



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:23 AM GMT

PDB ID : 3I8S
Title : Structure of the cytosolic domain of E. coli FeoB, nucleotide-free form
Authors : Petermann, N.; Hansen, G.; Hogg, T.; Hilgenfeld, R.
Deposited on : 2009-07-10
Resolution : 1.80 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

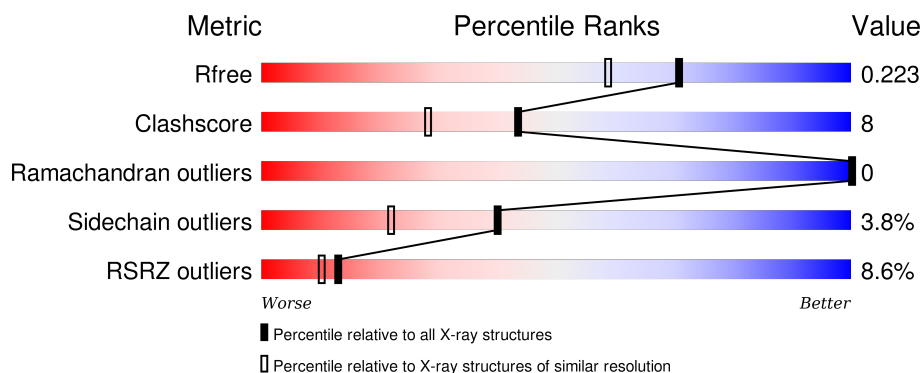
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

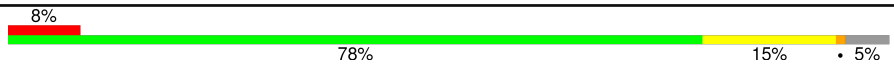
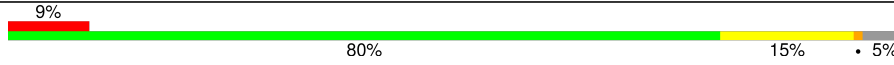

The reported resolution of this entry is 1.80 Å.

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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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	274	
1	B	274	
1	C	274	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6815 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
1	A	260	Total	C	N	O	S	0	14	0
			2084	1315	362	396	11			
1	B	261	Total	C	N	O	S	0	10	0
			2070	1302	363	393	12			
1	C	261	Total	C	N	O	S	0	8	0
			2057	1289	361	396	11			

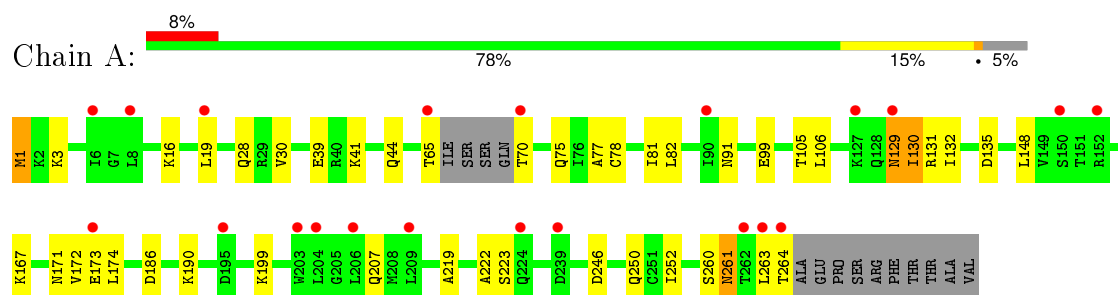
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	219	Total	O	0	0
			219	219		
2	B	203	Total	O	0	0
			203	203		
2	C	182	Total	O	0	0
			182	182		

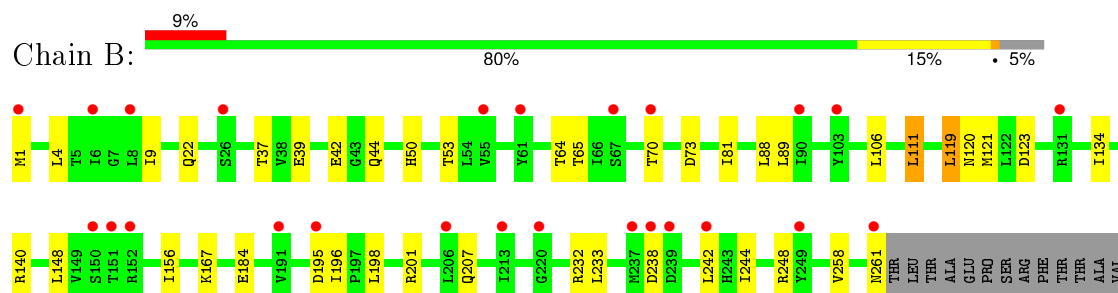
3 Residue-property plots

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.

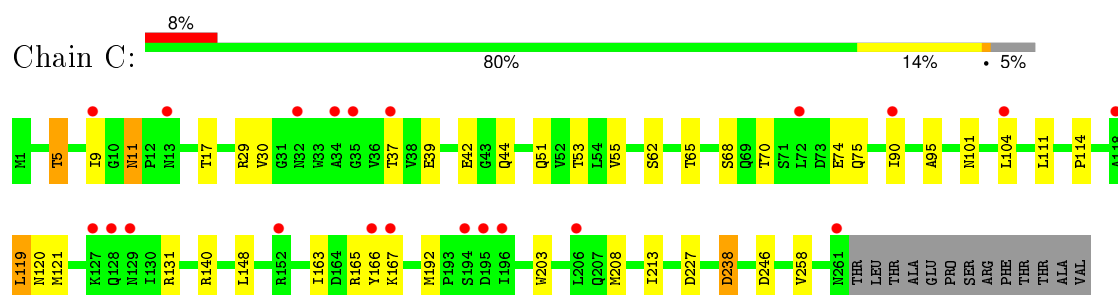
• Molecule 1: Ferrous iron transport protein B



• 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 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.40Å 56.28Å 90.62Å 90.00° 91.86° 90.00°	Depositor
Resolution (Å)	23.82 – 1.80 23.71 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (23.82-1.80) 99.9 (23.71-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.3.0037	Depositor
R, R_{free}	0.177 , 0.217 0.182 , 0.223	Depositor DCC
R_{free} test set	3513 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 51.1	EDS
Estimated twinning fraction	0.023 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 69594 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6815	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.10% 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	0.75	0/2144	0.80	0/2910
1	B	0.80	0/2125	0.87	5/2882 (0.2%)
1	C	0.79	1/2103 (0.0%)	0.82	1/2856 (0.0%)
All	All	0.78	1/6372 (0.0%)	0.83	6/8648 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	11	ASN	CB-CG	5.01	1.62	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	232	ARG	NE-CZ-NH1	5.77	123.19	120.30
1	B	111	LEU	CB-CG-CD1	5.73	120.74	111.00
1	B	73	ASP	CB-CG-OD1	5.36	123.13	118.30
1	B	119	LEU	CA-CB-CG	5.14	127.12	115.30
1	B	123	ASP	CB-CG-OD1	5.08	122.87	118.30
1	C	246	ASP	CB-CG-OD1	5.06	122.86	118.30

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	2084	0	2158	37	0
1	B	2070	0	2137	31	0
1	C	2057	0	2101	35	0
2	A	219	0	0	6	0
2	B	203	0	0	6	0
2	C	182	0	0	3	0
All	All	6815	0	6396	102	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 8.

All (102) 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:65:THR:HG21	1:B:70:THR:HG23	1.44	0.99
1:B:64:THR:HG21	1:B:207[A]:GLN:NE2	1.79	0.97
1:A:78:CYS:O	1:A:82[A]:LEU:HD13	1.67	0.93
1:B:44:GLN:OE1	1:B:53[B]:THR:HG22	1.69	0.93
1:A:106:LEU:HD22	1:A:252[A]:ILE:HD11	1.54	0.88
1:A:106:LEU:HD22	1:A:252[A]:ILE:CD1	2.08	0.83
1:C:140[A]:ARG:HH21	1:C:258:VAL:HG11	1.48	0.79
1:B:65:THR:CG2	1:B:70:THR:HG23	2.16	0.74
1:B:196:ILE:O	1:B:201:ARG:HD2	1.88	0.74
1:B:242:LEU:HD23	1:B:242:LEU:C	2.12	0.70
1:C:44:GLN:OE1	1:C:51:GLN:NE2	2.29	0.66
1:C:238:ASP:OD1	1:C:238:ASP:N	2.30	0.64
1:A:129:ASN:HD22	1:A:263:LEU:HB3	1.62	0.64
1:B:42:GLU:OE2	1:B:53[A]:THR:HG21	1.97	0.64
1:B:64:THR:HG21	1:B:207[A]:GLN:HE21	1.63	0.63
1:A:129:ASN:ND2	1:A:263:LEU:HB3	2.15	0.61
1:B:65:THR:CG2	1:B:70:THR:CG2	2.80	0.60
1:A:171:ASN:HB3	2:A:415:HOH:O	2.01	0.60
1:C:42:GLU:HG2	1:C:55:VAL:HG22	1.85	0.59
1:B:140[B]:ARG:CZ	1:B:258:VAL:HG11	2.33	0.58
1:C:5[A]:THR:HG23	2:C:330:HOH:O	2.02	0.58
1:B:64:THR:CG2	1:B:207[A]:GLN:NE2	2.63	0.58
1:C:192:MET:HE1	1:C:208:MET:CE	2.34	0.58
1:A:1:MET:HE2	1:A:3:LYS:HD3	1.85	0.57
1:A:246:ASP:O	1:A:250[B]:GLN:HG2	2.05	0.57
1:A:30:VAL:CG1	1:A:39:GLU:HG3	2.35	0.57
1:B:140[B]:ARG:NH2	1:B:258:VAL:HG11	2.20	0.56
1:C:44:GLN:NE2	1:C:53:THR:OG1	2.32	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:244:ILE:O	1:B:248[A]:ARG:HG3	2.05	0.56
1:A:99:GLU:HG3	1:A:252[A]:ILE:CG2	2.35	0.56
1:A:173:GLU:HG2	1:C:165:ARG:NE	2.21	0.56
1:B:44:GLN:OE1	1:B:53[B]:THR:CG2	2.49	0.55
1:A:30:VAL:HG13	1:A:39:GLU:HG3	1.87	0.55
1:B:42:GLU:OE2	1:B:53[A]:THR:CG2	2.55	0.55
1:A:30:VAL:HG22	1:A:41:LYS:HG2	1.88	0.55
1:B:248[C]:ARG:HD3	2:B:340:HOH:O	2.08	0.54
1:A:186:ASP:O	1:A:190:LYS:HG3	2.07	0.54
1:A:130:ILE:HD11	1:A:132:ILE:HD11	1.89	0.54
1:B:248[A]:ARG:HD2	2:B:411:HOH:O	2.08	0.54
1:B:9:ILE:HD12	2:B:317:HOH:O	2.08	0.54
1:A:172:VAL:HG23	1:A:174:LEU:HG	1.90	0.53
1:C:44:GLN:HB3	1:C:51:GLN:HE22	1.73	0.52
1:C:5[A]:THR:HG22	2:C:339:HOH:O	2.09	0.52
1:B:184:GLU:HG3	1:B:233:LEU:HG	1.92	0.52
1:C:120[B]:ASN:HB3	1:C:148:LEU:HD12	1.92	0.52
1:A:106:LEU:CD2	1:A:252[A]:ILE:CD1	2.84	0.52
1:C:213:ILE:H	1:C:213:ILE:HD12	1.74	0.52
1:C:30:VAL:CG1	1:C:39:GLU:HG3	2.41	0.51
1:C:114:PRO:HA	1:C:166[A]:TYR:OH	2.11	0.51
1:A:3:LYS:NZ	1:A:44:GLN:OE1	2.44	0.51
1:A:199:LYS:NZ	2:A:451:HOH:O	2.38	0.51
1:A:19:LEU:HD22	1:A:148:LEU:HD11	1.94	0.50
1:C:192:MET:CE	1:C:208:MET:HE1	2.42	0.49
1:C:70:THR:HG21	1:C:75:GLN:HE21	1.77	0.49
1:B:167[B]:LYS:NZ	2:B:387:HOH:O	2.34	0.49
1:A:78:CYS:O	1:A:82[A]:LEU:CD1	2.50	0.48
1:C:140[A]:ARG:NH2	1:C:258:VAL:HG11	2.23	0.48
1:C:192:MET:HE1	1:C:208:MET:HE2	1.95	0.48
1:A:106:LEU:HD22	1:A:252[B]:ILE:CG1	2.43	0.48
1:A:130:ILE:HD12	2:A:341:HOH:O	2.13	0.48
1:C:111:LEU:HD21	1:C:203:TRP:CE3	2.49	0.48
1:C:192:MET:HE1	1:C:208:MET:HE1	1.95	0.47
1:C:95:ALA:HB2	1:C:119:LEU:HD22	1.97	0.47
1:A:99:GLU:HG3	1:A:252[A]:ILE:HG22	1.96	0.47
1:C:192:MET:CE	1:C:208:MET:CE	2.93	0.47
1:B:120:ASN:HB3	1:B:148:LEU:HD12	1.97	0.47
1:A:65:THR:CG2	1:A:75:GLN:HE21	2.27	0.46
1:C:11:ASN:HD21	1:C:101:ASN:HA	1.80	0.46
1:C:70:THR:CG2	1:C:75:GLN:HE21	2.28	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5[A]:THR:CG2	2:C:330:HOH:O	2.64	0.45
1:C:166[A]:TYR:CG	1:C:167:LYS:N	2.83	0.45
1:B:81:ILE:HD12	1:B:89:LEU:HD11	1.98	0.45
1:A:207:GLN:HG2	2:A:441:HOH:O	2.15	0.45
1:B:119:LEU:HD12	1:B:134[B]:ILE:HD12	1.98	0.44
1:A:219:ALA:HB1	1:A:222:ALA:HB3	2.00	0.44
1:B:65:THR:HG21	1:B:70:THR:CG2	2.28	0.44
1:C:37:THR:HG22	1:C:37:THR:O	2.17	0.44
1:B:120:ASN:OD1	1:B:121[B]:MET:HG2	2.16	0.44
1:B:39:GLU:OE1	2:B:397:HOH:O	2.21	0.44
1:C:9:ILE:HG12	1:C:104:LEU:HD21	1.99	0.44
1:C:213:ILE:HD12	1:C:213:ILE:N	2.33	0.44
1:B:22:GLN:HB3	1:B:156:ILE:HD11	2.00	0.44
1:B:37:THR:O	2:B:301:HOH:O	2.21	0.44
1:C:65:THR:HG21	1:C:68:SER:OG	2.17	0.44
1:C:62:SER:O	1:C:74:GLU:HG2	2.18	0.43
1:C:114:PRO:HB3	1:C:166[A]:TYR:CE2	2.53	0.43
1:A:106:LEU:CD2	1:A:252[A]:ILE:HD13	2.47	0.43
1:A:106:LEU:CD2	1:A:252[A]:ILE:HD11	2.37	0.43
1:A:167:LYS:HE3	2:A:340:HOH:O	2.18	0.43
1:B:1:MET:O	1:B:50:HIS:CD2	2.72	0.43
1:A:131:ARG:HB2	1:A:263:LEU:HD13	2.01	0.42
1:C:90:ILE:HD11	1:C:163:ILE:HD11	2.01	0.42
1:B:81:ILE:HD12	1:B:89:LEU:CD1	2.48	0.42
1:A:77:ALA:O	1:A:81[A]:ILE:HG12	2.19	0.42
1:A:106:LEU:HD22	1:A:252[A]:ILE:HD13	1.96	0.42
1:B:22:GLN:HB3	1:B:156:ILE:CD1	2.51	0.41
1:A:91:ASN:HD22	1:A:105:THR:HG23	1.84	0.41
1:A:260:SER:C	1:A:261:ASN:HD22	2.24	0.41
1:C:166[A]:TYR:CD2	1:C:167:LYS:N	2.89	0.41
1:C:30:VAL:CG1	1:C:39:GLU:CG	2.99	0.41
1:A:16[B]:LYS:HE3	2:A:563:HOH:O	2.21	0.40
1:A:261:ASN:N	1:A:261:ASN:HD22	2.19	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	270/274 (98%)	266 (98%)	4 (2%)	0	100	100
1	B	269/274 (98%)	262 (97%)	7 (3%)	0	100	100
1	C	267/274 (97%)	260 (97%)	7 (3%)	0	100	100
All	All	806/822 (98%)	788 (98%)	18 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	233/231 (101%)	224 (96%)	9 (4%)	39	21
1	B	230/231 (100%)	222 (96%)	8 (4%)	43	25
1	C	228/231 (99%)	219 (96%)	9 (4%)	39	21
All	All	691/693 (100%)	665 (96%)	26 (4%)	40	22

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	28	GLN
1	A	70	THR
1	A	129	ASN
1	A	130	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	135	ASP
1	A	223	SER
1	A	261	ASN
1	A	264	THR
1	B	4	LEU
1	B	88	LEU
1	B	106	LEU
1	B	111	LEU
1	B	195	ASP
1	B	198	LEU
1	B	238	ASP
1	B	261	ASN
1	C	5[A]	THR
1	C	5[B]	THR
1	C	17	THR
1	C	29	ARG
1	C	119	LEU
1	C	121	MET
1	C	131	ARG
1	C	227	ASP
1	C	238	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	75	GLN
1	A	91	ASN
1	A	129	ASN
1	A	243	HIS
1	A	261	ASN
1	B	50	HIS
1	B	51	GLN
1	B	91	ASN
1	B	179	GLN
1	C	11	ASN
1	C	44	GLN
1	C	51	GLN
1	C	75	GLN
1	C	224	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	260/274 (94%)	0.65	21 (8%)	15 11	19, 26, 34, 40	0
1	B	261/274 (95%)	0.70	25 (9%)	10 7	14, 26, 36, 44	0
1	C	261/274 (95%)	0.69	21 (8%)	15 12	17, 26, 36, 40	0
All	All	782/822 (95%)	0.68	67 (8%)	13 10	14, 26, 36, 44	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	70	THR	5.9
1	B	152	ARG	5.6
1	A	264	THR	5.4
1	B	1	MET	5.2
1	C	194	SER	5.2
1	C	34	ALA	5.0
1	C	37	THR	5.0
1	B	238	ASP	4.9
1	A	263	LEU	4.8
1	C	195	ASP	4.6
1	C	129	ASN	4.4
1	B	151	THR	3.8
1	B	195	ASP	3.6
1	B	242	LEU	3.5
1	C	35	GLY	3.4
1	B	261	ASN	3.3
1	B	249	TYR	3.2
1	B	239	ASP	3.2
1	A	204	LEU	3.2
1	C	90	ILE	3.1
1	C	261	ASN	3.0
1	A	150	SER	3.0
1	B	67	SER	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	127	LYS	2.9
1	A	6	ILE	2.9
1	C	196	ILE	2.9
1	A	206	LEU	2.7
1	A	195	ASP	2.6
1	B	26	SER	2.6
1	C	118	ALA	2.6
1	A	262	THR	2.6
1	C	13	ASN	2.5
1	A	8	LEU	2.5
1	A	152	ARG	2.5
1	B	213	ILE	2.5
1	A	129	ASN	2.5
1	A	173	GLU	2.4
1	B	131	ARG	2.4
1	B	61	TYR	2.4
1	C	166[A]	TYR	2.4
1	B	150	SER	2.4
1	A	90	ILE	2.3
1	A	65	THR	2.3
1	A	203	TRP	2.2
1	B	220	GLY	2.2
1	A	224	GLN	2.2
1	C	167	LYS	2.2
1	C	72	LEU	2.2
1	C	206	LEU	2.2
1	B	70	THR	2.2
1	C	127	LYS	2.2
1	B	103[A]	TYR	2.2
1	A	239	ASP	2.2
1	C	152	ARG	2.2
1	C	128	GLN	2.2
1	A	209	LEU	2.2
1	B	8	LEU	2.1
1	C	104	LEU	2.1
1	B	237	MET	2.1
1	B	55	VAL	2.1
1	B	6	ILE	2.0
1	A	19	LEU	2.0
1	C	32	ASN	2.0
1	C	9	ILE	2.0
1	B	206	LEU	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	191	VAL	2.0
1	B	90	ILE	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.