



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 01:08 AM UTC

PDB ID : 2NRH / pdb\_00002nrh  
Title : Crystal structure of conserved putative Baf family transcriptional activator from *Campylobacter jejuni*  
Authors : Bonanno, J.B.; Dickey, M.; Bain, K.T.; Lau, C.; Wasserman, S.; Smith, D.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2006-11-02  
Resolution : 2.30 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](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  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (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

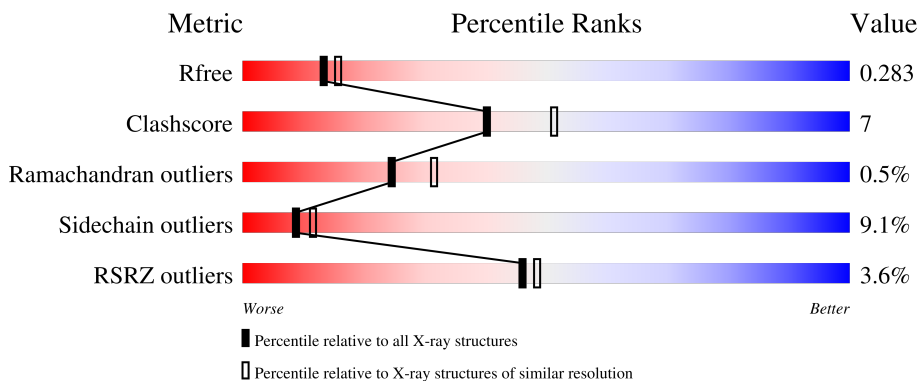
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 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	219	 4% 65% 19% 12%
1	B	219	 3% 70% 19% 9%

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 3258 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 Transcriptional activator, putative, Baf family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	192	Total	C	N	O	S	0	0	0
			1550	1022	245	279	4			
1	B	199	Total	C	N	O	S	0	0	0
			1607	1056	258	289	4			

There are 26 discrepancies between the modelled and reference sequences:

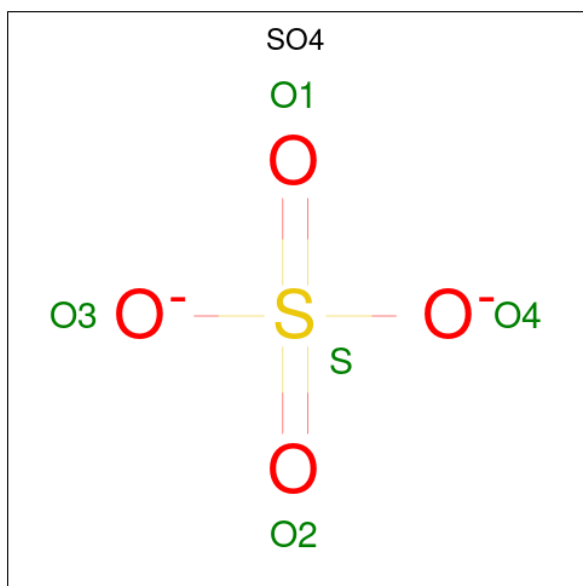
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	cloning artifact	UNP Q5HW73
A	0	SER	-	cloning artifact	UNP Q5HW73
A	1	LEU	-	cloning artifact	UNP Q5HW73
A	23	ASN	SER	engineered mutation	UNP Q5HW73
A	54	ASN	LYS	engineered mutation	UNP Q5HW73
A	210	GLU	-	expression tag	UNP Q5HW73
A	211	GLY	-	expression tag	UNP Q5HW73
A	212	HIS	-	expression tag	UNP Q5HW73
A	213	HIS	-	expression tag	UNP Q5HW73
A	214	HIS	-	expression tag	UNP Q5HW73
A	215	HIS	-	expression tag	UNP Q5HW73
A	216	HIS	-	expression tag	UNP Q5HW73
A	217	HIS	-	expression tag	UNP Q5HW73
B	-1	MET	-	cloning artifact	UNP Q5HW73
B	0	SER	-	cloning artifact	UNP Q5HW73
B	1	LEU	-	cloning artifact	UNP Q5HW73
B	23	ASN	SER	engineered mutation	UNP Q5HW73
B	54	ASN	LYS	engineered mutation	UNP Q5HW73
B	210	GLU	-	expression tag	UNP Q5HW73
B	211	GLY	-	expression tag	UNP Q5HW73
B	212	HIS	-	expression tag	UNP Q5HW73
B	213	HIS	-	expression tag	UNP Q5HW73
B	214	HIS	-	expression tag	UNP Q5HW73
B	215	HIS	-	expression tag	UNP Q5HW73
B	216	HIS	-	expression tag	UNP Q5HW73

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Chain	Residue	Modelled	Actual	Comment	Reference
B	217	HIS	-	expression tag	UNP Q5HW73

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S).



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

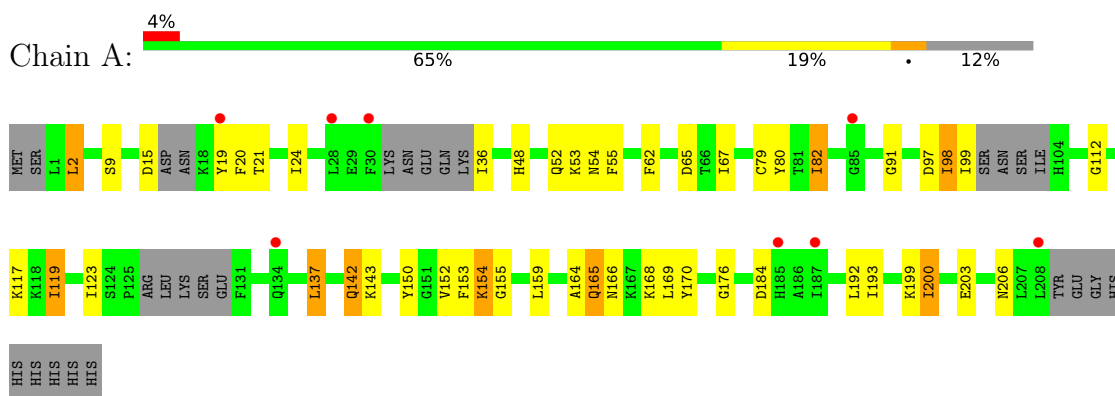
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	34	Total	O	0	0
			34	34		
3	B	52	Total	O	0	0
			52	52		

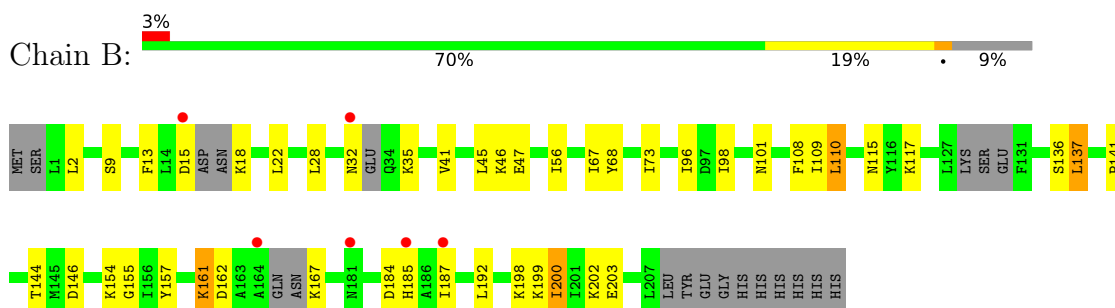
### 3 Residue-property plots [i](#)

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: Transcriptional activator, putative, Baf family



- Molecule 1: Transcriptional activator, putative, Baf family



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.29Å 73.97Å 116.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-2.30) 99.7 (20.00-2.30)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.82 (at 2.30Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.227 , 0.291 0.227 , 0.283	Depositor DCC
$R_{free}$ test set	1181 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.9	Xtrriage
Anisotropy	0.029	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 49.6	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	0.94	EDS
Total number of atoms	3258	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.15% 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](#)

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	1.05	2/1583 (0.1%)	1.14	6/2134 (0.3%)
1	B	1.14	2/1640 (0.1%)	1.13	4/2208 (0.2%)
All	All	1.10	4/3223 (0.1%)	1.14	10/4342 (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	A	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	82	ILE	CA-CB	6.24	1.62	1.54
1	B	109	ILE	C-O	5.66	1.29	1.24
1	A	2	LEU	CA-C	5.05	1.59	1.52
1	B	41	VAL	CA-CB	5.03	1.60	1.54

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	110	LEU	CA-C-N	-8.16	111.37	119.78
1	B	110	LEU	C-N-CA	-8.16	111.37	119.78
1	A	97	ASP	N-CA-C	6.17	119.45	109.40
1	A	98	ILE	N-CA-C	6.09	116.63	107.80
1	B	144	THR	N-CA-C	5.78	117.58	111.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	165	GLN	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	1550	0	1548	26	0
1	B	1607	0	1614	21	0
2	A	5	0	0	0	0
2	B	10	0	0	1	0
3	A	34	0	0	4	0
3	B	52	0	0	1	0
All	All	3258	0	3162	42	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 7.

The worst 5 of 42 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:82:ILE:HD11	1:A:192:LEU:HD11	1.37	1.07
1:B:137:LEU:HG	1:B:154:LYS:HG3	1.64	0.80
1:A:184:ASP:HA	3:A:226:HOH:O	1.88	0.71
1:A:200:ILE:HD11	3:A:227:HOH:O	1.94	0.68
1:B:9:SER:HB2	2:B:218:SO4:O2	2.02	0.59

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	182/219 (83%)	175 (96%)	5 (3%)	2 (1%)	11	13
1	B	189/219 (86%)	187 (99%)	2 (1%)	0	100	100
All	All	371/438 (85%)	362 (98%)	7 (2%)	2 (0%)	24	31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	ASN
1	A	152	VAL

### 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	166/192 (86%)	151 (91%)	15 (9%)	9	12
1	B	173/192 (90%)	157 (91%)	16 (9%)	8	11
All	All	339/384 (88%)	308 (91%)	31 (9%)	9	11

5 of 31 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	206	ASN
1	B	167	LYS
1	B	18	LYS
1	B	198	LYS
1	B	136	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	12	ASN

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Mol	Chain	Res	Type
1	B	134	GLN
1	B	177	GLN
1	A	122	HIS
1	A	10	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](#)

3 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	$\# Z  > 2$
2	SO4	B	218	-	4,4,4	0.27	0	6,6,6	0.25	0
2	SO4	B	219	-	4,4,4	0.14	0	6,6,6	0.28	0
2	SO4	A	218	-	4,4,4	0.30	0	6,6,6	0.90	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	218	SO4	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, 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	192/219 (87%)	0.64	8 (4%) 40 42	45, 66, 89, 103	0
1	B	199/219 (90%)	0.30	6 (3%) 52 54	44, 57, 74, 83	0
All	All	391/438 (89%)	0.47	14 (3%) 46 48	44, 61, 83, 103	0

The worst 5 of 14 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	85	GLY	3.7
1	A	30	PHE	3.1
1	A	19	TYR	3.1
1	A	28	LEU	2.8
1	A	208	LEU	2.8

### 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, 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	SO4	B	218	5/5	0.90	0.09	97,97,99,100	0
2	SO4	B	219	5/5	0.91	0.12	89,90,91,92	0
2	SO4	A	218	5/5	0.98	0.06	59,61,63,67	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.