



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 23, 2022 – 02:13 PM JST

PDB ID : 7WQL  
Title : Zinc binding strength of proteins dominants zinc uptake in Caco-2 cells  
Authors : Li, T.; Ma, J.; Zang, J.; Zhao, G.; Zhang, T.  
Deposited on : 2022-01-25  
Resolution : 2.00 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.

The types of validation reports are described at <http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	2.30
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)	:	2.30

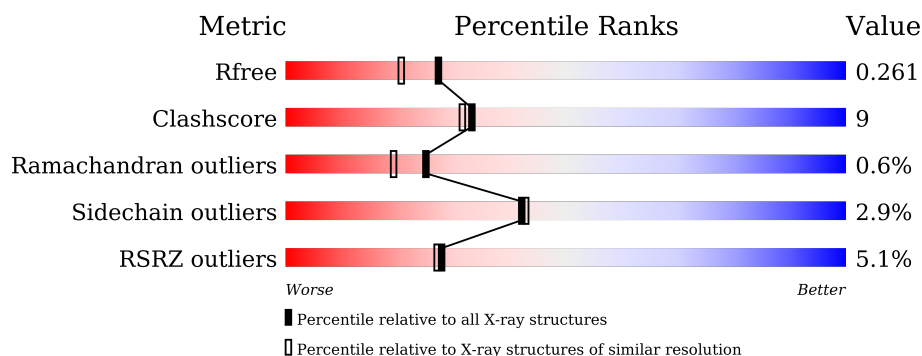
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	162	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 88%; height: 10px; background-color: green;"></div> <div style="width: 8%; height: 10px; background-color: yellow;"></div> <div style="width: 2%; height: 10px; background-color: orange;"></div> <div style="width: 2%; height: 10px; background-color: red;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div>
1	B	162	<div> <div style="width: 9%; height: 10px; background-color: red;"></div> <div style="width: 73%; height: 10px; background-color: green;"></div> <div style="width: 19%; height: 10px; background-color: yellow;"></div> <div style="width: 2%; height: 10px; background-color: orange;"></div> <div style="width: 2%; height: 10px; background-color: red;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5111 atoms, of which 2510 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 Beta-lactoglobulin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	157	Total	C	H	N	O	S	0	0	0
			2495	791	1255	200	240	9			
1	B	157	Total	C	H	N	O	S	0	0	0
			2496	791	1255	200	241	9			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	Zn	0	0
			5	5		
2	B	1	Total	Zn	0	0
			1	1		

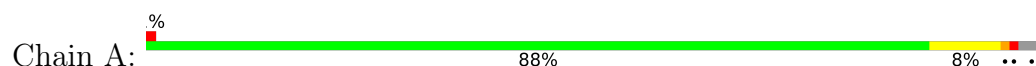
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	85	Total	O	0	0
			85	85		
3	B	29	Total	O	0	0
			29	29		

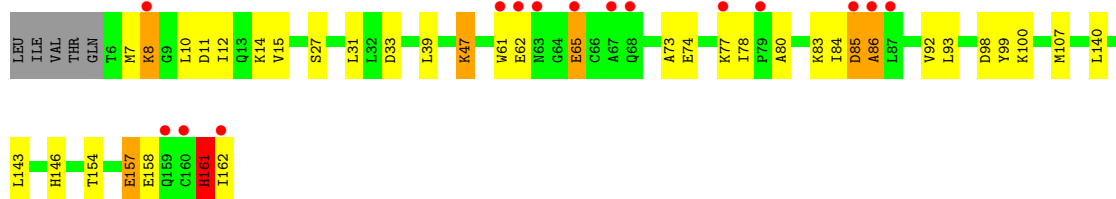
### 3 Residue-property plots [i](#)

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 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: Beta-lactoglobulin



#### • Molecule 1: Beta-lactoglobulin



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	132.15Å 42.46Å 70.04Å 90.00° 104.05° 90.00°	Depositor
Resolution (Å)	33.97 – 2.00 33.97 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.4 (33.97-2.00) 98.4 (33.97-2.00)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.96 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, $R_{free}$	0.207 , 0.261 0.207 , 0.261	Depositor DCC
$R_{free}$ test set	1997 reflections (7.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.8	Xtriage
Anisotropy	0.761	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 48.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5111	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.13% 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

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.63	0/1261	0.90	4/1706 (0.2%)
1	B	1.18	7/1262 (0.6%)	0.92	4/1706 (0.2%)
All	All	0.94	7/2523 (0.3%)	0.91	8/3412 (0.2%)

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

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	65	GLU	CG-CD	20.84	1.83	1.51
1	B	65	GLU	CB-CG	17.46	1.85	1.52
1	B	65	GLU	CD-OE1	17.17	1.44	1.25
1	B	65	GLU	CD-OE2	11.04	1.37	1.25
1	B	8	LYS	CE-NZ	9.43	1.72	1.49
1	B	8	LYS	CG-CD	-7.31	1.27	1.52
1	B	65	GLU	CA-CB	5.30	1.65	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	65	GLU	OE1-CD-OE2	-14.66	105.71	123.30
1	B	65	GLU	CA-CB-CG	12.81	141.59	113.40
1	A	148	ARG	NE-CZ-NH2	-11.45	114.58	120.30
1	A	40	ARG	NE-CZ-NH2	-10.58	115.01	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	148	ARG	NE-CZ-NH1	8.27	124.43	120.30
1	A	40	ARG	CG-CD-NE	-7.13	96.82	111.80
1	B	65	GLU	CB-CG-CD	5.35	128.65	114.20
1	B	65	GLU	N-CA-CB	5.12	119.82	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	85	ASP	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	1240	1255	1255	11	1
1	B	1241	1255	1255	33	0
2	A	5	0	0	0	1
2	B	1	0	0	0	0
3	A	85	0	0	4	3
3	B	29	0	0	1	0
All	All	2601	2510	2510	43	4

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 9.

All (43) 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:8:LYS:CE	1:B:8:LYS:NZ	1.72	1.52
1:B:65:GLU:CB	1:B:65:GLU:CG	1.85	1.51
1:B:65:GLU:CG	1:B:65:GLU:CD	1.83	1.46
1:A:131:GLU:OE1	3:A:301:HOH:O	2.02	0.77
1:B:84:ILE:HD12	1:B:86:ALA:HB2	1.74	0.69
1:A:138:LYS:NZ	3:A:303:HOH:O	2.26	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:ILE:HD13	1:B:92:VAL:HG23	1.77	0.67
1:A:148:ARG:NH2	3:A:302:HOH:O	2.23	0.64
1:B:12:ILE:O	1:B:15:VAL:HG22	1.99	0.62
1:B:161:HIS:O	1:B:162:ILE:HB	2.00	0.61
1:B:7:MET:HG2	1:B:8:LYS:N	2.18	0.58
1:B:47:LYS:NZ	3:B:302:HOH:O	2.39	0.56
1:B:93:LEU:N	1:B:93:LEU:HD12	2.22	0.55
1:B:7:MET:HG2	1:B:8:LYS:H	1.72	0.55
1:B:84:ILE:CD1	1:B:86:ALA:HB2	2.36	0.54
1:B:98:ASP:O	1:B:100:LYS:N	2.33	0.53
1:A:70:LYS:O	1:A:70:LYS:HG3	2.09	0.53
1:A:75:LYS:HE2	3:A:318:HOH:O	2.09	0.52
1:A:12:ILE:O	1:A:15:VAL:HG22	2.11	0.51
1:B:140:LEU:HD22	1:B:143:LEU:HD12	1.94	0.49
1:B:78:ILE:HG22	1:B:80:ALA:H	1.77	0.48
1:A:69:LYS:HE3	1:A:86:ALA:HB1	1.97	0.47
1:B:8:LYS:NZ	1:B:8:LYS:CD	2.68	0.47
1:B:39:LEU:HD22	1:B:107:MET:SD	2.55	0.47
1:B:7:MET:O	1:B:80:ALA:HB2	2.15	0.46
1:B:73:ALA:HA	1:B:84:ILE:HG22	1.98	0.46
1:B:100:LYS:HD3	1:B:100:LYS:HA	1.65	0.45
1:B:154:THR:O	1:B:158:GLU:HG2	2.16	0.45
1:A:33:ASP:HB2	1:B:33:ASP:O	2.17	0.45
1:B:14:LYS:HB3	1:B:99:TYR:CD1	2.52	0.44
1:B:154:THR:O	1:B:158:GLU:CG	2.66	0.44
1:B:61:TRP:HA	1:B:65:GLU:O	2.17	0.44
1:B:10:LEU:HG	1:B:11:ASP:N	2.33	0.43
1:A:93:LEU:HD12	1:A:93:LEU:N	2.34	0.43
1:B:31:LEU:HD22	1:B:39:LEU:HD12	2.00	0.43
1:B:84:ILE:HD13	1:B:92:VAL:CG2	2.48	0.43
1:B:157:GLU:OE2	1:B:157:GLU:N	2.38	0.43
1:B:74:GLU:HB3	1:B:83:LYS:HG3	2.02	0.42
1:B:7:MET:CG	1:B:8:LYS:H	2.31	0.42
1:B:7:MET:CG	1:B:8:LYS:N	2.81	0.42
1:B:27:SER:O	1:B:146:HIS:HD2	2.03	0.41
1:A:58:LEU:C	1:A:58:LEU:HD12	2.40	0.41
1:A:51:GLU:OE1	1:A:51:GLU:N	2.52	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:HIS:HD2	2:A:201:ZN:ZN[4_456]	1.11	0.49
3:A:372:HOH:O	3:A:380:HOH:O[4_446]	2.01	0.19
3:A:332:HOH:O	3:A:380:HOH:O[4_446]	2.06	0.14
3:A:323:HOH:O	3:A:364:HOH:O[4_446]	2.18	0.02

## 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	155/162 (96%)	146 (94%)	9 (6%)	0	100	100
1	B	155/162 (96%)	142 (92%)	11 (7%)	2 (1%)	12	6
All	All	310/324 (96%)	288 (93%)	20 (6%)	2 (1%)	25	19

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	86	ALA
1	B	161	HIS

### 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	138/143 (96%)	136 (99%)	2 (1%)	67	72
1	B	138/143 (96%)	132 (96%)	6 (4%)	29	26
All	All	276/286 (96%)	268 (97%)	8 (3%)	42	43

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ARG
1	A	148	ARG
1	B	47	LYS
1	B	62	GLU
1	B	77	LYS
1	B	85	ASP
1	B	157	GLU
1	B	161	HIS

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

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 6 are 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

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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	157/162 (96%)	0.03	1 (0%) 89 88	21, 34, 48, 57	0
1	B	157/162 (96%)	0.71	15 (9%) 8 7	21, 46, 66, 76	0
All	All	314/324 (96%)	0.37	16 (5%) 28 27	21, 39, 62, 76	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	162	ILE	5.9
1	B	67	ALA	4.6
1	B	61	TRP	4.4
1	B	79	PRO	4.1
1	B	87	LEU	3.6
1	B	86	ALA	3.5
1	B	8	LYS	3.4
1	B	63	ASN	3.2
1	B	159	GLN	3.0
1	B	77	LYS	2.8
1	B	62	GLU	2.6
1	B	160	CYS	2.4
1	B	85	ASP	2.4
1	A	100	LYS	2.2
1	B	68	GLN	2.1
1	B	65	GLU	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 monosaccharides 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, 95<sup>th</sup> 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	ZN	A	204	1/1	0.90	0.12	63,63,63,63	1
2	ZN	A	205	1/1	0.95	0.08	115,115,115,115	0
2	ZN	B	201	1/1	0.95	0.09	54,54,54,54	1
2	ZN	A	201	1/1	0.99	0.14	26,26,26,26	1
2	ZN	A	202	1/1	0.99	0.14	22,22,22,22	0
2	ZN	A	203	1/1	0.99	0.07	42,42,42,42	1

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

There are no such residues in this entry.