



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2016 – 08:14 PM EDT

PDB ID : 5IWO
Title : Bacterial sodium channel pore domain, low bromide
Authors : Shaya, D.; Findeisen, F.; Rohaim, A.; Minor, D.L.
Deposited on : 2016-03-22
Resolution : 3.33 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 : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0122
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107

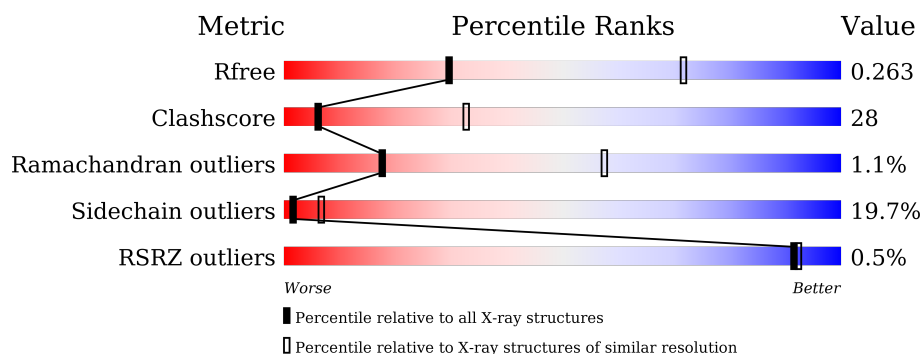
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.33 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	1004 (3.40-3.28)
Clashscore	102246	1072 (3.40-3.28)
Ramachandran outliers	100387	1055 (3.40-3.28)
Sidechain outliers	100360	1054 (3.40-3.28)
RSRZ outliers	91569	1009 (3.40-3.28)

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	152	<div> <div>%</div> <div> <div></div> <div>38%</div> <div>43%</div> <div>11%</div> <div>•</div> <div>9%</div> </div> </div>
1	B	152	<div> <div>37%</div> <div>41%</div> <div>13%</div> <div>•</div> <div>9%</div> </div>
1	C	152	<div> <div>%</div> <div> <div></div> <div>46%</div> <div>32%</div> <div>13%</div> <div>9%</div> </div> </div>
1	D	152	<div> <div>%</div> <div> <div></div> <div>40%</div> <div>43%</div> <div>7%</div> <div>•</div> <div>9%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BR	A	301	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4501 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 Ion transport protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	139	Total	C	N	O	S	0	0	0
			1124	742	181	195	6			
1	B	138	Total	C	N	O	S	0	1	0
			1126	745	180	195	6			
1	C	138	Total	C	N	O	S	0	0	0
			1122	742	178	196	6			
1	D	138	Total	C	N	O	S	0	0	0
			1128	745	181	196	6			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	137	GLY	-	expression tag	UNP Q0ABW0
A	138	PRO	-	expression tag	UNP Q0ABW0
A	139	SER	-	expression tag	UNP Q0ABW0
A	140	SER	-	expression tag	UNP Q0ABW0
A	141	PRO	-	expression tag	UNP Q0ABW0
A	142	SER	-	expression tag	UNP Q0ABW0
B	137	GLY	-	expression tag	UNP Q0ABW0
B	138	PRO	-	expression tag	UNP Q0ABW0
B	139	SER	-	expression tag	UNP Q0ABW0
B	140	SER	-	expression tag	UNP Q0ABW0
B	141	PRO	-	expression tag	UNP Q0ABW0
B	142	SER	-	expression tag	UNP Q0ABW0
C	137	GLY	-	expression tag	UNP Q0ABW0
C	138	PRO	-	expression tag	UNP Q0ABW0
C	139	SER	-	expression tag	UNP Q0ABW0
C	140	SER	-	expression tag	UNP Q0ABW0
C	141	PRO	-	expression tag	UNP Q0ABW0
C	142	SER	-	expression tag	UNP Q0ABW0
D	137	GLY	-	expression tag	UNP Q0ABW0
D	138	PRO	-	expression tag	UNP Q0ABW0
D	139	SER	-	expression tag	UNP Q0ABW0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	140	SER	-	expression tag	UNP Q0ABW0
D	141	PRO	-	expression tag	UNP Q0ABW0
D	142	SER	-	expression tag	UNP Q0ABW0

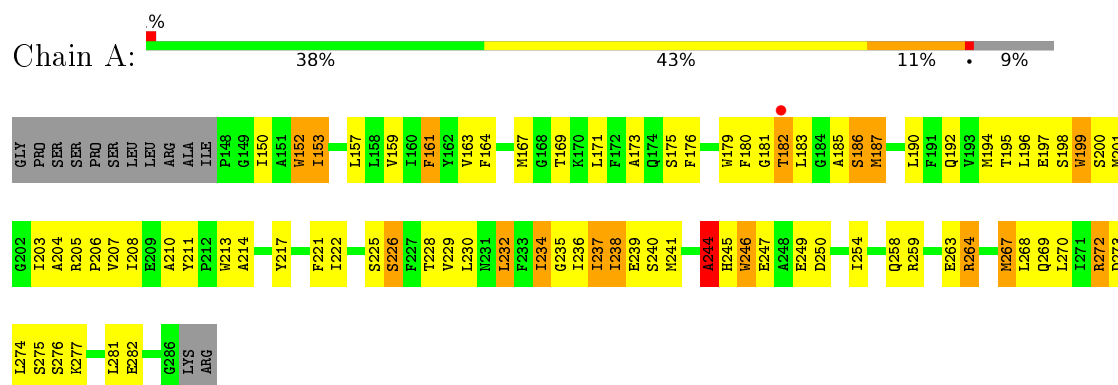
- Molecule 2 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Br 1 1	0	0

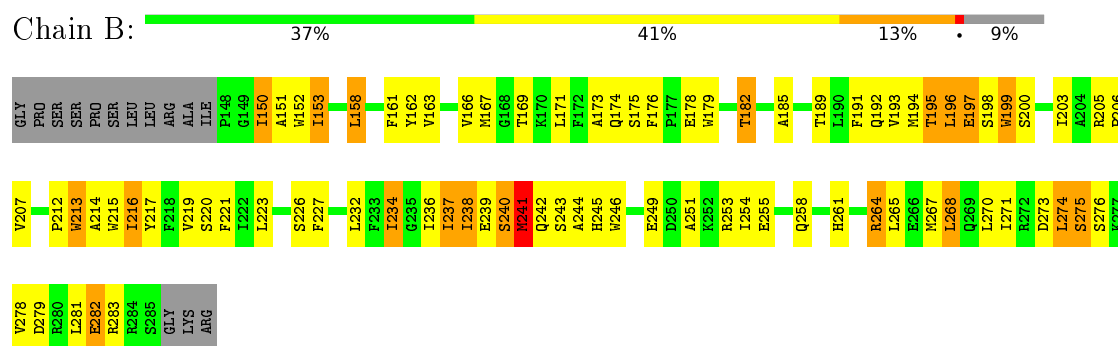
3 Residue-property plots

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.

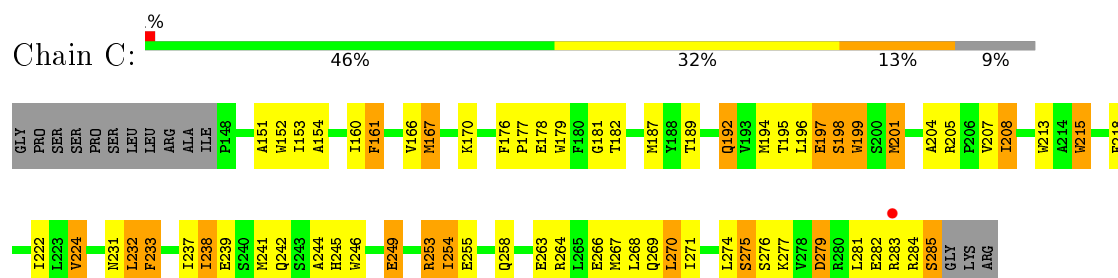
• Molecule 1: Ion transport protein



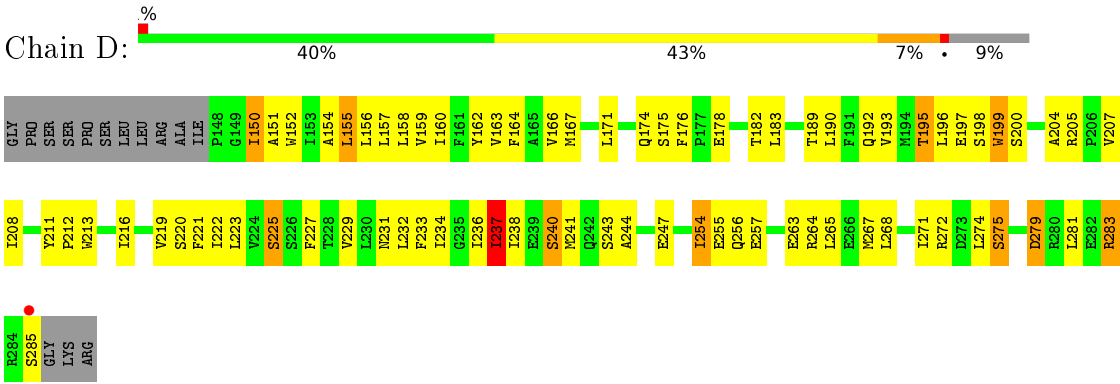
• Molecule 1: Ion transport protein



• Molecule 1: Ion transport protein



• Molecule 1: Ion transport protein



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	151.72Å 159.95Å 168.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.33 14.98 – 3.33	Depositor EDS
% Data completeness (in resolution range)	98.2 (15.00-3.33) 98.4 (14.98-3.33)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.23	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 3.32Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.209 , 0.262 0.209 , 0.263	Depositor DCC
R_{free} test set	1520 reflections (5.43%)	DCC
Wilson B-factor (Å ²)	103.9	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 85.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Outliers	0 of 29495 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4501	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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

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

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	0.84	3/1155 (0.3%)	0.96	4/1569 (0.3%)
1	B	0.75	4/1160 (0.3%)	0.94	4/1575 (0.3%)
1	C	0.81	4/1153 (0.3%)	0.93	2/1566 (0.1%)
1	D	0.81	3/1159 (0.3%)	0.89	1/1573 (0.1%)
All	All	0.80	14/4627 (0.3%)	0.93	11/6283 (0.2%)

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
1	B	0	1
All	All	0	2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	199	TRP	CD2-CE2	6.94	1.49	1.41
1	A	199	TRP	CD2-CE2	6.86	1.49	1.41
1	C	152	TRP	CD2-CE2	6.06	1.48	1.41
1	A	152	TRP	CD2-CE2	5.94	1.48	1.41
1	D	213	TRP	CD2-CE2	5.83	1.48	1.41
1	D	152	TRP	CD2-CE2	5.78	1.48	1.41
1	B	152	TRP	CD2-CE2	5.60	1.48	1.41
1	D	199	TRP	CD2-CE2	5.55	1.48	1.41
1	A	246	TRP	CD2-CE2	5.36	1.47	1.41
1	C	213	TRP	CD2-CE2	5.31	1.47	1.41
1	B	199	TRP	CD2-CE2	5.26	1.47	1.41
1	C	215	TRP	CD2-CE2	5.12	1.47	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	179	TRP	CD2-CE2	5.09	1.47	1.41
1	B	213	TRP	CD2-CE2	5.05	1.47	1.41

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	244	ALA	CB-CA-C	11.00	126.60	110.10
1	B	244	ALA	CB-CA-C	7.52	121.38	110.10
1	D	196	LEU	CB-CG-CD1	-6.48	99.98	111.00
1	B	158	LEU	CB-CG-CD2	-6.22	100.42	111.00
1	B	196	LEU	CB-CG-CD1	-6.15	100.55	111.00
1	C	244	ALA	CB-CA-C	5.68	118.63	110.10
1	A	183	LEU	CA-CB-CG	-5.63	102.34	115.30
1	A	244	ALA	N-CA-C	-5.21	96.92	111.00
1	A	264	ARG	NE-CZ-NH2	5.11	122.86	120.30
1	B	241	MET	CB-CA-C	-5.11	100.18	110.40
1	C	167	MET	CG-SD-CE	5.09	108.35	100.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	244	ALA	Peptide
1	B	241	MET	Peptide

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	1124	0	1111	91	0
1	B	1126	0	1119	71	0
1	C	1122	0	1112	72	0
1	D	1128	0	1123	76	0
2	A	1	0	0	7	0
All	All	4501	0	4465	253	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 28.

All (253) 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:264:ARG:HD3	2:A:301:BR:BR	1.67	1.48
2:A:301:BR:BR	1:C:264:ARG:HD3	1.88	1.27
1:A:204:ALA:O	1:A:208:ILE:HG13	1.49	1.11
1:A:264:ARG:CD	2:A:301:BR:BR	2.56	1.08
1:C:270:LEU:H	1:C:270:LEU:HD23	1.11	1.06
1:A:180:PHE:HD1	1:A:186:SER:HG	1.08	0.97
1:A:268:LEU:HD23	1:D:267:MET:HG2	1.44	0.96
1:A:246:TRP:O	1:A:250:ASP:HB2	1.66	0.94
1:A:221:PHE:O	1:A:225:SER:HB3	1.67	0.94
1:B:189:THR:O	1:B:193:VAL:HG23	1.71	0.91
2:A:301:BR:BR	1:C:264:ARG:CD	2.74	0.91
1:C:277:LYS:HD3	1:D:279:ASP:HB3	1.54	0.89
1:C:270:LEU:N	1:C:270:LEU:HD23	1.88	0.89
1:A:153:ILE:HG23	1:A:232:LEU:HD12	1.55	0.88
2:A:301:BR:BR	1:D:264:ARG:HD3	2.33	0.83
1:D:159:VAL:O	1:D:163:VAL:HG23	1.80	0.82
1:A:274:LEU:HD13	1:B:275:SER:HA	1.61	0.81
1:A:272:ARG:HG3	1:A:272:ARG:HH11	1.44	0.80
1:A:205:ARG:NH1	1:D:192:GLN:OE1	2.16	0.79
1:A:264:ARG:HB2	1:A:264:ARG:HH21	1.48	0.79
1:B:215:TRP:CE2	1:B:216:ILE:HD12	2.20	0.77
1:C:204:ALA:O	1:C:208:ILE:HG12	1.83	0.77
1:A:246:TRP:O	1:A:250:ASP:CB	2.33	0.76
1:A:200:SER:O	1:A:205:ARG:HB2	1.87	0.75
1:A:258:GLN:OE1	1:A:258:GLN:HA	1.88	0.74
1:B:193:VAL:HG13	1:B:199:TRP:HB2	1.70	0.73
1:C:270:LEU:CD2	1:C:270:LEU:H	1.95	0.72
1:C:160:ILE:HD13	1:C:224:VAL:CG1	2.20	0.72
1:C:233:PHE:HE1	1:C:237:ILE:HD12	1.54	0.72
1:A:159:VAL:O	1:A:163:VAL:HG23	1.90	0.71
1:A:167:MET:O	1:A:171:LEU:HB2	1.91	0.71
1:A:264:ARG:NH2	1:D:263:GLU:OE1	2.24	0.70
1:A:264:ARG:HB2	1:A:264:ARG:NH2	2.05	0.70
1:D:155:LEU:O	1:D:159:VAL:HG23	1.92	0.70
1:B:274:LEU:O	1:B:278:VAL:HG23	1.92	0.69
1:A:259:ARG:O	1:A:263:GLU:HG3	1.92	0.69
1:A:203:ILE:O	1:A:207:VAL:HG23	1.92	0.69
1:D:221:PHE:O	1:D:225:SER:HB3	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:154:ALA:O	1:D:158:LEU:HD12	1.93	0.68
1:C:177:PRO:HD2	1:C:178:GLU:OE1	1.94	0.68
1:A:268:LEU:HD11	1:A:272:ARG:NH2	2.09	0.68
1:A:281:LEU:HB2	1:B:282:GLU:HG3	1.76	0.68
1:C:270:LEU:CD2	1:C:270:LEU:N	2.57	0.67
1:C:263:GLU:OE1	1:D:264:ARG:NH2	2.27	0.67
1:C:161:PHE:HE1	1:C:194:MET:CE	2.08	0.67
1:B:242:GLN:HA	1:B:246:TRP:HB2	1.76	0.66
1:A:272:ARG:HG3	1:A:272:ARG:NH1	2.06	0.66
1:C:161:PHE:HE1	1:C:194:MET:HE3	1.58	0.66
1:D:157:LEU:HA	1:D:160:ILE:HD12	1.77	0.66
1:A:234:ILE:HD11	1:D:236:ILE:HG22	1.78	0.66
1:A:222:ILE:O	1:A:226:SER:HB2	1.97	0.65
1:A:264:ARG:CB	1:A:264:ARG:HH21	2.08	0.65
1:C:258:GLN:OE1	1:C:258:GLN:HA	1.96	0.65
1:D:241:MET:HA	1:D:244:ALA:HB3	1.78	0.65
1:B:196:LEU:N	1:B:196:LEU:HD23	2.11	0.65
1:B:162:TYR:O	1:B:166:VAL:HG23	1.96	0.65
1:B:197:GLU:HG3	1:C:199:TRP:CD1	2.32	0.64
1:A:192:GLN:HG3	1:B:200:SER:HB3	1.80	0.64
1:A:153:ILE:CG2	1:A:232:LEU:HD12	2.27	0.63
1:A:180:PHE:HD1	1:A:186:SER:OG	1.80	0.63
1:B:215:TRP:NE1	1:B:216:ILE:HD12	2.12	0.63
1:A:176:PHE:HZ	1:A:210:ALA:HB2	1.62	0.63
1:A:268:LEU:HD11	1:A:272:ARG:CZ	2.27	0.63
1:A:236:ILE:O	1:A:240:SER:HB2	1.98	0.63
1:A:274:LEU:HD11	1:B:278:VAL:HG21	1.80	0.63
1:C:160:ILE:HD13	1:C:224:VAL:HG12	1.81	0.63
1:A:192:GLN:HG3	1:B:200:SER:CB	2.30	0.62
1:B:216:ILE:O	1:B:220:SER:HB3	1.99	0.62
1:B:214:ALA:O	1:B:217:TYR:HB3	2.01	0.61
1:B:281:LEU:HD13	1:C:282:GLU:HG3	1.81	0.61
1:B:274:LEU:O	1:B:274:LEU:HD12	2.00	0.61
1:C:274:LEU:HD13	1:D:275:SER:HA	1.82	0.60
1:D:166:VAL:HG22	1:D:183:LEU:HD11	1.84	0.60
1:C:233:PHE:HZ	1:D:233:PHE:HE2	1.50	0.60
1:B:258:GLN:HA	1:B:258:GLN:OE1	2.01	0.59
1:B:276:SER:HA	1:B:279:ASP:HB2	1.84	0.59
1:B:245:HIS:CE1	1:C:246:TRP:HB3	2.38	0.59
1:A:230:LEU:HD11	1:D:233:PHE:CE1	2.38	0.58
1:C:151:ALA:HA	1:C:154:ALA:HB3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:192:GLN:HG3	1:D:200:SER:CB	2.34	0.58
1:C:277:LYS:CD	1:D:279:ASP:HB3	2.31	0.58
1:A:182:THR:HG23	1:A:185:ALA:HB2	1.86	0.57
1:B:163:VAL:O	1:B:167:MET:HB2	2.03	0.57
1:A:195:THR:HB	1:B:196:LEU:HD13	1.87	0.57
1:A:234:ILE:CD1	1:D:236:ILE:HG22	2.35	0.57
1:C:161:PHE:CE1	1:C:194:MET:HE3	2.39	0.57
1:C:271:ILE:HD11	1:D:271:ILE:HD11	1.86	0.57
1:A:277:LYS:HE2	1:B:279:ASP:OD1	2.04	0.57
1:A:267:MET:HG2	1:B:268:LEU:HD23	1.85	0.57
1:D:150:ILE:HG21	1:D:236:ILE:HG13	1.86	0.57
1:B:203:ILE:HG13	1:C:201:MET:CE	2.34	0.56
1:B:169:THR:O	1:B:173:ALA:HB2	2.05	0.56
1:B:241:MET:HA	1:C:238:ILE:HD11	1.86	0.56
1:D:205:ARG:HA	1:D:208:ILE:HD12	1.88	0.56
1:A:196:LEU:HD22	1:D:195:THR:CG2	2.36	0.56
1:A:199:TRP:CD1	1:D:197:GLU:HG3	2.41	0.56
1:A:150:ILE:O	1:A:153:ILE:N	2.39	0.56
1:B:176:PHE:CZ	1:B:207:VAL:HA	2.42	0.55
1:B:192:GLN:OE1	1:C:205:ARG:NH1	2.39	0.55
1:D:176:PHE:CZ	1:D:207:VAL:HA	2.42	0.55
1:B:261:HIS:CE1	1:B:265:LEU:HD13	2.42	0.54
1:C:277:LYS:HD3	1:D:279:ASP:CB	2.33	0.54
1:A:238:ILE:HG22	1:A:239:GLU:N	2.23	0.54
1:B:238:ILE:HG22	1:B:239:GLU:N	2.22	0.54
1:C:233:PHE:HZ	1:D:233:PHE:CE2	2.25	0.54
1:C:237:ILE:HD11	1:D:237:ILE:HD12	1.90	0.53
1:C:197:GLU:OE2	1:C:198:SER:HB2	2.08	0.53
1:C:241:MET:SD	1:D:241:MET:HE3	2.49	0.53
1:D:237:ILE:HG22	1:D:238:ILE:N	2.24	0.53
1:A:173:ALA:HB2	1:A:180:PHE:O	2.09	0.53
1:B:197:GLU:OE1	1:B:198:SER:HB2	2.09	0.52
1:C:192:GLN:OE1	1:D:205:ARG:NH2	2.35	0.52
1:B:251:ALA:O	1:B:254:ILE:HG22	2.08	0.52
1:B:178[A]:GLU:CD	1:B:178[A]:GLU:H	2.13	0.52
1:C:176:PHE:CZ	1:C:207:VAL:HA	2.45	0.52
1:D:236:ILE:O	1:D:240:SER:HB2	2.10	0.52
1:B:271:ILE:O	1:B:275:SER:HB2	2.10	0.52
1:A:281:LEU:HD23	1:D:281:LEU:HD11	1.92	0.52
1:B:215:TRP:CE2	1:B:216:ILE:CD1	2.92	0.51
1:C:197:GLU:HG3	1:D:199:TRP:CD1	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:283:ARG:C	1:C:285:SER:H	2.13	0.51
1:A:164:PHE:CD2	1:A:221:PHE:HD2	2.28	0.51
1:B:150:ILE:HG21	1:B:236:ILE:HG13	1.91	0.51
1:C:255:GLU:O	1:C:258:GLN:HB3	2.11	0.51
1:A:264:ARG:NE	2:A:301:BR:BR	2.99	0.51
1:D:166:VAL:HG22	1:D:183:LEU:CD1	2.41	0.51
1:A:205:ARG:N	1:A:206:PRO:CD	2.73	0.51
1:C:241:MET:SD	1:D:241:MET:CE	2.99	0.51
1:A:173:ALA:HB1	1:A:181:GLY:HA2	1.93	0.50
1:C:238:ILE:CG2	1:C:239:GLU:N	2.74	0.50
1:D:283:ARG:O	1:D:285:SER:N	2.44	0.50
1:D:200:SER:HA	1:D:204:ALA:HB3	1.94	0.50
1:A:161:PHE:HD2	1:A:187:MET:HE1	1.76	0.50
1:A:230:LEU:HD11	1:D:233:PHE:HE1	1.73	0.50
1:B:240:SER:O	1:B:243:SER:HB3	2.11	0.50
1:D:220:SER:OG	1:D:221:PHE:N	2.44	0.49
1:A:221:PHE:O	1:A:225:SER:CB	2.51	0.49
1:C:238:ILE:HG22	1:C:239:GLU:H	1.77	0.49
1:C:161:PHE:CE1	1:C:194:MET:CE	2.93	0.49
1:B:200:SER:O	1:B:205:ARG:HB2	2.12	0.49
1:D:271:ILE:O	1:D:275:SER:HB2	2.12	0.49
1:A:272:ARG:O	1:A:275:SER:HB2	2.12	0.49
1:D:156:LEU:HG	1:D:160:ILE:HD11	1.95	0.48
1:C:233:PHE:C	1:C:233:PHE:CD1	2.87	0.48
1:D:216:ILE:O	1:D:220:SER:HB3	2.12	0.48
1:B:237:ILE:O	1:B:240:SER:HB2	2.13	0.48
1:A:161:PHE:CD2	1:A:187:MET:CE	2.97	0.48
1:A:267:MET:HG2	1:B:268:LEU:CD2	2.42	0.48
1:C:192:GLN:NE2	1:C:197:GLU:O	2.45	0.47
1:C:208:ILE:HD12	1:C:215:TRP:HB3	1.95	0.47
1:C:275:SER:O	1:C:279:ASP:HB2	2.14	0.47
1:D:208:ILE:HA	1:D:211:TYR:O	2.13	0.47
1:A:263:GLU:OE1	1:B:264:ARG:NH1	2.45	0.47
1:B:215:TRP:CG	1:B:216:ILE:N	2.82	0.47
1:C:177:PRO:O	1:C:181:GLY:HA3	2.14	0.47
1:A:157:LEU:HD13	1:A:228:THR:HG22	1.96	0.47
1:A:176:PHE:CZ	1:A:210:ALA:HB2	2.48	0.47
1:B:196:LEU:HD23	1:B:196:LEU:H	1.80	0.47
1:C:176:PHE:CE2	1:C:207:VAL:HA	2.50	0.47
1:A:198:SER:HA	1:D:197:GLU:OE2	2.14	0.47
1:C:267:MET:HB3	1:D:268:LEU:CD2	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:SER:HA	1:D:274:LEU:HD13	1.97	0.46
1:B:189:THR:HG23	1:B:203:ILE:CD1	2.46	0.46
1:A:161:PHE:HD2	1:A:187:MET:CE	2.29	0.46
1:A:234:ILE:HG22	1:A:235:GLY:N	2.30	0.46
1:B:226:SER:O	1:B:227:PHE:C	2.55	0.46
1:B:274:LEU:HD21	1:C:274:LEU:HG	1.98	0.46
1:D:204:ALA:O	1:D:205:ARG:C	2.55	0.46
1:B:203:ILE:HG13	1:C:201:MET:HE3	1.98	0.46
2:A:301:BR:BR	1:D:264:ARG:CD	3.12	0.45
1:A:176:PHE:O	1:A:179:TRP:N	2.47	0.45
1:B:191:PHE:O	1:B:195:THR:HG23	2.16	0.45
1:B:238:ILE:HA	1:B:241:MET:HE2	1.98	0.45
1:C:258:GLN:OE1	1:C:258:GLN:CA	2.64	0.45
1:A:161:PHE:CD2	1:A:187:MET:HE1	2.52	0.45
1:C:179:TRP:O	1:C:189:THR:OG1	2.33	0.45
1:D:200:SER:O	1:D:205:ARG:N	2.38	0.45
1:A:194:MET:HG3	1:A:225:SER:OG	2.17	0.45
1:A:282:GLU:HA	1:D:281:LEU:HD21	1.98	0.45
1:B:169:THR:O	1:B:173:ALA:CB	2.65	0.45
1:C:266:GLU:O	1:C:269:GLN:HB2	2.17	0.45
1:B:167:MET:O	1:B:171:LEU:HG	2.17	0.44
1:A:269:GLN:OE1	1:A:269:GLN:HA	2.17	0.44
1:A:245:HIS:CE1	1:B:246:TRP:CG	3.05	0.44
1:C:237:ILE:HG13	1:D:234:ILE:HG12	2.00	0.44
1:A:190:LEU:HA	1:A:190:LEU:HD23	1.75	0.44
1:C:264:ARG:HB2	1:C:264:ARG:HH21	1.82	0.44
1:D:223:LEU:HD23	1:D:223:LEU:HA	1.43	0.44
1:B:215:TRP:CD2	1:B:216:ILE:N	2.86	0.44
1:C:266:GLU:O	1:C:270:LEU:HD23	2.18	0.44
1:B:205:ARG:N	1:B:206:PRO:CD	2.81	0.44
1:B:220:SER:OG	1:B:221:PHE:N	2.51	0.44
1:D:254:ILE:HG22	1:D:255:GLU:N	2.33	0.44
1:A:214:ALA:O	1:A:217:TYR:HB3	2.17	0.43
1:D:150:ILE:HG13	1:D:150:ILE:H	1.68	0.43
1:A:205:ARG:N	1:A:206:PRO:HD2	2.33	0.43
1:C:281:LEU:HD23	1:C:281:LEU:HA	1.42	0.43
1:D:189:THR:O	1:D:193:VAL:HG23	2.18	0.43
1:D:150:ILE:HD12	1:D:151:ALA:H	1.82	0.43
1:D:204:ALA:HB1	1:D:208:ILE:HD11	2.01	0.43
1:C:231:ASN:O	1:C:232:LEU:C	2.55	0.43
1:D:156:LEU:O	1:D:160:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:VAL:O	1:A:210:ALA:HB3	2.17	0.43
1:A:238:ILE:CG2	1:A:239:GLU:N	2.81	0.43
1:D:151:ALA:HA	1:D:154:ALA:HB3	1.99	0.43
1:D:190:LEU:HD22	1:D:221:PHE:CE2	2.54	0.43
1:D:254:ILE:CG2	1:D:255:GLU:N	2.81	0.43
1:A:182:THR:O	1:A:186:SER:HB2	2.19	0.43
1:C:253:ARG:NH2	1:C:254:ILE:HG13	2.34	0.43
1:D:272:ARG:CG	1:D:272:ARG:HH11	2.32	0.43
1:C:233:PHE:HD1	1:C:233:PHE:C	2.22	0.43
1:A:236:ILE:HG22	1:B:234:ILE:CD1	2.49	0.42
1:D:231:ASN:O	1:D:232:LEU:C	2.55	0.42
1:B:182:THR:O	1:B:185:ALA:HB3	2.19	0.42
1:B:273:ASP:N	1:B:273:ASP:OD1	2.51	0.42
1:C:264:ARG:NH2	1:C:264:ARG:HB2	2.34	0.42
1:B:213:TRP:CE3	1:B:214:ALA:HA	2.54	0.42
1:A:211:TYR:HB3	1:A:213:TRP:CE2	2.54	0.42
1:B:203:ILE:O	1:B:207:VAL:HG23	2.19	0.42
1:D:160:ILE:HG13	1:D:160:ILE:H	1.74	0.42
1:A:237:ILE:HA	1:B:234:ILE:HD11	2.01	0.42
1:B:193:VAL:HG13	1:B:199:TRP:CB	2.46	0.42
1:C:208:ILE:HG12	1:C:208:ILE:H	1.59	0.42
1:A:196:LEU:HD22	1:D:195:THR:HG22	2.02	0.42
1:B:205:ARG:N	1:B:206:PRO:HD2	2.34	0.42
1:C:238:ILE:HG22	1:C:239:GLU:N	2.34	0.42
1:C:245:HIS:NE2	1:C:249:GLU:HG2	2.35	0.42
1:D:204:ALA:O	1:D:207:VAL:N	2.52	0.42
1:A:182:THR:HG23	1:A:185:ALA:CB	2.49	0.41
1:A:238:ILE:HG22	1:A:239:GLU:H	1.85	0.41
1:A:241:MET:CE	1:D:241:MET:SD	3.08	0.41
1:B:153:ILE:HG13	1:B:153:ILE:H	1.53	0.41
1:D:225:SER:O	1:D:229:VAL:HG23	2.20	0.41
1:C:195:THR:O	1:C:196:LEU:HB2	2.19	0.41
1:B:150:ILE:O	1:B:153:ILE:HG13	2.20	0.41
1:A:241:MET:SD	1:B:241:MET:HE3	2.60	0.41
1:C:218:PHE:O	1:C:222:ILE:HG13	2.20	0.41
1:A:241:MET:HA	1:A:244:ALA:HB3	2.02	0.41
1:A:237:ILE:HG13	1:B:234:ILE:CG1	2.50	0.41
1:B:267:MET:SD	1:C:267:MET:HG3	2.61	0.41
1:A:161:PHE:CD2	1:A:187:MET:HE2	2.55	0.41
1:A:270:LEU:HA	1:A:273:ASP:HB2	2.02	0.41
1:C:153:ILE:HD13	1:C:231:ASN:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:SER:HB3	1:D:192:GLN:HG3	2.03	0.41
1:C:267:MET:HG2	1:C:267:MET:H	1.71	0.41
1:A:241:MET:HE3	1:D:241:MET:SD	2.60	0.41
1:D:162:TYR:CD1	1:D:162:TYR:C	2.94	0.40
1:B:219:VAL:O	1:B:223:LEU:HG	2.22	0.40
1:C:166:VAL:O	1:C:167:MET:C	2.59	0.40
1:D:190:LEU:HD23	1:D:190:LEU:HA	1.51	0.40
1:A:169:THR:HG23	1:A:180:PHE:O	2.21	0.40
1:D:164:PHE:C	1:D:166:VAL:N	2.74	0.40
1:D:257:GLU:O	1:D:257:GLU:HG3	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	137/152 (90%)	113 (82%)	24 (18%)	0	100	100
1	B	137/152 (90%)	116 (85%)	19 (14%)	2 (2%)	13	50
1	C	136/152 (90%)	110 (81%)	24 (18%)	2 (2%)	13	50
1	D	136/152 (90%)	110 (81%)	24 (18%)	2 (2%)	13	50
All	All	546/608 (90%)	449 (82%)	91 (17%)	6 (1%)	17	57

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	151	ALA
1	C	284	ARG
1	C	224	VAL
1	D	237	ILE
1	D	212	PRO

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Mol	Chain	Res	Type
1	B	212	PRO

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	117/131 (89%)	96 (82%)	21 (18%)	2	10
1	B	118/131 (90%)	92 (78%)	26 (22%)	1	4
1	C	118/131 (90%)	96 (81%)	22 (19%)	2	9
1	D	119/131 (91%)	95 (80%)	24 (20%)	1	6
All	All	472/524 (90%)	379 (80%)	93 (20%)	1	7

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	TRP
1	A	153	ILE
1	A	161	PHE
1	A	175	SER
1	A	182	THR
1	A	186	SER
1	A	187	MET
1	A	197	GLU
1	A	201	MET
1	A	226	SER
1	A	229	VAL
1	A	232	LEU
1	A	234	ILE
1	A	237	ILE
1	A	238	ILE
1	A	247	GLU
1	A	249	GLU
1	A	254	ILE
1	A	267	MET
1	A	272	ARG

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Mol	Chain	Res	Type
1	A	276	SER
1	B	150	ILE
1	B	153	ILE
1	B	158	LEU
1	B	161	PHE
1	B	174	GLN
1	B	175	SER
1	B	182	THR
1	B	194	MET
1	B	195	THR
1	B	197	GLU
1	B	216	ILE
1	B	232	LEU
1	B	234	ILE
1	B	237	ILE
1	B	238	ILE
1	B	240	SER
1	B	249	GLU
1	B	253	ARG
1	B	255	GLU
1	B	264	ARG
1	B	268	LEU
1	B	270	LEU
1	B	274	LEU
1	B	275	SER
1	B	282	GLU
1	B	283	ARG
1	C	161	PHE
1	C	170	LYS
1	C	182	THR
1	C	187	MET
1	C	192	GLN
1	C	197	GLU
1	C	198	SER
1	C	201	MET
1	C	208	ILE
1	C	232	LEU
1	C	233	PHE
1	C	238	ILE
1	C	242	GLN
1	C	249	GLU
1	C	253	ARG

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Mol	Chain	Res	Type
1	C	254	ILE
1	C	268	LEU
1	C	270	LEU
1	C	275	SER
1	C	276	SER
1	C	279	ASP
1	C	285	SER
1	D	150	ILE
1	D	155	LEU
1	D	167	MET
1	D	171	LEU
1	D	174	GLN
1	D	175	SER
1	D	178	GLU
1	D	182	THR
1	D	195	THR
1	D	198	SER
1	D	219	VAL
1	D	222	ILE
1	D	225	SER
1	D	227	PHE
1	D	237	ILE
1	D	240	SER
1	D	243	SER
1	D	247	GLU
1	D	254	ILE
1	D	256	GLN
1	D	265	LEU
1	D	275	SER
1	D	279	ASP
1	D	283	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	231	ASN
1	A	245	HIS
1	B	174	GLN
1	B	231	ASN
1	B	242	GLN
1	B	245	HIS
1	C	242	GLN

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Mol	Chain	Res	Type
1	C	245	HIS

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 1 ligands modelled in this entry, 1 is 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 ⓘ

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	139/152 (91%)	-0.57	1 (0%) 89 89	58, 111, 150, 178	0
1	B	138/152 (90%)	-0.55	0 100 100	59, 115, 152, 178	0
1	C	138/152 (90%)	-0.54	1 (0%) 89 89	71, 108, 158, 186	0
1	D	138/152 (90%)	-0.61	1 (0%) 89 89	68, 104, 152, 167	0
All	All	553/608 (90%)	-0.57	3 (0%) 91 92	58, 110, 154, 186	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	182	THR	2.8
1	D	285	SER	2.6
1	C	283	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	BR	A	301	1/1	0.94	0.18	-	170,170,170,170	0

6.5 Other polymers ⓘ

There are no such residues in this entry.