



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 05:14 AM UTC

PDB ID : 7NFT / pdb_00007nft
Title : Fujian capbinding domain in complex with Nb8208
Authors : Keown, J.R.; Grimes, J.M.; Fodor, E.
Deposited on : 2021-02-07
Resolution : 3.14 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
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

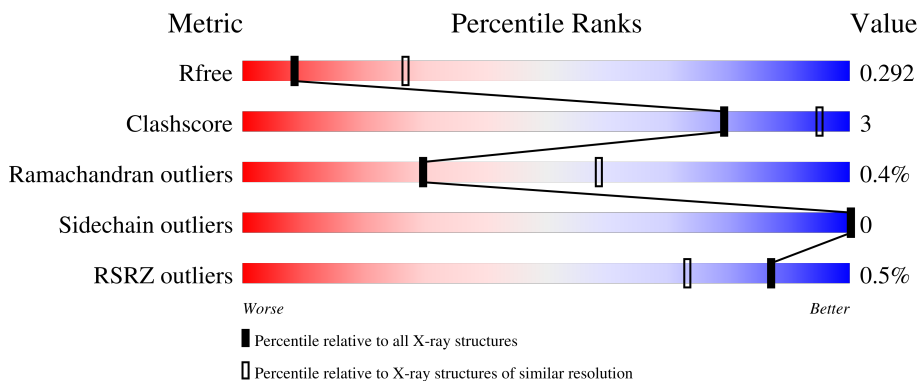
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.14 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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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 ≥ 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	304	 % 48% 5% 47%
1	B	304	 46% 6% 48%
2	C	132	 87% . 9%
2	D	132	 86% 5% 9%

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 8839 atoms, of which 4402 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 Polymerase basic protein 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	161	2587	802	1311	234	231	9	0	0	0
1	B	159	2556	790	1297	232	228	9	0	0	0

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	233	HIS	-	expression tag	UNP Q6E3N3
A	234	HIS	-	expression tag	UNP Q6E3N3
A	235	HIS	-	expression tag	UNP Q6E3N3
A	236	HIS	-	expression tag	UNP Q6E3N3
A	237	HIS	-	expression tag	UNP Q6E3N3
A	238	HIS	-	expression tag	UNP Q6E3N3
A	239	GLY	-	expression tag	UNP Q6E3N3
A	240	GLU	-	expression tag	UNP Q6E3N3
A	241	ASN	-	expression tag	UNP Q6E3N3
A	242	LEU	-	expression tag	UNP Q6E3N3
A	243	TYR	-	expression tag	UNP Q6E3N3
A	244	PHE	-	expression tag	UNP Q6E3N3
A	245	GLN	-	expression tag	UNP Q6E3N3
A	246	GLY	-	expression tag	UNP Q6E3N3
B	233	HIS	-	expression tag	UNP Q6E3N3
B	234	HIS	-	expression tag	UNP Q6E3N3
B	235	HIS	-	expression tag	UNP Q6E3N3
B	236	HIS	-	expression tag	UNP Q6E3N3
B	237	HIS	-	expression tag	UNP Q6E3N3
B	238	HIS	-	expression tag	UNP Q6E3N3
B	239	GLY	-	expression tag	UNP Q6E3N3
B	240	GLU	-	expression tag	UNP Q6E3N3
B	241	ASN	-	expression tag	UNP Q6E3N3
B	242	LEU	-	expression tag	UNP Q6E3N3
B	243	TYR	-	expression tag	UNP Q6E3N3

Continued on next page...

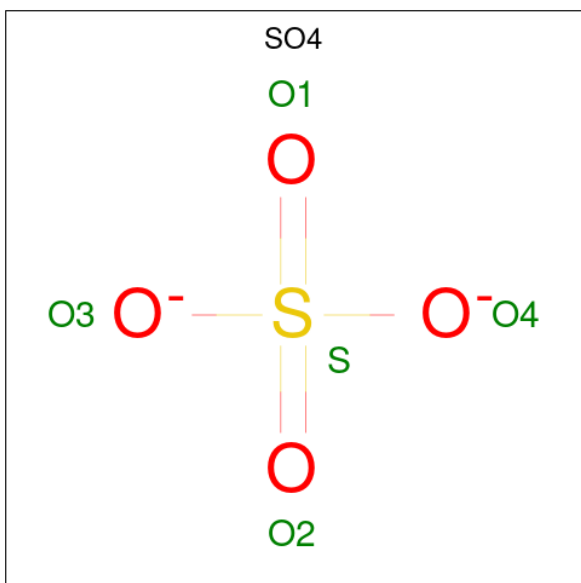
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	244	PHE	-	expression tag	UNP Q6E3N3
B	245	GLN	-	expression tag	UNP Q6E3N3
B	246	GLY	-	expression tag	UNP Q6E3N3

- Molecule 2 is a protein called Nb8208.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
2	C	120	Total	C	H	N	O	S	0	0	0
			1823	580	897	165	176	5			
2	D	120	Total	C	H	N	O	S	0	0	0
			1823	580	897	165	176	5			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total 5	O 4	S 1	0	0
3	B	1	Total 5	O 4	S 1	0	0
3	D	1	Total 5	O 4	S 1	0	0
3	D	1	Total 5	O 4	S 1	0	0

Q1	V2	G26	R38	E46	T91	T119	V120	SER	SER	HIS	HIS	HIS	HIS	HIS	HIS	HIS	PRO	GLU	GLU	ALA
----	----	-----	-----	-----	-----	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics i

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, α , β , γ	128.19Å 128.19Å 70.27Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	70.27 – 3.14 70.27 – 3.14	Depositor EDS
% Data completeness (in resolution range)	93.0 (70.27-3.14) 87.9 (70.27-3.14)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 3.13Å)	Xtrriage
Refinement program	PHENIX 1.19.1_4122	Depositor
R, R_{free}	0.244 , 0.291 0.248 , 0.292	Depositor DCC
R_{free} test set	582 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	60.8	Xtrriage
Anisotropy	0.095	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.077 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	8839	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.11	0/1294	0.27	0/1735
1	B	0.10	0/1276	0.27	0/1711
2	C	0.09	0/945	0.26	0/1275
2	D	0.09	0/945	0.30	0/1275
All	All	0.10	0/4460	0.27	0/5996

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	1276	1311	1311	8	2
1	B	1259	1297	1297	10	0
2	C	926	897	897	4	0
2	D	926	897	897	3	0
3	A	15	0	0	2	0
3	B	15	0	0	1	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
All	All	4437	4402	4402	24	2

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 3.

The worst 5 of 24 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:348:ASN:ND2	1:A:417:ASP:OD1	2.26	0.68
1:A:423:ARG:NH2	3:A:603:SO4:O2	2.28	0.66
1:B:423:ARG:NE	3:B:603:SO4:O3	2.30	0.63
1:B:403:VAL:HA	1:B:410:MET:HE2	1.87	0.55
1:B:348:ASN:ND2	1:B:417:ASP:OD1	2.39	0.54

All (2) 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:331:LYS:HZ1	1:A:455:ASP:OD2[2_545]	1.53	0.07
1:A:331:LYS:NZ	1:A:455:ASP:OD2[2_545]	2.14	0.06

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	159/304 (52%)	152 (96%)	6 (4%)	1 (1%)	21	50
1	B	157/304 (52%)	148 (94%)	8 (5%)	1 (1%)	21	50
2	C	118/132 (89%)	115 (98%)	3 (2%)	0	100	100
2	D	118/132 (89%)	115 (98%)	3 (2%)	0	100	100
All	All	552/872 (63%)	530 (96%)	20 (4%)	2 (0%)	30	59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	456	ASN
1	B	458	MET

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	140/266 (53%)	140 (100%)	0	100	100
1	B	138/266 (52%)	138 (100%)	0	100	100
2	C	94/105 (90%)	94 (100%)	0	100	100
2	D	94/105 (90%)	94 (100%)	0	100	100
All	All	466/742 (63%)	466 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	392	GLN
1	A	437	HIS
1	B	437	HIS
1	B	448	ASN

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](#)

10 ligands are modelled in this entry.

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
3	SO4	A	602	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	D	202	-	4,4,4	0.24	0	6,6,6	0.09	0
3	SO4	A	603	-	4,4,4	0.24	0	6,6,6	0.07	0
3	SO4	C	202	-	4,4,4	0.23	0	6,6,6	0.07	0
3	SO4	A	601	-	4,4,4	0.24	0	6,6,6	0.07	0
3	SO4	B	601	-	4,4,4	0.23	0	6,6,6	0.08	0
3	SO4	B	602	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	B	603	-	4,4,4	0.24	0	6,6,6	0.06	0
3	SO4	D	201	-	4,4,4	0.24	0	6,6,6	0.07	0
3	SO4	C	201	-	4,4,4	0.24	0	6,6,6	0.07	0

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.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	603	SO4	1	0
3	A	601	SO4	1	0
3	B	603	SO4	1	0

5.7 Other polymers [i](#)

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, 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	161/304 (52%)	0.12	2 (1%) 76 57	60, 72, 91, 102	0
1	B	159/304 (52%)	0.22	1 (0%) 85 70	61, 81, 99, 127	0
2	C	120/132 (90%)	0.20	0 100 100	63, 77, 94, 115	0
2	D	120/132 (90%)	0.26	0 100 100	63, 80, 111, 137	0
All	All	560/872 (64%)	0.20	3 (0%) 87 73	60, 78, 100, 137	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	357	HIS	2.4
1	A	337	SER	2.3
1	A	357	HIS	2.2

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](#)

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(\AA^2)	Q<0.9
3	SO4	D	202	5/5	0.80	0.14	72,73,80,81	0
3	SO4	D	201	5/5	0.86	0.14	56,60,61,65	0
3	SO4	C	201	5/5	0.89	0.10	49,49,52,59	0
3	SO4	B	601	5/5	0.91	0.13	58,61,65,66	0
3	SO4	C	202	5/5	0.92	0.12	68,72,78,80	0
3	SO4	B	602	5/5	0.92	0.11	60,64,67,72	0
3	SO4	A	603	5/5	0.93	0.08	55,61,65,66	0
3	SO4	A	601	5/5	0.94	0.11	55,56,59,65	0
3	SO4	B	603	5/5	0.94	0.10	47,47,49,53	0
3	SO4	A	602	5/5	0.95	0.14	55,61,66,68	0

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

There are no such residues in this entry.