

wwPDB X-ray Structure Validation Summary Report (i)

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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 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
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-RAY DIFFRACTION

The reported resolution of this entry is 1.25 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	Percentile Ranks	3	Value
Clashscore			0
Ramachandran outliers			0
Sidechain outliers			0
	Worse	Better	
	Percentile relative to all X-ray structures		
	Percentile relative to X-ray structures of similar resolution		

Metric	Whole archive	Similar resolution
Metric	$(\# {\rm Entries})$	$(\# { m Entries}, { m resolution} { m range}({ m \AA}))$
Clashscore	141614	1060 (1.28-1.24)
Ramachandran outliers	138981	1029 (1.28-1.24)
Sidechain outliers	138945	1028 (1.28-1.24)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain					
1	А	7	71%	29%				
2	В	7	71%	29%				
3	С	7	86%	14%				



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 175 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 Collagen like peptide.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
1	А	7	Total 42	C 26	N 7	O 9	0	0	0

• Molecule 2 is a protein called Collagen like peptide.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
2	В	7	Total 46	C 29	N 7	0 10	0	0	0

• Molecule 3 is a protein called Collagen like peptide.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
3	С	7	Total 45	C 29	N 7	0 9	0	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	14	Total O 14 14	0	0
4	В	15	Total O 15 15	0	0
4	С	13	Total O 13 13	0	0



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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 was not executed.

• Molecule 1: Collagen like peptide

Chain A:	71%	29%
6108 1008 6109 6109		
• Molecule 2: Colla	gen like peptide	
Chain B:	71%	29%
P202 P205 P205		
• Molecule 3: Colla	gen like peptide	
Chain C:	86%	14%



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	14.05Å 26.77Å 20.00Å	Depositor
a, b, c, α , β , γ	90.00° 106.76° 90.00°	Depositor
Resolution (Å)	10.00 - 1.25	Depositor
% Data completeness	89.1 (10.00-1.25)	Depositor
(in resolution range)	05.1 (10.00-1.25)	Depositor
R_{merge}	0.06	Depositor
R _{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.135 , 0.189	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	175	wwPDB-VP
Average B, all atoms $(Å^2)$	16.0	wwPDB-VP



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: HYP

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		
IVIOI	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.67	0/25	1.10	0/30	
2	В	0.71	0/22	0.85	0/28	
3	С	0.59	0/29	0.88	0/36	
All	All	0.65	0/76	0.95	0/94	

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	42	0	36	0	1
2	В	46	0	42	0	0
3	С	45	0	41	0	1
4	А	14	0	0	0	0
4	В	15	0	0	0	0
4	С	13	0	0	0	0
All	All	175	0	119	0	1

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 0.



There are no clashes within the asymmetric unit.

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)	
1:A:109:GLY:C	3:C:301:PRO:N[1_556]	1.34	0.86	

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	3/7~(43%)	3~(100%)	0	0	100	100
2	В	4/7~(57%)	4 (100%)	0	0	100	100
3	С	3/7~(43%)	3 (100%)	0	0	100	100
All	All	10/21~(48%)	10 (100%)	0	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 side chain 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	Analysed Rotameric		Percentiles		
1	А	2/2~(100%)	2~(100%)	0	100 100		
2	В	2/2~(100%)	2(100%)	0	100 100		
3	С	3/3~(100%)	3 (100%)	0	100 100		
All	All	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)

7 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).

Mol	ol Type Chain Res Li			Link	В	gths	Bond angles				
INIOI	Type	Unam	nes	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
2	HYP	В	202	2,3	$6,\!8,\!9$	0.94	0	$5,\!10,\!12$	1.16	1 (20%)	
2	HYP	В	205	2	6,8,9	1.10	1 (16%)	$5,\!10,\!12$	1.39	0	
1	HYP	А	108	1	6,8,9	0.94	0	5,10,12	1.20	0	
3	HYP	С	305	3	$6,\!8,\!9$	0.96	0	$5,\!10,\!12$	0.96	0	
1	HYP	А	105	1	$6,\!8,\!9$	1.03	0	$5,\!10,\!12$	1.14	1 (20%)	
3	HYP	С	302	3	6,8,9	0.74	0	5,10,12	0.77	0	
2	HYP	В	208	2,1	6,8,9	1.09	0	5,10,12	0.94	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.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HYP	В	202	2,3	-	0/0/11/13	0/1/1/1
2	HYP	В	205	2	-	0/0/11/13	0/1/1/1
1	HYP	А	108	1	-	0/0/11/13	0/1/1/1
3	HYP	С	305	3	-	0/0/11/13	0/1/1/1
1	HYP	А	105	1	-	0/0/11/13	0/1/1/1
3	HYP	С	302	3	-	0/0/11/13	0/1/1/1
2	HYP	В	208	2,1	-	0/0/11/13	0/1/1/1



All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	В	205	HYP	CB-CG	-2.26	1.48	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	В	202	HYP	OD1-CG-CD	-2.31	105.29	110.35
1	А	105	HYP	CB-CG-CD	2.17	105.93	103.27

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

