



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:09 PM GMT

PDB ID : 4R0Y
Title : Structure of Maltose-binding Protein Fusion with the C-terminal GH1 domain of Guanylate Kinase-associated Protein from Rattus norvegicus
Authors : Im, Y.J.; Tong, J.
Deposited on : 2014-08-03
Resolution : 2.00 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

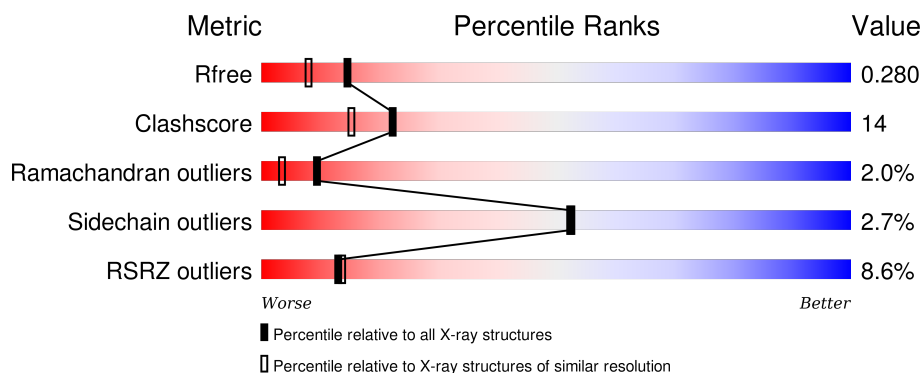
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 ≥ 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	501	
1	B	501	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7611 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 Maltose-binding periplasmic protein, Disks large-associated protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	469	Total	C	N	O	S	0	0	0
			3695	2367	613	701	14			
1	B	473	Total	C	N	O	S	0	0	0
			3735	2393	621	707	14			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	362	ALA	-	LINKER	UNP P0AEX9
A	363	MET	-	LINKER	UNP P0AEX9
A	917	VAL	-	LINKER	UNP P97836
A	918	ASP	-	LINKER	UNP P97836
B	362	ALA	-	LINKER	UNP P0AEX9
B	363	MET	-	LINKER	UNP P0AEX9
B	917	VAL	-	LINKER	UNP P97836
B	918	ASP	-	LINKER	UNP P97836

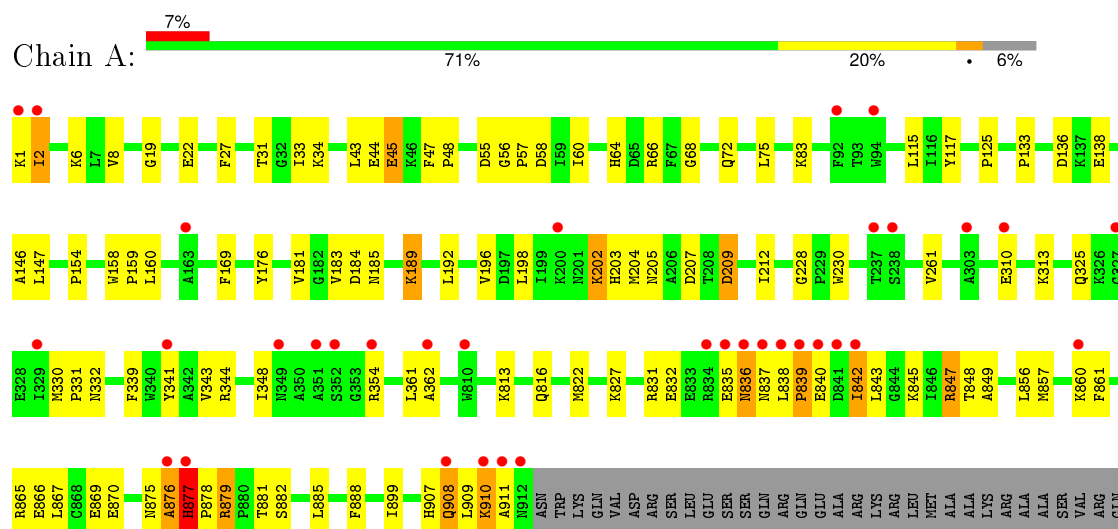
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	91	Total	O	0	0
			91	91		
2	B	90	Total	O	0	0
			90	90		

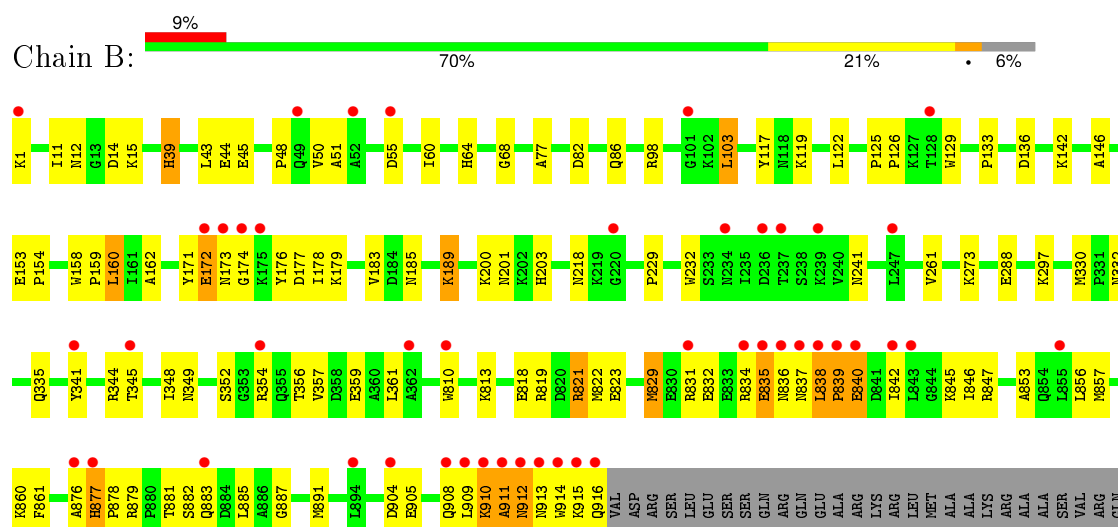
3 Residue-property plots [i](#)

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: Maltose-binding periplasmic protein, Disks large-associated protein 1



- Molecule 1: Maltose-binding periplasmic protein, Disks large-associated protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	99.11Å 158.68Å 65.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.52 – 2.00 39.52 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.7 (39.52-2.00) 96.8 (39.52-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.91 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.243 , 0.281 0.243 , 0.280	Depositor DCC
R_{free} test set	3452 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.583	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 67943 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7611	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.35	0/3783	0.56	0/5125
1	B	0.34	0/3825	0.56	0/5182
All	All	0.34	0/7608	0.56	0/10307

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	3695	0	3642	94	0
1	B	3735	0	3679	112	0
2	A	91	0	0	2	0
2	B	90	0	0	0	0
All	All	7611	0	7321	204	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 14.

All (204) 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:876:ALA:HB2	1:B:879:ARG:HB2	1.37	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:HIS:HD2	1:A:261:VAL:H	1.09	0.91
1:A:64:HIS:CD2	1:A:261:VAL:H	1.89	0.90
1:A:31:THR:HG23	1:A:33:ILE:H	1.41	0.84
1:B:43:LEU:HD13	1:B:60:ILE:HD11	1.61	0.81
1:B:172:GLU:CD	1:B:173:ASN:H	1.89	0.75
1:B:98:ARG:HG2	1:B:103:LEU:HD12	1.68	0.74
1:B:1:LYS:N	1:B:1:LYS:HD2	2.03	0.73
1:A:831:ARG:O	1:A:835:GLU:HG2	1.88	0.73
1:A:154:PRO:HG3	1:A:344:ARG:HA	1.72	0.71
1:A:837:ASN:C	1:A:838:LEU:HD23	2.12	0.70
1:A:44:GLU:HG2	1:A:45:GLU:OE1	1.91	0.70
1:A:816:GLN:CD	1:B:834:ARG:HH21	1.93	0.70
1:A:877:HIS:H	1:A:878:PRO:HA	1.57	0.69
1:A:66:ARG:HH21	1:A:66:ARG:HB3	1.57	0.69
1:A:66:ARG:NH2	1:A:66:ARG:HB3	2.07	0.69
1:B:876:ALA:CB	1:B:879:ARG:HB2	2.19	0.69
1:B:64:HIS:HD2	1:B:261:VAL:H	1.41	0.68
1:B:876:ALA:HB1	1:B:878:PRO:C	2.14	0.68
1:B:832:GLU:HG2	1:B:911:ALA:HB3	1.75	0.68
1:B:43:LEU:CD1	1:B:60:ILE:HD11	2.24	0.67
1:B:839:PRO:HB2	1:B:842:ILE:HB	1.75	0.67
1:B:154:PRO:HG3	1:B:344:ARG:HA	1.77	0.67
1:A:877:HIS:H	1:A:878:PRO:CA	2.08	0.67
1:A:27:PHE:O	1:A:31:THR:HG22	1.95	0.67
1:B:910:LYS:HD3	1:B:911:ALA:N	2.10	0.65
1:B:185:ASN:O	1:B:189:LYS:HE3	1.96	0.65
1:A:876:ALA:HA	1:A:879:ARG:HB2	1.77	0.65
1:A:192:LEU:O	1:A:196:VAL:HG23	1.97	0.65
1:A:907:HIS:C	1:A:909:LEU:H	2.01	0.64
1:B:877:HIS:H	1:B:878:PRO:CA	2.10	0.63
1:B:846:ILE:HD11	1:B:911:ALA:HB1	1.81	0.63
1:B:344:ARG:O	1:B:348:ILE:HG12	1.99	0.63
1:A:184:ASP:OD2	1:A:362:ALA:HA	1.99	0.62
1:A:6:LYS:HD2	1:A:34:LYS:HE2	1.80	0.62
1:B:178:ILE:HD12	1:B:335:GLN:HG2	1.80	0.62
1:B:881:THR:HG22	1:B:883:GLN:H	1.65	0.62
1:A:838:LEU:HD12	1:A:842:ILE:HG21	1.82	0.61
1:A:875:ASN:O	1:A:876:ALA:O	2.19	0.60
1:B:822:MET:HE1	1:B:853:ALA:HA	1.83	0.60
1:A:869:GLU:CD	1:B:847:ARG:HH12	2.05	0.59
1:A:198:LEU:HD13	1:A:204:MET:HE3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:TYR:CE2	1:A:813:LYS:HE2	2.37	0.59
1:B:910:LYS:O	1:B:912:ASN:N	2.36	0.58
1:B:839:PRO:O	1:B:840:GLU:CB	2.51	0.57
1:A:147:LEU:HD23	1:A:204:MET:HE2	1.85	0.57
1:A:185:ASN:O	1:A:189:LYS:HE3	2.04	0.57
1:A:68:GLY:HA3	1:A:332:ASN:O	2.03	0.57
1:B:877:HIS:H	1:B:878:PRO:C	2.07	0.57
1:A:822:MET:CE	1:A:899:ILE:HG21	2.35	0.57
1:A:64:HIS:HE1	1:A:330:MET:O	1.88	0.57
1:A:310:GLU:O	1:A:313:LYS:HG2	2.04	0.57
1:A:205:ASN:HB3	1:A:207:ASP:OD1	2.05	0.57
1:B:64:HIS:HE1	1:B:330:MET:O	1.87	0.56
1:A:837:ASN:C	1:A:839:PRO:HD3	2.26	0.56
1:A:909:LEU:O	1:A:910:LYS:CB	2.53	0.56
1:A:847:ARG:HG2	1:A:847:ARG:HH11	1.71	0.56
1:B:831:ARG:O	1:B:835:GLU:HB2	2.06	0.56
1:B:43:LEU:HD13	1:B:60:ILE:CD1	2.31	0.56
1:B:829:MET:HA	1:B:829:MET:HE3	1.87	0.56
1:A:847:ARG:NH1	1:A:847:ARG:HG2	2.21	0.55
1:B:834:ARG:O	1:B:835:GLU:C	2.45	0.55
1:A:198:LEU:HD13	1:A:204:MET:CE	2.37	0.55
1:A:838:LEU:HD13	1:A:911:ALA:HB1	1.89	0.55
1:B:832:GLU:HG2	1:B:911:ALA:CB	2.37	0.54
1:A:136:ASP:HA	1:A:146:ALA:HB2	1.90	0.54
1:A:838:LEU:O	1:A:842:ILE:HG22	2.06	0.54
1:A:83:LYS:HD3	1:A:83:LYS:O	2.08	0.54
1:A:909:LEU:O	1:A:910:LYS:HB3	2.07	0.54
1:B:818:GLU:OE1	1:B:821:ARG:HD3	2.07	0.54
1:B:881:THR:HG22	1:B:882:SER:N	2.23	0.54
1:A:344:ARG:O	1:A:348:ILE:HD13	2.07	0.54
1:B:189:LYS:HD3	1:B:361:LEU:HD12	1.89	0.54
1:B:200:LYS:HE3	1:B:201:ASN:HD21	1.73	0.53
1:B:178:ILE:CD1	1:B:335:GLN:HG2	2.39	0.53
1:B:1:LYS:H3	1:B:1:LYS:HD2	1.74	0.53
1:B:200:LYS:HG3	1:B:201:ASN:ND2	2.23	0.53
1:A:832:GLU:OE1	1:A:911:ALA:HB2	2.10	0.52
1:B:876:ALA:HB1	1:B:879:ARG:N	2.24	0.52
1:A:865:ARG:O	1:A:869:GLU:HG3	2.10	0.52
1:B:119:LYS:HB2	1:B:241:ASN:ND2	2.25	0.52
1:B:877:HIS:O	1:B:877:HIS:ND1	2.42	0.52
1:B:178:ILE:HD13	1:B:818:GLU:OE2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:PRO:HB3	1:A:203:HIS:CE1	2.44	0.52
1:B:229:PRO:HA	1:B:232:TRP:CE2	2.44	0.51
1:B:136:ASP:OD2	1:B:203:HIS:HD2	1.92	0.51
1:B:172:GLU:OE2	1:B:173:ASN:N	2.43	0.51
1:B:877:HIS:N	1:B:878:PRO:CA	2.74	0.51
1:A:838:LEU:N	1:A:839:PRO:HD3	2.25	0.51
1:B:189:LYS:HG3	1:B:357:VAL:HG12	1.93	0.51
1:B:341:TYR:CE2	1:B:813:LYS:HE2	2.44	0.51
1:A:845:LYS:O	1:A:848:THR:HG22	2.11	0.51
1:A:857:MET:O	1:A:861:PHE:HB2	2.10	0.51
1:B:12:ASN:ND2	1:B:14:ASP:OD1	2.38	0.51
1:B:288:GLU:H	1:B:288:GLU:CD	2.13	0.51
1:B:218:ASN:HD22	1:B:218:ASN:N	2.09	0.51
1:A:836:ASN:CB	1:A:838:LEU:HD22	2.41	0.50
1:A:325:GLN:NE2	2:A:1050:HOH:O	2.44	0.50
1:A:48:PRO:HA	1:A:75:LEU:HD13	1.94	0.50
1:A:183:VAL:CG1	1:A:361:LEU:HD22	2.42	0.50
1:B:43:LEU:C	1:B:43:LEU:HD12	2.33	0.49
1:A:8:VAL:HG13	1:A:57:PRO:HA	1.93	0.49
1:B:177:ASP:OD1	1:B:179:LYS:HB2	2.12	0.49
1:B:172:GLU:CD	1:B:173:ASN:N	2.64	0.49
1:B:172:GLU:OE2	1:B:173:ASN:ND2	2.45	0.49
1:B:904:ASP:O	1:B:908:GLN:HG3	2.12	0.49
1:A:64:HIS:HD2	1:A:261:VAL:N	1.93	0.49
1:A:842:ILE:HD13	1:A:842:ILE:O	2.12	0.49
1:B:356:THR:OG1	1:B:359:GLU:HG3	2.13	0.49
1:A:881:THR:O	1:A:885:LEU:HD13	2.12	0.49
1:B:158:TRP:CE2	1:B:162:ALA:HB2	2.47	0.49
1:A:907:HIS:C	1:A:909:LEU:N	2.65	0.49
1:B:838:LEU:HA	1:B:839:PRO:HD3	1.63	0.48
1:B:117:TYR:CE2	1:B:125:PRO:HD3	2.48	0.48
1:B:1:LYS:H1	1:B:1:LYS:HD2	1.74	0.48
1:B:183:VAL:HG21	1:B:810:TRP:CH2	2.48	0.48
1:A:341:TYR:CD2	1:A:813:LYS:HE2	2.49	0.48
1:B:842:ILE:HG21	1:B:911:ALA:HA	1.96	0.47
1:B:133:PRO:HB3	1:B:203:HIS:CE1	2.49	0.47
1:A:2:ILE:HG12	1:A:56:GLY:O	2.13	0.47
1:B:68:GLY:HA3	1:B:332:ASN:O	2.13	0.47
1:A:136:ASP:OD2	1:A:203:HIS:HD2	1.97	0.47
1:B:829:MET:HA	1:B:829:MET:CE	2.45	0.47
1:B:912:ASN:C	1:B:914:TRP:H	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:857:MET:O	1:B:861:PHE:HB2	2.15	0.47
1:A:908:GLN:O	1:A:909:LEU:HD23	2.15	0.47
1:B:39:HIS:N	1:B:39:HIS:CD2	2.83	0.46
1:B:171:TYR:CE1	1:B:174:GLY:HA2	2.50	0.46
1:A:907:HIS:O	1:A:909:LEU:N	2.48	0.46
1:A:867:LEU:HD13	1:A:888:PHE:CE2	2.51	0.46
1:A:83:LYS:C	1:A:83:LYS:HD3	2.36	0.46
1:A:1:LYS:HA	1:A:55:ASP:OD1	2.15	0.45
1:A:202:LYS:HB3	1:A:202:LYS:HZ2	1.81	0.45
1:B:45:GLU:O	1:B:48:PRO:HG2	2.17	0.45
1:A:910:LYS:HG3	1:A:911:ALA:N	2.32	0.45
1:A:848:THR:HG23	1:A:849:ALA:N	2.31	0.45
1:A:881:THR:HG22	1:A:882:SER:N	2.32	0.45
1:A:209:ASP:OD1	1:A:212:ILE:HG13	2.16	0.45
1:B:838:LEU:N	1:B:838:LEU:HD23	2.31	0.45
1:B:15:LYS:C	1:B:297:LYS:HD2	2.37	0.45
1:B:876:ALA:O	1:B:877:HIS:HB2	2.16	0.45
1:A:66:ARG:NH2	2:A:1072:HOH:O	2.49	0.45
1:A:19:GLY:O	1:A:22:GLU:HB2	2.16	0.45
1:A:47:PHE:CG	1:A:60:ILE:HD12	2.52	0.45
1:A:158:TRP:N	1:A:159:PRO:CD	2.80	0.45
1:B:887:GLY:O	1:B:891:MET:HG2	2.17	0.44
1:A:6:LYS:CE	1:A:34:LYS:HE2	2.47	0.44
1:A:154:PRO:HG3	1:A:344:ARG:CA	2.45	0.44
1:B:178:ILE:O	1:B:178:ILE:HG12	2.18	0.44
1:A:866:GLU:O	1:A:870:GLU:HG3	2.17	0.44
1:B:345:THR:HG22	1:B:349:ASN:ND2	2.32	0.44
1:A:176:TYR:CZ	1:A:331:PRO:HA	2.52	0.44
1:B:98:ARG:HG2	1:B:103:LEU:CD1	2.44	0.44
1:A:836:ASN:HB2	1:A:838:LEU:HD22	2.00	0.44
1:B:136:ASP:HA	1:B:146:ALA:HB2	1.98	0.44
1:B:856:LEU:HD12	1:B:860:LYS:CG	2.47	0.44
1:B:82:ASP:O	1:B:86:GLN:HG3	2.18	0.44
1:B:43:LEU:HD12	1:B:44:GLU:N	2.33	0.44
1:B:122:LEU:HD21	1:B:126:PRO:HD3	2.00	0.44
1:A:856:LEU:CD1	1:A:860:LYS:HD2	2.47	0.44
1:B:352:SER:OG	1:B:354:ARG:NH1	2.51	0.44
1:B:51:ALA:HA	1:B:55:ASP:O	2.17	0.44
1:A:202:LYS:HB3	1:A:202:LYS:NZ	2.33	0.43
1:A:31:THR:HG23	1:A:33:ILE:N	2.21	0.43
1:B:912:ASN:O	1:B:914:TRP:N	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:ARG:HH21	1:A:66:ARG:CB	2.27	0.43
1:B:153:GLU:HA	1:B:154:PRO:HD3	1.82	0.43
1:B:877:HIS:N	1:B:878:PRO:HA	2.32	0.43
1:B:50:VAL:HG23	1:B:51:ALA:N	2.33	0.43
1:B:173:ASN:OD1	1:B:173:ASN:O	2.37	0.43
1:B:909:LEU:O	1:B:910:LYS:O	2.37	0.43
1:B:835:GLU:HB3	1:B:836:ASN:H	1.64	0.43
1:B:171:TYR:HB2	1:B:176:TYR:CE1	2.53	0.43
1:B:64:HIS:CD2	1:B:261:VAL:H	2.28	0.43
1:A:339:PHE:CZ	1:A:343:VAL:HG21	2.54	0.43
1:B:356:THR:HG23	1:B:359:GLU:OE1	2.19	0.42
1:A:8:VAL:HG12	1:A:58:ASP:OD2	2.19	0.42
1:B:64:HIS:CE1	1:B:330:MET:O	2.69	0.42
1:B:914:TRP:CZ3	1:B:916:GLN:HB3	2.55	0.42
1:A:43:LEU:C	1:A:43:LEU:HD12	2.40	0.42
1:A:836:ASN:HB3	1:A:838:LEU:HD22	2.00	0.42
1:B:842:ILE:HG22	1:B:846:ILE:HG13	2.01	0.41
1:A:228:GLY:HA3	1:A:230:TRP:CH2	2.55	0.41
1:B:822:MET:HE1	1:B:853:ALA:CA	2.49	0.41
1:B:15:LYS:O	1:B:297:LYS:HD2	2.21	0.41
1:A:877:HIS:H	1:A:878:PRO:C	2.24	0.41
1:B:77:ALA:HB2	1:B:273:LYS:HE3	2.02	0.41
1:B:910:LYS:HD3	1:B:910:LYS:C	2.41	0.41
1:B:819:ARG:O	1:B:823:GLU:HG3	2.20	0.41
1:A:839:PRO:O	1:A:840:GLU:HB2	2.21	0.41
1:A:6:LYS:CD	1:A:34:LYS:HE2	2.46	0.41
1:A:877:HIS:N	1:A:878:PRO:CA	2.78	0.41
1:B:857:MET:HA	1:B:861:PHE:HB2	2.03	0.41
1:B:129:TRP:CE2	1:B:160:LEU:HG	2.55	0.41
1:B:845:LYS:HD3	1:B:905:GLU:OE2	2.21	0.41
1:B:11:ILE:O	1:B:39:HIS:HA	2.21	0.41
1:A:117:TYR:CE2	1:A:125:PRO:HD3	2.56	0.41
1:B:914:TRP:CE3	1:B:914:TRP:HA	2.55	0.40
1:B:158:TRP:N	1:B:159:PRO:CD	2.84	0.40
1:A:839:PRO:HA	1:A:843:LEU:CD1	2.51	0.40
1:B:881:THR:CG2	1:B:882:SER:N	2.83	0.40
1:B:142:LYS:HG2	1:B:142:LYS:O	2.21	0.40
1:B:136:ASP:OD2	1:B:203:HIS:CD2	2.73	0.40
1:A:169:PHE:CE1	1:A:181:VAL:HG22	2.56	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	467/501 (93%)	440 (94%)	20 (4%)	7 (2%)	13	5
1	B	471/501 (94%)	445 (94%)	14 (3%)	12 (2%)	7	2
All	All	938/1002 (94%)	885 (94%)	34 (4%)	19 (2%)	9	3

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	876	ALA
1	A	877	HIS
1	A	910	LYS
1	B	835	GLU
1	B	837	ASN
1	B	838	LEU
1	B	840	GLU
1	B	877	HIS
1	B	910	LYS
1	B	915	LYS
1	A	2	ILE
1	A	836	ASN
1	B	172	GLU
1	B	911	ALA
1	B	912	ASN
1	A	908	GLN
1	B	839	PRO
1	B	913	ASN
1	A	839	PRO

5.3.2 Protein sidechains [i](#)

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	384/411 (93%)	370 (96%)	14 (4%)	42	39
1	B	388/411 (94%)	381 (98%)	7 (2%)	66	69
All	All	772/822 (94%)	751 (97%)	21 (3%)	52	52

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	GLU
1	A	72	GLN
1	A	115	LEU
1	A	138	GLU
1	A	160	LEU
1	A	189	LYS
1	A	202	LYS
1	A	209	ASP
1	A	354	ARG
1	A	827	LYS
1	A	842	ILE
1	A	847	ARG
1	A	877	HIS
1	A	879	ARG
1	B	39	HIS
1	B	103	LEU
1	B	160	LEU
1	B	189	LYS
1	B	821	ARG
1	B	829	MET
1	B	885	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	HIS
1	A	86	GLN
1	A	201	ASN
1	A	203	HIS
1	A	218	ASN
1	A	325	GLN

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Mol	Chain	Res	Type
1	A	349	ASN
1	A	837	ASN
1	A	907	HIS
1	B	39	HIS
1	B	64	HIS
1	B	86	GLN
1	B	124	ASN
1	B	173	ASN
1	B	201	ASN
1	B	203	HIS
1	B	205	ASN
1	B	218	ASN
1	B	241	ASN
1	B	282	ASN
1	B	325	GLN
1	B	828	GLN
1	B	883	GLN
1	B	898	ASN
1	B	907	HIS
1	B	908	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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, 95th 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 ²)	Q<0.9
1	A	469/501 (93%)	0.74	35 (7%)	17 18	27, 45, 69, 95	0
1	B	473/501 (94%)	0.76	46 (9%)	10 10	30, 45, 78, 103	0
All	All	942/1002 (94%)	0.75	81 (8%)	13 14	27, 45, 74, 103	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	914	TRP	9.7
1	B	838	LEU	8.8
1	A	911	ALA	7.7
1	A	839	PRO	6.4
1	A	877	HIS	6.4
1	A	838	LEU	6.3
1	B	174	GLY	6.0
1	B	877	HIS	5.7
1	B	912	ASN	5.7
1	B	173	ASN	5.3
1	B	913	ASN	5.0
1	A	912	ASN	5.0
1	B	876	ALA	4.9
1	A	842	ILE	4.9
1	A	876	ALA	4.8
1	A	837	ASN	4.7
1	B	837	ASN	4.6
1	A	341	TYR	4.4
1	B	835	GLU	4.1
1	B	354	ARG	4.0
1	B	839	PRO	4.0
1	B	341	TYR	3.9
1	A	836	ASN	3.8
1	B	52	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	910	LYS	3.5
1	B	836	ASN	3.4
1	B	239	LYS	3.4
1	B	843	LEU	3.4
1	A	840	GLU	3.4
1	B	908	GLN	3.3
1	A	841	ASP	3.2
1	A	237	THR	3.2
1	A	1	LYS	3.2
1	B	916	GLN	3.2
1	B	911	ALA	3.1
1	A	2	ILE	3.1
1	B	840	GLU	3.0
1	B	49	GLN	2.9
1	B	842	ILE	2.9
1	A	352	SER	2.9
1	B	237	THR	2.8
1	B	915	LYS	2.8
1	B	345	THR	2.8
1	B	834	ARG	2.8
1	A	354	ARG	2.7
1	B	175	LYS	2.6
1	A	94	TRP	2.6
1	B	362	ALA	2.5
1	A	834	ARG	2.5
1	A	329	ILE	2.5
1	A	349	ASN	2.5
1	A	163	ALA	2.5
1	B	55	ASP	2.5
1	B	234	ASN	2.5
1	A	310	GLU	2.4
1	A	351	ALA	2.4
1	A	810	TRP	2.4
1	B	909	LEU	2.4
1	A	92	PHE	2.3
1	A	200	LYS	2.3
1	B	101	GLY	2.3
1	B	894	LEU	2.3
1	B	172	GLU	2.3
1	B	1	LYS	2.3
1	A	238	SER	2.3
1	A	835	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	303	ALA	2.2
1	B	810	TRP	2.2
1	A	327	GLY	2.2
1	B	236	ASP	2.1
1	A	362	ALA	2.1
1	B	910	LYS	2.1
1	B	855	LEU	2.1
1	B	883	GLN	2.1
1	A	860	LYS	2.1
1	B	247	LEU	2.1
1	B	831	ARG	2.1
1	B	904	ASP	2.1
1	B	128	THR	2.0
1	B	220	GLY	2.0
1	A	908	GLN	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.