

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

#### Aug 31, 2020 - 06:14 AM BST

PDB ID	:	5JPC
Title	:	Joint X-ray/neutron structure of MTAN complex with Formycin A
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Deposited on		
Resolution	:	2.50  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 following versions of software and data (see references (1)) were used in the production of this report:

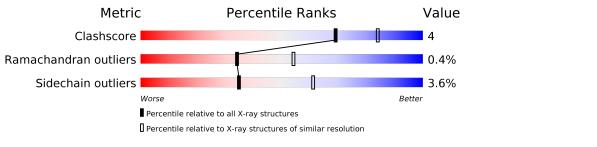
:	4.02b-467
:	1.8.5 (274361), CSD as541be (2020)
:	1.13
:	FAILED
:	1.1.7 (2018)
:	20191225.v01 (using entries in the PDB archive December 25th 2019)
:	Engh & Huber (2001)
:	Parkinson et al. (1996)
:	2.13
	:::::::::::::::::::::::::::::::::::::::

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: *NEUTRON DIFFRACTION, X-RAY DIFFRACTION* 

The reported resolution of this entry is 2.50 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
Clashscore	141614	$5346\ (2.50-2.50)$
Ramachandran outliers	138981	5231(2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

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 on 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%

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain	
1	٨	220		
	А	229	93%	6%



#### 5JPC

# 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3999 atoms, of which 1745 are hydrogens and 431 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 Aminodeoxyfutalosine nucleosidase.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace			
1	Δ	229	Total	С	D	Η	Ν	Ο	$\mathbf{S}$	0	192	0
	A	229	3808	1124	319	1738	286	333	8	0	192	0

• Molecule 2 is a ligand with the chemical component id FMC but its atom names do not match the existing wwPDB Chemical Component Dictionary definition for FMC. ERROR THIS SHOULD NOT HAPPEN FOLLOWING ANNOTATION.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
2	А	1	Total 32	C 10	D 6	11	N 5	0 4	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
3	А	53	Total 159	D 106	O 53	0	0

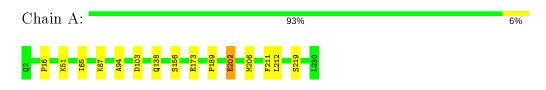


# 3 Residue-property plots (i)

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. 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. 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.

Note EDS failed to run properly.

• Molecule 1: Aminodeoxyfutalosine nucleosidase





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants	83.19Å $83.19$ Å $67.63$ Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor
Resolution (Å)	36.02 - 2.50	Depositor
% Data completeness	54.5(36.02-2.50)	Depositor
(in resolution range)	· · · · ·	-
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	$0.00 ~({\rm at}~2.10{ m \AA})$	Xtriage
Refinement program	nCNS 1.0.0	Depositor
$R, R_{free}$	0.282 , $0.266$	Depositor
Wilson B-factor $(Å^2)$	12.0	Xtriage
Anisotropy	0.474	Xtriage
L-test for twinning <sup>2</sup>	$< L >=0.46, < L^2>=0.29$	Xtriage
Estimated twinning fraction	0.094 for -h,-k,l	Xtriage
Total number of atoms	3999	wwPDB-VP
Average B, all atoms $(Å^2)$	33.0	wwPDB-VP

EDS failed to run properly - this section is therefore incomplete.

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.40% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<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: FMC, DOD

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		
	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.39	0/3266	0.61	0/4391	

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	А	2070	1738	301	7	0
2	А	25	7	13	1	0
3	А	159	0	0	0	0
All	All	2254	1745	314	8	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 (8) 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:202:GLU:CD	1:A:202:GLU:H	2.08	0.50	

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Atom-1	Atom-2	${f Interatomic}\ {f distance}\ ({ m \AA})$	Clash overlap (Å)
1:A:211:PHE:O	1:A:212:LEU:C	2.57	0.43
1:A:103:ASP:C	1:A:103:ASP:OD1	2.57	0.41
2:A:501:FMC:H2'	2:A:501:FMC:N3	2.29	0.41

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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	А	417/229 (182%)	389~(93%)	26~(6%)	2~(0%)	29 48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	51[A]	LYS
1	А	51[B]	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	А	352/192~(183%)	339~(96%)	13 (4%)	34 60

All (13) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	А	87[A]	LYS
1	А	87[B]	LYS
1	А	138[A]	GLN
1	А	138[B]	GLN
1	А	156[A]	SER
1	А	156[B]	SER
1	А	173[A]	GLU
1	А	173[B]	GLU
1	А	202	GLU
1	А	206[A]	MET
1	А	206[B]	MET
1	А	219[A]	SER
1	А	219[B]	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

#### 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 1 ligands modelled in this entry, 1 could not be matched to an existing wwPDB Chemical Component Dictionary definition at this stage - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

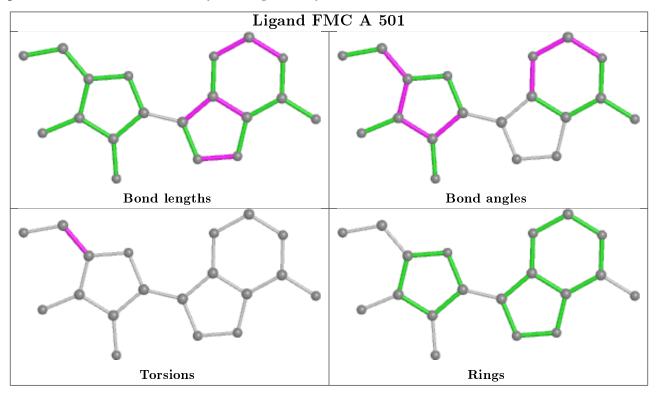
There are no torsion outliers.

There are no ring outliers.



No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and sufficient the outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



#### 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)

EDS failed to run properly - this section is therefore empty.

## 6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS failed to run properly - this section is therefore empty.

### 6.3 Carbohydrates (i)

EDS failed to run properly - this section is therefore empty.

### 6.4 Ligands (i)

EDS failed to run properly - this section is therefore empty.

### 6.5 Other polymers (i)

EDS failed to run properly - this section is therefore empty.

