



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:14 PM UTC

PDB ID : 6BWM / pdb_00006bwm
Title : Crystal structure of the TRPV2 ion channel
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Deposited on : 2017-12-15
Resolution : 3.90 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

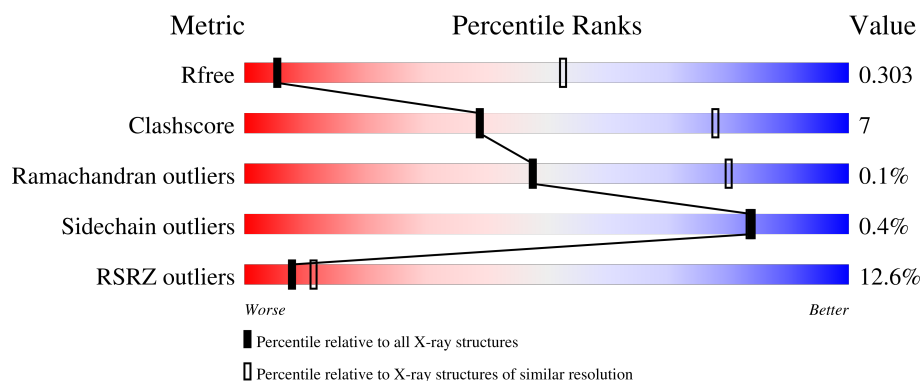
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.90 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}	180053	1270 (4.10-3.70)
Clashscore	190562	1034 (4.08-3.72)
Ramachandran outliers	187476	1251 (4.10-3.70)
Sidechain outliers	187428	1243 (4.10-3.70)
RSRZ outliers	180081	1269 (4.10-3.70)

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 $\leq 5\%$. 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	776	
1	B	776	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8512 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 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	606	Total	C	N	O	S	0	0	0
			4298	2758	742	776	22			
1	B	599	Total	C	N	O	S	0	0	0
			4213	2696	724	769	24			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	763	ASP	-	expression tag	UNP G1SNM3
A	764	TYR	-	expression tag	UNP G1SNM3
A	765	LYS	-	expression tag	UNP G1SNM3
A	766	ASP	-	expression tag	UNP G1SNM3
A	767	ASP	-	expression tag	UNP G1SNM3
A	768	ASP	-	expression tag	UNP G1SNM3
A	769	ASP	-	expression tag	UNP G1SNM3
A	770	LYS	-	expression tag	UNP G1SNM3
A	771	ALA	-	expression tag	UNP G1SNM3
A	772	HIS	-	expression tag	UNP G1SNM3
A	773	HIS	-	expression tag	UNP G1SNM3
A	774	HIS	-	expression tag	UNP G1SNM3
A	775	HIS	-	expression tag	UNP G1SNM3
A	776	HIS	-	expression tag	UNP G1SNM3
A	777	HIS	-	expression tag	UNP G1SNM3
B	763	ASP	-	expression tag	UNP G1SNM3
B	764	TYR	-	expression tag	UNP G1SNM3
B	765	LYS	-	expression tag	UNP G1SNM3
B	766	ASP	-	expression tag	UNP G1SNM3
B	767	ASP	-	expression tag	UNP G1SNM3
B	768	ASP	-	expression tag	UNP G1SNM3
B	769	ASP	-	expression tag	UNP G1SNM3
B	770	LYS	-	expression tag	UNP G1SNM3
B	771	ALA	-	expression tag	UNP G1SNM3

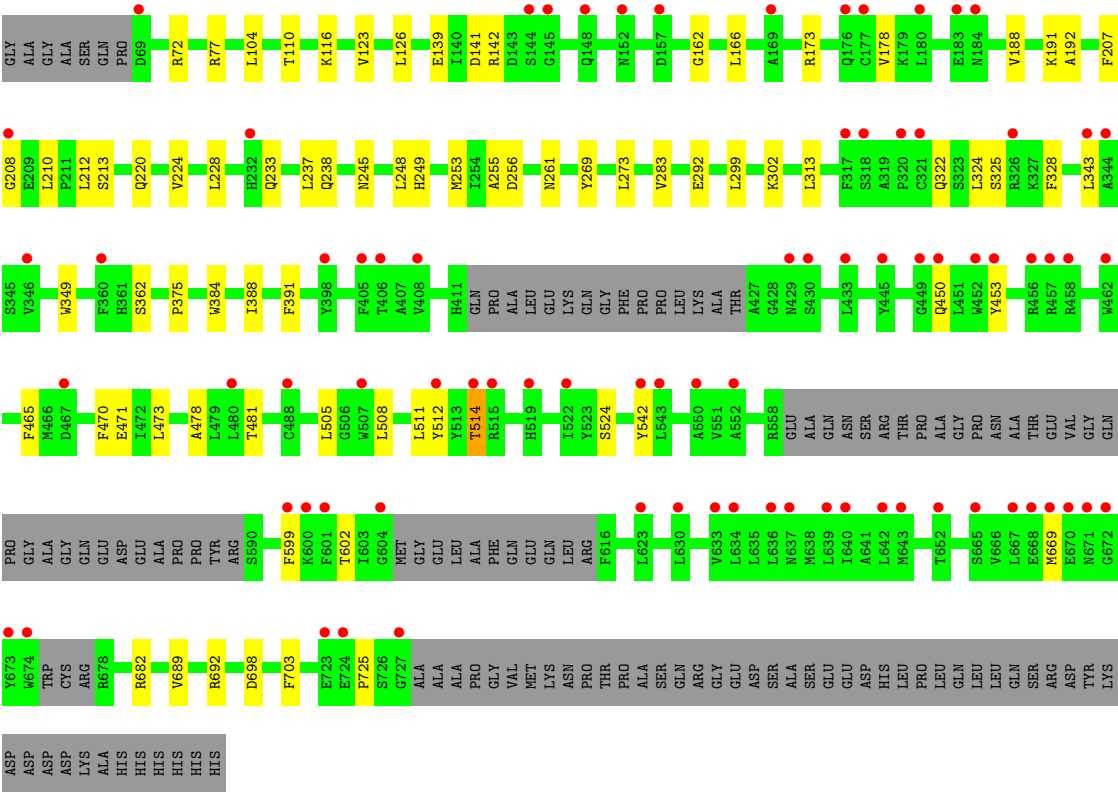
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Chain	Residue	Modelled	Actual	Comment	Reference
B	772	HIS	-	expression tag	UNP G1SNM3
B	773	HIS	-	expression tag	UNP G1SNM3
B	774	HIS	-	expression tag	UNP G1SNM3
B	775	HIS	-	expression tag	UNP G1SNM3
B	776	HIS	-	expression tag	UNP G1SNM3
B	777	HIS	-	expression tag	UNP G1SNM3

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	92.03Å 122.36Å 187.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.85 – 3.90 39.85 – 3.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.85-3.90) 99.7 (39.85-3.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 3.87Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.244 , 0.293 0.263 , 0.303	Depositor DCC
R_{free} test set	1058 reflections (5.31%)	wwPDB-VP
Wilson B-factor (Å ²)	89.7	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 71.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	8512	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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, 0.333 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: CA

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	0.15	0/4391	0.43	2/6013 (0.0%)
1	B	0.15	0/4301	0.40	0/5890
All	All	0.15	0/8692	0.42	2/11903 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	209	GLU	CA-C-N	6.09	129.94	120.68
1	A	209	GLU	C-N-CA	6.09	129.94	120.68

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	4298	0	3785	59	0
1	B	4213	0	3682	53	0
2	B	1	0	0	0	0
All	All	8512	0	7467	109	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 7.

All (109) 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:542:TYR:OH	1:B:602:THR:HG21	1.38	1.20
1:B:542:TYR:CZ	1:B:602:THR:HG21	2.04	0.91
1:B:542:TYR:OH	1:B:602:THR:CG2	2.21	0.88
1:B:207:PHE:CD1	1:B:208:GLY:O	2.26	0.88
1:B:207:PHE:HD1	1:B:208:GLY:O	1.62	0.82
1:A:625:LEU:HA	1:B:599:PHE:CZ	2.15	0.81
1:A:198:GLN:HA	1:A:206:TYR:HD1	1.48	0.78
1:B:542:TYR:CE1	1:B:602:THR:HG21	2.21	0.75
1:A:301:ALA:HA	1:A:370:MET:HE1	1.68	0.75
1:A:138:LEU:HD22	1:A:151:VAL:HG22	1.77	0.67
1:A:480:LEU:HD22	1:A:498:LEU:HD11	1.78	0.66
1:B:255:ALA:O	1:B:261:ASN:ND2	2.29	0.66
1:A:198:GLN:HA	1:A:206:TYR:CD1	2.32	0.63
1:A:470:PHE:HE2	1:A:512:TYR:HB2	1.65	0.61
1:A:511:LEU:O	1:A:514:THR:HG22	2.01	0.61
1:A:343:LEU:HD11	1:A:703:PHE:HB2	1.83	0.60
1:A:114:THR:HA	1:A:157:ASP:HB2	1.83	0.59
1:A:686:MET:HE3	1:A:699:GLU:HB3	1.84	0.59
1:B:343:LEU:HD11	1:B:703:PHE:HB2	1.84	0.59
1:A:255:ALA:HB3	1:A:305:LYS:HG2	1.84	0.59
1:A:69:ASP:O	1:A:96:TYR:HB2	2.03	0.59
1:A:70:LEU:O	1:A:99:ARG:NH1	2.37	0.58
1:B:212:LEU:HD21	1:B:237:LEU:HD23	1.86	0.58
1:B:453:TYR:CE2	1:B:471:GLU:HG2	2.39	0.57
1:A:173:ARG:HG2	1:A:220:GLN:NE2	2.19	0.56
1:A:447:LEU:HD11	1:A:479:LEU:HA	1.88	0.56
1:B:292:GLU:OE2	1:B:302:LYS:NZ	2.38	0.56
1:A:336:VAL:HG22	1:A:708:MET:HE2	1.88	0.56
1:B:508:LEU:O	1:B:511:LEU:HB3	2.06	0.55
1:B:207:PHE:CE1	1:B:208:GLY:O	2.60	0.55
1:B:245:ASN:HB3	1:B:249:HIS:HB2	1.88	0.55
1:B:72:ARG:O	1:B:77:ARG:NH1	2.39	0.54
1:A:698:ASP:OD2	1:A:700:ARG:NH2	2.42	0.53
1:B:173:ARG:HG2	1:B:220:GLN:NE2	2.24	0.52
1:A:323:SER:HB3	1:A:691:THR:O	2.10	0.51
1:A:82:VAL:HG11	1:A:137:LEU:HD11	1.92	0.50
1:B:692:ARG:HE	1:B:698:ASP:HB3	1.76	0.49
1:B:478:ALA:O	1:B:481:THR:OG1	2.28	0.49
1:A:252:VAL:HG22	1:A:308:ILE:HD13	1.94	0.49
1:A:336:VAL:CG2	1:A:708:MET:HE2	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:628:VAL:CB	1:B:599:PHE:HZ	2.26	0.49
1:B:191:LYS:HG2	1:B:210:LEU:HD23	1.94	0.48
1:B:450:GLN:HE22	1:B:471:GLU:HG3	1.77	0.48
1:A:515:ARG:HA	1:A:521:GLY:HA2	1.95	0.48
1:B:110:THR:HA	1:B:116:LYS:O	2.14	0.48
1:B:256:ASP:HB3	1:B:261:ASN:HD22	1.79	0.48
1:B:473:LEU:HD11	1:B:505:LEU:HD11	1.95	0.47
1:A:256:ASP:CG	1:A:261:ASN:HD22	2.22	0.47
1:B:384:TRP:O	1:B:388:ILE:HG13	2.14	0.47
1:B:166:LEU:HD11	1:B:178:VAL:HG13	1.97	0.47
1:B:514:THR:HG23	1:B:524:SER:HB2	1.96	0.47
1:A:248:LEU:HB3	1:A:269:TYR:HE1	1.80	0.47
1:A:303:GLU:HB3	1:A:305:LYS:NZ	2.30	0.47
1:A:202:ASP:OD1	1:A:202:ASP:N	2.48	0.46
1:A:505:LEU:HD23	1:A:508:LEU:HD12	1.98	0.46
1:A:599:PHE:O	1:A:602:THR:HG22	2.14	0.46
1:B:104:LEU:HD12	1:B:141:ASP:HB2	1.96	0.46
1:B:248:LEU:HB3	1:B:269:TYR:HE1	1.79	0.46
1:B:349:TRP:CG	1:B:682:ARG:HE	2.34	0.46
1:B:188:VAL:HG22	1:B:233:GLN:NE2	2.30	0.46
1:B:349:TRP:CE3	1:B:682:ARG:HG3	2.50	0.46
1:A:328:PHE:HE2	1:A:700:ARG:HD2	1.81	0.46
1:B:302:LYS:HA	1:B:362:SER:HB3	1.98	0.46
1:A:315:ARG:NH1	1:A:324:LEU:O	2.29	0.45
1:B:139:GLU:OE1	1:B:142:ARG:NH2	2.43	0.45
1:B:208:GLY:O	1:B:213:SER:HB2	2.17	0.45
1:A:119:LEU:HD22	1:A:138:LEU:HD21	1.99	0.45
1:A:406:THR:OG1	1:A:507:TRP:NE1	2.44	0.45
1:B:470:PHE:CE2	1:B:512:TYR:HB2	2.52	0.45
1:A:111:GLU:HB2	1:A:116:LYS:HB3	1.99	0.45
1:A:364:SER:O	1:A:367:ARG:HG3	2.17	0.45
1:A:126:LEU:HD23	1:A:131:ASN:HB2	1.99	0.44
1:A:206:TYR:OH	1:A:245:ASN:ND2	2.42	0.44
1:B:123:VAL:O	1:B:126:LEU:HG	2.17	0.44
1:A:182:VAL:HG11	1:A:230:ASN:HD22	1.82	0.44
1:A:222:ASP:OD1	1:A:222:ASP:N	2.49	0.44
1:B:542:TYR:OH	1:B:602:THR:CB	2.66	0.44
1:A:152:ASN:OD1	1:A:187:ASN:N	2.51	0.44
1:B:511:LEU:O	1:B:514:THR:HG22	2.18	0.44
1:B:470:PHE:CZ	1:B:512:TYR:HB2	2.53	0.43
1:B:238:GLN:NE2	1:B:283:VAL:HG21	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:GLY:HA3	1:B:192:ALA:HA	2.00	0.43
1:A:182:VAL:HG11	1:A:230:ASN:ND2	2.33	0.43
1:A:380:LEU:HD13	1:A:663:ALA:HB2	2.00	0.42
1:A:672:GLY:O	1:A:673:TYR:HB3	2.19	0.42
1:B:328:PHE:CZ	1:B:689:VAL:HG11	2.54	0.42
1:A:160:TYR:CE1	1:B:725:PRO:HG2	2.54	0.42
1:B:313:LEU:HB3	1:B:375:PRO:HG3	2.02	0.42
1:B:322:GLN:O	1:B:325:SER:OG	2.38	0.42
1:A:319:ALA:HB1	1:A:320:PRO:HD2	2.02	0.41
1:B:253:MET:SD	1:B:299:LEU:HD21	2.60	0.41
1:A:93:LEU:HD23	1:A:140:ILE:HD13	2.02	0.41
1:A:685:VAL:HG21	1:A:704:ARG:HH12	1.84	0.41
1:B:349:TRP:CD1	1:B:682:ARG:HH21	2.39	0.41
1:A:245:ASN:HB3	1:A:249:HIS:HB2	2.01	0.41
1:A:297:LEU:HD22	1:A:354:VAL:HG21	2.03	0.41
1:A:173:ARG:HG2	1:A:220:GLN:HE21	1.86	0.41
1:A:238:GLN:HE21	1:A:283:VAL:HG11	1.84	0.41
1:A:238:GLN:HE21	1:A:283:VAL:HG21	1.85	0.41
1:A:331:TRP:CE3	1:A:725:PRO:HB3	2.56	0.41
1:A:462:TRP:O	1:A:462:TRP:CG	2.74	0.41
1:A:694:ASP:OD1	1:A:695:GLY:N	2.53	0.41
1:B:450:GLN:NE2	1:B:471:GLU:HG3	2.35	0.41
1:A:129:GLY:HA3	1:A:176:GLN:NE2	2.36	0.40
1:A:209:GLU:OE2	1:A:242:SER:OG	2.33	0.40
1:A:248:LEU:HB3	1:A:269:TYR:CE1	2.57	0.40
1:B:273:LEU:HB3	1:B:324:LEU:HD13	2.03	0.40
1:B:391:PHE:HB2	1:B:669:MET:HE1	2.02	0.40
1:B:224:VAL:O	1:B:228:LEU:HG	2.22	0.40

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	598/776 (77%)	582 (97%)	16 (3%)	0	100	100
1	B	589/776 (76%)	569 (97%)	19 (3%)	1 (0%)	43	74
All	All	1187/1552 (76%)	1151 (97%)	35 (3%)	1 (0%)	48	80

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	465	PHE

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	367/666 (55%)	365 (100%)	2 (0%)	81	82
1	B	361/666 (54%)	360 (100%)	1 (0%)	86	85
All	All	728/1332 (55%)	725 (100%)	3 (0%)	84	83

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	514	THR
1	A	602	THR
1	B	514	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	ASN
1	A	238	GLN
1	B	238	GLN
1	B	261	ASN
1	B	381	GLN
1	B	450	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 oligosaccharides 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	606/776 (78%)	0.89	73 (12%)	9 12	36, 91, 186, 217	0
1	B	599/776 (77%)	0.90	79 (13%)	7 10	41, 101, 183, 213	0
All	All	1205/1552 (77%)	0.89	152 (12%)	8 11	36, 96, 185, 217	0

All (152) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	605	MET	8.7
1	B	456	ARG	8.6
1	A	668	GLU	8.3
1	B	318	SER	8.2
1	A	430	SER	7.7
1	B	452	TRP	7.6
1	A	604	GLY	7.2
1	A	452	TRP	6.8
1	A	723	GLU	6.6
1	B	69	ASP	6.3
1	B	672	GLY	6.1
1	B	639	LEU	5.7
1	A	398	TYR	5.4
1	B	640	ILE	5.4
1	A	429	ASN	5.3
1	B	148	GLN	5.2
1	B	643	MET	5.2
1	A	665	SER	5.2
1	B	453	TYR	5.0
1	A	664	ILE	4.7
1	A	453	TYR	4.7
1	A	669	MET	4.5
1	A	427	ALA	4.5
1	B	321	CYS	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	601	PHE	4.3
1	A	512	TYR	4.2
1	A	643	MET	4.1
1	B	669	MET	4.1
1	A	671	ASN	3.9
1	B	429	ASN	3.9
1	B	642	LEU	3.8
1	A	509	ASN	3.7
1	A	602	THR	3.7
1	B	320	PRO	3.7
1	B	667	LEU	3.6
1	B	346	VAL	3.6
1	B	671	ASN	3.5
1	A	727	GLY	3.4
1	B	450	GLN	3.3
1	A	504	VAL	3.3
1	B	652	THR	3.3
1	A	318	SER	3.3
1	B	183	GLU	3.3
1	B	674	TRP	3.2
1	A	449	GLY	3.2
1	B	552	ALA	3.2
1	B	433	LEU	3.2
1	A	69	ASP	3.1
1	B	398	TYR	3.0
1	B	670	GLU	3.0
1	A	548	GLY	3.0
1	B	169	ALA	3.0
1	A	71	ASN	2.9
1	B	634	LEU	2.9
1	B	406	THR	2.9
1	B	144	SER	2.9
1	B	604	GLY	2.9
1	B	668	GLU	2.8
1	A	189	HIS	2.8
1	B	727	GLY	2.8
1	B	633	VAL	2.8
1	B	724	GLU	2.8
1	A	157	ASP	2.8
1	A	320	PRO	2.8
1	B	623	LEU	2.8
1	A	400	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	595	SER	2.8
1	A	456	ARG	2.8
1	A	601	PHE	2.8
1	B	449	GLY	2.8
1	B	180	LEU	2.7
1	A	551	VAL	2.7
1	A	450	GLN	2.7
1	A	600	LYS	2.7
1	A	676	CYS	2.7
1	B	176	GLN	2.7
1	A	433	LEU	2.7
1	B	522	ILE	2.7
1	A	541	VAL	2.6
1	B	600	LYS	2.6
1	A	638	MET	2.6
1	A	599	PHE	2.6
1	A	616	PHE	2.6
1	A	200	ASN	2.6
1	A	448	LEU	2.6
1	B	458	ARG	2.6
1	A	203	THR	2.5
1	A	646	THR	2.5
1	B	514	THR	2.5
1	A	598	LEU	2.5
1	A	293	GLY	2.5
1	B	515	ARG	2.5
1	A	399	LEU	2.5
1	A	590	SER	2.5
1	A	360	PHE	2.4
1	B	457	ARG	2.4
1	B	488	CYS	2.4
1	A	675	TRP	2.4
1	A	515	ARG	2.4
1	B	177	CYS	2.4
1	B	636	LEU	2.4
1	B	360	PHE	2.4
1	B	344	ALA	2.4
1	B	507	TRP	2.4
1	B	343	LEU	2.4
1	B	480	LEU	2.4
1	B	467	ASP	2.4
1	A	634	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	408	VAL	2.3
1	A	238	GLN	2.3
1	B	317	PHE	2.3
1	A	404	ILE	2.3
1	B	519	HIS	2.3
1	A	603	ILE	2.3
1	B	599	PHE	2.3
1	A	724	GLU	2.3
1	A	127	GLN	2.3
1	B	723	GLU	2.3
1	A	690	GLY	2.3
1	B	637	ASN	2.3
1	A	474	PHE	2.3
1	A	467	ASP	2.3
1	A	635	LEU	2.2
1	A	366	HIS	2.2
1	B	462	TRP	2.2
1	A	673	TYR	2.2
1	B	542	TYR	2.2
1	B	665	SER	2.2
1	A	645	GLU	2.2
1	A	550	ALA	2.1
1	A	501	SER	2.1
1	A	401	TYR	2.1
1	B	445	TYR	2.1
1	A	505	LEU	2.1
1	B	543	LEU	2.1
1	B	157	ASP	2.1
1	B	326	ARG	2.1
1	A	99	ARG	2.1
1	B	550	ALA	2.1
1	B	630	LEU	2.1
1	B	232	HIS	2.1
1	B	673	TYR	2.1
1	A	111	GLU	2.1
1	B	152	ASN	2.1
1	B	184	ASN	2.1
1	B	208	GLY	2.0
1	B	430	SER	2.0
1	A	722	CYS	2.0
1	B	512	TYR	2.0
1	B	405	PHE	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	145	GLY	2.0
1	A	667	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
2	CA	B	801	1/1	0.85	0.11	145,145,145,145	1

6.5 Other polymers [i](#)

There are no such residues in this entry.