



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 22, 2024 – 04:18 PM EDT

PDB ID : 6UV7  
Title : Crystal structure of alr1298, a pentapeptide repeat protein from Nostoc Pcc 7120, determined at 2.3 Angstrom resolution  
Authors : Kennedy, M.A.; Zhang, R.  
Deposited on : 2019-11-01  
Resolution : 2.27 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

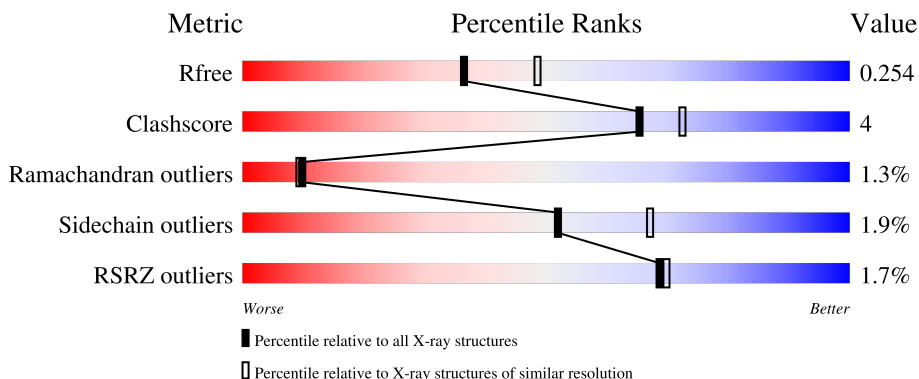
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.27 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))
$R_{free}$	164625	8487 (2.30-2.26)
Clashscore	180529	9437 (2.30-2.26)
Ramachandran outliers	177936	9341 (2.30-2.26)
Sidechain outliers	177891	9342 (2.30-2.26)
RSRZ outliers	164620	8487 (2.30-2.26)

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	187	 2% 75% 5% • 19%
1	B	187	 % 72% 10% • 18%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4843 atoms, of which 2413 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 Alr1298 protein.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	152	Total	C	H	N	O	S	Se	0	0	0
			2392	746	1196	215	230	2	3			
1	B	154	Total	C	H	N	O	S	Se	0	0	0
			2434	760	1217	218	234	2	3			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	initiating methionine	UNP Q8YXB7
A	-18	GLY	-	expression tag	UNP Q8YXB7
A	-17	SER	-	expression tag	UNP Q8YXB7
A	-16	SER	-	expression tag	UNP Q8YXB7
A	-15	HIS	-	expression tag	UNP Q8YXB7
A	-14	HIS	-	expression tag	UNP Q8YXB7
A	-13	HIS	-	expression tag	UNP Q8YXB7
A	-12	HIS	-	expression tag	UNP Q8YXB7
A	-11	HIS	-	expression tag	UNP Q8YXB7
A	-10	HIS	-	expression tag	UNP Q8YXB7
A	-9	SER	-	expression tag	UNP Q8YXB7
A	-8	SER	-	expression tag	UNP Q8YXB7
A	-7	GLY	-	expression tag	UNP Q8YXB7
A	-6	LEU	-	expression tag	UNP Q8YXB7
A	-5	VAL	-	expression tag	UNP Q8YXB7
A	-4	PRO	-	expression tag	UNP Q8YXB7
A	-3	ARG	-	expression tag	UNP Q8YXB7
A	-2	GLY	-	expression tag	UNP Q8YXB7
A	-1	SER	-	expression tag	UNP Q8YXB7
A	0	HIS	-	expression tag	UNP Q8YXB7
A	89	MSE	LEU	conflict	UNP Q8YXB7
A	124	MSE	LEU	conflict	UNP Q8YXB7
A	159	GLU	LYS	conflict	UNP Q8YXB7
B	-19	MSE	-	initiating methionine	UNP Q8YXB7
B	-18	GLY	-	expression tag	UNP Q8YXB7

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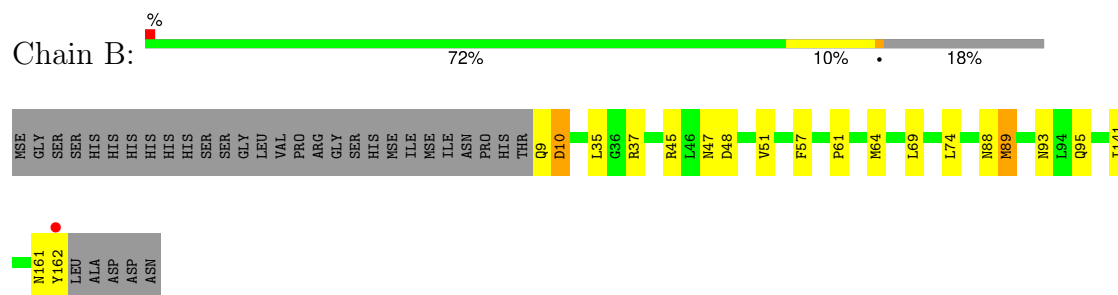
Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	SER	-	expression tag	UNP Q8YXB7
B	-16	SER	-	expression tag	UNP Q8YXB7
B	-15	HIS	-	expression tag	UNP Q8YXB7
B	-14	HIS	-	expression tag	UNP Q8YXB7
B	-13	HIS	-	expression tag	UNP Q8YXB7
B	-12	HIS	-	expression tag	UNP Q8YXB7
B	-11	HIS	-	expression tag	UNP Q8YXB7
B	-10	HIS	-	expression tag	UNP Q8YXB7
B	-9	SER	-	expression tag	UNP Q8YXB7
B	-8	SER	-	expression tag	UNP Q8YXB7
B	-7	GLY	-	expression tag	UNP Q8YXB7
B	-6	LEU	-	expression tag	UNP Q8YXB7
B	-5	VAL	-	expression tag	UNP Q8YXB7
B	-4	PRO	-	expression tag	UNP Q8YXB7
B	-3	ARG	-	expression tag	UNP Q8YXB7
B	-2	GLY	-	expression tag	UNP Q8YXB7
B	-1	SER	-	expression tag	UNP Q8YXB7
B	0	HIS	-	expression tag	UNP Q8YXB7
B	89	MSE	LEU	conflict	UNP Q8YXB7
B	124	MSE	LEU	conflict	UNP Q8YXB7
B	159	GLU	LYS	conflict	UNP Q8YXB7

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	8	Total O 8 8	0	0
2	B	9	Total O 9 9	0	0



- Molecule 1: Alr1298 protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.84Å 86.36Å 96.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.02 – 2.27 30.02 – 2.27	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.02-2.27) 99.8 (30.02-2.27)	Depositor EDS
$R_{merge}$	0.03	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.60 (at 2.27Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, $R_{free}$	0.199 , 0.251 0.206 , 0.254	Depositor DCC
$R_{free}$ test set	16321 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	55.3	Xtriage
Anisotropy	0.515	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 42.8	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	0.96	EDS
Total number of atoms	4843	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.96% 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 [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	0.63	0/1207	0.74	0/1617
1	B	0.70	0/1229	0.77	0/1647
All	All	0.66	0/2436	0.76	0/3264

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	1196	1196	1200	6	0
1	B	1217	1217	1217	12	0
2	A	8	0	0	0	0
2	B	9	0	0	2	0
All	All	2430	2413	2417	18	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.

The worst 5 of 18 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:85:ILE:HD11	1:A:103:ASP:OD2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:LEU:HB2	1:B:89:MSE:HE3	1.88	0.54
1:B:9:GLN:O	1:B:10:ASP:CB	2.57	0.53
1:B:35:LEU:HD21	1:B:61:PRO:HA	1.91	0.52
1:A:48:ASP:OD1	1:A:49:SER:N	2.43	0.50

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	150/187 (80%)	139 (93%)	8 (5%)	3 (2%)	6	4
1	B	152/187 (81%)	149 (98%)	2 (1%)	1 (1%)	19	22
All	All	302/374 (81%)	288 (95%)	10 (3%)	4 (1%)	10	9

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	B	10	ASP
1	A	143	ASN
1	A	65	LYS

### 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	133/158 (84%)	131 (98%)	2 (2%)	60	74
1	B	135/158 (85%)	132 (98%)	3 (2%)	47	62
All	All	268/316 (85%)	263 (98%)	5 (2%)	52	67

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	CYS
1	A	143	ASN
1	B	37	ARG
1	B	48	ASP
1	B	89	MSE

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 oligosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

## 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	149/187 (79%)	0.13	4 (2%) 56 57	55, 72, 99, 130	0
1	B	151/187 (80%)	-0.10	1 (0%) 84 85	47, 63, 86, 100	0
All	All	300/374 (80%)	0.02	5 (1%) 69 70	47, 68, 94, 130	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	161	ASN	3.4
1	B	162	TYR	3.1
1	A	11	ILE	2.7
1	A	141	ILE	2.4
1	A	62	GLY	2.3

### 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](#)

There are no ligands in this entry.

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

There are no such residues in this entry.