Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:57 AM GMT

PDB ID	:	3CRA
Title	:	Crystal Structure of Escherichia coli MazG, the Regulator of Nutritional Stress Response
Authors	:	Lee, S.; Kim, M.H.; Kang, B.S.; Kim, J.S.; Kim, Y.G.; Kim, K.J.
Deposited on	:	2008-04-05
Resolution	:	2.10 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:

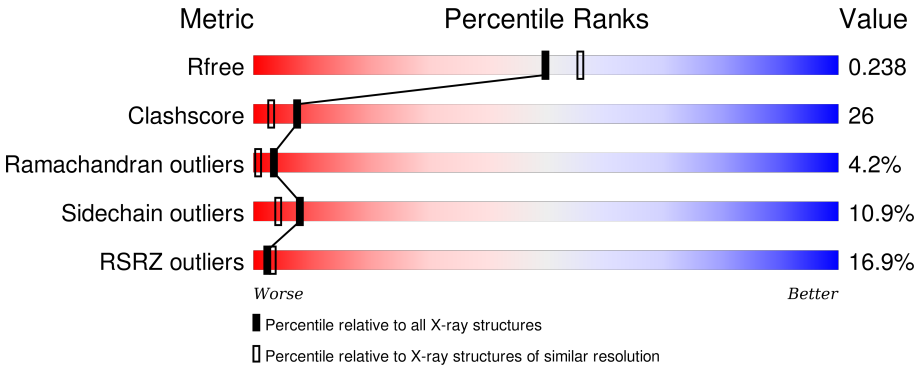
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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 <=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	265	20% 59% 20% 8% • 10%
1	B	265	9% 52% 23% 6% • 16%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3870 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 Protein mazG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	239	Total	C	N	O	S	0	0	0
			1956	1223	345	378	10			
1	B	222	Total	C	N	O	S	0	0	0
			1806	1135	312	350	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP P0AEY3
A	0	HIS	-	EXPRESSION TAG	UNP P0AEY3
B	-1	GLY	-	EXPRESSION TAG	UNP P0AEY3
B	0	HIS	-	EXPRESSION TAG	UNP P0AEY3

- Molecule 2 is water.

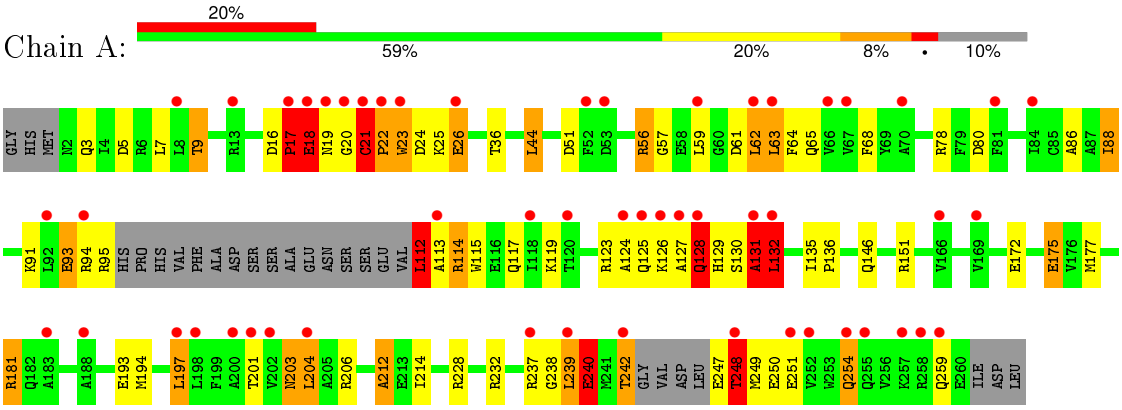
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		
2	B	64	Total	O	0	0
			64	64		

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.

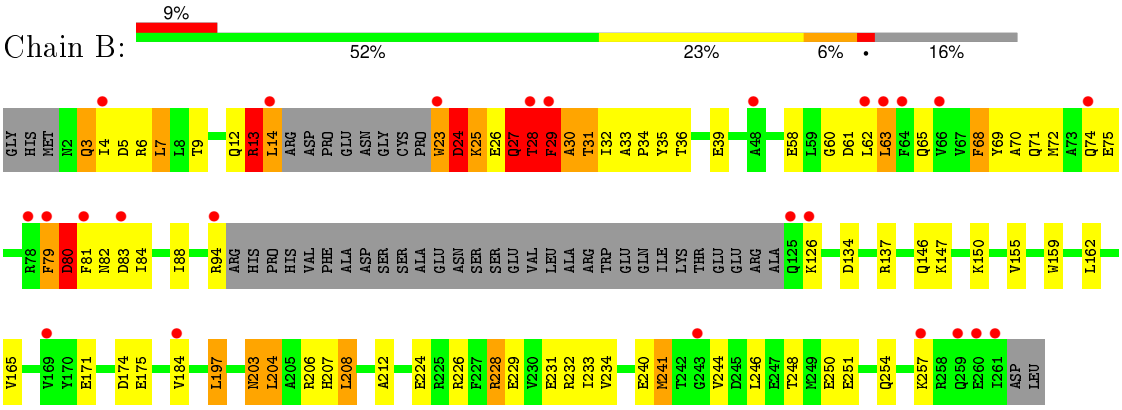
- Molecule 1: Protein mazG

Chain A:



- Molecule 1: Protein mazG

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.79Å 67.57Å 140.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.47 – 2.10 32.85 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.2 (28.47-2.10) 93.4 (32.85-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.69 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.226 , 0.284 0.238 , 0.238	Depositor DCC
R_{free} test set	1702 reflections (5.28%)	DCC
Wilson B-factor (Å^2)	35.1	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 53.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	1 of 38465 reflections (0.003%)	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	3870	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.73% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	1.12	3/1984 (0.2%)	1.12	17/2671 (0.6%)
1	B	1.11	1/1830 (0.1%)	1.09	13/2464 (0.5%)
All	All	1.11	4/3814 (0.1%)	1.10	30/5135 (0.6%)

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	5
1	B	0	6
All	All	0	11

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	175	GLU	CB-CG	-11.57	1.30	1.52
1	A	212	ALA	CA-CB	6.06	1.65	1.52
1	A	172	GLU	CD-OE2	-5.51	1.19	1.25
1	B	147	LYS	CD-CE	5.50	1.65	1.51

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	228	ARG	NE-CZ-NH1	9.75	125.17	120.30
1	A	228	ARG	NE-CZ-NH2	-8.37	116.11	120.30
1	A	21	CYS	C-N-CD	-8.34	102.24	120.60
1	B	174	ASP	CB-CG-OD1	7.19	124.77	118.30
1	B	197	LEU	CB-CG-CD2	7.14	123.14	111.00
1	B	228	ARG	NE-CZ-NH1	-7.02	116.79	120.30
1	A	21	CYS	C-N-CA	6.89	150.96	122.00
1	A	131	ALA	O-C-N	-6.82	111.78	122.70

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	228	ARG	NE-CZ-NH2	6.82	123.71	120.30
1	B	134	ASP	CB-CG-OD1	-6.74	112.23	118.30
1	A	131	ALA	C-N-CA	6.73	138.52	121.70
1	B	241	MET	CG-SD-CE	6.36	110.38	100.20
1	A	131	ALA	CA-C-N	6.22	130.88	117.20
1	B	80	ASP	N-CA-C	5.95	127.07	111.00
1	B	134	ASP	CB-CG-OD2	5.88	123.59	118.30
1	A	194	MET	CG-SD-CE	-5.85	90.84	100.20
1	B	162	LEU	CB-CG-CD2	-5.83	101.08	111.00
1	A	132	LEU	CB-CG-CD1	-5.77	101.19	111.00
1	A	131	ALA	N-CA-C	5.57	126.03	111.00
1	B	7	LEU	CA-CB-CG	5.42	127.77	115.30
1	B	232	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	132	LEU	CB-CG-CD2	5.37	120.12	111.00
1	A	80	ASP	CB-CG-OD2	5.31	123.08	118.30
1	A	181	ARG	NE-CZ-NH2	-5.25	117.68	120.30
1	A	44	LEU	CA-CB-CG	5.23	127.32	115.30
1	A	248	THR	N-CA-C	5.22	125.10	111.00
1	B	28	THR	N-CA-C	-5.22	96.92	111.00
1	A	78	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	B	80	ASP	CB-CG-OD1	-5.02	113.78	118.30
1	A	204	LEU	CA-CB-CG	5.01	126.83	115.30

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	112	LEU	Peptide
1	A	131	ALA	Peptide
1	A	17	PRO	Peptide
1	A	240	GLU	Peptide
1	A	248	THR	Peptide
1	B	23	TRP	Peptide
1	B	24	ASP	Peptide
1	B	25	LYS	Peptide
1	B	27	GLN	Peptide
1	B	29	PHE	Peptide
1	B	79	PHE	Peptide

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	1956	0	1919	109	5
1	B	1806	0	1779	113	5
2	A	44	0	0	4	0
2	B	64	0	0	9	0
All	All	3870	0	3698	191	5

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 26.

All (191) 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:112:LEU:HB3	1:B:14:LEU:CD2	1.12	1.54
1:A:112:LEU:CB	1:B:14:LEU:HD21	1.33	1.50
1:B:74:GLN:CG	1:B:80:ASP:HB2	1.49	1.40
1:A:112:LEU:CB	1:B:14:LEU:CD2	1.95	1.35
1:A:132:LEU:HD22	2:A:307:HOH:O	1.27	1.29
1:B:74:GLN:HG2	1:B:80:ASP:CB	1.62	1.28
1:A:22:PRO:CB	1:A:23:TRP:HB3	1.65	1.26
1:B:74:GLN:NE2	1:B:80:ASP:OD1	1.73	1.20
1:A:112:LEU:HB3	1:B:14:LEU:HD22	1.17	1.15
1:B:12:GLN:O	1:B:13:ARG:HB3	1.51	1.10
1:A:21:CYS:HB3	1:A:22:PRO:O	1.50	1.09
1:A:112:LEU:HB2	1:B:14:LEU:HD21	1.36	1.07
1:A:22:PRO:HB3	1:A:23:TRP:CB	1.88	1.04
1:A:22:PRO:HB2	1:A:24:ASP:H	1.25	0.99
1:A:22:PRO:HB3	1:A:23:TRP:HB3	0.99	0.98
1:B:74:GLN:HG2	1:B:80:ASP:HB2	0.96	0.96
1:B:203:ASN:ND2	1:B:206:ARG:HH21	1.63	0.96
1:A:112:LEU:HB3	1:B:14:LEU:HD23	1.46	0.95
1:A:115:TRP:HH2	1:B:61:ASP:HB3	1.31	0.95
1:B:12:GLN:O	1:B:13:ARG:CB	2.18	0.91
1:B:204:LEU:HD22	1:B:208:LEU:HD22	1.51	0.91
1:B:74:GLN:HG3	1:B:80:ASP:HB2	1.53	0.90
1:A:132:LEU:HG	1:B:228:ARG:NH1	1.86	0.90

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:ASN:HD21	1:B:206:ARG:NH2	1.70	0.90
1:A:57:GLY:O	1:A:61:ASP:OD2	1.90	0.89
1:B:203:ASN:HD21	1:B:206:ARG:HH21	0.88	0.86
1:A:112:LEU:CG	1:B:14:LEU:HD21	2.06	0.86
1:A:112:LEU:HD13	1:A:113:ALA:HA	1.56	0.85
1:B:36:THR:HA	1:B:65:GLN:HE22	1.42	0.85
1:A:113:ALA:HB3	1:A:114:ARG:HA	1.61	0.83
1:A:132:LEU:C	1:A:132:LEU:HD12	1.99	0.82
1:A:132:LEU:O	1:A:132:LEU:HD12	1.78	0.82
1:A:95:ARG:C	1:A:112:LEU:HD21	2.00	0.81
1:A:22:PRO:CB	1:A:23:TRP:CB	2.50	0.81
1:B:74:GLN:HE21	1:B:81:PHE:H	1.28	0.81
1:A:22:PRO:HB2	1:A:24:ASP:N	1.96	0.80
1:B:28:THR:O	1:B:29:PHE:CD2	2.34	0.80
1:A:146:GLN:HE22	1:A:212:ALA:H	1.29	0.78
1:A:21:CYS:SG	1:A:22:PRO:C	2.62	0.78
1:A:21:CYS:CB	1:A:22:PRO:O	2.32	0.78
1:A:91:LYS:O	1:A:95:ARG:HG3	1.85	0.76
1:A:88:ILE:HG12	1:B:60:GLY:CA	2.16	0.75
1:B:146:GLN:HE22	1:B:212:ALA:H	1.32	0.74
1:A:115:TRP:CH2	1:B:61:ASP:HB3	2.21	0.74
1:A:132:LEU:CD1	1:B:228:ARG:HH12	1.99	0.74
1:A:22:PRO:HB2	1:A:23:TRP:HB3	1.68	0.73
1:A:132:LEU:CG	1:B:228:ARG:HH12	2.01	0.73
1:A:113:ALA:HB1	1:A:114:ARG:O	1.88	0.72
1:B:23:TRP:C	1:B:25:LYS:H	1.92	0.72
1:A:127:ALA:HB1	1:A:128:GLN:O	1.89	0.72
1:A:247:GLU:HB2	1:A:249:MET:H	1.54	0.71
1:B:240:GLU:O	1:B:244:VAL:HG23	1.91	0.71
1:B:30:ALA:HA	1:B:32:ILE:N	2.07	0.70
1:A:114:ARG:HG2	1:A:117:GLN:HG3	1.73	0.70
1:A:132:LEU:CG	1:B:228:ARG:NH1	2.54	0.70
1:A:113:ALA:CB	1:A:114:ARG:CA	2.70	0.70
1:A:113:ALA:CB	1:A:114:ARG:C	2.61	0.69
1:B:74:GLN:CD	1:B:80:ASP:OD1	2.30	0.69
1:A:86:ALA:HA	1:B:3:GLN:HG2	1.74	0.69
1:B:234:VAL:HG21	1:B:241:MET:CE	2.23	0.68
1:A:112:LEU:N	1:B:24:ASP:OD2	2.26	0.68
1:B:74:GLN:CG	1:B:80:ASP:CB	2.40	0.68
1:A:44:LEU:HD22	1:B:33:ALA:HB1	1.75	0.68
1:B:171:GLU:O	1:B:175:GLU:HG3	1.94	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:CYS:HB2	1:A:25:LYS:HD3	1.75	0.67
1:A:62:LEU:O	1:A:65:GLN:HB2	1.94	0.67
1:B:4:ILE:H	1:B:4:ILE:HD12	1.60	0.67
1:A:113:ALA:HB3	1:A:114:ARG:CA	2.25	0.66
1:A:113:ALA:HB1	1:A:114:ARG:C	2.16	0.66
1:B:244:VAL:HG12	1:B:248:THR:HB	1.77	0.66
1:B:74:GLN:CD	2:B:267:HOH:O	2.33	0.66
1:B:68:PHE:CE2	1:B:72:MET:HE1	2.31	0.65
1:A:63:LEU:O	1:A:64:PHE:C	2.34	0.65
1:B:74:GLN:HG2	1:B:80:ASP:HB3	1.75	0.64
1:A:214:ILE:HD11	1:B:34:PRO:HB3	1.81	0.62
1:A:175:GLU:HG2	1:A:193:GLU:OE2	1.98	0.62
1:A:3:GLN:HB2	1:B:82:ASN:OD1	1.99	0.62
1:B:28:THR:O	1:B:29:PHE:HD2	1.83	0.60
1:B:39:GLU:OE2	1:B:58:GLU:OE2	2.19	0.60
1:B:30:ALA:CA	1:B:32:ILE:N	2.64	0.60
1:A:130:SER:OG	1:A:132:LEU:HB3	2.01	0.60
1:A:247:GLU:N	1:A:251:GLU:OE2	2.35	0.60
2:A:269:HOH:O	1:B:137:ARG:HD3	2.01	0.59
1:B:30:ALA:CA	1:B:32:ILE:H	2.15	0.59
1:B:234:VAL:HG21	1:B:241:MET:HE3	1.85	0.59
1:B:71:GLN:O	1:B:75:GLU:HG3	2.01	0.59
1:A:21:CYS:HB3	1:A:22:PRO:C	2.19	0.59
1:B:23:TRP:N	1:B:24:ASP:HB3	2.17	0.59
1:B:224:GLU:OE2	1:B:228:ARG:NH1	2.36	0.59
1:B:159:TRP:HZ3	2:B:295:HOH:O	1.86	0.59
1:A:132:LEU:HD12	1:B:228:ARG:HH12	1.68	0.59
1:A:127:ALA:HA	1:A:128:GLN:HB3	1.84	0.58
1:B:23:TRP:N	1:B:24:ASP:CB	2.67	0.58
1:B:204:LEU:HD22	1:B:208:LEU:CD2	2.31	0.58
1:B:234:VAL:HG21	1:B:241:MET:HE1	1.86	0.58
1:A:61:ASP:O	1:A:65:GLN:HG2	2.03	0.58
1:A:113:ALA:HB3	1:A:114:ARG:C	2.24	0.58
1:A:88:ILE:HG12	1:B:60:GLY:HA3	1.86	0.57
1:B:68:PHE:HE2	1:B:72:MET:CE	2.19	0.56
1:A:59:LEU:HD11	1:B:69:TYR:HB3	1.87	0.56
1:A:214:ILE:CD1	1:B:34:PRO:HB3	2.36	0.55
1:A:132:LEU:CD1	1:A:132:LEU:O	2.54	0.55
1:B:146:GLN:NE2	1:B:212:ALA:H	2.03	0.55
1:A:21:CYS:SG	1:A:25:LYS:HB2	2.47	0.55
1:A:238:GLY:O	1:A:240:GLU:N	2.40	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:ILE:O	1:B:35:TYR:HB2	2.07	0.54
1:A:5:ASP:O	1:A:9:THR:HG23	2.08	0.54
1:A:113:ALA:CB	1:A:114:ARG:HA	2.28	0.54
1:B:27:GLN:HA	1:B:28:THR:OG1	2.07	0.54
1:A:7:LEU:CD2	1:A:63:LEU:HD13	2.38	0.54
1:A:201:THR:HG21	2:B:304:HOH:O	2.08	0.54
1:B:30:ALA:C	1:B:32:ILE:H	2.11	0.54
1:B:68:PHE:HE2	1:B:72:MET:HE1	1.71	0.53
1:B:226:ARG:CZ	1:B:257:LYS:HG2	2.38	0.53
1:B:23:TRP:C	1:B:25:LYS:N	2.60	0.52
1:B:79:PHE:O	1:B:83:ASP:HB2	2.09	0.52
1:B:70:ALA:O	1:B:74:GLN:HG3	2.10	0.52
1:B:68:PHE:CE2	1:B:72:MET:CE	2.93	0.52
1:A:64:PHE:HB2	1:B:88:ILE:HG12	1.92	0.52
1:B:27:GLN:O	1:B:72:MET:HE1	2.10	0.51
1:A:16:ASP:OD1	1:A:18:GLU:O	2.27	0.51
1:A:127:ALA:HA	1:A:128:GLN:CB	2.42	0.50
1:B:80:ASP:N	2:B:300:HOH:O	2.44	0.50
1:B:26:GLU:O	1:B:27:GLN:HG3	2.11	0.50
1:B:4:ILE:H	1:B:4:ILE:CD1	2.25	0.50
1:B:80:ASP:CA	2:B:300:HOH:O	2.60	0.49
1:A:95:ARG:C	1:A:112:LEU:CD2	2.76	0.49
1:B:74:GLN:HE21	1:B:80:ASP:CG	2.16	0.49
1:B:203:ASN:ND2	1:B:206:ARG:NH2	2.43	0.49
1:B:9:THR:O	1:B:12:GLN:O	2.31	0.48
1:A:123:ARG:HD3	1:A:151:ARG:CZ	2.43	0.48
1:A:132:LEU:HD13	1:A:132:LEU:HA	1.10	0.48
1:A:203:ASN:ND2	1:A:206:ARG:HH21	2.12	0.48
1:A:21:CYS:CB	1:A:22:PRO:C	2.78	0.48
1:B:36:THR:CA	1:B:65:GLN:HE22	2.20	0.48
1:A:119:LYS:NZ	1:B:39:GLU:OE1	2.33	0.48
1:A:23:TRP:HA	1:A:26:GLU:CG	2.44	0.47
1:A:18:GLU:CB	1:A:19:ASN:ND2	2.77	0.47
1:B:5:ASP:O	1:B:9:THR:HG23	2.14	0.47
1:A:93:GLU:OE1	1:B:6:ARG:NH1	2.47	0.47
1:B:246:LEU:O	1:B:250:GLU:HG3	2.15	0.47
1:A:135:ILE:O	1:A:136:PRO:C	2.50	0.47
1:A:22:PRO:CB	1:A:23:TRP:CA	2.93	0.47
1:A:125:GLN:HA	1:A:126:LYS:CG	2.46	0.46
1:A:203:ASN:HD22	1:A:206:ARG:HH21	1.62	0.46
1:A:112:LEU:CG	1:B:14:LEU:CD2	2.79	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:GLU:HG3	2:B:281:HOH:O	2.15	0.46
1:A:112:LEU:CB	1:B:14:LEU:HD23	2.21	0.46
1:B:30:ALA:HA	1:B:31:THR:C	2.33	0.46
1:B:248:THR:O	1:B:251:GLU:HG2	2.16	0.46
1:A:177:MET:O	1:A:181:ARG:HD2	2.16	0.46
1:A:197:LEU:O	1:A:201:THR:HG23	2.16	0.45
1:A:132:LEU:HG	1:B:228:ARG:HH11	1.76	0.45
1:A:62:LEU:O	1:A:65:GLN:CB	2.65	0.45
1:A:36:THR:OG1	1:A:65:GLN:NE2	2.48	0.45
1:A:201:THR:CG2	2:B:304:HOH:O	2.65	0.45
1:B:23:TRP:N	1:B:24:ASP:CA	2.79	0.45
1:B:240:GLU:O	1:B:244:VAL:CG2	2.62	0.45
1:A:23:TRP:O	1:A:26:GLU:HG3	2.17	0.45
1:A:131:ALA:O	1:A:151:ARG:HB3	2.18	0.44
1:B:28:THR:N	1:B:30:ALA:O	2.50	0.44
1:A:22:PRO:HB2	1:A:23:TRP:CA	2.48	0.44
1:B:74:GLN:NE2	1:B:81:PHE:H	2.06	0.44
1:B:74:GLN:NE2	1:B:80:ASP:CG	2.63	0.43
1:B:234:VAL:HG11	1:B:241:MET:HE3	2.01	0.43
1:A:130:SER:OG	1:B:231:GLU:OE1	2.22	0.43
1:B:4:ILE:HD12	1:B:4:ILE:N	2.32	0.43
1:A:23:TRP:HA	1:A:26:GLU:HG3	2.01	0.43
1:A:249:MET:HE2	2:A:301:HOH:O	2.18	0.42
1:A:112:LEU:HA	1:A:113:ALA:O	2.19	0.42
1:A:112:LEU:HA	1:A:112:LEU:HD22	1.76	0.42
1:A:119:LYS:HA	1:A:119:LYS:HD3	1.73	0.42
1:B:30:ALA:C	1:B:32:ILE:N	2.72	0.42
1:A:16:ASP:HA	1:A:17:PRO:HD3	1.89	0.41
1:B:229:GLU:O	1:B:233:ILE:HG13	2.21	0.41
1:B:159:TRP:CE3	1:B:165:VAL:HG22	2.56	0.41
1:B:207:HIS:HD2	2:B:280:HOH:O	2.04	0.41
1:A:5:ASP:O	1:A:9:THR:CG2	2.69	0.41
1:A:132:LEU:CD2	2:A:307:HOH:O	2.14	0.41
1:A:21:CYS:HG	1:A:25:LYS:H	1.65	0.41
1:A:250:GLU:OE2	1:A:254:GLN:NE2	2.48	0.41
1:B:81:PHE:O	1:B:84:ILE:HB	2.20	0.41
1:A:7:LEU:HD21	1:A:63:LEU:HD13	2.03	0.41
1:A:19:ASN:HB3	1:A:20:GLY:O	2.21	0.41
1:B:203:ASN:HD22	1:B:203:ASN:HA	1.74	0.40
1:B:28:THR:O	1:B:29:PHE:CB	2.69	0.40
1:B:159:TRP:CZ3	2:B:295:HOH:O	2.57	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:LEU:HB2	1:B:14:LEU:CD2	2.14	0.40
1:B:23:TRP:N	1:B:24:ASP:C	2.74	0.40
1:A:56:ARG:HG3	1:A:57:GLY:N	2.35	0.40
1:B:63:LEU:HD22	1:B:63:LEU:O	2.22	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:LEU:CD2	1:B:254:GLN:NE2[4_456]	1.24	0.96
1:A:239:LEU:CD2	1:B:254:GLN:CD[4_456]	1.55	0.65
1:A:239:LEU:CG	1:B:254:GLN:NE2[4_456]	2.08	0.12
1:A:242:THR:OG1	1:B:250:GLU:OE1[4_456]	2.10	0.10
1:A:239:LEU:CD2	1:B:254:GLN:OE1[4_456]	2.13	0.07

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	233/265 (88%)	206 (88%)	14 (6%)	13 (6%)	2 0
1	B	216/265 (82%)	203 (94%)	7 (3%)	6 (3%)	6 2
All	All	449/530 (85%)	409 (91%)	21 (5%)	19 (4%)	3 1

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	PRO
1	A	18	GLU
1	A	21	CYS
1	A	22	PRO
1	A	239	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	29	PHE
1	A	23	TRP
1	A	94	ARG
1	A	132	LEU
1	A	240	GLU
1	B	13	ARG
1	B	24	ASP
1	B	28	THR
1	B	31	THR
1	A	124	ALA
1	A	128	GLN
1	A	248	THR
1	B	30	ALA
1	A	259	GLN

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	205/227 (90%)	182 (89%)	23 (11%)	7 4
1	B	190/227 (84%)	170 (90%)	20 (10%)	8 5
All	All	395/454 (87%)	352 (89%)	43 (11%)	8 4

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	THR
1	A	18	GLU
1	A	26	GLU
1	A	51	ASP
1	A	56	ARG
1	A	62	LEU
1	A	63	LEU
1	A	68	PHE
1	A	88	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	93	GLU
1	A	112	LEU
1	A	114	ARG
1	A	128	GLN
1	A	129	HIS
1	A	132	LEU
1	A	197	LEU
1	A	203	ASN
1	A	204	LEU
1	A	232	ARG
1	A	237	ARG
1	A	240	GLU
1	A	242	THR
1	A	254	GLN
1	B	3	GLN
1	B	7	LEU
1	B	13	ARG
1	B	14	LEU
1	B	24	ASP
1	B	27	GLN
1	B	28	THR
1	B	62	LEU
1	B	63	LEU
1	B	68	PHE
1	B	80	ASP
1	B	94	ARG
1	B	126	LYS
1	B	150	LYS
1	B	155	VAL
1	B	184	VAL
1	B	197	LEU
1	B	203	ASN
1	B	204	LEU
1	B	208	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	65	GLN
1	A	74	GLN
1	A	146	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	203	ASN
1	A	259	GLN
1	B	65	GLN
1	B	74	GLN
1	B	129	HIS
1	B	146	GLN
1	B	203	ASN
1	B	259	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

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	239/265 (90%)	1.11	53 (22%)	1	1	22, 44, 78, 89	0
1	B	222/265 (83%)	0.78	25 (11%)	7	9	20, 41, 73, 89	0
All	All	461/530 (86%)	0.95	78 (16%)	2	3	20, 42, 76, 89	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	79	PHE	9.7
1	B	125	GLN	9.2
1	A	239	LEU	8.5
1	B	261	ILE	7.3
1	A	124	ALA	7.1
1	A	19	ASN	6.0
1	A	23	TRP	5.7
1	A	66	VAL	5.2
1	B	14	LEU	5.0
1	A	21	CYS	4.7
1	B	63	LEU	4.6
1	A	248	THR	4.6
1	A	18	GLU	4.6
1	A	126	LYS	4.4
1	A	118	ILE	4.3
1	A	63	LEU	4.2
1	A	62	LEU	3.9
1	A	22	PRO	3.9
1	A	258	ARG	3.8
1	B	257	LYS	3.8
1	B	74	GLN	3.7
1	A	94	ARG	3.6
1	A	84	ILE	3.5
1	B	83	ASP	3.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	242	THR	3.4
1	A	59	LEU	3.4
1	A	257	LYS	3.3
1	A	237	ARG	3.3
1	A	125	GLN	3.3
1	B	259	GLN	3.3
1	A	128	GLN	3.2
1	A	255	GLN	3.2
1	A	17	PRO	3.2
1	A	198	LEU	3.1
1	B	4	ILE	3.0
1	A	259	GLN	3.0
1	A	70	ALA	3.0
1	A	67	VAL	3.0
1	A	81	PHE	2.9
1	B	29	PHE	2.9
1	B	94	ARG	2.9
1	A	26	GLU	2.7
1	A	254	GLN	2.7
1	B	62	LEU	2.7
1	A	20	GLY	2.7
1	A	251	GLU	2.6
1	B	260	GLU	2.6
1	A	120	THR	2.6
1	A	252	VAL	2.6
1	A	169	VAL	2.5
1	B	66	VAL	2.5
1	A	52	PHE	2.5
1	A	53	ASP	2.4
1	B	64	PHE	2.4
1	A	201	THR	2.3
1	B	169	VAL	2.3
1	A	131	ALA	2.3
1	B	81	PHE	2.2
1	A	132	LEU	2.2
1	A	13	ARG	2.2
1	A	202	VAL	2.2
1	A	113	ALA	2.2
1	A	8	LEU	2.1
1	A	92	LEU	2.1
1	A	197	LEU	2.1
1	A	127	ALA	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	183	ALA	2.1
1	B	184	VAL	2.1
1	B	126	LYS	2.1
1	A	204	LEU	2.1
1	B	243	GLY	2.1
1	B	23	TRP	2.1
1	A	166	VAL	2.1
1	A	188	ALA	2.1
1	B	48	ALA	2.0
1	B	28	THR	2.0
1	A	200	ALA	2.0
1	B	78	ARG	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.