



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 02:15 am BST

PDB ID : 6BX7
Title : Crystal Structure of Human Protocadherin-1 EC1-4
Authors : Modak, D.; Sotomayor, M.
Deposited on : 2017-12-17
Resolution : 2.85 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

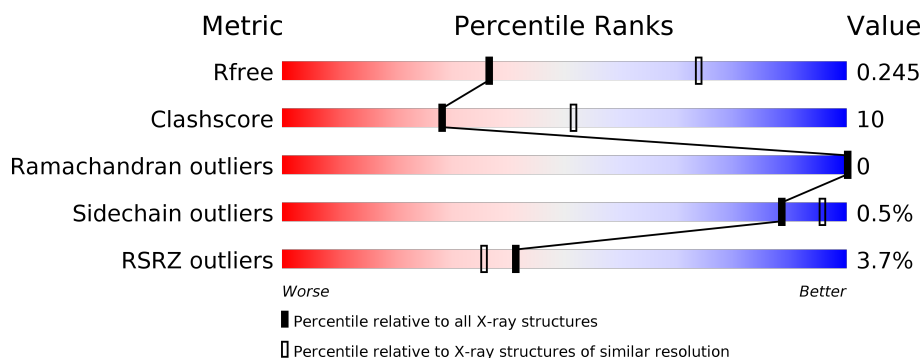
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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 $\leq 5\%$. 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	455	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>16%</div> <div>• 6%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3339 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 Protocadherin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	S	0	0	0
			3288	2039	563	680	6			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q08174
A	447	LEU	-	expression tag	UNP Q08174
A	448	GLU	-	expression tag	UNP Q08174
A	449	HIS	-	expression tag	UNP Q08174
A	450	HIS	-	expression tag	UNP Q08174
A	451	HIS	-	expression tag	UNP Q08174
A	452	HIS	-	expression tag	UNP Q08174
A	453	HIS	-	expression tag	UNP Q08174
A	454	HIS	-	expression tag	UNP Q08174

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	Ca	0	0
			11	11		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total	O	0	0
			25	25		

- Molecule 1: Protocadherin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	146.98 Å 146.98 Å 147.69 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.26 – 2.85 48.21 – 2.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.26-2.85) 99.9 (48.21-2.85)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.86 Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.221 , 0.243 0.224 , 0.245	Depositor DCC
R_{free} test set	1120 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3339	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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, EPE

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	0.56	0/3338	0.79	2/4548 (0.0%)

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	A	0	5

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	423	ASP	CB-CA-C	-7.73	94.94	110.40
1	A	277	ARG	CB-CA-C	-5.66	99.08	110.40

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	185	ARG	Sidechain
1	A	210	ARG	Sidechain
1	A	277	ARG	Sidechain
1	A	38	GLY	Peptide
1	A	60	ARG	Sidechain

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	3288	0	3224	64	0
2	A	15	0	18	1	0
3	A	11	0	0	0	0
4	A	25	0	0	0	0
All	All	3339	0	3242	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 10.

All (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:A:349:GLU:OE1	1:A:409:TYR:HB2	1.53	1.08
1:A:16:LEU:HA	1:A:52:PHE:CD1	1.92	1.03
1:A:349:GLU:OE1	1:A:441:ASP:HA	1.75	0.84
1:A:409:TYR:HA	1:A:412:VAL:CG2	2.07	0.84
1:A:124:THR:HG23	1:A:177:VAL:HG21	1.64	0.77
1:A:347:ILE:O	1:A:439:VAL:HA	1.90	0.71
1:A:423:ASP:HB3	1:A:425:GLY:H	1.56	0.71
1:A:349:GLU:OE1	1:A:409:TYR:CB	2.37	0.71
1:A:354:GLU:HB2	1:A:402:GLN:HE21	1.57	0.69
1:A:333:ARG:NH2	1:A:359:LEU:HD23	2.10	0.67
1:A:16:LEU:HA	1:A:52:PHE:CE1	2.31	0.66
1:A:60:ARG:NH1	1:A:61:GLU:OE2	2.33	0.61
1:A:415:TYR:HB2	1:A:437:VAL:HG22	1.82	0.60
1:A:16:LEU:HB2	1:A:52:PHE:CE1	2.36	0.60
1:A:347:ILE:HD11	1:A:351:VAL:HG11	1.84	0.60
1:A:409:TYR:HA	1:A:412:VAL:HG21	1.84	0.59
1:A:166:GLU:OE1	1:A:171:LYS:HG2	2.03	0.58
1:A:94:LEU:HD12	1:A:94:LEU:C	2.24	0.58
1:A:168:GLN:O	1:A:169:GLU:HG2	2.02	0.58
1:A:19:SER:HB3	1:A:50:ASP:HB3	1.87	0.57
1:A:83:ILE:HD13	1:A:95:LEU:HD23	1.87	0.56
1:A:409:TYR:CG	1:A:441:ASP:HB2	2.42	0.54
1:A:353:GLU:O	1:A:354:GLU:HG2	2.08	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:GLU:HG3	2:A:501:EPE:H62	1.92	0.52
1:A:51:ILE:O	1:A:52:PHE:CD1	2.63	0.52
1:A:77:LEU:C	1:A:77:LEU:HD23	2.30	0.52
1:A:349:GLU:CD	1:A:442:VAL:H	2.13	0.51
1:A:163:GLN:HB2	1:A:176:ILE:HD11	1.93	0.51
1:A:354:GLU:HB2	1:A:402:GLN:NE2	2.26	0.51
1:A:77:LEU:HD23	1:A:78:GLU:N	2.25	0.51
1:A:274:ARG:HG2	1:A:274:ARG:HH21	1.74	0.51
1:A:423:ASP:OD2	1:A:429:LEU:HB2	2.11	0.51
1:A:335:ILE:CG2	1:A:359:LEU:HD22	2.41	0.51
1:A:51:ILE:O	1:A:52:PHE:HD1	1.94	0.50
1:A:79:PHE:CZ	1:A:97:GLY:HA3	2.47	0.50
1:A:20:LEU:HD22	1:A:33:TYR:CE1	2.48	0.49
1:A:412:VAL:HG23	1:A:439:VAL:HG21	1.94	0.49
1:A:245:VAL:HG21	1:A:300:VAL:HG11	1.94	0.48
1:A:409:TYR:HA	1:A:412:VAL:HG22	1.95	0.48
1:A:135:ALA:HB3	1:A:195:VAL:HG21	1.95	0.48
1:A:19:SER:HB3	1:A:50:ASP:CB	2.43	0.48
1:A:16:LEU:HA	1:A:52:PHE:HD1	1.69	0.47
1:A:383:PHE:CE1	1:A:403:THR:HG22	2.49	0.47
1:A:259:GLU:OE2	1:A:305:ARG:NH2	2.47	0.47
1:A:51:ILE:C	1:A:52:PHE:CD1	2.88	0.47
1:A:60:ARG:NH2	1:A:140:ALA:HB2	2.30	0.47
1:A:409:TYR:CD2	1:A:441:ASP:HB2	2.51	0.46
1:A:166:GLU:OE2	1:A:171:LYS:HE3	2.16	0.46
1:A:33:TYR:CE1	1:A:83:ILE:HD11	2.51	0.45
1:A:3:VAL:HG22	1:A:4:VAL:H	1.81	0.45
1:A:412:VAL:CG2	1:A:439:VAL:HG21	2.47	0.45
1:A:335:ILE:HG23	1:A:359:LEU:HB2	2.00	0.44
1:A:257:GLU:O	1:A:304:ASP:HA	2.18	0.43
1:A:8:PRO:HA	1:A:102:GLN:HB2	2.01	0.43
1:A:17:ILE:HG12	1:A:51:ILE:O	2.18	0.42
1:A:16:LEU:CB	1:A:52:PHE:CE1	3.01	0.42
1:A:221:PRO:HA	1:A:248:ASN:O	2.20	0.42
1:A:19:SER:HB3	1:A:50:ASP:HA	2.01	0.42
1:A:413:LYS:N	1:A:439:VAL:HG22	2.35	0.41
1:A:79:PHE:CE1	1:A:97:GLY:HA3	2.55	0.41
1:A:12:PRO:O	1:A:15:THR:OG1	2.37	0.41
1:A:362:VAL:CG1	1:A:373:VAL:HG11	2.51	0.41
1:A:408:ASP:C	1:A:410:GLU:H	2.24	0.41
1:A:420:VAL:HG22	1:A:432:THR:HG22	2.03	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	423/455 (93%)	390 (92%)	33 (8%)	0	100	100

There are no Ramachandran outliers to report.

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	370/392 (94%)	368 (100%)	2 (0%)	88	96

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	271	ARG
1	A	349	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

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 ⓘ

Of 12 ligands modelled in this entry, 11 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	EPE	A	501	-	15,15,15	1.94	1 (6%)	18,20,20	1.63	2 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EPE	A	501	-	-	2/9/19/19	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	EPE	C10-S	-6.99	1.67	1.77

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	EPE	O2S-S-C10	4.59	112.44	106.92
2	A	501	EPE	O3S-S-O2S	-2.14	106.04	111.27

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	EPE	C8-C7-N4-C3
2	A	501	EPE	C9-C10-S-O2S

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	EPE	1	0

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	A	429/455 (94%)	0.16	16 (3%) 41 36	39, 67, 136, 163	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	337	LEU	4.2
1	A	84	THR	4.0
1	A	350	ASP	3.1
1	A	83	ILE	2.8
1	A	5	TYR	2.8
1	A	17	ILE	2.5
1	A	52	PHE	2.4
1	A	16	LEU	2.4
1	A	407	LEU	2.4
1	A	95	LEU	2.4
1	A	47	LYS	2.2
1	A	91	SER	2.2
1	A	64	ARG	2.1
1	A	414	ASP	2.1
1	A	32	LEU	2.0
1	A	35	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 carbohydrates 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(Å ²)	Q<0.9
3	CA	A	511	1/1	0.69	0.08	110,110,110,110	0
2	EPE	A	501	15/15	0.93	0.27	97,106,112,112	0
3	CA	A	512	1/1	0.94	0.17	62,62,62,62	0
3	CA	A	502	1/1	0.96	0.17	75,75,75,75	0
3	CA	A	508	1/1	0.97	0.15	66,66,66,66	0
3	CA	A	510	1/1	0.98	0.14	53,53,53,53	0
3	CA	A	506	1/1	0.99	0.17	45,45,45,45	0
3	CA	A	505	1/1	0.99	0.14	53,53,53,53	0
3	CA	A	503	1/1	0.99	0.17	60,60,60,60	0
3	CA	A	504	1/1	0.99	0.19	50,50,50,50	0
3	CA	A	509	1/1	1.00	0.11	49,49,49,49	0
3	CA	A	507	1/1	1.00	0.14	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.