wwPDB X-ray Structure Validation Summary Report (i)

May 31, 2022 – 07:09 pm BST

PDB ID	:	7PGF
Title	:	Calcium-selective Sp1 channel pore domain only
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Deposited on	:	2021-08-13
Resolution	:	3.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at *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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

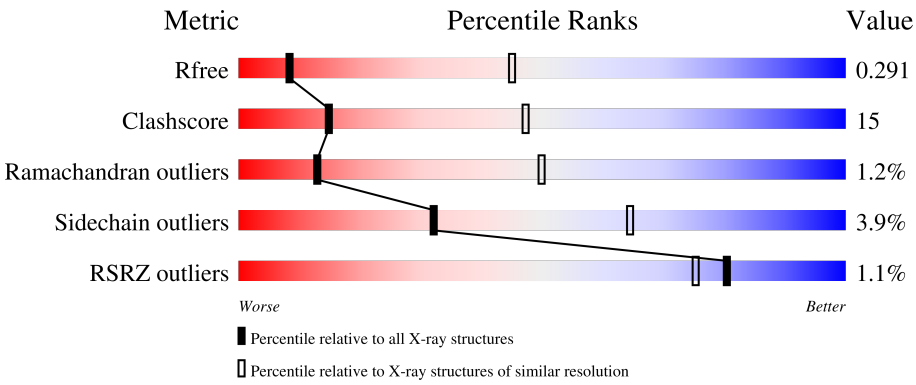
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.28.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.28.1

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 3.50 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	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

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	C	138	58% 33% • • 6%
1	D	138	2% 56% 36% • • •

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2031 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 transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	C	130	1008	666	158	178	6	0	0	0
			Total	C	N	O	S			
1	D	134	1023	674	157	186	6	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

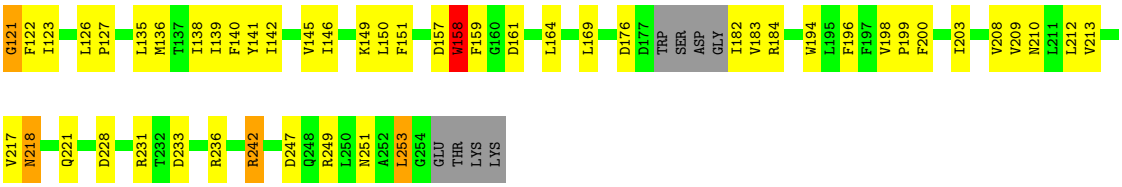
Chain	Residue	Modelled	Actual	Comment	Reference
C	176	ASP	GLU	engineered mutation	UNP A0A7Y7EFG5
C	177	ASP	SER	engineered mutation	UNP A0A7Y7EFG5
C	180	ASP	MET	engineered mutation	UNP A0A7Y7EFG5
D	176	ASP	GLU	engineered mutation	UNP A0A7Y7EFG5
D	177	ASP	SER	engineered mutation	UNP A0A7Y7EFG5
D	180	ASP	MET	engineered mutation	UNP A0A7Y7EFG5

3 Residue-property plots (i)

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.

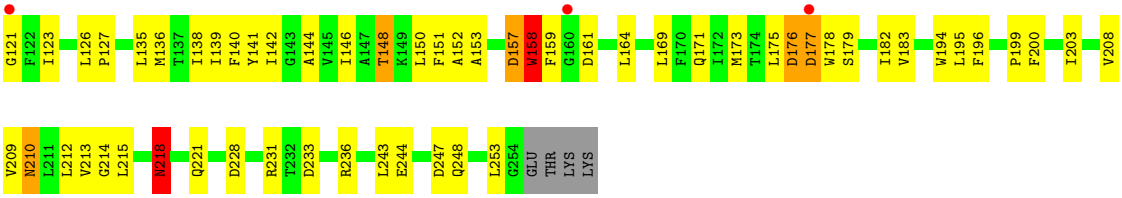
- Molecule 1: Ion transporter

Chain C:



- Molecule 1: Ion transporter

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	133.68Å 133.68Å 130.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	14.93 – 3.50 12.48 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (14.93-3.50) 100.0 (12.48-3.50)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.61 (at 3.53Å)	Xtriage
Refinement program	REFMAC 5.8.0, PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.265 , 0.290 0.273 , 0.291	Depositor DCC
R_{free} test set	420 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å^2)	134.6	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.42$, $< L^2 > = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	2031	wwPDB-VP
Average B, all atoms (Å^2)	153.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% of the height of the origin peak. No significant pseudotranslation is detected.*

[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 (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	C	0.48	3/1032 (0.3%)	0.80	6/1408 (0.4%)
1	D	0.46	3/1048 (0.3%)	0.75	6/1433 (0.4%)
All	All	0.47	6/2080 (0.3%)	0.78	12/2841 (0.4%)

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	C	0	3
1	D	0	2
All	All	0	5

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	210	ASN	CG-ND2	-7.07	1.15	1.32
1	C	210	ASN	CG-ND2	-6.65	1.16	1.32
1	C	210	ASN	CG-OD1	-6.11	1.10	1.24
1	D	210	ASN	CG-OD1	-6.10	1.10	1.24
1	C	218	ASN	CG-OD1	-5.75	1.11	1.24

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	242	ARG	CB-CG-CD	7.47	131.04	111.60
1	D	121	GLY	C-N-CA	7.33	140.02	121.70
1	C	242	ARG	CA-CB-CG	7.18	129.20	113.40
1	C	253	LEU	CA-CB-CG	6.95	131.29	115.30
1	C	161	ASP	CB-CG-OD1	6.34	124.01	118.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	121	GLY	Peptide
1	C	158	TRP	Peptide
1	C	218	ASN	Sidechain
1	D	158	TRP	Peptide
1	D	218	ASN	Sidechain

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	C	1008	0	992	33	1
1	D	1023	0	991	36	1
All	All	2031	0	1983	61	1

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 15.

The worst 5 of 61 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:C:198:VAL:HG21	1:D:195:LEU:HD22	1.49	0.91
1:D:175:LEU:O	1:D:177:ASP:N	2.18	0.76
1:D:176:ASP:O	1:D:178:TRP:N	2.21	0.72
1:C:249:ARG:O	1:C:253:LEU:HD23	1.90	0.72
1:D:148:THR:HG22	1:D:152:ALA:HB2	1.75	0.68

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:249:ARG:NH2	1:D:153:ALA:O[7_455]	2.19	0.01

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	C	126/138 (91%)	116 (92%)	9 (7%)	1 (1%)	19	58
1	D	132/138 (96%)	122 (92%)	8 (6%)	2 (2%)	10	45
All	All	258/276 (94%)	238 (92%)	17 (7%)	3 (1%)	13	50

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	176	ASP
1	C	122	PHE
1	D	177	ASP

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	C	103/115 (90%)	101 (98%)	2 (2%)	57	80
1	D	104/115 (90%)	98 (94%)	6 (6%)	20	53
All	All	207/230 (90%)	199 (96%)	8 (4%)	32	64

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	247	ASP
1	D	218	ASN
1	D	158	TRP

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Mol	Chain	Res	Type
1	D	157	ASP
1	D	210	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	210	ASN
1	D	218	ASN
1	D	221	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 monosaccharides 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 (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, 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	C	130/138 (94%)	-0.60	0	100	100	99, 149, 218, 263	0
1	D	134/138 (97%)	-0.48	3 (2%)	62	56	97, 148, 224, 281	0
All	All	264/276 (95%)	-0.54	3 (1%)	80	75	97, 149, 222, 281	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	160	GLY	2.8
1	D	121	GLY	2.4
1	D	177	ASP	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 monosaccharides 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.