B:56:MET:O	1:B:60:ARG:HG3	2.08	0.53
1:A:386:MET:HE3	1:A:404:THR:HB	1.90	0.53
1:B:140:LEU:HD21	1:B:305:THR:HA	1.90	0.53
1:A:71:PRO:HD2	1:A:73:LEU:H	1.73	0.53
1:A:187:ALA:HB1	1:A:406:GLU:OE1	2.08	0.53
1:A:278:LEU:HB3	1:A:279:PRO:C	2.28	0.53
1:B:285:TRP:HA	1:B:285:TRP:CE3	2.44	0.53
1:A:255:ILE:HD13	1:A:427:VAL:HG21	1.91	0.52
1:B:46:GLY:HA3	1:B:183:THR:HG21	1.91	0.52
1:B:395:ARG:HG3	1:B:429:GLU:OE2	2.10	0.52
1:A:96:PHE:O	1:A:122:ARG:HD2	2.10	0.51
1:A:350:ALA:HB3	1:A:351:PRO:HD3	1.93	0.51
1:B:366:HIS:O	1:B:369:LEU:HB3	2.11	0.51
1:B:400:ALA:O	1:B:404:THR:HG23	2.11	0.51
1:B:28:HIS:NE2	1:B:30:ARG:HB2	2.26	0.51
1:A:396:ALA:HB1	1:A:399:THR:HG23	1.93	0.50
1:A:38:VAL:O	1:A:42:THR:HG22	2.11	0.50
1:B:144:GLU:CG	1:B:349:PHE:HB2	2.40	0.50
1:B:37:VAL:O	1:B:41:ILE:HG12	2.12	0.50
1:B:40:LEU:HG	1:B:219:MET:HE2	1.94	0.50
1:B:316:PHE:CD2	1:B:316:PHE:N	2.77	0.50
1:B:208:LEU:HD12	1:B:211:HIS:CE1	2.46	0.50
1:B:447:LYS:C	1:B:449:ASN:H	2.16	0.50
1:B:140:LEU:N	1:B:140:LEU:HD23	2.26	0.49
1:B:303:PRO:O	1:B:332:ARG:HD2	2.12	0.49
1:A:116:LEU:HD23	1:A:117:PRO:HD2	1.93	0.49
1:A:46:GLY:HA3	1:A:183:THR:HG21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:PHE:CZ	1:A:416:PRO:HB3	2.47	0.49
1:B:81:ILE:O	1:B:85:MET:HG3	2.12	0.49
1:A:242:VAL:HG22	1:A:380:VAL:HG21	1.95	0.49
1:B:140:LEU:CD2	1:B:305:THR:HA	2.43	0.49
1:A:40:LEU:HD23	1:A:223:ILE:HG13	1.94	0.49
1:B:411:TYR:CE1	1:B:414:ILE:HD12	2.47	0.49
1:B:404:THR:OG1	1:B:417:LEU:HD11	2.13	0.49
1:A:119:VAL:HG23	1:A:122:ARG:HB2	1.95	0.49
1:B:41:ILE:HG22	1:B:154:ILE:HD13	1.94	0.48
1:A:40:LEU:O	1:A:44:VAL:HG23	2.13	0.48
1:B:36:ILE:O	1:B:40:LEU:HB2	2.13	0.48
1:B:415:LEU:HB2	1:B:416:PRO:HD3	1.94	0.48
1:A:354:GLY:O	1:A:358:ILE:HG13	2.14	0.48
1:A:144:GLU:CG	1:A:349:PHE:HB2	2.38	0.48
1:B:124:LEU:HB3	1:B:125:PRO:HD3	1.96	0.48
1:A:204:ARG:HA	1:A:205:SER:HA	1.50	0.47
1:B:367:PHE:CD2	1:B:381:MET:HE1	2.49	0.47
1:A:369:LEU:HD13	1:A:369:LEU:C	2.34	0.47
1:A:41:ILE:HG22	1:A:154:ILE:HD13	1.96	0.47
1:A:94:LYS:HB3	1:A:98:PRO:HG3	1.95	0.47
1:A:380:VAL:HG13	1:A:413:VAL:HG13	1.95	0.47
1:B:401:ILE:HG23	1:B:417:LEU:HD13	1.97	0.47
1:A:438:THR:O	1:A:442:GLU:HG3	2.15	0.47
1:B:246:SER:HB2	1:B:249:MET:HE3	1.96	0.46
1:A:303:PRO:O	1:A:332:ARG:HD2	2.15	0.46
1:B:249:MET:HG2	1:B:367:PHE:CE1	2.49	0.46
1:A:95:ARG:NH2	1:A:450:ARG:HH12	2.14	0.46
1:B:82:SER:HB3	1:B:291:GLY:O	2.16	0.46
1:A:304:LEU:HD23	1:A:304:LEU:H	1.80	0.46
1:B:250:PHE:CZ	1:B:416:PRO:HB3	2.50	0.46
1:A:199:MET:O	1:A:199:MET:HG2	2.15	0.46
1:B:204:ARG:HA	1:B:205:SER:HA	1.51	0.46
1:B:140:LEU:HB3	1:B:336:THR:HG22	1.97	0.46
1:B:441:LEU:HD22	1:B:441:LEU:O	2.15	0.46
1:A:81:ILE:O	1:A:85:MET:HG3	2.15	0.46
1:A:92:LEU:O	1:A:96:PHE:HB2	2.15	0.46
1:B:332:ARG:O	1:B:336:THR:HG23	2.16	0.45
1:A:140:LEU:HB3	1:A:336:THR:HG22	1.97	0.45
1:A:436:ILE:O	1:A:440:LEU:HB2	2.16	0.45
1:A:65:GLN:HA	1:A:68:ALA:HB2	1.98	0.45
1:B:40:LEU:HD23	1:B:223:ILE:HG13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:447:LYS:HB3	1:B:450:ARG:HG2	1.99	0.45
1:A:384:ALA:HB1	1:A:420:THR:HG21	1.99	0.45
1:A:312:VAL:C	1:A:314:TRP:H	2.19	0.45
1:A:413:VAL:O	1:A:417:LEU:HB2	2.17	0.45
1:B:449:ASN:O	1:B:450:ARG:C	2.55	0.45
1:B:242:VAL:HG11	1:B:416:PRO:CG	2.46	0.45
1:A:246:SER:HB2	1:A:249:MET:HE3	1.99	0.45
1:B:395:ARG:HD3	1:B:435:PRO:HG3	2.00	0.44
1:B:121:GLN:H	1:B:121:GLN:CD	2.19	0.44
1:B:369:LEU:HD13	1:B:369:LEU:C	2.38	0.44
1:B:71:PRO:HD2	1:B:73:LEU:H	1.83	0.44
1:A:119:VAL:O	1:A:119:VAL:HG23	2.18	0.44
1:A:225:ARG:HA	1:A:229:GLY:CA	2.47	0.44
1:B:234:ILE:HG23	1:B:234:ILE:O	2.17	0.44
1:A:107:GLN:CB	1:A:149:GLN:HE21	2.31	0.44
1:A:332:ARG:O	1:A:336:THR:HG23	2.18	0.44
1:B:356:ALA:HB2	1:B:389:LEU:HB2	2.00	0.44
1:B:67:LEU:C	1:B:69:PRO:HD2	2.38	0.44
1:A:105:ILE:HB	1:A:106:PRO:HD3	2.00	0.43
1:A:403:LEU:O	1:A:407:MET:HG2	2.18	0.43
1:B:285:TRP:HA	1:B:285:TRP:HE3	1.83	0.43
1:A:68:ALA:N	1:A:69:PRO:CD	2.80	0.43
1:B:303:PRO:HG3	1:B:314:TRP:CD2	2.52	0.43
1:A:297:LEU:HD13	1:A:335:LEU:HD11	2.01	0.43
1:A:140:LEU:HD12	1:A:336:THR:HG23	2.01	0.43
1:A:197:GLU:OE1	1:A:399:THR:HG22	2.18	0.43
1:A:63:LEU:O	1:A:63:LEU:HD22	2.19	0.43
1:B:68:ALA:N	1:B:69:PRO:CD	2.81	0.43
1:B:386:MET:HB3	1:B:386:MET:HE2	1.75	0.43
1:B:229:GLY:HA3	1:B:230:GLN:HA	1.78	0.43
1:A:250:PHE:CE1	1:A:416:PRO:HB3	2.54	0.43
1:B:71:PRO:HB2	1:B:72:PRO:HD3	1.97	0.43
1:B:316:PHE:HA	1:B:317:ASN:HA	1.74	0.43
1:B:369:LEU:O	1:B:369:LEU:HD13	2.19	0.43
1:A:148:ILE:HD11	1:A:181:LEU:HB2	2.01	0.43
1:B:208:LEU:HD12	1:B:211:HIS:ND1	2.34	0.43
1:A:197:GLU:OE2	1:A:397:PRO:HD2	2.19	0.42
1:B:241:ARG:HD2	1:B:376:PRO:HG2	1.99	0.42
1:A:200:HIS:HA	1:A:201:PRO:HD3	1.86	0.42
1:B:198:GLU:O	1:B:199:MET:HB3	2.19	0.42
1:B:436:ILE:O	1:B:440:LEU:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:277:ARG:O	1:B:278:LEU:C	2.58	0.42
1:A:415:LEU:HB2	1:A:416:PRO:HD3	2.00	0.42
1:B:119:VAL:O	1:B:119:VAL:HG23	2.19	0.42
1:B:200:HIS:HA	1:B:201:PRO:HD3	1.91	0.42
1:B:225:ARG:HA	1:B:229:GLY:CA	2.49	0.42
1:A:169:GLN:O	1:A:173:ILE:HG13	2.19	0.42
1:A:56:MET:O	1:A:60:ARG:HG3	2.20	0.42
1:B:108:ILE:O	1:B:112:LEU:HB2	2.19	0.42
1:B:413:VAL:O	1:B:417:LEU:HB2	2.20	0.42
1:A:28:HIS:CE1	1:A:30:ARG:HB2	2.55	0.42
1:B:105:ILE:N	1:B:106:PRO:CD	2.82	0.42
1:B:116:LEU:HD23	1:B:117:PRO:CD	2.46	0.42
1:B:65:GLN:HA	1:B:68:ALA:HB2	2.02	0.42
1:A:410:ASN:OD1	1:A:413:VAL:HG23	2.19	0.42
1:A:100:THR:HG21	1:A:126:ILE:HG22	2.01	0.42
1:A:49:PHE:O	1:A:53:VAL:HG23	2.19	0.42
1:A:111:HIS:CD2	1:A:111:HIS:C	2.92	0.42
1:A:94:LYS:HA	1:A:95:ARG:HA	1.75	0.42
1:B:181:LEU:HD22	1:B:194:LEU:HD22	2.02	0.41
1:B:88:LEU:HD23	1:B:88:LEU:O	2.19	0.41
1:A:217:CYS:O	1:A:221:THR:HG23	2.19	0.41
1:B:105:ILE:HB	1:B:106:PRO:HD3	2.01	0.41
1:A:402:LEU:O	1:A:406:GLU:HG3	2.20	0.41
1:A:411:TYR:CE1	1:B:406:GLU:HG2	2.56	0.41
1:A:415:LEU:HD23	1:A:415:LEU:HA	1.88	0.41
1:B:305:THR:O	1:B:306:ASP:HB2	2.20	0.41
1:B:386:MET:HE2	1:B:404:THR:HB	2.02	0.41
1:B:70:ILE:HA	1:B:71:PRO:HA	1.88	0.41
1:B:242:VAL:HG13	1:B:380:VAL:HG11	2.02	0.41
1:A:285:TRP:CE3	1:A:285:TRP:HA	2.55	0.41
1:A:447:LYS:O	1:A:450:ARG:HG3	2.21	0.41
1:A:127:LYS:CD	1:A:146:PRO:HA	2.50	0.41
1:A:104:GLY:HA2	1:A:149:GLN:HG3	2.01	0.41
1:A:27:LEU:HD23	1:A:211:HIS:CG	2.56	0.41
1:A:70:ILE:O	1:A:70:ILE:HG23	2.21	0.41
1:B:119:VAL:HG23	1:B:122:ARG:HB2	2.03	0.41
1:A:28:HIS:HA	1:A:29:PRO:HD3	1.73	0.41
1:B:96:PHE:HB3	1:B:126:ILE:HG12	2.02	0.41
1:A:316:PHE:HA	1:A:317:ASN:HA	1.76	0.41
1:B:127:LYS:CD	1:B:146:PRO:HA	2.51	0.41
1:B:313:LEU:C	1:B:316:PHE:O	2.58	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:VAL:O	1:B:42:THR:HG23	2.21	0.41
1:A:90:PHE:HD2	1:A:287:GLY:HA2	1.86	0.41
1:A:303:PRO:HG3	1:A:314:TRP:CD2	2.56	0.41
1:A:422:LEU:HD11	1:B:215:PHE:CZ	2.56	0.41
1:A:193:ALA:HB2	1:A:402:LEU:HD12	2.03	0.41
1:A:88:LEU:O	1:A:88:LEU:HD23	2.21	0.41
1:B:315:ALA:HB1	1:B:361:VAL:HG21	2.02	0.41
1:A:124:LEU:HB3	1:A:125:PRO:HD3	2.04	0.40
1:A:234:ILE:HG23	1:A:234:ILE:O	2.21	0.40
1:A:250:PHE:HB3	1:A:420:THR:HB	2.03	0.40
1:B:94:LYS:HA	1:B:95:ARG:HA	1.80	0.40
1:A:290:LEU:HD21	1:A:338:ILE:HG22	2.03	0.40
1:A:389:LEU:HA	1:A:389:LEU:HD23	1.89	0.40
1:A:41:ILE:HB	1:A:154:ILE:HG21	2.03	0.40
1:A:257:PHE:HA	1:A:260:MET:HB2	2.03	0.40
1:B:386:MET:CE	1:B:404:THR:HB	2.52	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	423/466 (91%)	365 (86%)	43 (10%)	15 (4%)	4	28
1	B	423/466 (91%)	375 (89%)	36 (8%)	12 (3%)	6	34
All	All	846/932 (91%)	740 (88%)	79 (9%)	27 (3%)	5	30

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	71	PRO
1	A	103	SER

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Mol	Chain	Res	Type
1	A	449	ASN
1	B	71	PRO
1	B	96	PHE
1	B	103	SER
1	B	278	LEU
1	A	96	PHE
1	A	278	LEU
1	A	318	SER
1	B	449	ASN
1	A	104	GLY
1	A	142	GLY
1	B	318	SER
1	A	116	LEU
1	A	197	GLU
1	A	203	PHE
1	A	315	ALA
1	B	116	LEU
1	B	142	GLY
1	B	315	ALA
1	A	434	LYS
1	B	203	PHE
1	B	234	ILE
1	A	188	PRO
1	A	234	ILE
1	B	345	ILE

### 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	334/368 (91%)	293 (88%)	41 (12%)	5	25
1	B	334/368 (91%)	300 (90%)	34 (10%)	8	34
All	All	668/736 (91%)	593 (89%)	75 (11%)	7	29

All (75) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	A	27	LEU
1	A	28	HIS
1	A	30	ARG
1	A	40	LEU
1	A	42	THR
1	A	63	LEU
1	A	95	ARG
1	A	96	PHE
1	A	112	LEU
1	A	115	LYS
1	A	116	LEU
1	A	120	TRP
1	A	121	GLN
1	A	124	LEU
1	A	140	LEU
1	A	149	GLN
1	A	165	THR
1	A	166	GLN
1	A	169	GLN
1	A	170	ARG
1	A	215	PHE
1	A	228	ARG
1	A	236	LEU
1	A	256	LEU
1	A	263	THR
1	A	268	LEU
1	A	274	TRP
1	A	304	LEU
1	A	316	PHE
1	A	317	ASN
1	A	321	HIS
1	A	349	PHE
1	A	353	LEU
1	A	375	ILE
1	A	377	GLU
1	A	389	LEU
1	A	399	THR
1	A	403	LEU
1	A	417	LEU
1	A	440	LEU
1	A	448	GLN
1	B	27	LEU
1	B	40	LEU

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Mol	Chain	Res	Type
1	B	42	THR
1	B	63	LEU
1	B	73	LEU
1	B	95	ARG
1	B	96	PHE
1	B	115	LYS
1	B	116	LEU
1	B	120	TRP
1	B	121	GLN
1	B	124	LEU
1	B	140	LEU
1	B	149	GLN
1	B	165	THR
1	B	166	GLN
1	B	169	GLN
1	B	215	PHE
1	B	228	ARG
1	B	268	LEU
1	B	274	TRP
1	B	304	LEU
1	B	316	PHE
1	B	317	ASN
1	B	349	PHE
1	B	353	LEU
1	B	375	ILE
1	B	389	LEU
1	B	399	THR
1	B	403	LEU
1	B	417	LEU
1	B	440	LEU
1	B	441	LEU
1	B	448	GLN

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

Mol	Chain	Res	Type
1	B	230	GLN

### 5.3.3 RNA ⓘ

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 2 ligands modelled in this entry, 2 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 [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	425/466 (91%)	-0.04	7 (1%)	72 59	39, 67, 108, 198	0
1	B	425/466 (91%)	-0.02	7 (1%)	72 59	37, 69, 115, 253	0
All	All	850/932 (91%)	-0.03	14 (1%)	72 59	37, 69, 112, 253	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	70	ILE	4.7
1	B	205	SER	4.5
1	B	204	ARG	4.3
1	A	205	SER	3.9
1	B	70	ILE	3.5
1	A	141	ALA	3.1
1	B	450	ARG	2.7
1	A	204	ARG	2.7
1	B	299	LEU	2.3
1	B	71	PRO	2.2
1	A	140	LEU	2.1
1	B	449	ASN	2.1
1	A	301	PRO	2.0
1	A	69	PRO	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	CL	B	467	1/1	0.96	0.38	1.36	88,88,88,88	0
2	CL	A	467	1/1	0.94	0.25	-0.28	79,79,79,79	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.