



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 01:35 am BST

PDB ID : 2WIA  
Title : Crystal Structures of the N-terminal Intracellular Domain of FeoB from *Klebsiella Pneumoniae* in Apo Form  
Authors : Hung, K.-W.; Chang, Y.-W.; Chen, J.-H.; Chen, Y.-C.; Sun, Y.-J.; Hsiao, C.-D.; Huang, T.-H.  
Deposited on : 2009-05-09  
Resolution : 2.45 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

<https://www.wwpdb.org/validation/2017/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) : 1.13  
EDS : 2.11  
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) : 2.11

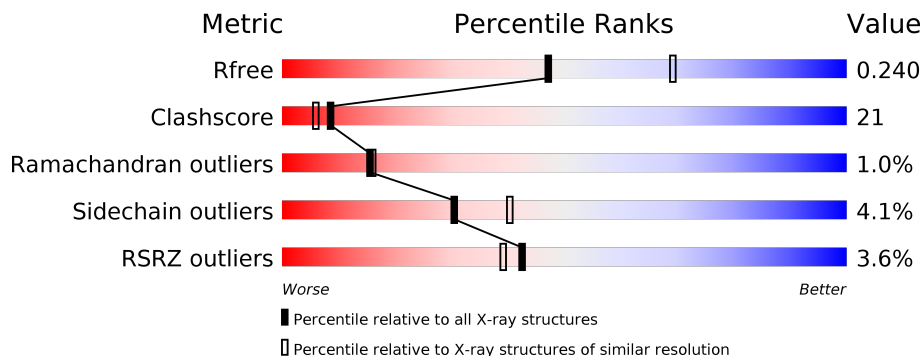
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 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  $\leq 5\%$ . 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	267	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 57%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 33%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">2%      57%      33%      • • • 5%</p>
1	B	267	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 56%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 34%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">4%      56%      34%      • • • 6%</p>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3889 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 FERROUS IRON TRANSPORT PROTEIN B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	253	1864	1170	331	354	9	0	0	0
1	B	252	1825	1142	323	351	9	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	129	GLN	LYS	conflict	UNP A6TF32
B	129	GLN	LYS	conflict	UNP A6TF32

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

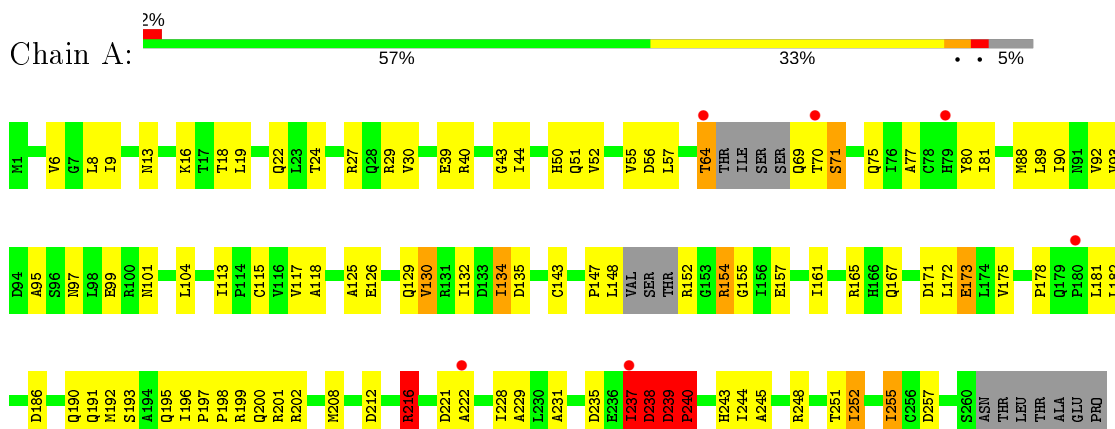
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	103	103	103	0	0
3	B	95	95	95	0	0

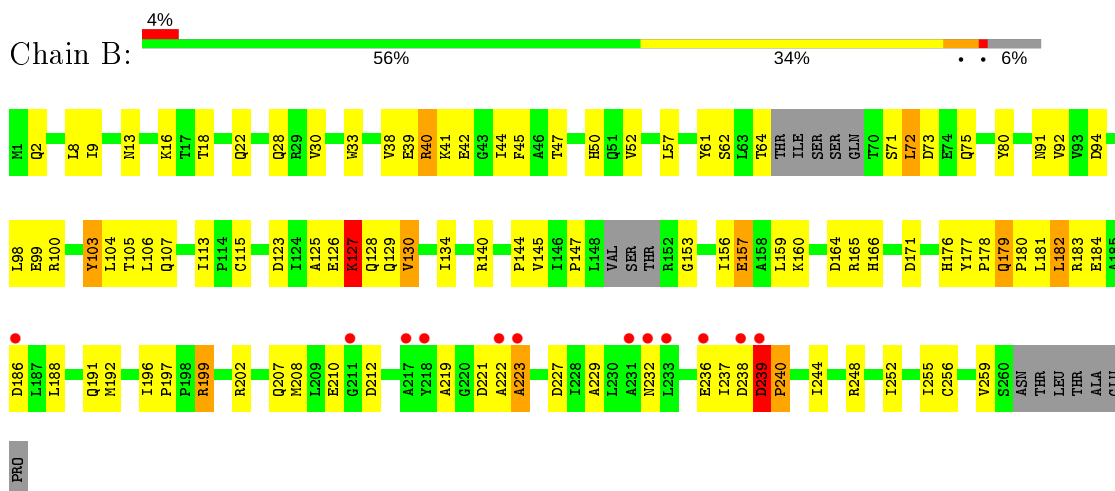
### 3 Residue-property plots [i](#)

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: FERROUS IRON TRANSPORT PROTEIN B



#### • Molecule 1: FERROUS IRON TRANSPORT PROTEIN B



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.20 Å 84.20 Å 122.46 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.13 – 2.45 25.13 – 2.45	Depositor EDS
% Data completeness (in resolution range)	93.3 (25.13-2.45) 97.7 (25.13-2.45)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.38 (at 2.44 Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.209 , 0.230 0.215 , 0.240	Depositor DCC
$R_{free}$ test set	876 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.9	Xtrriage
Anisotropy	0.552	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 22.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.216 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	3889	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.60% 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 [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	1.41	9/1889 (0.5%)	1.36	15/2576 (0.6%)
1	B	1.43	13/1849 (0.7%)	1.43	26/2525 (1.0%)
All	All	1.42	22/3738 (0.6%)	1.39	41/5101 (0.8%)

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	8
1	B	0	3
All	All	0	11

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	236	GLU	C-N	-12.70	1.04	1.34
1	A	240	PRO	N-CD	7.34	1.58	1.47
1	B	103	TYR	C-O	7.17	1.36	1.23
1	A	252	ILE	CB-CG2	7.17	1.75	1.52
1	A	55	VAL	CB-CG2	6.55	1.66	1.52
1	B	191	GLN	C-O	6.02	1.34	1.23
1	B	180	PRO	N-CD	5.98	1.56	1.47
1	B	239	ASP	C-N	5.93	1.45	1.34
1	B	259	VAL	CB-CG2	-5.92	1.40	1.52
1	B	157	GLU	CD-OE2	5.83	1.32	1.25
1	B	127	LYS	CA-CB	5.80	1.66	1.53
1	B	61	TYR	CD2-CE2	5.55	1.47	1.39
1	A	175	VAL	CB-CG2	5.53	1.64	1.52
1	B	61	TYR	CZ-OH	5.50	1.47	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	115	CYS	CB-SG	-5.48	1.72	1.81
1	B	179	GLN	C-N	5.28	1.44	1.34
1	A	40	ARG	CZ-NH2	5.16	1.39	1.33
1	A	237	ILE	C-N	5.16	1.46	1.34
1	A	117	VAL	CB-CG1	5.13	1.63	1.52
1	A	93	VAL	CB-CG1	5.13	1.63	1.52
1	A	134	ILE	CA-CB	5.08	1.66	1.54
1	B	145	VAL	CB-CG1	5.04	1.63	1.52

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	238	ASP	C-N-CA	12.88	153.91	121.70
1	B	164	ASP	CB-CG-OD2	-12.71	106.86	118.30
1	B	239	ASP	C-N-CD	11.62	152.80	128.40
1	A	240	PRO	CA-N-CD	-10.33	97.04	111.50
1	B	176	HIS	CA-C-O	8.74	138.46	120.10
1	A	71	SER	O-C-N	-7.84	110.15	122.70
1	A	239	ASP	C-N-CD	7.82	144.83	128.40
1	A	40	ARG	NE-CZ-NH1	-7.76	116.42	120.30
1	A	237	ILE	CA-C-N	7.57	133.86	117.20
1	B	164	ASP	CB-CG-OD1	7.50	125.05	118.30
1	A	239	ASP	O-C-N	-7.45	106.94	121.10
1	B	179	GLN	C-N-CD	7.02	143.13	128.40
1	B	237	ILE	C-N-CA	-6.89	104.47	121.70
1	B	94	ASP	CB-CG-OD1	6.84	124.46	118.30
1	B	240	PRO	CA-N-CD	-6.77	102.02	111.50
1	A	237	ILE	O-C-N	-6.76	111.89	122.70
1	B	221	ASP	CB-CG-OD2	6.69	124.32	118.30
1	B	73	ASP	CB-CA-C	-6.57	97.25	110.40
1	B	72	LEU	CB-CG-CD2	-6.31	100.27	111.00
1	B	94	ASP	CB-CG-OD2	-6.24	112.69	118.30
1	A	199	ARG	NE-CZ-NH2	-6.03	117.29	120.30
1	B	106	LEU	CB-CG-CD1	-5.89	100.98	111.00
1	B	199	ARG	NE-CZ-NH2	-5.78	117.41	120.30
1	B	180	PRO	CA-N-CD	-5.75	103.45	111.50
1	B	182	LEU	CB-CG-CD1	-5.70	101.30	111.00
1	A	248	ARG	NE-CZ-NH1	-5.55	117.52	120.30
1	A	135	ASP	CB-CG-OD1	5.51	123.26	118.30
1	A	216	ARG	CB-CA-C	-5.49	99.43	110.40
1	B	239	ASP	O-C-N	-5.46	110.72	121.10
1	A	239	ASP	N-CA-C	5.42	125.63	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	223	ALA	CA-C-N	-5.41	105.31	117.20
1	B	248	ARG	NE-CZ-NH1	-5.40	117.60	120.30
1	B	100	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	A	51	GLN	N-CA-C	-5.34	96.57	111.00
1	A	238	ASP	CB-CA-C	5.31	121.02	110.40
1	B	40	ARG	NE-CZ-NH1	-5.29	117.65	120.30
1	B	103	TYR	CB-CA-C	5.29	120.99	110.40
1	B	40	ARG	CB-CG-CD	-5.24	97.97	111.60
1	B	165	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	B	176	HIS	CA-C-N	-5.21	105.74	117.20
1	B	38	VAL	CB-CA-C	-5.07	101.78	111.40

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	237	ILE	Mainchain,Peptide
1	A	238	ASP	Mainchain,Peptide
1	A	239	ASP	Mainchain
1	A	240	PRO	Mainchain
1	A	69	GLN	Mainchain
1	A	70	THR	Mainchain
1	B	103	TYR	Mainchain
1	B	223	ALA	Mainchain
1	B	239	ASP	Mainchain

## 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	1864	0	1822	90	0
1	B	1825	0	1750	61	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	103	0	0	8	0
3	B	95	0	0	8	0
All	All	3889	0	3572	150	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 21.

All (150) 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:252:ILE:CB	1:A:252:ILE:CG2	1.75	1.57
1:A:228:ILE:O	1:A:231:ALA:HB3	1.62	0.99
1:B:64:THR:HG21	1:B:207:GLN:HE21	1.28	0.98
1:B:140:ARG:HD2	1:B:255:ILE:HD13	1.49	0.93
1:A:216:ARG:CG	1:A:216:ARG:HH21	1.85	0.90
1:B:184:GLU:OE1	1:B:232:ASN:HB2	1.72	0.89
1:A:237:ILE:HB	1:A:238:ASP:HA	1.54	0.88
1:A:216:ARG:HB2	1:A:216:ARG:NH2	1.88	0.88
1:B:64:THR:HG21	1:B:207:GLN:NE2	1.89	0.87
1:B:47:THR:HG22	1:B:160:LYS:HB3	1.58	0.86
1:B:178:PRO:HD2	1:B:181:LEU:HD12	1.58	0.84
1:A:251:THR:O	1:A:255:ILE:HD13	1.77	0.82
1:A:157:GLU:O	1:A:161:ILE:HD13	1.80	0.82
1:A:216:ARG:CB	1:A:216:ARG:HH21	1.94	0.80
1:B:30:VAL:CG1	1:B:39:GLU:HG3	2.12	0.79
1:A:216:ARG:HH21	1:A:216:ARG:HG2	1.47	0.79
1:A:216:ARG:HB2	1:A:216:ARG:HH21	1.46	0.79
1:B:140:ARG:HB3	1:B:255:ILE:HD12	1.64	0.79
1:A:134:ILE:HD13	1:A:147:PRO:HG3	1.65	0.78
1:A:252:ILE:HB	1:A:252:ILE:CG2	2.12	0.78
1:A:237:ILE:HG13	1:A:238:ASP:HB3	1.68	0.75
1:B:239:ASP:O	1:B:240:PRO:C	2.25	0.74
1:A:27:ARG:HH11	1:A:27:ARG:HG3	1.50	0.73
1:A:178:PRO:HD2	1:A:181:LEU:HD12	1.69	0.73
1:A:30:VAL:HG13	1:A:39:GLU:HG3	1.70	0.72
1:A:192:MET:CE	1:A:208:MET:CE	2.68	0.72
1:A:196:ILE:O	1:A:201:ARG:NH1	2.22	0.71
1:B:179:GLN:HG2	3:B:2078:HOH:O	1.91	0.70
1:A:192:MET:HE1	1:A:208:MET:CE	2.24	0.68
1:B:210:GLU:HG3	1:B:244:ILE:HB	1.75	0.67
1:A:252:ILE:CA	1:A:252:ILE:CG2	2.71	0.66
1:B:199:ARG:HB2	3:B:2081:HOH:O	1.95	0.66
1:B:40:ARG:NH1	1:B:42:GLU:OE2	2.29	0.66
1:A:228:ILE:O	1:A:231:ALA:CB	2.41	0.65
1:A:64:THR:HB	1:A:212:ASP:OD2	1.97	0.65
1:B:126:GLU:O	1:B:129:GLN:N	2.30	0.64
1:B:64:THR:HB	1:B:212:ASP:OD2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:ASP:HB2	3:B:2074:HOH:O	1.95	0.64
1:B:181:LEU:CD2	1:B:240:PRO:HB3	2.28	0.64
1:A:191:GLN:O	1:A:221:ASP:HB3	1.97	0.63
1:A:237:ILE:HG13	1:A:238:ASP:CB	2.29	0.63
1:A:29:ARG:HB3	3:A:2021:HOH:O	1.98	0.63
1:B:123:ASP:N	1:B:123:ASP:OD1	2.31	0.62
1:A:192:MET:CE	1:A:208:MET:HE1	2.29	0.61
1:A:238:ASP:HB2	1:A:240:PRO:HG3	1.80	0.61
1:B:140:ARG:HB3	1:B:255:ILE:CD1	2.31	0.60
1:B:153:GLY:O	1:B:157:GLU:HG2	2.00	0.60
1:A:8:LEU:HD21	1:A:16:LYS:HG3	1.83	0.60
1:B:210:GLU:HA	1:B:244:ILE:HD12	1.83	0.60
1:B:140:ARG:CD	1:B:255:ILE:HD13	2.27	0.60
1:B:9:ILE:HA	3:B:2002:HOH:O	2.02	0.60
1:B:188:LEU:HD13	1:B:208:MET:HE3	1.84	0.60
1:A:6:VAL:HG11	1:A:90:ILE:HG13	1.84	0.59
1:A:237:ILE:HB	1:A:238:ASP:CA	2.31	0.59
1:A:50:HIS:CE1	3:A:2027:HOH:O	2.56	0.58
1:A:44:ILE:HA	1:A:52:VAL:O	2.03	0.58
1:A:192:MET:CE	1:A:208:MET:HE2	2.33	0.57
1:A:71:SER:O	1:A:75:GLN:HG3	2.04	0.57
1:A:81:ILE:HG23	1:A:113:ILE:HD11	1.86	0.57
1:A:77:ALA:O	1:A:81:ILE:HG12	2.05	0.56
1:A:181:LEU:HD13	1:A:244:ILE:HG13	1.88	0.56
1:B:30:VAL:HG12	1:B:39:GLU:HG3	1.88	0.56
1:B:99:GLU:HG3	1:B:252:ILE:HG22	1.86	0.56
1:A:216:ARG:CB	1:A:216:ARG:NH2	2.59	0.56
1:A:6:VAL:CG1	1:A:90:ILE:HG13	2.36	0.56
1:A:192:MET:HE2	1:A:208:MET:CE	2.35	0.55
1:A:50:HIS:HE1	3:A:2027:HOH:O	1.89	0.55
1:A:27:ARG:NH1	1:A:27:ARG:HG3	2.22	0.55
1:B:210:GLU:CA	1:B:244:ILE:HD12	2.37	0.55
1:A:229:ALA:HA	3:A:2097:HOH:O	2.07	0.54
1:A:81:ILE:HG23	1:A:113:ILE:CD1	2.37	0.54
1:A:134:ILE:CD1	1:A:147:PRO:HG3	2.35	0.54
1:A:64:THR:HG21	3:A:2091:HOH:O	2.08	0.54
1:B:57:LEU:HD21	1:B:80:TYR:CD2	2.44	0.54
1:B:153:GLY:O	1:B:157:GLU:CG	2.57	0.53
1:A:182:LEU:HD22	1:A:202:ARG:HH21	1.73	0.53
1:A:192:MET:HE2	1:A:208:MET:HE2	1.90	0.53
1:B:30:VAL:CG1	1:B:39:GLU:CG	2.85	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:LEU:O	1:B:107:GLN:HB2	2.08	0.53
1:A:18:THR:O	1:A:22:GLN:HG3	2.10	0.52
1:A:132:ILE:HD12	1:A:132:ILE:N	2.25	0.51
1:A:8:LEU:CD2	1:A:16:LYS:HG3	2.40	0.51
1:B:44:ILE:N	1:B:44:ILE:HD12	2.25	0.51
1:B:8:LEU:HG	1:B:16:LYS:HG3	1.93	0.50
1:B:9:ILE:HG13	1:B:104:LEU:HD21	1.94	0.50
1:A:81:ILE:HD12	1:A:89:LEU:HD11	1.93	0.50
1:B:126:GLU:O	1:B:127:LYS:C	2.49	0.50
1:A:24:THR:HG21	1:A:43:GLY:HA3	1.94	0.50
1:A:197:PRO:O	1:A:201:ARG:HG3	2.11	0.49
1:A:244:ILE:HG22	1:A:245:ALA:N	2.28	0.49
1:A:89:LEU:HD11	1:A:113:ILE:HD13	1.93	0.49
1:B:18:THR:O	1:B:22:GLN:HG3	2.12	0.49
1:B:238:ASP:CB	3:B:2087:HOH:O	2.60	0.48
1:B:219:ALA:HB1	1:B:222:ALA:O	2.13	0.48
1:A:115:CYS:O	1:A:143:CYS:HB2	2.14	0.47
1:B:156:ILE:O	1:B:159:LEU:HB3	2.14	0.47
1:A:118:ALA:HB1	1:A:148:LEU:HD11	1.96	0.47
1:A:197:PRO:HG2	1:A:200:GLN:HG3	1.97	0.47
1:A:8:LEU:HD11	1:A:92:VAL:CG2	2.45	0.47
1:B:182:LEU:O	1:B:183:ARG:C	2.52	0.47
1:B:186:ASP:OD2	1:B:202:ARG:NH2	2.48	0.47
1:B:8:LEU:HD11	1:B:92:VAL:CG2	2.45	0.47
1:A:155:GLY:N	3:A:2074:HOH:O	2.46	0.46
1:A:30:VAL:CG1	1:A:39:GLU:HG3	2.44	0.46
1:B:240:PRO:O	1:B:244:ILE:HG13	2.15	0.46
1:A:239:ASP:HB2	1:A:243:HIS:HE1	1.80	0.46
1:A:161:ILE:O	1:A:165:ARG:HG3	2.16	0.46
1:A:125:ALA:HB1	1:A:130:VAL:HG13	1.97	0.46
1:A:255:ILE:N	1:A:255:ILE:CD1	2.80	0.45
1:B:144:PRO:HD3	1:B:166:HIS:CE1	2.52	0.45
1:B:140:ARG:HD2	1:B:255:ILE:CD1	2.32	0.45
1:B:256:CYS:HB3	3:B:2093:HOH:O	2.17	0.45
1:A:57:LEU:HD21	1:A:80:TYR:CD2	2.52	0.45
1:A:216:ARG:HG2	1:A:216:ARG:NH2	2.23	0.45
1:B:177:TYR:HB2	1:B:182:LEU:HD11	1.99	0.45
1:A:193:SER:OG	1:A:195:GLN:HB2	2.17	0.44
1:B:28:GLN:HA	1:B:42:GLU:O	2.17	0.44
1:B:45:PHE:CZ	1:B:52:VAL:HG11	2.53	0.44
1:A:126:GLU:O	1:A:129:GLN:N	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:ASN:OD1	1:B:71:SER:HB2	2.17	0.44
1:A:64:THR:CB	1:A:212:ASP:OD2	2.65	0.44
1:A:81:ILE:CG2	1:A:113:ILE:HD11	2.48	0.44
1:B:9:ILE:HD12	3:B:2002:HOH:O	2.16	0.44
1:B:134:ILE:HG12	1:B:147:PRO:HG3	1.99	0.44
1:A:6:VAL:HG23	1:A:52:VAL:HG13	2.00	0.43
1:A:152:ARG:C	1:A:154:ARG:H	2.20	0.43
1:A:154:ARG:O	1:A:155:GLY:C	2.55	0.43
1:B:33:TRP:HA	3:B:2015:HOH:O	2.19	0.43
1:A:257:ASP:HA	3:A:2101:HOH:O	2.19	0.43
1:B:2:GLN:HB2	1:B:50:HIS:CD2	2.54	0.43
1:B:196:ILE:HA	1:B:197:PRO:HD2	1.84	0.42
1:A:167:GLN:OE1	1:A:167:GLN:HA	2.19	0.42
1:A:6:VAL:HA	1:A:88:MET:O	2.19	0.42
1:A:19:LEU:HD23	1:A:92:VAL:HG22	2.01	0.42
1:A:56:ASP:OD2	3:A:2034:HOH:O	2.22	0.42
1:A:95:ALA:HB1	1:A:132:ILE:HD11	2.03	0.41
1:B:125:ALA:O	1:B:130:VAL:HG13	2.20	0.41
1:B:91:ASN:HD22	1:B:105:THR:HG23	1.85	0.41
1:A:191:GLN:HB2	1:A:222:ALA:HB2	2.03	0.41
1:B:126:GLU:O	1:B:128:GLN:N	2.54	0.41
1:A:197:PRO:HG2	1:A:200:GLN:CD	2.41	0.41
1:A:99:GLU:HG3	1:A:252:ILE:HG22	2.02	0.41
1:A:192:MET:HE1	1:A:208:MET:HE2	1.97	0.41
1:A:197:PRO:HG2	1:A:200:GLN:CG	2.50	0.41
1:A:9:ILE:HG12	1:A:104:LEU:HD21	2.02	0.41
1:B:98:LEU:HA	1:B:98:LEU:HD12	1.73	0.41
1:A:97:ASN:O	1:A:101:ASN:ND2	2.46	0.40
1:A:172:LEU:O	1:A:173:GLU:C	2.59	0.40
1:B:184:GLU:OE1	1:B:229:ALA:HA	2.21	0.40
1:B:72:LEU:O	1:B:75:GLN:HB2	2.21	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	247/267 (92%)	226 (92%)	18 (7%)	3 (1%)	13	12
1	B	246/267 (92%)	228 (93%)	16 (6%)	2 (1%)	19	22
All	All	493/534 (92%)	454 (92%)	34 (7%)	5 (1%)	15	16

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	227	ASP
1	A	154	ARG
1	A	235	ASP
1	B	127	LYS
1	A	173	GLU

### 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	186/221 (84%)	178 (96%)	8 (4%)	29	38
1	B	178/221 (80%)	171 (96%)	7 (4%)	32	42
All	All	364/442 (82%)	349 (96%)	15 (4%)	30	40

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	THR
1	A	130	VAL
1	A	171	ASP
1	A	186	ASP
1	A	190	GLN
1	A	198	PRO
1	A	216	ARG
1	A	255	ILE

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Mol	Chain	Res	Type
1	B	13	ASN
1	B	41	LYS
1	B	62	SER
1	B	113	ILE
1	B	130	VAL
1	B	192	MET
1	B	239	ASP

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

Mol	Chain	Res	Type
1	A	28	GLN
1	A	97	ASN
1	A	179	GLN
1	B	28	GLN
1	B	91	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	236:GLU	C	237:ILE	N	1.04

## 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	253/267 (94%)	-0.02	6 (2%) 59 54	11, 27, 53, 64	0
1	B	252/267 (94%)	0.01	12 (4%) 30 28	12, 27, 56, 63	0
All	All	505/534 (94%)	-0.01	18 (3%) 42 39	11, 27, 55, 64	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	222	ALA	5.2
1	A	222	ALA	5.1
1	B	233	LEU	4.7
1	B	217	ALA	4.3
1	B	231	ALA	3.5
1	B	238	ASP	3.2
1	B	232	ASN	3.0
1	B	223	ALA	2.6
1	B	218	TYR	2.5
1	B	236	GLU	2.5
1	A	79	HIS	2.3
1	B	186	ASP	2.3
1	A	64	THR	2.2
1	B	211	GLY	2.2
1	A	180	PRO	2.2
1	A	237	ILE	2.2
1	A	70	THR	2.2
1	B	239	ASP	2.1

### 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. 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	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	B	1262	1/1	0.62	0.27	34,34,34,34	1
2	MG	A	1262	1/1	0.88	0.07	19,19,19,19	1

### 6.5 Other polymers [i](#)

There are no such residues in this entry.