



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:54 PM GMT

PDB ID : 4N5Q
Title : Crystal structure of the N-terminal ankyrin repeat domain of TRPV3
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Deposited on : 2013-10-10
Resolution : 1.95 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

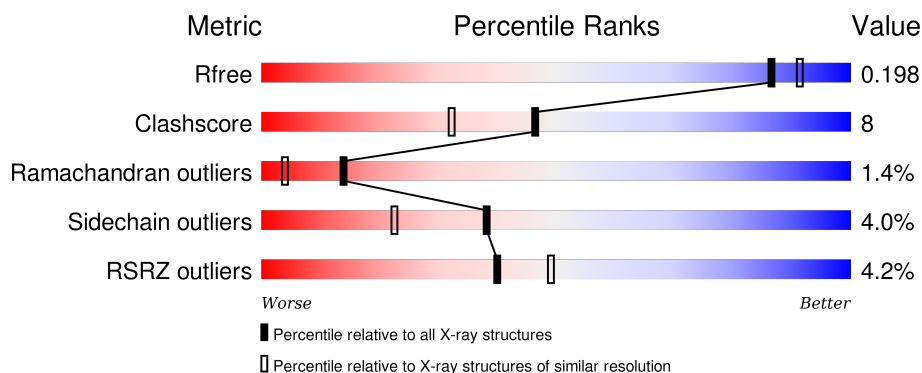
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	2910 (1.96-1.92)
Clashscore	102246	3095 (1.96-1.92)
Ramachandran outliers	100387	3062 (1.96-1.92)
Sidechain outliers	100360	3062 (1.96-1.92)
RSRZ outliers	91569	2915 (1.96-1.92)

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	257	<div> <div>3%</div> <div>82%</div> <div>12%</div> <div>• •</div> </div>
1	B	257	<div> <div>5%</div> <div>82%</div> <div>13%</div> <div>• • •</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4545 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 Transient receptor potential cation channel subfamily V member 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	0	0
			1978	1243	346	377	12			
1	B	254	Total	C	N	O	S	0	1	0
			2029	1274	358	384	13			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	117	MET	-	EXPRESSION TAG	UNP Q8K424
A	368	HIS	-	EXPRESSION TAG	UNP Q8K424
A	369	HIS	-	EXPRESSION TAG	UNP Q8K424
A	370	HIS	-	EXPRESSION TAG	UNP Q8K424
A	371	HIS	-	EXPRESSION TAG	UNP Q8K424
A	372	HIS	-	EXPRESSION TAG	UNP Q8K424
A	373	HIS	-	EXPRESSION TAG	UNP Q8K424
B	117	MET	-	EXPRESSION TAG	UNP Q8K424
B	368	HIS	-	EXPRESSION TAG	UNP Q8K424
B	369	HIS	-	EXPRESSION TAG	UNP Q8K424
B	370	HIS	-	EXPRESSION TAG	UNP Q8K424
B	371	HIS	-	EXPRESSION TAG	UNP Q8K424
B	372	HIS	-	EXPRESSION TAG	UNP Q8K424
B	373	HIS	-	EXPRESSION TAG	UNP Q8K424

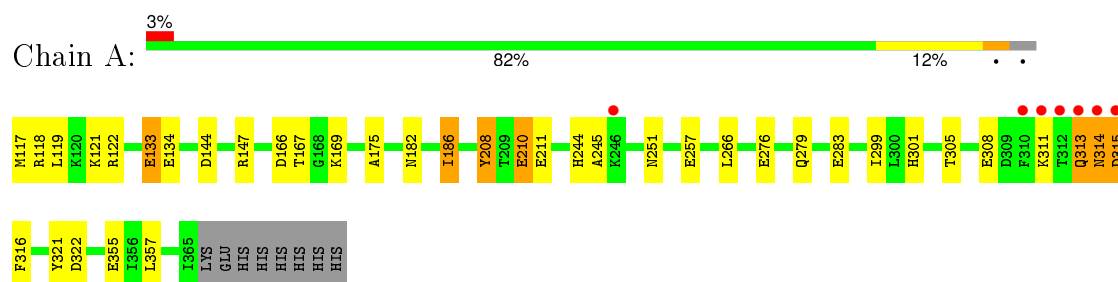
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	275	Total	O	0	0
			275	275		
2	B	263	Total	O	0	0
			263	263		

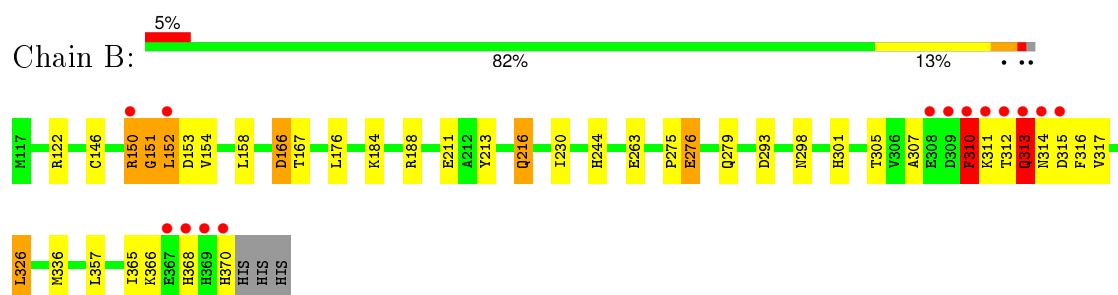
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Transient receptor potential cation channel subfamily V member 3



- Molecule 1: Transient receptor potential cation channel subfamily V member 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.52Å 78.56Å 82.97Å 90.00° 93.30° 90.00°	Depositor
Resolution (Å)	41.40 – 1.95 41.42 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.3 (41.40-1.95) 93.0 (41.42-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 1.95Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.169 , 0.201 0.165 , 0.198	Depositor DCC
R_{free} test set	2718 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 57524 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4545	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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	0.73	0/2005	0.76	0/2705
1	B	0.73	0/2062	0.77	3/2781 (0.1%)
All	All	0.73	0/4067	0.76	3/5486 (0.1%)

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

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	166	ASP	N-CA-CB	-5.64	100.45	110.60
1	B	326	LEU	CA-CB-CG	-5.22	103.29	115.30
1	B	151	GLY	N-CA-C	5.05	125.72	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	310	PHE	Peptide

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	1978	0	1989	25	0
1	B	2029	0	2034	37	0
2	A	275	0	0	11	0
2	B	263	0	0	9	0
All	All	4545	0	4023	61	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 8.

All (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:A:321:TYR:OH	2:A:554:HOH:O	1.87	0.92
1:B:279:GLN:OE1	2:B:504:HOH:O	1.96	0.84
1:A:210:GLU:OE1	2:A:644:HOH:O	1.95	0.83
1:B:370:HIS:ND1	2:B:557:HOH:O	2.10	0.82
1:B:153:ASP:OD2	2:B:645:HOH:O	2.00	0.79
1:A:257:GLU:OE2	2:A:514:HOH:O	2.01	0.78
1:A:313:GLN:O	1:A:315:ASP:N	2.20	0.72
1:B:314:ASN:O	1:B:316:PHE:N	2.22	0.72
1:B:370:HIS:CG	2:B:557:HOH:O	2.45	0.70
1:B:311:LYS:HG2	1:B:313:GLN:HG3	1.76	0.68
1:B:314:ASN:C	1:B:316:PHE:H	1.96	0.68
1:A:322:ASP:OD2	2:A:565:HOH:O	2.12	0.68
1:A:134:GLU:OE1	2:A:557:HOH:O	2.12	0.67
1:B:150:ARG:HA	1:B:154:VAL:HG23	1.78	0.66
1:B:151:GLY:O	1:B:152:LEU:HB2	1.96	0.65
1:B:313:GLN:CD	1:B:314:ASN:H	2.01	0.64
1:A:144:ASP:OD1	2:A:588:HOH:O	2.13	0.64
1:B:298:ASN:HD22	1:B:336:MET:HE2	1.62	0.64
1:B:310:PHE:HD2	1:B:311:LYS:N	1.98	0.61
1:B:211:GLU:H	1:B:211:GLU:CD	2.04	0.60
1:A:314:ASN:ND2	1:A:355:GLU:OE1	2.36	0.59
1:B:370:HIS:HB3	2:B:557:HOH:O	2.05	0.56
1:A:313:GLN:NE2	1:A:313:GLN:O	2.40	0.55
1:B:276:GLU:H	1:B:276:GLU:CD	2.09	0.54
1:A:118:ARG:HH21	1:B:122:ARG:NH1	2.06	0.53
1:B:314:ASN:C	1:B:316:PHE:N	2.63	0.52
1:A:279:GLN:O	1:A:283:GLU:HG3	2.10	0.51
1:A:175:ALA:HB2	1:A:186:ILE:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:ARG:HD2	2:B:533:HOH:O	2.10	0.51
1:B:365:ILE:O	1:B:368:HIS:HB3	2.10	0.51
1:B:275:PRO:O	1:B:279:GLN:HG3	2.12	0.49
1:B:301:HIS:O	1:B:305:THR:HG23	2.12	0.49
1:B:244:HIS:HD2	2:B:549:HOH:O	1.96	0.47
1:B:311:LYS:HB3	1:B:313:GLN:O	2.15	0.47
1:B:216:GLN:HG3	2:B:426:HOH:O	2.14	0.46
1:A:244:HIS:CD2	2:A:529:HOH:O	2.69	0.46
1:B:176:LEU:HD21	1:B:230:ILE:HG21	1.96	0.46
1:A:251:ASN:ND2	2:A:560:HOH:O	2.38	0.45
1:A:121:LYS:NZ	2:A:581:HOH:O	2.41	0.45
1:A:133:GLU:H	1:A:133:GLU:CD	2.17	0.45
1:B:166:ASP:HB3	1:B:167:THR:HG23	1.97	0.45
1:A:166:ASP:HB3	1:A:167:THR:HG23	1.99	0.45
1:B:312:THR:OG1	1:B:313:GLN:N	2.50	0.44
1:B:146:CYS:SG	1:B:158:LEU:HD21	2.58	0.44
1:A:169:LYS:HE2	1:A:208:TYR:OH	2.17	0.44
1:A:301:HIS:O	1:A:305:THR:HG23	2.17	0.44
1:B:326:LEU:HD11	1:B:366:LYS:NZ	2.33	0.44
1:B:263:GLU:OE1	1:B:293:ASP:HB2	2.17	0.44
1:A:117:MET:HG3	1:A:119:LEU:H	1.81	0.44
1:A:266:LEU:HG	1:A:299:ILE:HD12	2.01	0.43
1:B:307:ALA:HB2	1:B:317:VAL:HG11	2.00	0.43
1:B:184:LYS:O	1:B:188:ARG:HG3	2.19	0.43
1:B:310:PHE:HD2	1:B:311:LYS:H	1.66	0.42
1:B:310:PHE:CD2	1:B:311:LYS:N	2.84	0.42
1:A:122:ARG:HD2	2:A:557:HOH:O	2.18	0.42
1:A:245:ALA:HB3	2:A:550:HOH:O	2.19	0.41
1:A:182:ASN:O	1:A:186:ILE:HG23	2.21	0.41
1:B:150:ARG:HA	1:B:154:VAL:CG2	2.48	0.41
1:B:275:PRO:HD2	1:B:276:GLU:OE1	2.21	0.41
1:B:213:TYR:HE1	2:B:558:HOH:O	2.03	0.41
1:A:313:GLN:C	1:A:315:ASP:H	2.16	0.41

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	247/257 (96%)	239 (97%)	6 (2%)	2 (1%)	24	10
1	B	253/257 (98%)	238 (94%)	10 (4%)	5 (2%)	9	2
All	All	500/514 (97%)	477 (95%)	16 (3%)	7 (1%)	14	4

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	311	LYS
1	A	314	ASN
1	B	150	ARG
1	B	152	LEU
1	B	310	PHE
1	B	315	ASP
1	B	313	GLN

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	A	212/220 (96%)	200 (94%)	12 (6%)	25	10
1	B	218/220 (99%)	213 (98%)	5 (2%)	58	48
All	All	430/440 (98%)	413 (96%)	17 (4%)	38	22

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	GLU
1	A	147	ARG
1	A	186	ILE
1	A	208	TYR
1	A	210	GLU
1	A	211	GLU
1	A	276	GLU
1	A	308	GLU
1	A	313	GLN
1	A	315	ASP
1	A	316	PHE
1	A	357	LEU
1	B	216	GLN
1	B	276	GLU
1	B	310	PHE
1	B	313	GLN
1	B	357	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	279	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 carbohydrates 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

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	249/257 (96%)	0.02	7 (2%) 56 65	17, 29, 63, 124	0
1	B	254/257 (98%)	0.23	14 (5%) 29 37	17, 28, 76, 122	0
All	All	503/514 (97%)	0.12	21 (4%) 40 49	17, 29, 69, 124	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	310	PHE	11.6
1	A	314	ASN	9.7
1	B	312	THR	9.5
1	B	152	LEU	7.2
1	B	313	GLN	6.1
1	B	370	HIS	5.8
1	B	150	ARG	5.8
1	A	312	THR	5.6
1	A	310	PHE	5.5
1	A	315	ASP	5.4
1	B	309	ASP	4.6
1	B	368	HIS	4.1
1	B	311	LYS	3.8
1	B	314	ASN	3.5
1	B	315	ASP	3.3
1	A	313	GLN	3.2
1	B	367	GLU	3.0
1	A	311	LYS	3.0
1	B	308	GLU	2.8
1	B	369	HIS	2.8
1	A	246	LYS	2.6

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.