



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 30, 2020 – 04:01 PM EST

PDB ID : 7KP4  
Title : Crystal structure of human claudin-4 in complex with Clostridium perfringens enterotoxin C-terminal domain  
Authors : Vecchio, A.J.; Stroud, R.M.  
Deposited on : 2020-11-10  
Resolution : 3.37 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](mailto: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.

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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	:	2.14.6
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)	:	2.14.6

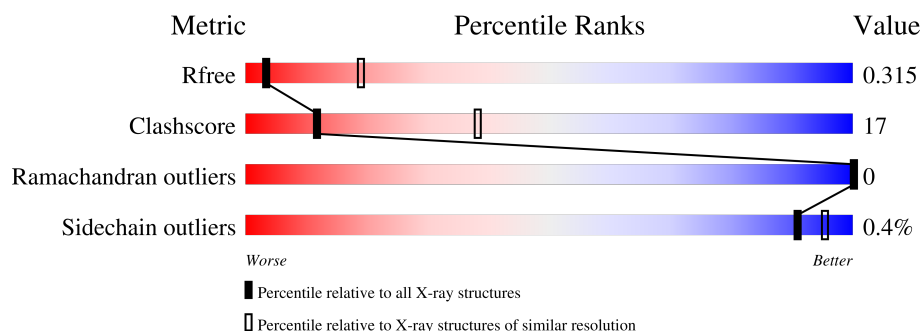
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.37 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	1691 (3.46-3.30)
Clashscore	141614	1762 (3.46-3.30)
Ramachandran outliers	138981	1732 (3.46-3.30)
Sidechain outliers	138945	1731 (3.46-3.30)

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  $\geq 3$ , 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  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	214	
2	B	134	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2375 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 Claudin-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	182	Total	C	N	O	S	0	0	0
			1318	854	216	231	17			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	210	GLY	-	expression tag	UNP O14493
A	211	LEU	-	expression tag	UNP O14493
A	212	VAL	-	expression tag	UNP O14493
A	213	PRO	-	expression tag	UNP O14493
A	214	ARG	-	expression tag	UNP O14493

- Molecule 2 is a protein called Heat-labile enterotoxin B chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	134	Total	C	N	O	S	0	0	0
			1057	673	174	208	2			

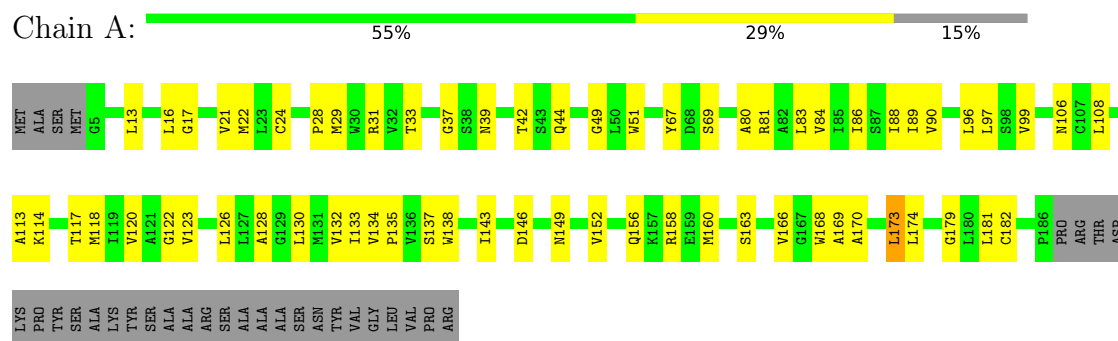
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	191	MET	-	initiating methionine	UNP P01558
B	320	GLY	-	expression tag	UNP P01558
B	321	LEU	-	expression tag	UNP P01558
B	322	VAL	-	expression tag	UNP P01558
B	323	PRO	-	expression tag	UNP P01558
B	324	ARG	-	expression tag	UNP P01558

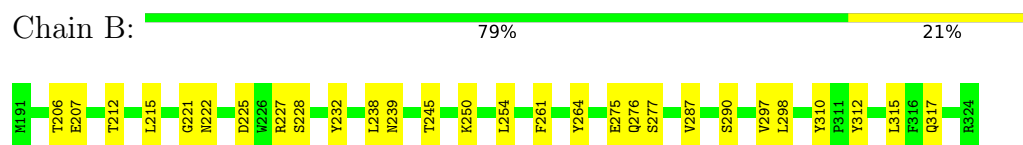
### 3 Residue-property plots

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. 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. 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: Claudin-4



#### • Molecule 2: Heat-labile enterotoxin B chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.36Å 116.23Å 118.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.81 – 3.37 59.02 – 3.25	Depositor EDS
% Data completeness (in resolution range)	91.8 (44.81-3.37) 74.7 (59.02-3.25)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.11 (at 3.26Å)	Xtriage
Refinement program	PHENIX (1.14 _ 3260: ???)	Depositor
R, $R_{free}$	0.294 , 0.296 0.307 , 0.315	Depositor DCC
$R_{free}$ test set	1185 reflections (7.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	111.1	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.19 , 55.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.054 for -h,l,k	Xtriage
$F_o, F_c$ correlation	0.79	EDS
Total number of atoms	2375	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	170.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.96% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

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	0.30	0/1341	0.53	0/1832
2	B	0.31	0/1079	0.50	0/1467
All	All	0.30	0/2420	0.52	0/3299

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	1318	0	1357	69	0
2	B	1057	0	1026	18	0
All	All	2375	0	2383	83	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 17.

All (83) 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:17:GLY:N	1:A:173:LEU:HD11	1.44	1.32
1:A:17:GLY:HA2	1:A:173:LEU:HD21	1.20	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:GLY:N	1:A:173:LEU:CD1	2.24	0.99
1:A:16:LEU:C	1:A:173:LEU:HD11	1.86	0.96
1:A:17:GLY:HA2	1:A:173:LEU:CD2	1.98	0.93
1:A:170:ALA:HA	1:A:173:LEU:CD2	2.00	0.91
1:A:17:GLY:CA	1:A:173:LEU:HD11	2.06	0.85
1:A:17:GLY:CA	1:A:173:LEU:HD21	2.11	0.71
1:A:17:GLY:CA	1:A:173:LEU:CG	2.69	0.71
1:A:170:ALA:O	1:A:173:LEU:HD23	1.91	0.71
1:A:170:ALA:HA	1:A:173:LEU:HD21	1.77	0.67
1:A:122:GLY:HA3	1:A:182:CYS:HB2	1.78	0.65
1:A:17:GLY:HA3	1:A:173:LEU:HG	1.80	0.64
1:A:17:GLY:CA	1:A:173:LEU:CD1	2.72	0.62
1:A:17:GLY:CA	1:A:173:LEU:CD2	2.76	0.62
1:A:170:ALA:HA	1:A:173:LEU:HD23	1.82	0.61
1:A:146:ASP:OD1	2:B:227:ARG:NH1	2.31	0.61
1:A:130:LEU:HA	1:A:133:ILE:HG22	1.83	0.60
1:A:21:VAL:HG22	1:A:132:VAL:HG11	1.84	0.60
2:B:264:TYR:HB2	2:B:297:VAL:HG12	1.84	0.59
1:A:24:CYS:SG	1:A:163:SER:O	2.56	0.58
1:A:17:GLY:HA2	1:A:173:LEU:CG	2.34	0.57
1:A:17:GLY:CA	1:A:173:LEU:HG	2.34	0.56
1:A:170:ALA:CA	1:A:173:LEU:HD23	2.36	0.56
1:A:120:VAL:HA	1:A:123:VAL:HG12	1.87	0.56
1:A:128:ALA:O	1:A:132:VAL:HG23	2.06	0.55
2:B:261:PHE:CE1	2:B:277:SER:HB3	2.41	0.55
2:B:275:GLU:HG2	2:B:287:VAL:HG21	1.88	0.55
1:A:123:VAL:HA	1:A:126:LEU:HD12	1.90	0.54
1:A:83:LEU:HD23	1:A:135:PRO:HB3	1.90	0.54
1:A:22:MET:HG2	1:A:88:ILE:HD12	1.90	0.53
1:A:86:ILE:HD13	1:A:89:ILE:HD11	1.91	0.53
1:A:170:ALA:C	1:A:173:LEU:HD23	2.29	0.52
1:A:132:VAL:O	1:A:135:PRO:HD2	2.10	0.52
1:A:42:THR:HG21	2:B:222:ASN:ND2	2.25	0.51
1:A:90:VAL:HG12	1:A:128:ALA:HB2	1.91	0.51
1:A:39:ASN:OD1	2:B:317:GLN:NE2	2.36	0.51
1:A:84:VAL:O	1:A:88:ILE:HG12	2.10	0.51
1:A:113:ALA:O	1:A:117:THR:HG23	2.11	0.50
1:A:21:VAL:HG21	1:A:132:VAL:HG21	1.94	0.49
1:A:17:GLY:O	1:A:21:VAL:HG23	2.13	0.49
1:A:134:VAL:O	1:A:138:TRP:HB2	2.13	0.49
1:A:42:THR:OG1	1:A:44:GLN:NE2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:PRO:HA	1:A:49:GLY:HA3	1.96	0.48
2:B:250:LYS:HA	2:B:290:SER:HA	1.96	0.47
2:B:212:THR:HG21	2:B:245:THR:HB	1.96	0.47
2:B:207:GLU:HG3	2:B:238:LEU:HD11	1.95	0.47
1:A:114:LYS:O	1:A:118:MET:HG2	2.14	0.47
1:A:29:MET:HB2	1:A:163:SER:H	1.80	0.46
2:B:232:TYR:N	2:B:310:TYR:O	2.43	0.46
1:A:96:LEU:O	1:A:99:VAL:HG12	2.15	0.46
1:A:31:ARG:HD3	1:A:158:ARG:NE	2.32	0.45
1:A:108:LEU:HD21	1:A:113:ALA:HB3	1.99	0.45
1:A:13:LEU:O	1:A:173:LEU:HD12	2.17	0.45
2:B:254:LEU:HB3	2:B:315:LEU:HB3	1.98	0.45
1:A:24:CYS:HB2	1:A:166:VAL:HG12	1.98	0.45
1:A:169:ALA:O	1:A:173:LEU:HD23	2.17	0.45
1:A:80:ALA:O	1:A:84:VAL:HG12	2.18	0.44
2:B:275:GLU:HG3	2:B:276:GLN:N	2.33	0.44
1:A:170:ALA:O	1:A:173:LEU:CD2	2.65	0.43
1:A:33:THR:HG22	1:A:158:ARG:HG2	2.01	0.43
1:A:21:VAL:HG23	1:A:170:ALA:HB1	2.00	0.43
1:A:67:TYR:O	1:A:69:SER:N	2.52	0.43
1:A:173:LEU:HG	1:A:174:LEU:N	2.32	0.43
2:B:225:ASP:OD2	2:B:227:ARG:NH2	2.51	0.43
1:A:17:GLY:HA2	1:A:173:LEU:HD11	1.93	0.43
1:A:97:LEU:HD11	1:A:120:VAL:HG13	2.01	0.43
1:A:17:GLY:H	1:A:173:LEU:CD1	2.24	0.42
1:A:149:ASN:O	1:A:152:VAL:HG12	2.19	0.42
1:A:83:LEU:HD12	1:A:83:LEU:HA	1.85	0.42
2:B:215:LEU:O	2:B:221:GLY:HA2	2.20	0.42
1:A:179:GLY:C	1:A:181:LEU:H	2.24	0.41
1:A:117:THR:HA	1:A:120:VAL:HG12	2.03	0.41
1:A:51:TRP:NE1	1:A:81:ARG:HD2	2.36	0.41
2:B:261:PHE:CD2	2:B:298:LEU:HD11	2.56	0.41
1:A:156:GLN:OE1	2:B:225:ASP:HB2	2.21	0.41
1:A:134:VAL:HG13	1:A:135:PRO:HD3	2.02	0.41
1:A:37:GLY:C	1:A:39:ASN:H	2.24	0.41
2:B:228:SER:OG	2:B:312:TYR:N	2.53	0.41
1:A:137:SER:N	1:A:168:TRP:HE1	2.18	0.40
1:A:143:ILE:HD11	1:A:160:MET:HA	2.03	0.40
1:A:106:ASN:N	1:A:106:ASN:OD1	2.55	0.40
2:B:206:THR:HA	2:B:239:ASN:O	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	180/214 (84%)	168 (93%)	12 (7%)	0	100	100
2	B	132/134 (98%)	131 (99%)	1 (1%)	0	100	100
All	All	312/348 (90%)	299 (96%)	13 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	139/168 (83%)	138 (99%)	1 (1%)	84	92
2	B	116/118 (98%)	116 (100%)	0	100	100
All	All	255/286 (89%)	254 (100%)	1 (0%)	91	95

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	173	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	A	44	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 monosaccharides 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.