wwPDB X-ray Structure Validation Summary Report (i)

Sep 20, 2023 – 09:49 PM EDT

PDB ID	:	5IWN
Title	:	Bacterial sodium channel pore domain, high bromide
Authors	:	Shaya, D.; Findeisen, F.; Rohaim, A.; Minor, D.L.
Deposited on	:	2016-03-22
Resolution	:	3.75 Å(reported)

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

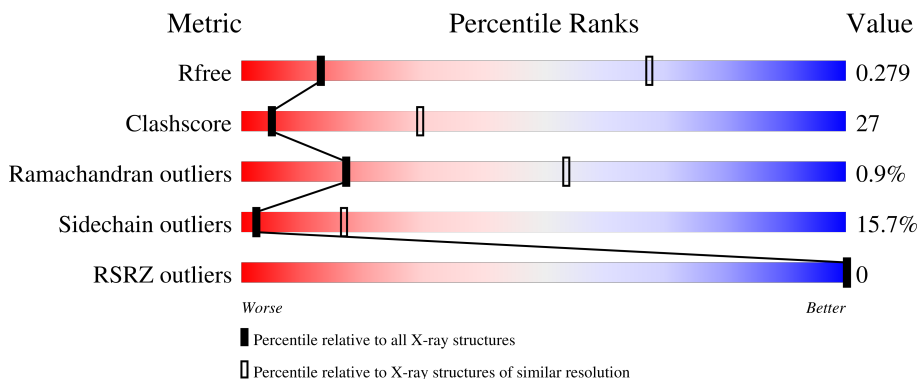|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation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

1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.75 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	Percentile Ranks	Value
Rfree		0.279
Clashscore		27
Ramachandran outliers		0.9%
Sidechain outliers		15.7%
RSRZ outliers		0

Worse   *Better*

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1039 (3.94-3.58)
Clashscore	141614	1051 (3.92-3.60)
Ramachandran outliers	138981	1015 (3.92-3.60)
Sidechain outliers	138945	1011 (3.92-3.60)
RSRZ outliers	127900	1050 (3.96-3.56)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	152	47% 35% 9% 9%
1	B	152	40% 41% 9% 9%
1	C	152	39% 44% 8% 9%
1	D	152	36% 43% 12% • 9%

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 criteria:

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

2 Entity composition ⓘ

There are 2 unique types of molecules in this entry. The entry contains 4502 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	2	Total 2	Br 2	0	0

3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.
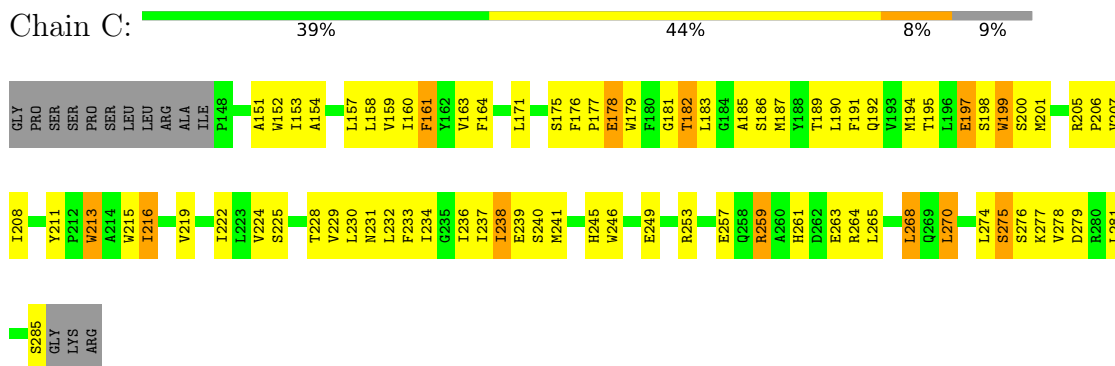
• Molecule 1: Ion transport protein
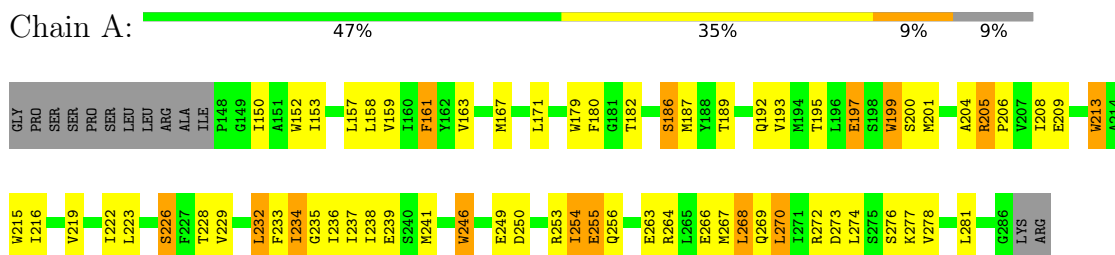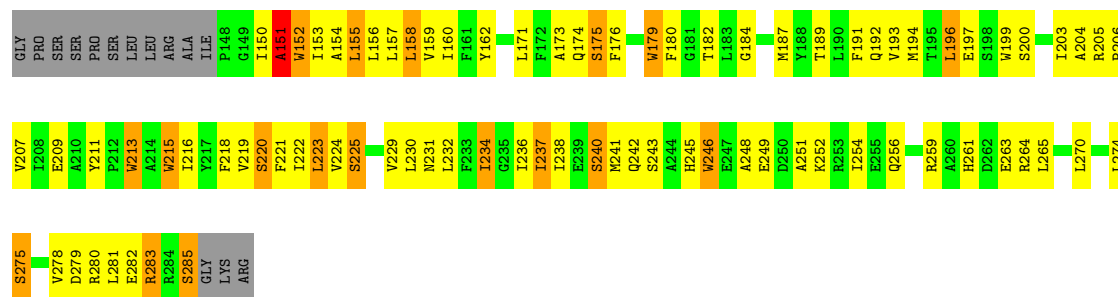
Chain A:

• Molecule 1: Ion transport protein
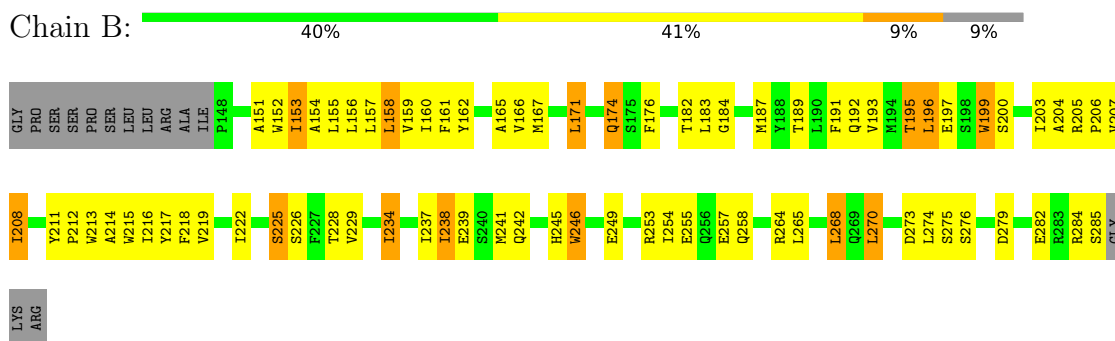
Chain B:

• Molecule 1: Ion transport protein

Chain C:

• Molecule 1: Ion transport protein

Chain D:

4   Data and refinement statistics   ⓘ

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	157.40Å 160.87Å 167.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.75 14.98 – 3.75	Depositor EDS
% Data completeness (in resolution range)	99.0 (15.00-3.75) 99.5 (14.98-3.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.26	Depositor
$< I/\sigma(I) >$ [1]	1.54 (at 3.77Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.228 , 0.280 0.222 , 0.279	Depositor DCC
R_{free} test set	1126 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å^2)	144.1	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.26 , 104.8	EDS
L-test for twinning[2]	$< |L| > = 0.35, < L^2 > = 0.18$	Xtriage
Estimated twinning fraction	0.129 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4502	wwPDB-VP
Average B, all atoms (Å^2)	169.0	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.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

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	0.70	3/1155 (0.3%)	0.78	1/1569 (0.1%)
1	B	0.70	3/1160 (0.3%)	0.77	1/1575 (0.1%)
1	C	0.76	4/1153 (0.3%)	0.81	0/1566
1	D	0.72	5/1159 (0.4%)	0.80	3/1573 (0.2%)
All	All	0.72	15/4627 (0.3%)	0.79	5/6283 (0.1%)

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	199	TRP	CD2-CE2	6.69	1.49	1.41
1	C	152	TRP	CD2-CE2	5.74	1.48	1.41
1	D	179	TRP	CD2-CE2	5.50	1.48	1.41
1	A	152	TRP	CD2-CE2	5.50	1.48	1.41
1	C	213	TRP	CD2-CE2	5.43	1.47	1.41

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	196	LEU	CB-CG-CD1	-5.77	101.19	111.00
1	D	196	LEU	CB-CG-CD1	-5.73	101.26	111.00
1	D	223	LEU	CA-CB-CG	-5.49	102.67	115.30
1	D	151	ALA	CB-CA-C	5.29	118.04	110.10
1	A	270	LEU	CB-CG-CD1	-5.02	102.47	111.00

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	1124	0	1111	71	0
1	B	1126	0	1119	61	0
1	C	1122	0	1112	85	0
1	D	1128	0	1123	78	0
2	A	2	0	0	6	0
All	All	4502	0	4465	244	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 27.

The worst 5 of 244 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)
2:A:301:BR:BR	1:C:264:ARG:HD3	1.81	1.33
1:A:264:ARG:HD3	2:A:301:BR:BR	1.87	1.28
1:A:264:ARG:NE	2:A:301:BR:BR	2.25	1.23
1:A:264:ARG:CD	2:A:301:BR:BR	2.44	1.21
1:B:208:ILE:HD11	1:B:215:TRP:HB3	1.25	1.17

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	137/152 (90%)	117 (85%)	19 (14%)	1 (1%)	22	59

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	137/152 (90%)	119 (87%)	16 (12%)	2 (2%)	10	45
1	C	136/152 (90%)	114 (84%)	21 (15%)	1 (1%)	22	59
1	D	136/152 (90%)	115 (85%)	20 (15%)	1 (1%)	22	59
All	All	546/608 (90%)	465 (85%)	76 (14%)	5 (1%)	17	54

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	151	ALA
1	A	199	TRP
1	B	151	ALA
1	B	212	PRO
1	C	216	ILE

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%)	101 (86%)	16 (14%)	3	22
1	B	118/131 (90%)	98 (83%)	20 (17%)	2	14
1	C	118/131 (90%)	100 (85%)	18 (15%)	2	17
1	D	119/131 (91%)	99 (83%)	20 (17%)	2	14
All	All	472/524 (90%)	398 (84%)	74 (16%)	2	17

5 of 74 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	175	SER
1	D	275	SER
1	D	209	GLU
1	D	242	GLN
1	B	225	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	242	GLN
1	C	242	GLN
1	D	245	HIS
1	C	245	HIS
1	B	231	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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

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 ⓘ

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(Å²)	Q<0.9
1	A	139/152 (91%)	-0.81	0	100	100	125, 169, 207, 233	0
1	B	138/152 (90%)	-0.79	0	100	100	130, 171, 213, 242	0
1	C	138/152 (90%)	-0.75	0	100	100	117, 165, 214, 258	0
1	D	138/152 (90%)	-0.78	0	100	100	115, 160, 203, 246	0
All	All	553/608 (90%)	-0.78	0	100	100	115, 166, 211, 258	0

There are no RSRZ outliers to report.

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 monosaccharides 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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å²)	Q<0.9
2	BR	A	301	1/1	0.92	0.49	178,178,178,178	0
2	BR	A	302	1/1	0.92	0.30	195,195,195,195	0

6.5 Other polymers (i)

There are no such residues in this entry.