Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 09:12 pm BST

PDB ID	:	2Z6H
Title	:	Crystal Structure of Beta-Catenin Armadillo Repeat Region and Its C-Terminal domain
Authors	:	Xing, Y.; Takemaru, K.; Liu, J.; Zheng, J.; Moon, R.; Xu, W.
Deposited on	:	2007-08-01
Resolution	:	2.20 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 following versions of software and data (see references ⓘ) were used in the production of this report:

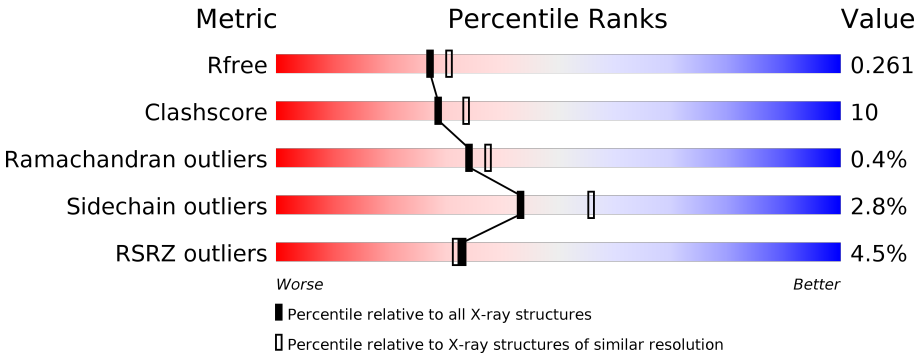
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 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	644	4% 65% 17% • 17%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4058 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 Catenin beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	533	4030	2537	728	739	26	0	0	0

- Molecule 2 is water.

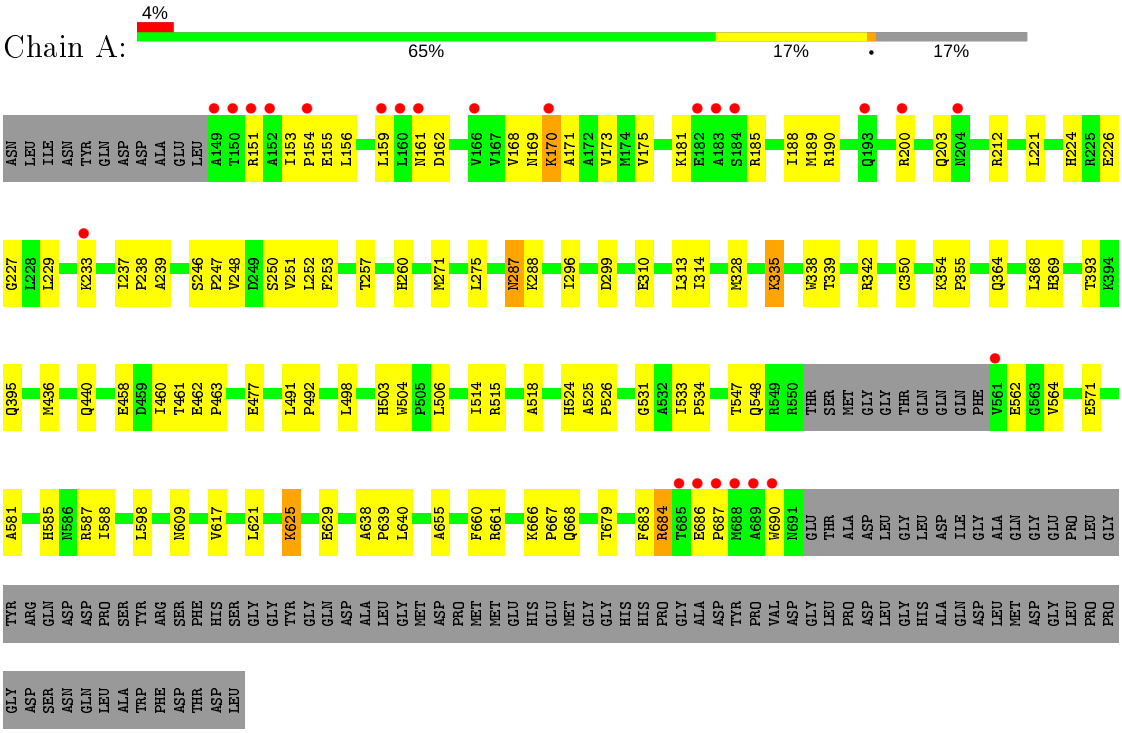
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	28	28	28	0	0

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 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: Catenin beta-1

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α, β, γ	104.23Å 104.23Å 259.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.69 – 2.20 48.69 – 2.20	Depositor EDS
% Data completeness (in resolution range)	80.9 (48.69-2.20) 88.3 (48.69-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	2.46 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.213 , 0.249 0.224 , 0.261	Depositor DCC
R_{free} test set	2854 reflections (8.01%)	wwPDB-VP
Wilson B-factor (Å^2)	30.8	Xtriage
Anisotropy	0.285	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 42.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	4058	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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

5.1 Standard geometry

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	0.33	0/4089	0.54	0/5556

There are no bond length outliers.

There are no bond angle outliers.

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	4030	0	4140	79	0
2	A	28	0	0	0	0
All	All	4058	0	4140	79	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 10.

All (79) 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:161:ASN:HD21	1:A:200:ARG:HH21	1.16	0.88
1:A:161:ASN:ND2	1:A:200:ARG:HH21	1.72	0.88
1:A:189:MET:HB3	1:A:221:LEU:HD22	1.54	0.87
1:A:229:LEU:HG	1:A:233:LYS:HE3	1.55	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ARG:HH11	1:A:224:HIS:HE1	1.25	0.81
1:A:161:ASN:HD21	1:A:200:ARG:NH2	1.82	0.78
1:A:169:ASN:HD22	1:A:170:LYS:HE3	1.53	0.74
1:A:354:LYS:HB3	1:A:355:PRO:HD3	1.70	0.73
1:A:687:PRO:HG2	1:A:690:TRP:HB2	1.72	0.72
1:A:364:GLN:HE22	1:A:395:GLN:HE21	1.38	0.71
1:A:503:HIS:HB3	1:A:562:GLU:OE1	1.93	0.69
1:A:151:ARG:HD3	1:A:151:ARG:O	1.94	0.68
1:A:237:ILE:HB	1:A:238:PRO:HD3	1.77	0.66
1:A:328:MET:O	1:A:369:HIS:HE1	1.78	0.66
1:A:666:LYS:HB3	1:A:667:PRO:HD2	1.78	0.66
1:A:260:HIS:HE1	1:A:299:ASP:OD1	1.81	0.64
1:A:169:ASN:HD22	1:A:170:LYS:CE	2.16	0.59
1:A:156:LEU:HD23	1:A:159:LEU:HD12	1.84	0.58
1:A:525:ALA:HB3	1:A:526:PRO:HD3	1.86	0.58
1:A:185:ARG:HH11	1:A:224:HIS:CE1	2.13	0.57
1:A:625:LYS:HE2	1:A:629:GLU:HG3	1.87	0.57
1:A:625:LYS:HE2	1:A:629:GLU:CG	2.35	0.57
1:A:504:TRP:HB3	1:A:564:VAL:HG21	1.85	0.56
1:A:328:MET:O	1:A:369:HIS:CE1	2.57	0.56
1:A:169:ASN:ND2	1:A:170:LYS:HE3	2.20	0.56
1:A:153:ILE:N	1:A:154:PRO:HD2	2.20	0.55
1:A:287:ASN:HD22	1:A:288:LYS:N	2.05	0.55
1:A:684:ARG:HD2	1:A:684:ARG:N	2.22	0.54
1:A:156:LEU:HA	1:A:159:LEU:HD12	1.89	0.54
1:A:212:ARG:HD3	1:A:250:SER:OG	2.07	0.54
1:A:189:MET:CE	1:A:226:GLU:HB2	2.38	0.54
1:A:229:LEU:CG	1:A:233:LYS:HE3	2.33	0.54
1:A:221:LEU:O	1:A:227:GLY:HA3	2.09	0.53
1:A:257:THR:HG22	1:A:296:ILE:HD13	1.89	0.53
1:A:462:GLU:HB3	1:A:463:PRO:HD3	1.91	0.53
1:A:460:ILE:O	1:A:463:PRO:HD2	2.09	0.52
1:A:625:LYS:O	1:A:629:GLU:HG3	2.10	0.52
1:A:533:ILE:HB	1:A:534:PRO:HD3	1.92	0.51
1:A:185:ARG:NH1	1:A:224:HIS:HE1	2.02	0.51
1:A:436:MET:SD	1:A:477:GLU:HG2	2.51	0.51
1:A:436:MET:O	1:A:440:GLN:HG3	2.10	0.51
1:A:547:THR:HG22	1:A:548:GLN:HE21	1.76	0.50
1:A:679:THR:HG23	1:A:683:PHE:CE2	2.46	0.50
1:A:504:TRP:HB3	1:A:564:VAL:CG2	2.42	0.50
1:A:571:GLU:OE1	1:A:609:ASN:HB3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:MET:HE2	1:A:226:GLU:HB2	1.94	0.48
1:A:188:ILE:C	1:A:190:ARG:H	2.17	0.48
1:A:271:MET:O	1:A:275:LEU:HG	2.13	0.48
1:A:338:TRP:CZ2	1:A:342:ARG:HD2	2.49	0.47
1:A:364:GLN:O	1:A:368:LEU:HD13	2.16	0.47
1:A:169:ASN:O	1:A:173:VAL:HG23	2.15	0.46
1:A:364:GLN:HE22	1:A:395:GLN:NE2	2.12	0.46
1:A:458:GLU:HA	1:A:461:THR:OG1	2.16	0.46
1:A:518:ALA:O	1:A:524:HIS:HE1	1.99	0.46
1:A:491:LEU:HB2	1:A:492:PRO:HD3	1.98	0.45
1:A:335:LYS:O	1:A:339:THR:HG23	2.16	0.45
1:A:458:GLU:HG2	1:A:506:LEU:HD22	1.98	0.45
1:A:287:ASN:C	1:A:287:ASN:HD22	2.19	0.44
1:A:171:ALA:O	1:A:175:VAL:HG23	2.18	0.44
1:A:625:LYS:HE3	1:A:625:LYS:HA	1.98	0.44
1:A:585:HIS:O	1:A:588:ILE:HG22	2.18	0.44
1:A:181:LYS:HB3	1:A:181:LYS:HE3	1.83	0.44
1:A:515:ARG:C	1:A:515:ARG:HD3	2.39	0.43
1:A:393:THR:HG22	1:A:393:THR:O	2.19	0.43
1:A:640:LEU:HD22	1:A:655:ALA:HA	2.01	0.43
1:A:155:GLU:O	1:A:159:LEU:HG	2.18	0.42
1:A:203:GLN:NE2	1:A:239:ALA:HA	2.34	0.42
1:A:260:HIS:CE1	1:A:299:ASP:OD1	2.67	0.42
1:A:498:LEU:HD11	1:A:514:ILE:CD1	2.50	0.41
1:A:248:VAL:HG13	1:A:251:VAL:HB	2.02	0.41
1:A:310:GLU:O	1:A:314:ILE:HG13	2.20	0.41
1:A:581:ALA:O	1:A:587:ARG:CG	2.68	0.41
1:A:617:VAL:O	1:A:621:LEU:HG	2.21	0.41
1:A:168:VAL:HG13	1:A:169:ASN:N	2.36	0.41
1:A:638:ALA:HB3	1:A:639:PRO:CD	2.50	0.41
1:A:338:TRP:O	1:A:342:ARG:HG3	2.20	0.41
1:A:246:SER:HA	1:A:247:PRO:HD3	1.82	0.40
1:A:531:GLY:O	1:A:534:PRO:HD2	2.20	0.40
1:A:253:PHE:O	1:A:257:THR:HG23	2.20	0.40

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	529/644 (82%)	508 (96%)	19 (4%)	2 (0%)	34 37

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	686	GLU
1	A	162	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	A	432/532 (81%)	420 (97%)	12 (3%)	43 56

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	170	LYS
1	A	252	LEU
1	A	287	ASN
1	A	313	LEU
1	A	335	LYS
1	A	350	CYS
1	A	598	LEU
1	A	625	LYS
1	A	660	PHE

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Mol	Chain	Res	Type
1	A	661	ARG
1	A	668	GLN
1	A	684	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	161	ASN
1	A	169	ASN
1	A	186	HIS
1	A	193	GLN
1	A	203	GLN
1	A	224	HIS
1	A	260	HIS
1	A	287	ASN
1	A	326	ASN
1	A	369	HIS
1	A	395	GLN
1	A	407	GLN
1	A	440	GLN
1	A	524	HIS
1	A	530	GLN
1	A	538	GLN
1	A	548	GLN
1	A	594	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 carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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, 95^{th} 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	533/644 (82%)	0.08	24 (4%) 33 32	17, 37, 71, 85	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	686	GLU	7.1
1	A	149	ALA	7.1
1	A	152	ALA	6.1
1	A	561	VAL	5.3
1	A	687	PRO	5.1
1	A	150	THR	4.7
1	A	151	ARG	4.0
1	A	689	ALA	3.5
1	A	690	TRP	3.0
1	A	184	SER	3.0
1	A	204	ASN	2.9
1	A	161	ASN	2.8
1	A	166	VAL	2.8
1	A	183	ALA	2.7
1	A	685	THR	2.5
1	A	688	MET	2.5
1	A	160	LEU	2.3
1	A	233	LYS	2.3
1	A	193	GLN	2.2
1	A	154	PRO	2.2
1	A	159	LEU	2.1
1	A	182	GLU	2.0
1	A	200	ARG	2.0
1	A	170	LYS	2.0

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 carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.