



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 03:25 PM UTC

PDB ID : 7FBC / pdb\_00007fbc  
Title : De novo design protein D22 with MBP tag  
Authors : Bin, H.  
Deposited on : 2021-07-09  
Resolution : 1.85 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](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 : 4-5-2 with Phenix2.0  
Xtrriage (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

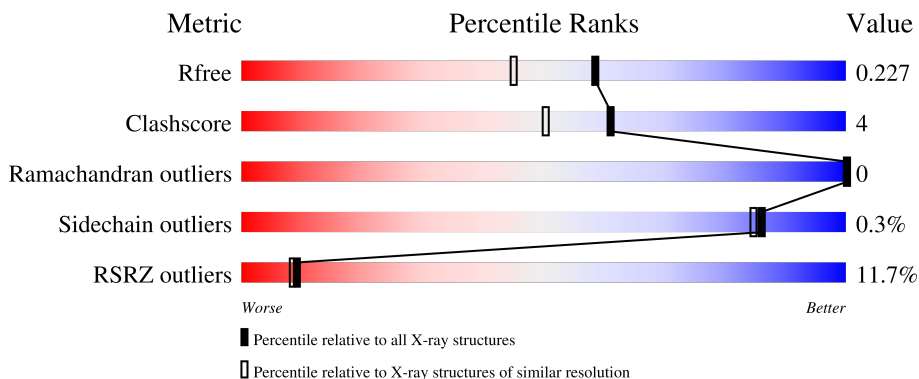
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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))
$R_{free}$	180053	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	461	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3612 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 Maltodextrin-binding protein,De novo design protein D22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	461	3496	2241	581	665	9	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	83	ALA	ASP	engineered mutation	UNP A0A4P1LXE0
A	84	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
A	173	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
A	174	ALA	ASN	engineered mutation	UNP A0A4P1LXE0
A	240	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
A	360	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
A	363	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
A	364	ALA	ASP	engineered mutation	UNP A0A4P1LXE0

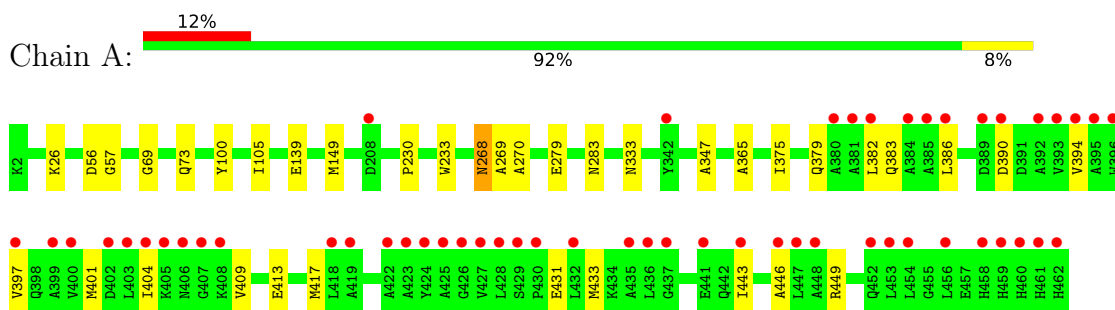
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	116	Total	O	0	0
			116	116		

### 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: Maltodextrin-binding protein, De novo design protein D22



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.63Å 77.14Å 94.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.07 – 1.85 48.07 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.07-1.85) 99.9 (48.07-1.85)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.89 (at 1.86Å)	Xtrriage
Refinement program	PHENIX 1.10	Depositor
R, $R_{free}$	0.184 , 0.215 (Not available) , 0.227	Depositor DCC
$R_{free}$ test set	1939 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.8	Xtrriage
Anisotropy	0.247	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 33.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3612	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.64% 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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.40	0/3573	0.55	0/4860

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 [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	3496	0	3465	29	0
2	A	116	0	0	0	0
All	All	3612	0	3465	29	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 4.

All (29) 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:413:GLU:CD	1:A:449:ARG:HH22	1.91	0.78
1:A:413:GLU:OE1	1:A:449:ARG:NH2	2.29	0.66
1:A:73:GLN:HG2	1:A:100:TYR:OH	1.98	0.63
1:A:268:ASN:ND2	1:A:270:ALA:H	2.01	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:GLY:O	1:A:73:GLN:HG3	2.04	0.58
1:A:268:ASN:HD22	1:A:270:ALA:H	1.50	0.57
1:A:383:GLN:HG3	1:A:386:LEU:HD12	1.88	0.54
1:A:69:GLY:HA3	1:A:333:ASN:O	2.10	0.51
1:A:431:GLU:H	1:A:431:GLU:CD	2.19	0.50
1:A:404:ILE:HG12	1:A:409:VAL:HB	1.93	0.50
1:A:100:TYR:HB2	1:A:105:ILE:HD13	1.94	0.50
1:A:279:GLU:OE1	1:A:283:ASN:ND2	2.44	0.49
1:A:268:ASN:HD22	1:A:269:ALA:N	2.13	0.46
1:A:268:ASN:HD22	1:A:268:ASN:C	2.24	0.46
1:A:417:MET:HE2	1:A:446:ALA:HB1	1.98	0.46
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.52	0.45
1:A:73:GLN:HG2	1:A:100:TYR:CZ	2.51	0.45
1:A:279:GLU:CD	1:A:283:ASN:HD22	2.24	0.45
1:A:347:ALA:HB2	1:A:365:ALA:HB2	2.00	0.44
1:A:26:LYS:HA	1:A:26:LYS:HD2	1.71	0.44
1:A:382:LEU:HD23	1:A:382:LEU:HA	1.89	0.43
1:A:397:VAL:O	1:A:401:MET:HG2	2.19	0.42
1:A:433:MET:HG3	1:A:443:ILE:HD11	2.02	0.42
1:A:56:ASP:CG	1:A:57:GLY:H	2.27	0.42
1:A:382:LEU:O	1:A:386:LEU:HG	2.20	0.41
1:A:390:ASP:O	1:A:394:VAL:HG23	2.20	0.41
1:A:139:GLU:OE1	1:A:139:GLU:HA	2.19	0.41
1:A:375:ILE:O	1:A:379:GLN:HB2	2.20	0.41
1:A:149:MET:HE2	1:A:149:MET:HB3	1.84	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	459/461 (100%)	452 (98%)	7 (2%)	0	<a href="#">100</a> <a href="#">100</a>

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	347/352 (99%)	346 (100%)	1 (0%)	86 85

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

Mol	Chain	Res	Type
1	A	268	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	202	ASN
1	A	268	ASN
1	A	326	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](#)

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

### 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<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	461/461 (100%)	0.25	54 (11%) <b>9</b> <b>8</b>	13, 23, 63, 82	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	386	LEU	4.7
1	A	460	HIS	4.4
1	A	456	LEU	4.3
1	A	402	ASP	4.2
1	A	458	HIS	4.1
1	A	427	VAL	3.9
1	A	399	ALA	3.8
1	A	406	ASN	3.7
1	A	394	VAL	3.7
1	A	403	LEU	3.6
1	A	424	TYR	3.5
1	A	430	PRO	3.4
1	A	404	ILE	3.3
1	A	380	ALA	3.2
1	A	441	GLU	3.2
1	A	423	ALA	3.1
1	A	446	ALA	3.1
1	A	461	HIS	3.1
1	A	396	TRP	3.1
1	A	432	LEU	3.1
1	A	405	LYS	3.0
1	A	462	HIS	3.0
1	A	407	GLY	3.0
1	A	382	LEU	2.8
1	A	385	ALA	2.8
1	A	459	HIS	2.8
1	A	392	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	429	SER	2.7
1	A	454	LEU	2.7
1	A	381	ALA	2.7
1	A	342	TYR	2.7
1	A	428	LEU	2.7
1	A	393	VAL	2.6
1	A	400	VAL	2.6
1	A	422	ALA	2.6
1	A	435	ALA	2.6
1	A	447	LEU	2.5
1	A	448	ALA	2.5
1	A	426	GLY	2.5
1	A	453	LEU	2.5
1	A	395	ALA	2.5
1	A	397	VAL	2.5
1	A	389	ASP	2.4
1	A	436	LEU	2.4
1	A	425	ALA	2.4
1	A	437	GLY	2.3
1	A	408	LYS	2.3
1	A	443	ILE	2.2
1	A	390	ASP	2.2
1	A	384	ALA	2.2
1	A	419	ALA	2.1
1	A	208	ASP	2.1
1	A	418	LEU	2.0
1	A	452	GLN	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.