

# Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 10:47 AM EDT

PDB ID : 8T6H

Title: X-ray crystal structure of PfA-M1(E319A)

Authors: Webb, C.T.; McGowan, S.

Deposited on : 2023-06-16

Resolution : 1.98 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
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

MolProbity : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.37.1

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

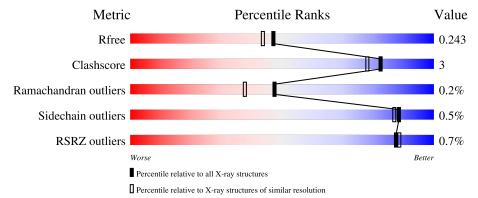
Validation Pipeline (wwPDB-VP) : 2.37.1

## 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$ 

The reported resolution of this entry is 1.98 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}({\rm \AA})) \end{array}$
$R_{free}$	130704	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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 <=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		
			<u>%</u>		
1	A	898	91%	7% •	



# 2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 7932 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 Aminopeptidase N.

Mol	Chain	Residues		A	toms			ZeroOcc	AltConf	Trace
1	A	889	Total 7224	C 4660	N 1162	O 1373	S 29	0	12	0

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	194	MET	-	initiating methionine	UNP O96935
A	213	GLN	ASN	conflict	UNP O96935
A	223	GLN	ASN	conflict	UNP O96935
A	319	ALA	GLU	conflict	UNP O96935
A	378	PRO	HIS	conflict	UNP O96935
A	501	GLN	ASN	conflict	UNP O96935
A	745	GLN	ASN	conflict	UNP O96935
A	795	GLN	ASN	conflict	UNP O96935
A	1069	GLN	ASN	conflict	UNP O96935
A	1086	HIS	-	expression tag	UNP O96935
A	1087	HIS	-	expression tag	UNP O96935
A	1088	HIS	-	expression tag	UNP O96935
A	1089	HIS	-	expression tag	UNP O96935
A	1090	HIS	-	expression tag	UNP O96935
A	1091	HIS	-	expression tag	UNP O96935

• Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

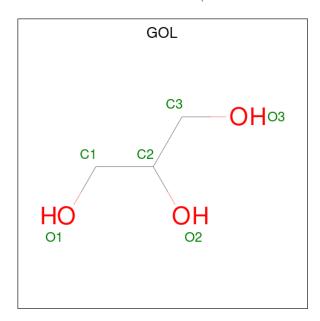
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0

• Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	3	Total Mg 3 3	0	0

 $\bullet$  Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $\mathrm{C_3H_8O_3}).$ 



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0

• Molecule 5 is water.

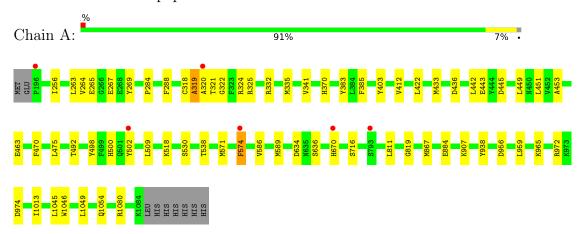
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	680	Total O 680 680	0	0



# 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: Aminopeptidase N





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	74.96Å 109.15Å 117.16Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	61.79 - 1.98	Depositor
Resolution (A)	61.79 - 1.98	EDS
% Data completeness	99.8 (61.79-1.98)	Depositor
(in resolution range)	99.8 (61.79-1.98)	EDS
$R_{merge}$	0.21	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.21 (at 1.98Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
D D	0.193 , 0.243	Depositor
$R, R_{free}$	0.195 , $0.243$	DCC
$R_{free}$ test set	3418 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.6	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.35, 50.1	EDS
L-test for twinning <sup>2</sup>	$ < L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7932	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.41% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of <|L|>,  $<L^2>$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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: ZN, GOL, MG

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).

Mal	Chain		lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.26	0/7415	0.42	1/10043 (0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}(^{o})$
1	A	574	PHE	CD1-CE1-CZ	5.06	126.18	120.10

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	7224	0	7055	36	0
2	A	1	0	0	0	0
3	A	3	0	0	0	0
4	A	24	0	32	0	0
5	A	680	0	0	6	0
All	All	7932	0	7087	36	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 3.



All (36) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
		distance (Å)	overlap (Å)
1:A:670[B]:HIS:HE1	1:A:716:SER:HB2	1.58	0.69
1:A:884:GLU:OE1	5:A:1201:HOH:O	2.12	0.67
1:A:574:PHE:CG	1:A:574:PHE:CE2	2.91	0.58
1:A:492:THR:HG22	1:A:530:SER:OG	2.04	0.57
1:A:1080:ARG:NH1	5:A:1215:HOH:O	2.38	0.56
1:A:319:ALA:HA	1:A:383:TYR:HB2	1.86	0.55
1:A:538:THR:HG21	1:A:819:GLY:HA3	1.89	0.54
1:A:509:LEU:HD21	1:A:518:LYS:HB3	1.90	0.52
1:A:451:LEU:HG	1:A:470:PHE:CE2	2.46	0.51
1:A:586:VAL:HA	1:A:589:MET:HE2	1.93	0.50
1:A:956:ASP:HB3	1:A:959:LEU:HB3	1.94	0.49
1:A:318:CYS:HA	1:A:322:GLY:HA3	1.95	0.49
1:A:325:ARG:NH1	5:A:1228:HOH:O	2.44	0.49
1:A:445:ASP:OD1	5:A:1202:HOH:O	2.20	0.49
1:A:463:GLU:HG3	1:A:500:HIS:HB3	1.94	0.49
1:A:403:TYR:N	1:A:443:GLU:OE2	2.36	0.47
1:A:264:VAL:HG23	1:A:267:GLU:HB3	1.99	0.45
1:A:341:VAL:HG21	1:A:385:PHE:HE1	1.81	0.44
1:A:907:LYS:HG2	1:A:938:TYR:CZ	2.54	0.43
1:A:284:PRO:HG2	1:A:288:PHE:CD1	2.53	0.43
1:A:453:ALA:HB2	1:A:475:LEU:HD23	2.01	0.43
1:A:634:ASP:OD2	1:A:636:SER:OG	2.21	0.43
1:A:811:LEU:HB3	1:A:867:MET:SD	2.59	0.42
1:A:332:ARG:HG3	1:A:335:MET:HG2	2.01	0.42
1:A:412:VAL:HB	1:A:422:LEU:HD22	2.02	0.42
1:A:965:LYS:NZ	5:A:1254:HOH:O	2.52	0.41
1:A:1045:LEU:HD13	1:A:1049:LEU:HD21	2.02	0.41
1:A:1046:TRP:O	1:A:1054:GLN:HG2	2.21	0.41
1:A:433:MET:HG2	1:A:498:TYR:CZ	2.56	0.41
1:A:442:LEU:HD22	1:A:502:TYR:CE1	2.55	0.41
1:A:972:ARG:HB3	1:A:974:ASP:OD1	2.21	0.41
1:A:256:ILE:HG13	1:A:263:LEU:HB2	2.02	0.41
1:A:321:THR:O	1:A:324:ARG:HG2	2.20	0.41
1:A:1013:ILE:HD12	1:A:1013:ILE:HA	1.90	0.40
1:A:256:ILE:HD11	1:A:269:TYR:CD1	2.56	0.40
1:A:265:GLU:OE2	5:A:1203:HOH:O	2.22	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	898/898 (100%)	880 (98%)	16 (2%)	2 (0%)	47 38

#### All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	319	ALA
1	A	320	ALA

#### 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 Out		Outliers	Percentiles
1	A	790/829 (95%)	786 (100%)	4 (0%)	88 87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	370	HIS
1	A	436	ASP
1	A	449	LEU
1	A	571	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	317	GLN
1	A	464	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 monosaccharides in this entry.

### 5.6 Ligand geometry (i)

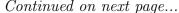
Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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).

Mal	Acl Trans Chair Box Lin		Timle	В	Bond lengths			Bond angles		
Mol	Type	Chain	Res	Link	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	A	1103	-	5,5,5	1.11	0	5,5,5	0.38	0
4	GOL	A	1104	-	5,5,5	1.10	1 (20%)	5,5,5	0.26	0
4	GOL	A	1105	-	5,5,5	1.12	1 (20%)	5,5,5	0.29	0
4	GOL	A	1106	-	5,5,5	1.10	1 (20%)	5,5,5	0.24	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

$\mathbf{Mol}$	Type	Chain	$\operatorname{Res}$	Link	Chirals	Torsions	Rings
4	GOL	A	1103	-	-	0/4/4/4	-





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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	1104	-	-	2/4/4/4	_
4	GOL	A	1105	-	-	0/4/4/4	-
4	GOL	A	1106	-	-	2/4/4/4	_

All (3) bond length outliers are listed below:

$\mathbf{Mol}$	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}( ext{\AA})$
4	A	1104	GOL	C3-C2	2.05	1.60	1.51
4	A	1105	GOL	C3-C2	2.05	1.60	1.51
4	A	1106	GOL	C3-C2	2.04	1.60	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1104	GOL	O1-C1-C2-C3
4	A	1106	GOL	O1-C1-C2-C3
4	A	1104	GOL	O1-C1-C2-O2
4	A	1106	GOL	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

## 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,  $95^{th}$  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>	$\# \mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	889/898 (98%)	0.02	6 (0%) 87 88	13, 25, 43, 55	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ	
1	A 320 ALA		3.1		
1	A	502	TYR	2.5	
1	A	670[A]	HIS	2.4	
1	A	574	PHE	2.3	
1	A	793[A]	SER	2.2	
1	A	196	PRO	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 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,  $95^{th}$  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	${f B-factors}({f \AA}^2)$	Q < 0.9
4	GOL	A	1106	6/6	0.82	0.18	36,37,52,54	0

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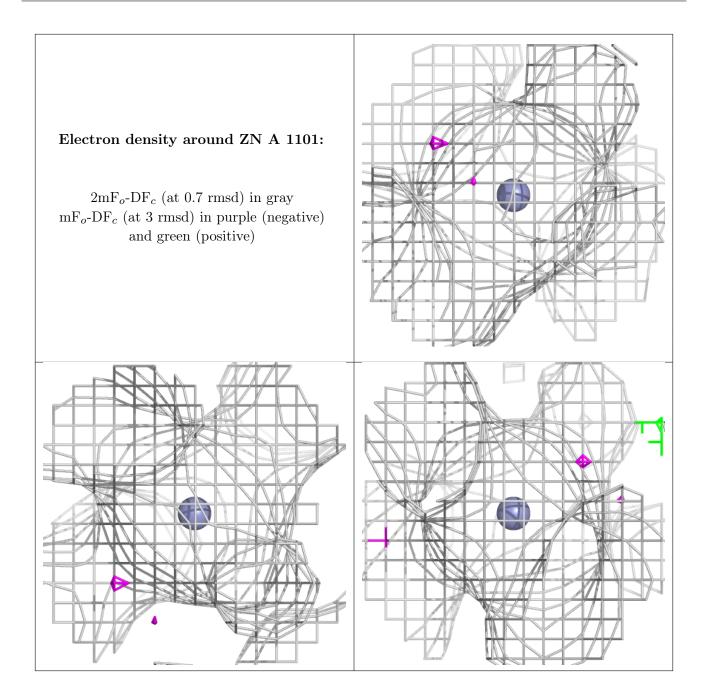


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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	GOL	A	1105	6/6	0.90	0.21	36,37,55,55	0
3	MG	A	1102	1/1	0.94	0.06	32,32,32,32	0
4	GOL	A	1104	6/6	0.95	0.13	18,22,42,42	0
4	GOL	A	1103	6/6	0.97	0.13	20,23,24,24	0
3	MG	A	1108	1/1	0.97	0.13	34,34,34,34	0
3	MG	A	1107	1/1	0.99	0.10	39,39,39,39	0
2	ZN	A	1101	1/1	1.00	0.08	27,27,27,27	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





# 6.5 Other polymers (i)

There are no such residues in this entry.

