W O R L D W I D E

wwPDB

P R O T E I N   D A T A   B A N K

wwPDB X-ray Structure Validation Summary Report (i)

Oct 13, 2024 – 10:26 AM EDT

PDB ID	:	3A2A
Title	:	The structure of the carboxyl-terminal domain of the human voltage-gated proton channel Hv1
Authors	:	Li, S.J.; Unno, H.; Zhou, Q.; Zhao, Q.; Zhai, Y.; Sun, F.
Deposited on	:	2009-05-08
Resolution	:	2.00 Å(reported)

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

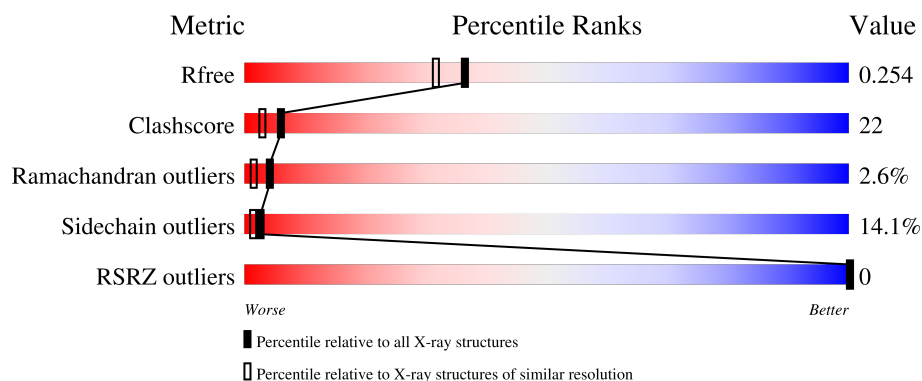|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 2.00 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		0.254
Clashscore		22
Ramachandran outliers		2.6%
Sidechain outliers		14.1%
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}	164625	9409 (2.00-2.00)
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (2.00-2.00)

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	58	53% 12% 5% 29%
1	B	58	36% 21% 10% • 29%
1	C	58	50% 14% 7% 29%
1	D	58	36% 22% 9% • 29%

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2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 1466 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 Voltage-gated hydrogen channel 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	41	Total	C	N	O	S	0	0	0
			349	217	69	61	2			
1	B	41	Total	C	N	O	S	0	0	0
			349	217	69	61	2			
1	C	41	Total	C	N	O	S	0	0	0
			349	217	69	61	2			
1	D	41	Total	C	N	O	S	0	0	0
			349	217	69	61	2			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	216	GLY	-	expression tag	UNP Q96D96
A	217	PRO	-	expression tag	UNP Q96D96
A	218	LEU	-	expression tag	UNP Q96D96
A	219	GLY	-	expression tag	UNP Q96D96
A	220	SER	-	expression tag	UNP Q96D96
B	216	GLY	-	expression tag	UNP Q96D96
B	217	PRO	-	expression tag	UNP Q96D96
B	218	LEU	-	expression tag	UNP Q96D96
B	219	GLY	-	expression tag	UNP Q96D96
B	220	SER	-	expression tag	UNP Q96D96
C	216	GLY	-	expression tag	UNP Q96D96
C	217	PRO	-	expression tag	UNP Q96D96
C	218	LEU	-	expression tag	UNP Q96D96
C	219	GLY	-	expression tag	UNP Q96D96
C	220	SER	-	expression tag	UNP Q96D96
D	216	GLY	-	expression tag	UNP Q96D96
D	217	PRO	-	expression tag	UNP Q96D96
D	218	LEU	-	expression tag	UNP Q96D96
D	219	GLY	-	expression tag	UNP Q96D96
D	220	SER	-	expression tag	UNP Q96D96

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• Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Cl 1	0	0
2	C	1	Total 1	Cl 1	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	14	Total 14	O 14	0	0
3	B	20	Total 20	O 20	0	0
3	C	14	Total 14	O 14	0	0
3	D	20	Total 20	O 20	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: Voltage-gated hydrogen channel 1

Chain A:

- Molecule 1: Voltage-gated hydrogen channel 1

Chain B:

- Molecule 1: Voltage-gated hydrogen channel 1

Chain C:

- Molecule 1: Voltage-gated hydrogen channel 1

Chain D:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α, β, γ	37.65Å 37.65Å 137.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.65 – 2.00 37.65 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (37.65-2.00) 100.0 (37.65-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.86 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.252 , 0.266 0.251 , 0.254	Depositor DCC
R_{free} test set	628 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å^2)	31.8	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 63.7	EDS
L-test for twinning[2]	$< |L| > = 0.50, < L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.487 for h,-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	1466	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.27% 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: CL

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	1.45	1/351 (0.3%)	1.17	2/465 (0.4%)
1	B	1.20	1/351 (0.3%)	1.18	4/465 (0.9%)
1	C	1.47	2/351 (0.6%)	1.19	2/465 (0.4%)
1	D	1.23	1/351 (0.3%)	1.18	3/465 (0.6%)
All	All	1.34	5/1404 (0.4%)	1.18	11/1860 (0.6%)

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	B	0	1
1	D	0	1
All	All	0	2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	249	CYS	CB-SG	9.00	1.97	1.82
1	C	249	CYS	CB-SG	8.72	1.97	1.82
1	B	246	GLU	CD-OE1	-5.76	1.19	1.25
1	D	246	GLU	CD-OE1	-5.62	1.19	1.25
1	C	233	GLN	CA-CB	-5.24	1.42	1.53

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	263	LEU	CA-CB-CG	8.50	134.84	115.30
1	D	263	LEU	CA-CB-CG	8.02	133.73	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	230	ARG	NE-CZ-NH2	-7.29	116.65	120.30
1	A	230	ARG	NE-CZ-NH2	-6.72	116.94	120.30
1	C	230	ARG	NE-CZ-NH1	5.77	123.18	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	229	LEU	Peptide
1	D	229	LEU	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	349	0	368	7	0
1	B	349	0	368	26	0
1	C	349	0	368	12	0
1	D	349	0	368	26	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0
3	A	14	0	0	1	0
3	B	20	0	0	6	0
3	C	14	0	0	2	0
3	D	20	0	0	7	0
All	All	1466	0	1472	64	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 22.

The worst 5 of 64 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:264:ARG:HD3	3:B:32:HOH:O	1.48	1.11
1:D:264:ARG:HD3	3:D:66:HOH:O	1.52	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:ARG:HG3	1:D:233:GLN:HE22	1.35	0.92
1:D:264:ARG:NH1	3:D:50:HOH:O	2.02	0.92
1:B:264:ARG:NH1	3:B:16:HOH:O	2.07	0.86

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	39/58 (67%)	39 (100%)	0	0	100	100
1	B	39/58 (67%)	36 (92%)	1 (3%)	2 (5%)	1	0
1	C	39/58 (67%)	39 (100%)	0	0	100	100
1	D	39/58 (67%)	36 (92%)	1 (3%)	2 (5%)	1	0
All	All	156/232 (67%)	150 (96%)	2 (1%)	4 (3%)	4	1

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	230	ARG
1	D	230	ARG
1	B	229	LEU
1	D	229	LEU

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	39/52 (75%)	35 (90%)	4 (10%)	6	3
1	B	39/52 (75%)	32 (82%)	7 (18%)	1	1
1	C	39/52 (75%)	35 (90%)	4 (10%)	6	3
1	D	39/52 (75%)	32 (82%)	7 (18%)	1	1
All	All	156/208 (75%)	134 (86%)	22 (14%)	3	1

5 of 22 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	262	LEU
1	D	234	MET
1	D	229	LEU
1	D	250	SER
1	B	234	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	260	ASN
1	C	260	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 oligosaccharides 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 ⓘ

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	41/58 (70%)	-1.12	0	100	100	20, 30, 50, 60	0
1	B	41/58 (70%)	-1.15	0	100	100	22, 32, 57, 68	0
1	C	41/58 (70%)	-1.11	0	100	100	20, 30, 50, 60	0
1	D	41/58 (70%)	-1.16	0	100	100	23, 32, 57, 68	0
All	All	164/232 (70%)	-1.14	0	100	100	20, 31, 57, 68	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	CL	A	1	1/1	1.00	0.01	23,23,23,23	1
2	CL	C	2	1/1	1.00	0.01	23,23,23,23	1

6.5 Other polymers ⓘ

There are no such residues in this entry.