

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

#### Jan 14, 2024 - 08:42 am GMT

PDB ID	:	6Y1S
Title	:	Bicyclic peptide bp70 in I4132 at 1.0 Angstrom resolution
Authors	:	Baeriswyl, S.; Stocker, A.; Reymond, JL.
Deposited on		
Resolution	:	1.04 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:

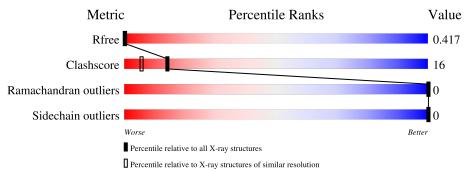
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as $541$ be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

# 1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.04 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$		
R <sub>free</sub>	130704	1596 (1.10-0.98)		
Clashscore	141614	1677 (1.10-0.98)		
Ramachandran outliers	138981	1591 (1.10-0.98)		
Sidechain outliers	138945	1589 (1.10-0.98)		

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%

Mol	Chain	Length		Quality of chain						
1	Δ	12	220/	620/	150/					
1	А	10	23%	62%	15%					



#### 6Y1S

# 2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 209 atoms, of which 104 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 bp70.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	А	13	Total 209	C 70	Н 104	N 19	0 14	S 2	0	0	1



# 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.

• Molecule 1: bp70

Chain A:	23%	62%	15%
0651 H2 A3 L4 A6 A6 C7	18 710 711 M1213 M1213		



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 41 3 2	Depositor
Cell constants	55.46Å $55.46$ Å $55.46$ Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	39.24 - 1.04	Depositor
	39.22 - 1.04	EDS
% Data completeness	99.9 (39.24-1.04)	Depositor
(in resolution range)	94.9 (39.22-1.04)	EDS
R <sub>merge</sub>	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	$0.55 (at 1.04 \text{\AA})$	Xtriage
Refinement program	SHELX	Depositor
D D.	(Not available), $0.234$	Depositor
$R, R_{free}$	0.437 , $0.417$	DCC
$R_{free}$ test set	363 reflections $(5.01%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	16.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.48, 39.8	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	0.83	EDS
Total number of atoms	209	wwPDB-VP
Average B, all atoms $(Å^2)$	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.54% 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: NH2, ORN, DAB, O65, NLE

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	Bond	lengths	Bond angles		
Mol (	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.79	0/70	2.40	6/90~(6.7%)	

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	2	HIS	CA-CB-CG	7.90	127.03	113.60
1	А	9	ARG	NE-CZ-NH1	7.06	123.83	120.30
1	А	9	ARG	NH1-CZ-NH2	-5.81	113.01	119.40
1	А	9	ARG	NE-CZ-NH2	5.73	123.16	120.30
1	А	11	TYR	CB-CG-CD1	5.09	124.06	121.00
1	А	11	TYR	CA-C-O	-5.04	109.51	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	А	105	104	87	3	0
All	All	105	104	87	3	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 16.



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)	
1:A:4:NLE:O	1:A:8:ILE:HD12	2.02	0.59	
1:A:7:CYS:O	1:A:10:CYS:HB2	2.17	0.44	
1:A:1:O65:O3	1:A:6:DAB:HB2	2.21	0.40	

All (3) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

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	Perce	ntiles
1	А	7/13~(54%)	5 (71%)	2(29%)	0	100	100

There are no Ramachandran outliers to report.

#### 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	7/7~(100%)	7~(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. 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)

3 non-standard protein/DNA/RNA residues 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).

Mal	Mol Type Chain Res		Link	B	Bond lengths			Bond angles		
IVIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2
1	NLE	А	4	1	6,7,8	5.65	4 (66%)	2,7,9	1.33	0
1	ORN	А	3	1	6,7,8	1.47	2 (33%)	2,7,9	1.80	1 (50%)
1	DAB	А	6	1	$5,\!6,\!7$	2.76	3 (60%)	1,6,8	4.74	1 (100%)

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.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	NLE	А	4	1	-	2/5/6/8	-
1	ORN	А	3	1	-	1/5/6/8	-
1	DAB	А	6	1	-	2/4/5/7	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	А	4	NLE	CB-CA	12.81	1.70	1.53
1	А	4	NLE	CA-N	-4.14	1.35	1.48
1	А	6	DAB	O-C	-3.90	1.04	1.19
1	А	6	DAB	CA-N	-3.84	1.36	1.48
1	А	6	DAB	CB-CA	2.52	1.56	1.53
1	А	4	NLE	CG-CB	-2.44	1.41	1.52
1	А	3	ORN	O-C	2.27	1.29	1.19
1	А	3	ORN	CB-CA	2.14	1.56	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	А	4	NLE	CE-CD	-2.06	1.33	1.49

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	6	DAB	CB-CG-ND	4.74	147.07	112.82
1	А	3	ORN	CG-CD-NE	-2.49	94.95	112.78

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	А	3	ORN	O-C-CA-CB
1	А	6	DAB	N-CA-CB-CG
1	А	6	DAB	CA-CB-CG-ND
1	А	4	NLE	CA-CB-CG-CD
1	А	4	NLE	CE-CD-CG-CB

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	А	4	NLE	1	0
1	А	6	DAB	1	0

## 5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry (i)

There are no ligands in this entry.

## 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)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

## 6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

## 6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

## 6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

