Full wwPDB X-ray Structure Validation Report (i)

Oct 8, 2023 – 03:13 PM EDT

PDB ID	:	6M8R
Title	:	Crystal structure of the KCTD16 BTB domain in complex with GABAB2 peptide
Authors	:	Zheng, S.; Kruse, A.C.
Deposited on	:	2018-08-22
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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

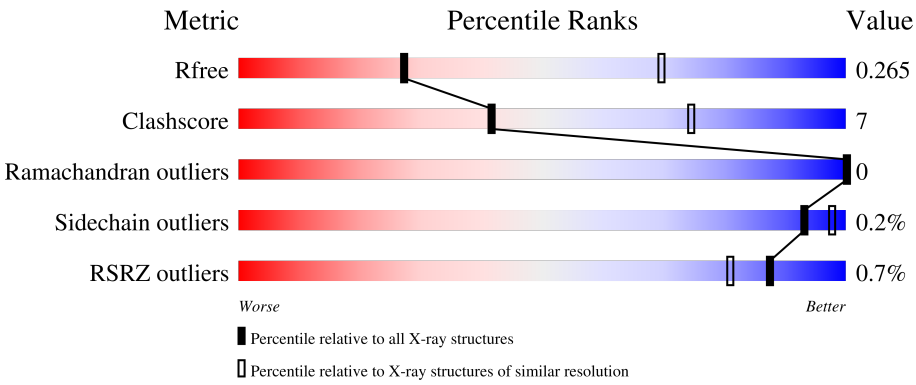
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 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}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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	103	2%; 73%; 17%; •; 8%
1	B	103	77%; 17%; 7%
1	C	103	78%; 17%; 6%
1	D	103	%; 82%; 14%; •; •
1	E	103	2%; 83%; 11%; 7%

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Mol	Chain	Length	Quality of chain
1	F	103	78% 16% 7%
1	G	103	81% 17%
1	H	103	81% 12% 8%
1	I	103	88% 9%
1	J	103	% 82% 13% 6%
2	K	41	66% 15% 20%
2	L	41	2% 61% 17% 22%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8578 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 BTB/POZ domain-containing protein KCTD16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	G	100	824	542	140	140	2	0	0	0
1	H	95	800	528	136	135	1	0	0	0
1	I	100	831	545	140	144	2	0	0	0
1	J	97	805	531	137	135	2	0	0	0
1	B	96	801	527	136	136	2	0	0	0
1	C	97	806	530	136	139	1	0	0	0
1	D	99	826	543	140	141	2	0	0	0
1	E	96	790	521	134	134	1	0	0	0
1	A	95	795	524	133	137	1	0	0	0
1	F	96	799	526	136	136	1	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	22	MET	-	initiating methionine	UNP Q68DU8
H	22	MET	-	initiating methionine	UNP Q68DU8
I	22	MET	-	initiating methionine	UNP Q68DU8
J	22	MET	-	initiating methionine	UNP Q68DU8
B	22	MET	-	initiating methionine	UNP Q68DU8
C	22	MET	-	initiating methionine	UNP Q68DU8
D	22	MET	-	initiating methionine	UNP Q68DU8
E	22	MET	-	initiating methionine	UNP Q68DU8
A	22	MET	-	initiating methionine	UNP Q68DU8

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Chain	Residue	Modelled	Actual	Comment	Reference
F	22	MET	-	initiating methionine	UNP Q68DU8

- Molecule 2 is a protein called Gamma-aminobutyric acid type B receptor subunit 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	K	33	254	161	48	45	0	0	0
			Total	C	N	O			
2	L	32	245	155	47	43	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	873	GLY	-	expression tag	UNP O75899
K	874	PRO	-	expression tag	UNP O75899
K	875	GLU	-	expression tag	UNP O75899
L	873	GLY	-	expression tag	UNP O75899
L	874	PRO	-	expression tag	UNP O75899
L	875	GLU	-	expression tag	UNP O75899

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

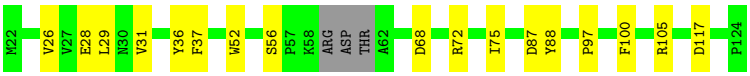
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
3	A	1	1	1	0	0
			Total	Mg		
3	F	1	1	1	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

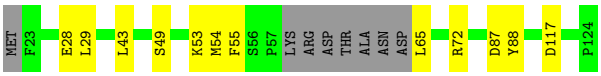
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain G:



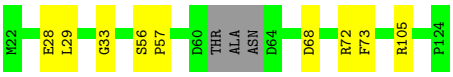
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain H:



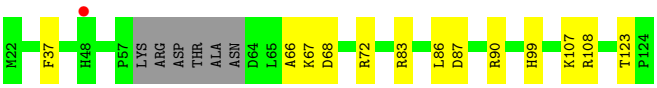
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain I:



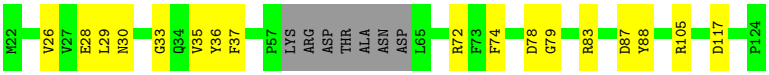
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain J:



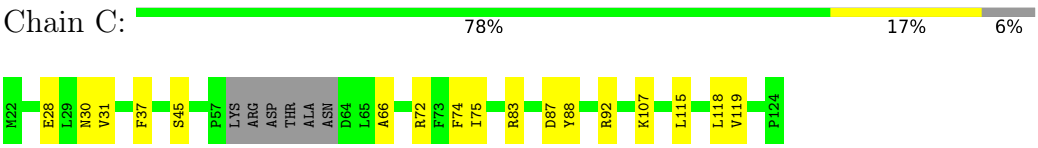
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain B:



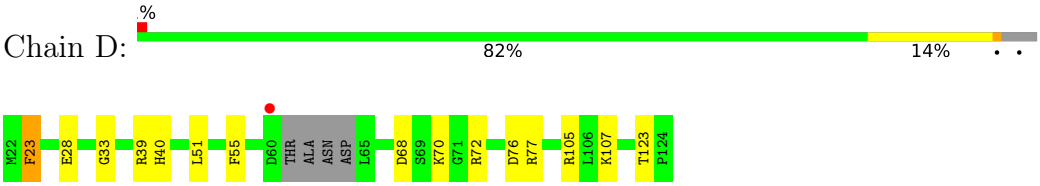
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain C:



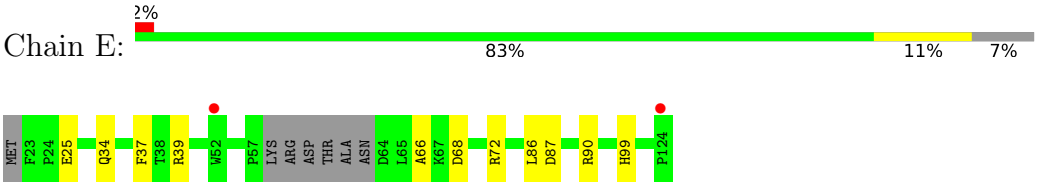
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain D:



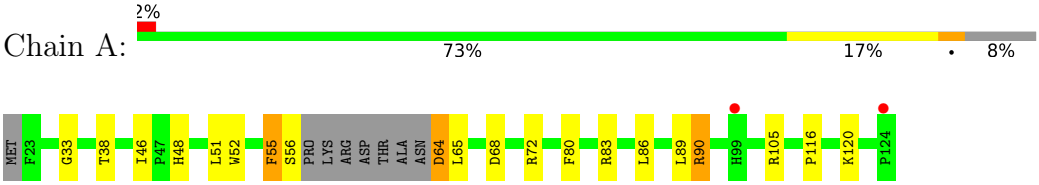
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain E:



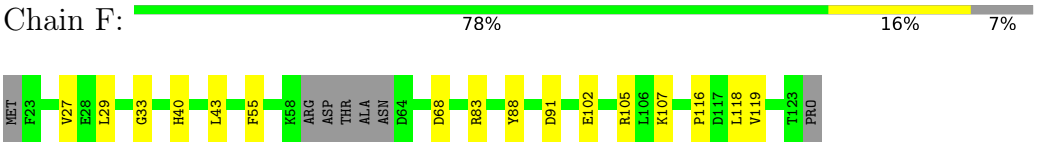
• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain A:



• Molecule 1: BTB/POZ domain-containing protein KCTD16

Chain F:

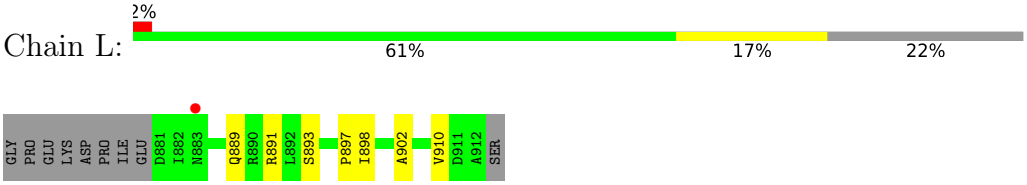


• Molecule 2: Gamma-aminobutyric acid type B receptor subunit 2

Chain K:

• Molecule 2: Gamma-aminobutyric acid type B receptor subunit 2

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	91.96Å 64.94Å 114.14Å 90.00° 99.79° 90.00°	Depositor
Resolution (Å)	38.64 – 3.20 38.64 – 3.20	Depositor EDS
% Data completeness (in resolution range)	98.9 (38.64-3.20) 98.9 (38.64-3.20)	Depositor EDS
R_{merge}	0.33	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.29 (at 3.18Å)	Xtriage
Refinement program	PHENIX (1.14_3211: ???)	Depositor
R, R_{free}	0.217 , 0.265 0.217 , 0.265	Depositor DCC
R_{free} test set	1686 reflections (7.66%)	wwPDB-VP
Wilson B-factor (Å^2)	60.0	Xtriage
Anisotropy	0.625	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 40.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	8578	wwPDB-VP
Average B, all atoms (Å^2)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.02% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.53	1/817 (0.1%)	0.75	3/1103 (0.3%)
1	B	0.26	0/824	0.47	0/1113
1	C	0.26	0/829	0.45	0/1122
1	D	0.26	0/849	0.45	0/1146
1	E	0.25	0/813	0.42	0/1102
1	F	0.25	0/821	0.44	0/1109
1	G	0.25	0/847	0.44	0/1145
1	H	0.25	0/823	0.44	0/1111
1	I	0.25	0/854	0.43	0/1154
1	J	0.25	0/828	0.45	0/1119
2	K	0.27	0/260	0.50	0/354
2	L	0.25	0/251	0.47	0/342
All	All	0.29	1/8816 (0.0%)	0.48	3/11920 (0.0%)

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	64	ASP	N-CA	12.27	1.70	1.46

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	64	ASP	N-CA-CB	-13.99	85.41	110.60
1	A	64	ASP	N-CA-C	9.74	137.30	111.00
1	A	55	PHE	CB-CA-C	6.44	123.28	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	90	ARG	Sidechain

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	795	0	792	35	0
1	B	801	0	795	15	0
1	C	806	0	792	12	0
1	D	826	0	823	16	0
1	E	790	0	773	11	0
1	F	799	0	792	12	0
1	G	824	0	815	12	0
1	H	800	0	806	8	0
1	I	831	0	818	6	0
1	J	805	0	802	10	0
2	K	254	0	254	6	0
2	L	245	0	240	7	0
3	A	1	0	0	0	0
3	F	1	0	0	0	0
All	All	8578	0	8502	119	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 7.

All (119) 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:64:ASP:N	1:A:64:ASP:CA	1.70	1.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:HIS:N	1:A:52:TRP:CE3	2.06	1.23
1:A:64:ASP:N	1:A:64:ASP:CB	2.20	1.04
1:A:48:HIS:N	1:A:52:TRP:CZ3	2.26	1.00
1:A:48:HIS:HA	1:A:52:TRP:HZ3	1.30	0.97
1:A:48:HIS:HA	1:A:52:TRP:CZ3	2.00	0.96
1:A:48:HIS:CA	1:A:52:TRP:CZ3	2.55	0.89
1:A:86:LEU:HD21	1:A:90:ARG:HH11	1.38	0.87
1:D:23:PHE:HE2	1:D:55:PHE:HD1	1.26	0.83
1:B:26:VAL:HG21	1:A:68:ASP:HA	1.59	0.82
1:A:64:ASP:OD1	1:A:65:LEU:N	2.13	0.80
1:B:87:ASP:OD2	1:A:105:ARG:NH2	2.15	0.79
1:A:48:HIS:CA	1:A:52:TRP:CE3	2.67	0.77
1:G:87:ASP:OD2	1:F:105:ARG:NH2	2.19	0.76
1:A:64:ASP:CG	1:A:65:LEU:H	1.86	0.76
1:A:86:LEU:HD23	1:A:90:ARG:HD2	1.70	0.73
1:A:48:HIS:O	1:A:48:HIS:ND1	2.22	0.71
1:J:67:LYS:H	2:K:883:ASN:HD21	1.39	0.71
1:A:64:ASP:N	1:A:64:ASP:HB2	2.05	0.70
1:A:86:LEU:CD2	1:A:90:ARG:HH11	2.03	0.70
1:A:48:HIS:N	1:A:52:TRP:HE3	1.87	0.68
1:B:35:VAL:HG12	2:L:910:VAL:O	1.97	0.65
2:L:897:PRO:HB3	2:L:902:ALA:HA	1.80	0.64
1:A:48:HIS:CA	1:A:52:TRP:HZ3	2.01	0.64
1:D:105:ARG:NH2	1:E:87:ASP:OD2	2.32	0.62
1:B:105:ARG:NH2	1:C:87:ASP:OD2	2.33	0.62
1:I:105:ARG:NH2	1:J:87:ASP:OD2	2.34	0.61
1:A:90:ARG:HH12	2:L:889:GLN:NE2	2.00	0.60
1:B:28:GLU:HG2	1:B:37:PHE:CZ	2.36	0.60
1:D:70:LYS:HD2	1:D:72:ARG:HH12	1.66	0.59
1:D:23:PHE:HD2	1:D:40:HIS:CD2	2.21	0.59
1:A:64:ASP:CG	1:A:65:LEU:N	2.56	0.58
1:A:46:ILE:HD12	1:A:89:LEU:HD23	1.86	0.57
1:D:23:PHE:CE2	1:D:55:PHE:HD1	2.15	0.57
1:G:105:ARG:NH2	1:H:87:ASP:OD2	2.36	0.57
1:A:86:LEU:O	1:A:90:ARG:HG3	2.05	0.57
1:G:26:VAL:HG21	1:F:68:ASP:HA	1.86	0.56
1:J:99:HIS:ND1	1:J:99:HIS:O	2.38	0.56
1:J:66:ALA:HB2	2:K:882:ILE:HD12	1.87	0.56
1:H:28:GLU:HG2	1:H:72:ARG:HG2	1.87	0.55
1:D:68:ASP:OD1	1:D:72:ARG:N	2.40	0.55
1:E:66:ALA:HB1	2:L:891:ARG:HH12	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ASP:OD1	1:A:72:ARG:N	2.40	0.54
1:A:116:PRO:O	1:A:120:LYS:HG3	2.06	0.54
1:G:52:TRP:O	1:G:56:SER:HB2	2.08	0.54
1:A:48:HIS:HA	1:A:52:TRP:CE3	2.39	0.54
1:A:86:LEU:O	1:A:90:ARG:CG	2.56	0.54
1:A:38:THR:HB	1:A:86:LEU:HD11	1.89	0.53
1:F:27:VAL:HG11	1:F:55:PHE:HE2	1.72	0.53
1:E:66:ALA:HB1	2:L:891:ARG:NH1	2.24	0.53
1:D:23:PHE:HE2	1:D:55:PHE:CD1	2.17	0.53
1:F:107:LYS:HG2	1:F:119:VAL:HG23	1.91	0.52
1:E:34:GLN:HE21	2:L:898:ILE:HD12	1.75	0.51
1:C:28:GLU:HG2	1:C:72:ARG:HG2	1.92	0.51
1:F:29:LEU:HD11	1:F:43:LEU:HD21	1.92	0.51
1:A:80:PHE:O	1:A:83:ARG:HG2	2.11	0.51
1:G:28:GLU:HG2	1:G:37:PHE:CE2	2.46	0.50
1:E:68:ASP:OD1	1:E:72:ARG:N	2.44	0.50
1:J:108:ARG:HD3	1:E:99:HIS:CE1	2.47	0.50
1:C:107:LYS:HG3	1:C:119:VAL:HG13	1.94	0.50
1:A:51:LEU:O	1:A:55:PHE:HD1	1.94	0.50
1:A:86:LEU:HD21	1:A:90:ARG:NH1	2.18	0.49
1:E:25:GLU:OE2	1:E:39:ARG:NH2	2.46	0.49
1:D:51:LEU:O	1:D:55:PHE:HD2	1.96	0.49
1:E:72:ARG:NH2	2:L:893:SER:OG	2.45	0.49
1:G:97:PRO:HD2	1:G:100:PHE:CD1	2.47	0.48
1:H:29:LEU:HD21	1:H:55:PHE:HZ	1.78	0.48
1:D:23:PHE:HD2	1:D:40:HIS:HD2	1.61	0.48
1:B:33:GLY:HA2	1:C:37:PHE:CD1	2.48	0.48
1:I:33:GLY:HA2	1:J:37:PHE:CD1	2.49	0.48
1:C:31:VAL:HG13	1:C:75:ILE:HB	1.96	0.48
1:F:116:PRO:O	1:F:119:VAL:HG12	2.15	0.47
1:C:66:ALA:HB2	1:D:39:ARG:HH12	1.79	0.47
1:F:91:ASP:OD2	2:K:890:ARG:NH1	2.48	0.47
1:H:88:TYR:OH	1:H:117:ASP:HB2	2.15	0.46
1:C:115:LEU:O	1:C:119:VAL:HG23	2.15	0.46
1:F:40:HIS:NE2	1:F:55:PHE:O	2.43	0.46
1:D:70:LYS:HD2	1:D:72:ARG:NH1	2.31	0.46
1:J:68:ASP:OD1	1:J:72:ARG:N	2.49	0.45
1:B:78:ASP:OD2	1:C:83:ARG:NE	2.48	0.45
1:D:28:GLU:HG2	1:D:72:ARG:HG2	1.98	0.45
1:A:52:TRP:CD1	1:A:52:TRP:O	2.69	0.45
1:J:107:LYS:HE3	1:J:123:THR:OG1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:88:TYR:OH	1:G:117:ASP:HB2	2.17	0.44
1:D:76:ASP:OD2	1:E:39:ARG:HG3	2.17	0.44
1:I:56:SER:OG	1:I:57:PRO:HD3	2.17	0.44
1:I:68:ASP:OD2	1:I:72:ARG:HB2	2.17	0.44
1:G:36:TYR:OH	2:K:908:GLY:O	2.24	0.44
1:B:28:GLU:HB3	1:B:72:ARG:HG2	2.00	0.44
1:I:28:GLU:HG2	1:I:72:ARG:HG2	1.98	0.44
1:H:43:LEU:HD13	1:H:55:PHE:CE2	2.52	0.44
1:B:30:ASN:O	1:B:74:PHE:HA	2.18	0.44
1:B:105:ARG:HH12	1:C:83:ARG:CZ	2.31	0.43
1:G:31:VAL:HG22	1:G:75:ILE:HB	1.98	0.43
1:J:83:ARG:HH11	1:J:83:ARG:HG3	1.84	0.43
1:C:88:TYR:CG	1:C:118:LEU:HD22	2.54	0.43
1:C:45:SER:HB3	1:C:92:ARG:HH22	1.84	0.43
1:J:86:LEU:HG	1:J:90:ARG:NH1	2.34	0.42
1:B:29:LEU:HD23	1:B:29:LEU:HA	1.89	0.42
1:G:29:LEU:HB2	1:G:36:TYR:HB2	2.00	0.42
1:H:54:MET:HE3	1:H:65:LEU:HD11	2.01	0.42
1:I:29:LEU:HD23	1:I:73:PHE:HB2	2.00	0.42
1:C:30:ASN:O	1:C:74:PHE:HA	2.20	0.42
1:B:37:PHE:HB2	1:A:33:GLY:N	2.35	0.41
1:F:83:ARG:NE	2:K:892:LEU:HD13	2.35	0.41
1:D:33:GLY:HA2	1:E:37:PHE:CD1	2.56	0.41
1:D:107:LYS:NZ	1:D:123:THR:HG22	2.35	0.41
1:F:33:GLY:O	2:K:910:VAL:HG21	2.19	0.41
1:F:88:TYR:CD2	1:F:118:LEU:HD12	2.55	0.41
1:D:76:ASP:O	1:D:77:ARG:HD2	2.21	0.41
1:H:49:SER:O	1:H:53:LYS:HG3	2.20	0.41
1:B:36:TYR:OH	1:B:79:GLY:O	2.26	0.41
1:G:97:PRO:HB3	1:F:102:GLU:HG2	2.03	0.41
1:B:83:ARG:HH21	1:A:80:PHE:HB3	1.86	0.41
1:B:88:TYR:OH	1:B:117:ASP:HB2	2.20	0.41
1:E:86:LEU:HG	1:E:90:ARG:NH1	2.35	0.41
1:H:29:LEU:HD21	1:H:55:PHE:CZ	2.56	0.41
1:G:68:ASP:OD1	1:G:72:ARG:N	2.54	0.40
1:A:48:HIS:CA	1:A:52:TRP:HE3	2.26	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	91/103 (88%)	89 (98%)	2 (2%)	0	100	100
1	B	92/103 (89%)	90 (98%)	2 (2%)	0	100	100
1	C	93/103 (90%)	92 (99%)	1 (1%)	0	100	100
1	D	95/103 (92%)	92 (97%)	3 (3%)	0	100	100
1	E	92/103 (89%)	89 (97%)	3 (3%)	0	100	100
1	F	92/103 (89%)	87 (95%)	5 (5%)	0	100	100
1	G	96/103 (93%)	92 (96%)	4 (4%)	0	100	100
1	H	91/103 (88%)	88 (97%)	3 (3%)	0	100	100
1	I	96/103 (93%)	93 (97%)	3 (3%)	0	100	100
1	J	93/103 (90%)	90 (97%)	3 (3%)	0	100	100
2	K	31/41 (76%)	29 (94%)	2 (6%)	0	100	100
2	L	30/41 (73%)	28 (93%)	2 (7%)	0	100	100
All	All	992/1112 (89%)	959 (97%)	33 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	87/95 (92%)	86 (99%)	1 (1%)	73	88
1	B	87/95 (92%)	87 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	87/95 (92%)	87 (100%)	0	100	100
1	D	90/95 (95%)	89 (99%)	1 (1%)	73	88
1	E	84/95 (88%)	84 (100%)	0	100	100
1	F	86/95 (90%)	86 (100%)	0	100	100
1	G	88/95 (93%)	88 (100%)	0	100	100
1	H	88/95 (93%)	88 (100%)	0	100	100
1	I	90/95 (95%)	90 (100%)	0	100	100
1	J	87/95 (92%)	87 (100%)	0	100	100
2	K	28/36 (78%)	28 (100%)	0	100	100
2	L	26/36 (72%)	26 (100%)	0	100	100
All	All	928/1022 (91%)	926 (100%)	2 (0%)	93	98

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	23	PHE
1	A	56	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	I	99	HIS
1	D	40	HIS
1	E	99	HIS
2	K	883	ASN
2	L	889	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 monosaccharides 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, 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	95/103 (92%)	-0.12	2 (2%) 63 49	39, 68, 95, 118	0
1	B	96/103 (93%)	-0.39	0 100 100	25, 44, 73, 97	0
1	C	97/103 (94%)	-0.43	0 100 100	27, 41, 69, 86	0
1	D	99/103 (96%)	-0.34	1 (1%) 82 72	25, 47, 78, 93	0
1	E	96/103 (93%)	0.06	2 (2%) 63 49	43, 72, 106, 127	0
1	F	96/103 (93%)	-0.16	0 100 100	37, 60, 82, 95	0
1	G	100/103 (97%)	-0.41	0 100 100	28, 46, 78, 87	0
1	H	95/103 (92%)	-0.43	0 100 100	23, 41, 65, 80	0
1	I	100/103 (97%)	-0.40	0 100 100	25, 44, 69, 79	0
1	J	97/103 (94%)	-0.18	1 (1%) 82 72	34, 56, 84, 105	0
2	K	33/41 (80%)	-0.03	0 100 100	35, 56, 80, 87	0
2	L	32/41 (78%)	0.02	1 (3%) 49 32	39, 55, 96, 97	0
All	All	1036/1112 (93%)	-0.26	7 (0%) 87 81	23, 51, 88, 127	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	124	PRO	4.3
2	L	883	ASN	3.3
1	A	124	PRO	2.6
1	E	52	TRP	2.4
1	D	60	ASP	2.0
1	J	48	HIS	2.0
1	A	99	HIS	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 monosaccharides 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. The B-factors column lists the minimum, median, 95th 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	B-factors(Å^2)	Q<0.9
3	MG	F	201	1/1	0.89	0.08	59,59,59,59	0
3	MG	A	201	1/1	0.93	0.20	85,85,85,85	0

6.5 Other polymers (i)

There are no such residues in this entry.