



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:00 PM GMT

PDB ID : 4B7O
Title : THE FrpB IRON TRANSPORTER FROM NEISSERIA MENINGITIDIS
(F5-1 VARIANT) APOPROTEIN FORM
Authors : Saleem, M.; Prince, S.M.; Derrick, J.P.
Deposited on : 2012-08-21
Resolution : 2.32 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

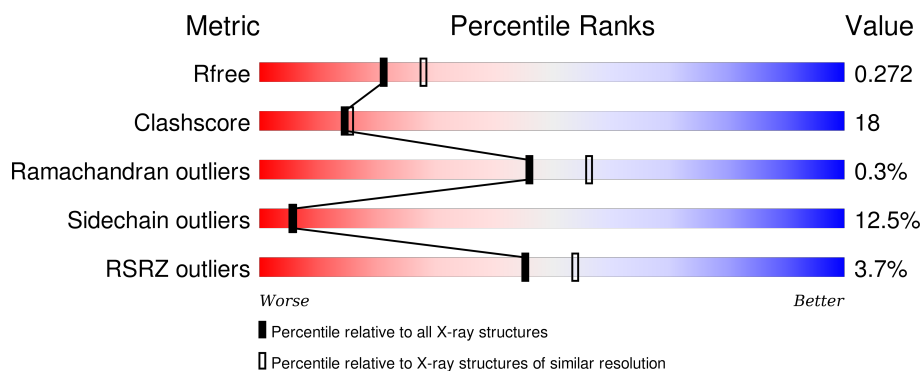
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.32 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	4425 (2.34-2.30)
Clashscore	102246	5057 (2.34-2.30)
Ramachandran outliers	100387	5008 (2.34-2.30)
Sidechain outliers	100360	5007 (2.34-2.30)
RSRZ outliers	91569	4432 (2.34-2.30)

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	745	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5182 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 IRON-REGULATED OUTER MEMBRANE PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	649	Total	C	N	O	S	0	0	0
			5103	3197	922	979	5			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP Q841A2
A	2	HIS	-	EXPRESSION TAG	UNP Q841A2
A	3	HIS	-	EXPRESSION TAG	UNP Q841A2
A	4	HIS	-	EXPRESSION TAG	UNP Q841A2
A	5	HIS	-	EXPRESSION TAG	UNP Q841A2
A	6	HIS	-	EXPRESSION TAG	UNP Q841A2
A	7	HIS	-	EXPRESSION TAG	UNP Q841A2
A	8	SER	-	EXPRESSION TAG	UNP Q841A2
A	9	SER	-	EXPRESSION TAG	UNP Q841A2
A	10	GLY	-	EXPRESSION TAG	UNP Q841A2
A	11	LEU	-	EXPRESSION TAG	UNP Q841A2
A	12	VAL	-	EXPRESSION TAG	UNP Q841A2
A	13	PRO	-	EXPRESSION TAG	UNP Q841A2
A	14	ARG	-	EXPRESSION TAG	UNP Q841A2
A	15	GLY	-	EXPRESSION TAG	UNP Q841A2
A	16	SER	-	EXPRESSION TAG	UNP Q841A2
A	17	GLY	-	EXPRESSION TAG	UNP Q841A2
A	18	MET	-	EXPRESSION TAG	UNP Q841A2
A	19	LYS	-	EXPRESSION TAG	UNP Q841A2
A	20	GLU	-	EXPRESSION TAG	UNP Q841A2
A	21	THR	-	EXPRESSION TAG	UNP Q841A2
A	22	ALA	-	EXPRESSION TAG	UNP Q841A2
A	23	ALA	-	EXPRESSION TAG	UNP Q841A2
A	24	ALA	-	EXPRESSION TAG	UNP Q841A2
A	25	LYS	-	EXPRESSION TAG	UNP Q841A2
A	26	PHE	-	EXPRESSION TAG	UNP Q841A2
A	27	GLU	-	EXPRESSION TAG	UNP Q841A2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ARG	-	EXPRESSION TAG	UNP Q841A2
A	29	GLN	-	EXPRESSION TAG	UNP Q841A2
A	30	HIS	-	EXPRESSION TAG	UNP Q841A2
A	31	MET	-	EXPRESSION TAG	UNP Q841A2
A	32	ASP	-	EXPRESSION TAG	UNP Q841A2
A	33	SER	-	EXPRESSION TAG	UNP Q841A2
A	34	PRO	-	EXPRESSION TAG	UNP Q841A2
A	35	ASP	-	EXPRESSION TAG	UNP Q841A2
A	36	LEU	-	EXPRESSION TAG	UNP Q841A2
A	37	GLY	-	EXPRESSION TAG	UNP Q841A2
A	38	THR	-	EXPRESSION TAG	UNP Q841A2
A	39	ASP	-	EXPRESSION TAG	UNP Q841A2
A	40	ASP	-	EXPRESSION TAG	UNP Q841A2
A	41	ASP	-	EXPRESSION TAG	UNP Q841A2
A	42	ASP	-	EXPRESSION TAG	UNP Q841A2
A	43	LYS	-	EXPRESSION TAG	UNP Q841A2
A	44	MET	-	EXPRESSION TAG	UNP Q841A2
A	45	ALA	-	EXPRESSION TAG	UNP Q841A2
A	46	GLU	-	EXPRESSION TAG	UNP Q841A2
A	47	ASN	-	EXPRESSION TAG	UNP Q841A2
A	48	ASN	-	EXPRESSION TAG	UNP Q841A2
A	49	ALA	-	EXPRESSION TAG	UNP Q841A2
A	50	LYS	-	EXPRESSION TAG	UNP Q841A2
A	51	VAL	-	EXPRESSION TAG	UNP Q841A2
A	724	GLN	-	EXPRESSION TAG	UNP Q841A2
A	725	ARG	-	EXPRESSION TAG	UNP Q841A2
A	726	TRP	-	EXPRESSION TAG	UNP Q841A2
A	727	THR	-	EXPRESSION TAG	UNP Q841A2
A	728	ASN	-	EXPRESSION TAG	UNP Q841A2
A	729	THR	-	EXPRESSION TAG	UNP Q841A2
A	730	LEU	-	EXPRESSION TAG	UNP Q841A2
A	731	PRO	-	EXPRESSION TAG	UNP Q841A2
A	732	GLY	-	EXPRESSION TAG	UNP Q841A2
A	733	VAL	-	EXPRESSION TAG	UNP Q841A2
A	734	GLY	-	EXPRESSION TAG	UNP Q841A2
A	735	ARG	-	EXPRESSION TAG	UNP Q841A2
A	736	ASP	-	EXPRESSION TAG	UNP Q841A2
A	737	VAL	-	EXPRESSION TAG	UNP Q841A2
A	738	ARG	-	EXPRESSION TAG	UNP Q841A2
A	739	LEU	-	EXPRESSION TAG	UNP Q841A2
A	740	GLY	-	EXPRESSION TAG	UNP Q841A2
A	741	VAL	-	EXPRESSION TAG	UNP Q841A2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	742	ASN	-	EXPRESSION TAG	UNP Q841A2
A	743	TYR	-	EXPRESSION TAG	UNP Q841A2
A	744	LYS	-	EXPRESSION TAG	UNP Q841A2
A	745	PHE	-	EXPRESSION TAG	UNP Q841A2
A	272	GLU	LYS	SEE REMARK 999	UNP Q841A2
A	302	PRO	SER	SEE REMARK 999	UNP Q841A2
A	638	ALA	THR	SEE REMARK 999	UNP Q841A2

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	79	Total O 79 79	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	175.21 Å 79.10 Å 74.75 Å 90.00° 97.58° 90.00°	Depositor
Resolution (Å)	37.05 – 2.32 37.05 – 2.32	Depositor EDS
% Data completeness (in resolution range)	98.0 (37.05-2.32) 98.0 (37.05-2.32)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.31 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.233 , 0.281 0.228 , 0.272	Depositor DCC
R_{free} test set	2172 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 33.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 43114 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5182	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 4.50% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.91	3/5210 (0.1%)	0.95	7/7027 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	697	TRP	CD2-CE2	7.64	1.50	1.41
1	A	185	TRP	CD2-CE2	6.30	1.49	1.41
1	A	92	GLU	CD-OE2	5.82	1.32	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	685	ARG	NE-CZ-NH1	6.77	123.69	120.30
1	A	385	ARG	NE-CZ-NH2	-6.20	117.20	120.30
1	A	685	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	A	697	TRP	CA-CB-CG	5.89	124.89	113.70
1	A	646	ARG	NE-CZ-NH2	-5.50	117.55	120.30
1	A	337	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	A	335	LYS	CD-CE-NZ	5.10	123.43	111.70

There are no chirality outliers.

There are no planarity outliers.

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	5103	0	4945	180	0
2	A	79	0	0	22	0
All	All	5182	0	4945	180	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 18.

All (180) 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:615:ARG:HH12	1:A:634:GLN:NE2	1.38	1.19
1:A:316:THR:HB	2:A:2041:HOH:O	1.40	1.18
1:A:427:LYS:HB2	2:A:2052:HOH:O	1.49	1.10
1:A:265:ARG:HB3	2:A:2041:HOH:O	1.56	1.06
1:A:81:THR:HG21	1:A:710:SER:OG	1.55	1.05
1:A:211:ASN:HD22	1:A:211:ASN:N	1.56	1.03
1:A:419:ASP:HB3	1:A:422:LYS:HE2	1.41	1.02
1:A:615:ARG:HH12	1:A:634:GLN:HE21	1.06	1.00
1:A:656:TRP:HE1	1:A:691:ASN:HD22	1.07	0.98
1:A:211:ASN:H	1:A:211:ASN:ND2	1.61	0.97
1:A:706:ASN:HD22	1:A:742:ASN:HD21	1.05	0.96
1:A:582:VAL:HG12	2:A:2059:HOH:O	1.71	0.89
1:A:71:ILE:CD1	1:A:149:VAL:HG23	2.03	0.88
1:A:599:ARG:HD2	2:A:2063:HOH:O	1.72	0.87
1:A:265:ARG:HH11	1:A:265:ARG:HG3	1.38	0.87
1:A:615:ARG:NH1	1:A:634:GLN:NE2	2.22	0.86
1:A:603:LEU:HD11	1:A:643:LEU:HD11	1.56	0.86
1:A:242:PRO:HD2	1:A:286:THR:HG21	1.57	0.85
1:A:145:LYS:HD3	1:A:176:ASP:OD2	1.76	0.85
1:A:615:ARG:NH1	1:A:634:GLN:HE21	1.74	0.83
1:A:472:LYS:HE2	1:A:473:THR:O	1.79	0.82
1:A:211:ASN:ND2	1:A:211:ASN:N	2.17	0.81
1:A:287:VAL:C	2:A:2044:HOH:O	2.20	0.79
1:A:478:THR:O	1:A:479:VAL:HG23	1.83	0.78
1:A:150:GLN:O	1:A:165:GLY:HA3	1.84	0.77
1:A:99:GLY:H	1:A:724:GLN:HE22	1.32	0.76
1:A:656:TRP:HE1	1:A:691:ASN:ND2	1.81	0.76
1:A:211:ASN:H	1:A:211:ASN:HD22	0.82	0.74
1:A:99:GLY:H	1:A:724:GLN:NE2	1.86	0.73
1:A:341:ASP:OD2	1:A:343:LYS:HE3	1.89	0.73
1:A:373:ALA:HB3	1:A:376:THR:CG2	2.18	0.72
1:A:71:ILE:HD12	1:A:149:VAL:HG23	1.71	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:LEU:HD23	1:A:490:VAL:HG12	1.71	0.72
1:A:570:GLN:CD	1:A:582:VAL:HG11	2.10	0.71
1:A:242:PRO:CD	1:A:286:THR:HG21	2.19	0.71
1:A:440:PRO:HG3	1:A:474:HIS:HB3	1.73	0.70
1:A:373:ALA:HB3	1:A:376:THR:HG23	1.73	0.69
1:A:518:GLN:NE2	2:A:2057:HOH:O	1.61	0.69
1:A:81:THR:HG21	1:A:710:SER:HG	1.54	0.69
1:A:116:SER:HB2	1:A:164:ASN:HD22	1.58	0.69
1:A:691:ASN:O	1:A:714:VAL:HG23	1.93	0.69
1:A:208:LYS:NZ	2:A:2034:HOH:O	2.26	0.68
1:A:242:PRO:O	1:A:286:THR:CG2	2.43	0.67
1:A:472:LYS:CE	1:A:473:THR:O	2.43	0.66
1:A:599:ARG:CD	2:A:2063:HOH:O	2.35	0.65
1:A:603:LEU:HD11	1:A:643:LEU:CD1	2.25	0.65
1:A:535:GLU:HG2	1:A:565:ALA:HA	1.78	0.65
1:A:242:PRO:O	1:A:286:THR:HG23	1.96	0.65
1:A:255:ILE:CD1	1:A:268:LEU:HB3	2.27	0.64
1:A:265:ARG:HH11	1:A:265:ARG:CG	2.10	0.64
1:A:234:ASN:HB2	1:A:238:GLY:N	2.13	0.64
1:A:255:ILE:HD12	1:A:268:LEU:HB3	1.80	0.63
1:A:569:PRO:HB2	1:A:579:ARG:HB3	1.80	0.62
1:A:421:GLU:C	1:A:421:GLU:OE1	2.37	0.62
1:A:63:GLN:NE2	1:A:66:LYS:HD3	2.14	0.62
1:A:472:LYS:HE3	1:A:476:GLY:HA2	1.80	0.62
1:A:461:THR:HG22	2:A:2053:HOH:O	2.00	0.62
1:A:145:LYS:HB3	1:A:172:VAL:HA	1.83	0.61
1:A:369:ASP:OD1	1:A:379:LYS:HG2	2.01	0.60
1:A:177:LEU:HD12	1:A:215:LEU:HD12	1.83	0.59
1:A:478:THR:O	1:A:479:VAL:CG2	2.50	0.59
1:A:478:THR:HG22	1:A:479:VAL:N	2.17	0.59
1:A:529:ALA:HA	1:A:580:GLU:OE1	2.03	0.58
1:A:247:ASP:OD2	1:A:249:ARG:NH2	2.37	0.58
1:A:101:ASN:O	1:A:104:SER:HB2	2.04	0.58
1:A:529:ALA:O	1:A:532:THR:HG23	2.03	0.58
1:A:116:SER:CB	1:A:164:ASN:HD22	2.16	0.57
1:A:264:HIS:HD2	1:A:315:TYR:OH	1.87	0.57
1:A:372:LEU:HD21	1:A:378:LEU:HD22	1.86	0.57
1:A:650:PRO:HB2	1:A:697:TRP:CZ3	2.39	0.57
1:A:265:ARG:NH1	1:A:265:ARG:HG3	2.17	0.57
1:A:430:ASP:O	1:A:434:SER:OG	2.23	0.56
1:A:648:GLN:HG3	2:A:2070:HOH:O	2.05	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:ASP:HB3	1:A:706:ASN:ND2	2.21	0.56
1:A:177:LEU:HD12	1:A:215:LEU:CD1	2.35	0.56
1:A:535:GLU:HG3	1:A:566:LEU:HD12	1.88	0.56
1:A:272:GLU:CD	2:A:2040:HOH:O	2.45	0.55
1:A:115:ASN:HB2	1:A:630:GLU:HG2	1.88	0.55
1:A:234:ASN:HD22	1:A:238:GLY:HA3	1.71	0.55
1:A:71:ILE:HD11	1:A:149:VAL:HG23	1.86	0.55
1:A:65:SER:CB	1:A:593:GLU:HG3	2.37	0.55
1:A:706:ASN:HD22	1:A:742:ASN:ND2	1.88	0.54
1:A:276:ARG:HB3	1:A:305:GLU:HG2	1.88	0.54
1:A:61:ASP:OD1	1:A:69:THR:HG23	2.07	0.54
1:A:562:ILE:HB	1:A:587:ILE:HB	1.90	0.54
1:A:324:GLU:HB2	2:A:2046:HOH:O	2.08	0.53
1:A:421:GLU:OE1	1:A:421:GLU:O	2.26	0.53
1:A:150:GLN:O	1:A:165:GLY:CA	2.56	0.53
1:A:328:ALA:HA	1:A:365:ASN:O	2.09	0.53
1:A:428:ASP:O	1:A:431:LEU:N	2.41	0.53
1:A:622:LYS:O	1:A:624:LEU:HG	2.10	0.52
1:A:264:HIS:CD2	1:A:315:TYR:OH	2.62	0.52
1:A:265:ARG:NH1	1:A:265:ARG:CG	2.71	0.52
1:A:65:SER:HB3	1:A:593:GLU:HG3	1.91	0.52
1:A:526:ILE:HD11	1:A:579:ARG:HH21	1.75	0.51
1:A:435:TYR:CD2	1:A:525:ILE:HG13	2.46	0.51
1:A:478:THR:CG2	1:A:479:VAL:N	2.73	0.51
1:A:81:THR:HG23	1:A:738:ARG:HD3	1.93	0.51
1:A:615:ARG:HH12	1:A:634:GLN:HE22	1.45	0.51
1:A:278:ILE:HD13	1:A:278:ILE:N	2.26	0.51
1:A:598:TYR:CE2	1:A:600:THR:HG22	2.46	0.50
1:A:282:ARG:NH1	2:A:2043:HOH:O	2.31	0.50
1:A:448:TYR:HB3	1:A:465:ARG:HG2	1.93	0.50
1:A:514:TYR:CD2	1:A:526:ILE:HG13	2.47	0.50
1:A:153:ALA:HB3	1:A:539:ASN:HB3	1.93	0.50
1:A:685:ARG:NH2	1:A:724:GLN:O	2.44	0.50
1:A:388:GLU:HG3	1:A:441:THR:HG22	1.94	0.50
1:A:302:PRO:HD2	2:A:2045:HOH:O	2.11	0.49
1:A:274:GLN:OE1	1:A:276:ARG:NH1	2.45	0.49
1:A:322:PHE:O	1:A:370:SER:HB3	2.11	0.49
1:A:529:ALA:HB3	1:A:532:THR:CG2	2.42	0.49
1:A:258:THR:HG23	1:A:264:HIS:O	2.12	0.49
1:A:506:TYR:OH	1:A:538:ARG:HD3	2.12	0.49
1:A:67:ILE:HG12	1:A:150:GLN:HG3	1.95	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:653:GLU:O	1:A:695:ALA:HA	2.13	0.48
1:A:569:PRO:HA	1:A:580:GLU:O	2.14	0.48
1:A:396:ASN:CG	1:A:437:LEU:HD22	2.34	0.48
1:A:380:TYR:CD1	1:A:380:TYR:N	2.82	0.47
1:A:426:LYS:HZ3	1:A:426:LYS:HB3	1.80	0.47
1:A:675:ASP:OD1	1:A:676:ARG:N	2.48	0.47
1:A:590:HIS:HE1	2:A:2060:HOH:O	1.96	0.47
1:A:143:LEU:O	1:A:172:VAL:HG12	2.15	0.47
1:A:234:ASN:HD22	1:A:238:GLY:CA	2.27	0.46
1:A:517:LEU:O	1:A:517:LEU:HG	2.15	0.46
1:A:548:ASP:C	1:A:548:ASP:OD1	2.52	0.46
1:A:243:TYR:HE2	1:A:287:VAL:HG21	1.80	0.46
1:A:243:TYR:HE2	1:A:287:VAL:CG2	2.29	0.46
1:A:263:ASP:HA	1:A:318:LYS:HD2	1.97	0.46
1:A:615:ARG:CZ	1:A:634:GLN:HE21	2.29	0.46
1:A:423:ALA:HA	1:A:426:LYS:HZ3	1.81	0.46
1:A:421:GLU:HG2	2:A:2051:HOH:O	2.15	0.46
1:A:99:GLY:N	1:A:721:PRO:HG2	2.31	0.46
1:A:697:TRP:CH2	1:A:699:PRO:HA	2.50	0.46
1:A:217:SER:O	1:A:251:TYR:HA	2.17	0.45
1:A:242:PRO:N	1:A:286:THR:HG21	2.31	0.45
1:A:426:LYS:HB3	1:A:426:LYS:NZ	2.31	0.45
1:A:626:SER:HB3	1:A:631:PHE:CE2	2.51	0.45
1:A:423:ALA:HB2	2:A:2051:HOH:O	2.16	0.45
1:A:140:ASP:HA	1:A:141:PRO:HD2	1.86	0.45
1:A:255:ILE:HD11	1:A:268:LEU:HB3	1.98	0.45
1:A:71:ILE:CD1	1:A:149:VAL:CG2	2.87	0.45
1:A:439:ASN:HA	1:A:440:PRO:HD3	1.84	0.45
1:A:242:PRO:HB2	1:A:286:THR:HG22	1.99	0.44
1:A:590:HIS:CE1	2:A:2060:HOH:O	2.70	0.44
1:A:396:ASN:N	1:A:396:ASN:OD1	2.47	0.44
1:A:674:LYS:HB3	1:A:674:LYS:NZ	2.32	0.44
1:A:349:GLY:HA2	1:A:522:LYS:HD3	2.00	0.44
1:A:225:GLU:HG2	1:A:244:SER:H	1.82	0.44
1:A:423:ALA:CB	2:A:2051:HOH:O	2.65	0.44
1:A:615:ARG:HD3	1:A:615:ARG:HA	1.81	0.44
1:A:225:GLU:OE2	1:A:243:TYR:N	2.35	0.43
1:A:369:ASP:HB3	1:A:377:LEU:HD11	2.00	0.43
1:A:97:PHE:CE2	1:A:107:LEU:HD12	2.53	0.43
1:A:143:LEU:HD22	1:A:252:LEU:HB2	2.00	0.43
1:A:327:ASP:O	1:A:327:ASP:OD1	2.36	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:THR:CG2	1:A:738:ARG:HD3	2.49	0.43
1:A:557:TYR:CD2	1:A:592:TYR:HB3	2.52	0.43
1:A:114:GLN:HB3	1:A:136:ARG:HD2	2.01	0.43
1:A:233:ARG:HA	1:A:233:ARG:HD3	1.78	0.42
1:A:211:ASN:CB	1:A:258:THR:O	2.68	0.42
1:A:446:GLY:HA2	1:A:466:TYR:O	2.19	0.42
1:A:139:VAL:HG22	1:A:140:ASP:N	2.35	0.42
1:A:255:ILE:HD11	1:A:268:LEU:HD23	2.00	0.42
1:A:652:LEU:HD12	1:A:697:TRP:HB2	2.01	0.42
1:A:442:LYS:HE3	1:A:444:ASP:OD2	2.19	0.42
1:A:224:LYS:HD3	1:A:224:LYS:N	2.33	0.42
1:A:251:TYR:HE2	2:A:2040:HOH:O	2.01	0.42
1:A:342:ASP:OD2	1:A:355:ASN:HB3	2.20	0.41
1:A:423:ALA:HA	1:A:426:LYS:NZ	2.35	0.41
1:A:107:LEU:C	1:A:107:LEU:HD23	2.39	0.41
1:A:376:THR:HA	1:A:452:ILE:O	2.19	0.41
1:A:435:TYR:HB3	1:A:437:LEU:HD13	2.03	0.41
1:A:316:THR:CB	2:A:2041:HOH:O	2.24	0.41
1:A:502:SER:HA	1:A:541:GLU:O	2.20	0.41
1:A:569:PRO:HB3	1:A:580:GLU:C	2.41	0.41
1:A:263:ASP:HA	1:A:318:LYS:CD	2.51	0.41
1:A:579:ARG:HB2	1:A:579:ARG:HE	1.55	0.41
1:A:526:ILE:CD1	1:A:579:ARG:HH21	2.35	0.40
1:A:211:ASN:HB3	1:A:258:THR:O	2.22	0.40
1:A:729:THR:OG1	1:A:730:LEU:N	2.54	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	639/745 (86%)	613 (96%)	24 (4%)	2 (0%)	46	56

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	234	ASN
1	A	675	ASP

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	535/615 (87%)	468 (88%)	67 (12%)	6 5

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	THR
1	A	73	THR
1	A	81	THR
1	A	116	SER
1	A	117	VAL
1	A	146	VAL
1	A	150	GLN
1	A	211	ASN
1	A	212	PHE
1	A	215	LEU
1	A	217	SER
1	A	239	LYS
1	A	255	ILE
1	A	261	ASP
1	A	265	ARG
1	A	276	ARG
1	A	286	THR
1	A	318	LYS
1	A	320	LEU
1	A	327	ASP
1	A	333	LEU
1	A	335	LYS
1	A	343	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	357	THR
1	A	364	MET
1	A	370	SER
1	A	372	LEU
1	A	376	THR
1	A	396	ASN
1	A	402	SER
1	A	418	THR
1	A	420	GLU
1	A	421	GLU
1	A	426	LYS
1	A	427	LYS
1	A	430	ASP
1	A	431	LEU
1	A	434	SER
1	A	435	TYR
1	A	437	LEU
1	A	445	THR
1	A	460	LEU
1	A	470	LYS
1	A	472	LYS
1	A	473	THR
1	A	474	HIS
1	A	481	SER
1	A	488	PHE
1	A	503	SER
1	A	532	THR
1	A	533	LYS
1	A	540	THR
1	A	560	GLN
1	A	580	GLU
1	A	582	VAL
1	A	593	GLU
1	A	600	THR
1	A	607	VAL
1	A	611	HIS
1	A	648	GLN
1	A	674	LYS
1	A	676	ARG
1	A	679	LYS
1	A	697	TRP
1	A	702	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	737	VAL
1	A	742	ASN

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	63	GLN
1	A	164	ASN
1	A	211	ASN
1	A	264	HIS
1	A	387	GLN
1	A	560	GLN
1	A	611	HIS
1	A	634	GLN
1	A	691	ASN
1	A	724	GLN
1	A	728	ASN
1	A	742	ASN

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, 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	649/745 (87%)	0.26	24 (3%) 45 54	29, 43, 61, 88	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	680	LEU	4.8
1	A	621	PRO	3.9
1	A	320	LEU	3.8
1	A	322	PHE	3.6
1	A	117	VAL	3.5
1	A	182	ASP	3.2
1	A	119	ILE	3.2
1	A	352	VAL	2.9
1	A	326	LEU	2.9
1	A	679	LYS	2.8
1	A	622	LYS	2.4
1	A	600	THR	2.4
1	A	419	ASP	2.3
1	A	607	VAL	2.3
1	A	674	LYS	2.3
1	A	676	ARG	2.2
1	A	598	TYR	2.2
1	A	420	GLU	2.2
1	A	166	ALA	2.2
1	A	97	PHE	2.1
1	A	138	ILE	2.1
1	A	401	ILE	2.1
1	A	672	GLY	2.1
1	A	343	LYS	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.