

wwPDB X-ray Structure Validation Summary Report (i)

Sep 16, 2021 – 02:06 pm BST

PDB ID	:	70NW
Title	:	Crystal structure of PBP3 from E. coli in complex with AIC499
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Deposited on		
$\operatorname{Resolution}$:	2.70 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 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:

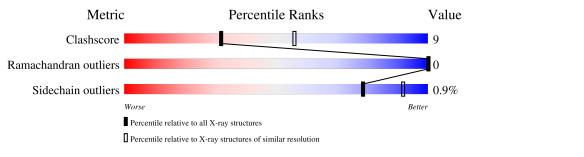
Xtriage (Phenix) EDS buster-report Percentile statistics Ideal geometry (proteins)	: : : :	20191225.v01 (using entries in the PDB archive December 25th 2019) Engh & Huber (2001)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)	:	Parkinson et al. (1996) 2.23.1

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 2.70 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	3122 (2.70-2.70)
Ramachandran outliers	138981	3069(2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)

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	564	68%	18%	14%



$70\mathrm{NW}$

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3593 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 Peptidoglycan D,D-transpeptidase FtsI.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Λ	485	Total	С	Ν	Ο	\mathbf{S}	0	1	0
	A	400	3516	2220	621	664	11	0	1	0

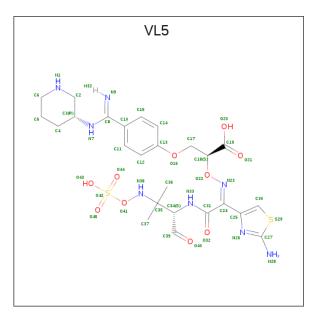
Chain	Residue	Modelled	Actual	Comment	Reference
А	25	MET	-	initiating methionine	UNP P0AD68
А	26	GLY	-	expression tag	UNP P0AD68
А	27	SER	-	expression tag	UNP P0AD68
A	28	SER	-	expression tag	UNP P0AD68
A	29	HIS	-	expression tag	UNP P0AD68
A	30	HIS	-	expression tag	UNP P0AD68
A	31	HIS	-	expression tag	UNP P0AD68
A	32	HIS	-	expression tag	UNP P0AD68
A	33	HIS	-	expression tag	UNP P0AD68
A	34	HIS	-	expression tag	UNP P0AD68
A	35	SER	-	expression tag	UNP P0AD68
A	36	SER	-	expression tag	UNP P0AD68
A	37	GLY	-	expression tag	UNP P0AD68
A	38	LEU	-	expression tag	UNP P0AD68
A	39	VAL	-	expression tag	UNP P0AD68
A	40	PRO	-	expression tag	UNP P0AD68
A	41	ARG	-	expression tag	UNP P0AD68
A	42	GLY	-	expression tag	UNP P0AD68
A	43	SER	-	expression tag	UNP P0AD68
А	44	HIS	-	expression tag	UNP P0AD68
A	45	MET	-	expression tag	UNP P0AD68
A	46	ALA	-	expression tag	UNP P0AD68
A	47	SER	-	expression tag	UNP P0AD68
А	48	MET	-	expression tag	UNP P0AD68

There are 24 discrepancies between the modelled and reference sequences:

• Molecule 2 is (2S)-2-[(Z)-[1-(2-azanyl-1,3-thiazol-4-yl)-2-[[(2S)-3-methyl-1-oxidanylidene-3-(sulfooxyamino)butan-2-yl]amino]-2-oxidanylidene-ethylidene]amino]oxy-3-[4-[N-[(3R)

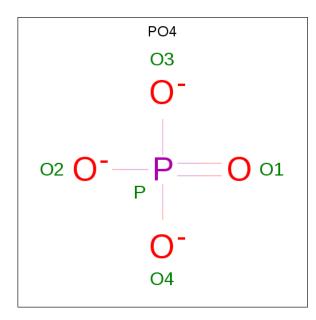


-piperidin-3-yl]carbamimidoyl]phenoxy]propanoic acid (three-letter code: VL5) (formula: $C_{25}H_{34}N_8O_{10}S_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues		Ato	\mathbf{pms}			ZeroOcc	AltConf
0	Λ	1	Total	С	Ν	Ο	S	0	0
	А	T	45	25	8	10	2	0	0

• Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Ato	\mathbf{ms}		Zero	Occ	AltConf
3	А	1	Total 5	O 4	Р 1	0		0

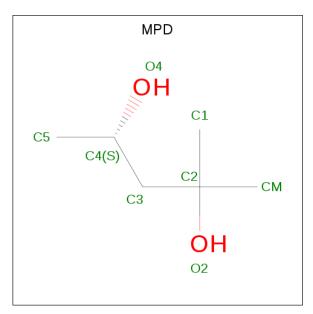
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Mol	Chain	Residues	Ate	\mathbf{pms}		ZeroOcc	AltConf
3	А	1	Total 5	O 4	Р 1	0	0

• Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0
4	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	6	Total O 6 6	0	0

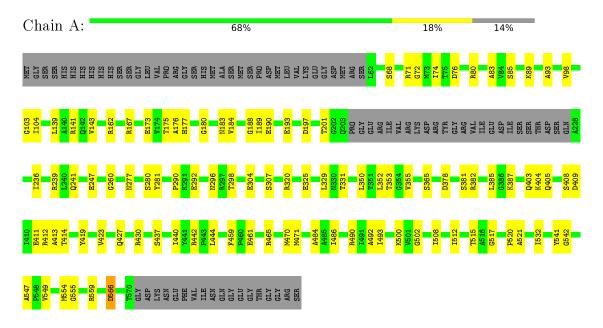


3 Residue-property plots (i)

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 was not executed.

• Molecule 1: Peptidoglycan D,D-transpeptidase FtsI





4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants	106.90Å 106.90 Å 285.83 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.64 - 2.70	Depositor
% Data completeness	51.1 (48.64-2.70)	Depositor
(in resolution range)	51.1 (40.04-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PHENIX 1.19.1	Depositor
R, R_{free}	0.234 , 0.271	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3593	wwPDB-VP
Average B, all atoms $(Å^2)$	82.0	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: VL5, PO4, MPD

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	Bond angles		
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.24	0/3586	0.48	0/4895

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	3516	0	3364	62	0
2	А	45	0	0	2	0
3	А	10	0	0	1	0
4	А	16	0	28	0	0
5	А	6	0	0	1	0
All	All	3593	0	3392	62	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 9.

The worst 5 of 62 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:385:LEU:HD22	1:A:423:VAL:HG11	1.70	0.73
1:A:381:SER:O	1:A:387:LYS:NZ	2.26	0.69
1:A:304:GLU:OE1	2:A:601:VL5:N28	2.27	0.67
1:A:71:ARG:HB3	1:A:167:ARG:HH12	1.59	0.67
1:A:85:SER:OG	1:A:167:ARG:NH1	2.30	0.65

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	Percentile	s
1	А	482/564~(86%)	458 (95%)	24~(5%)	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	A	341/460~(74%)	338~(99%)	3~(1%)	78 92

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	459	PHE
1	А	541	TYR

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Mol	Chain	\mathbf{Res}	Type
1	А	566	ASP

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)

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)

5 ligands 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	Trees	Chain	nain Res Link Bond lengths				Bond angles			
INIOI	Type	Chain	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
2	VL5	А	601	1	$35,\!47,\!47$	1.43	3 (8%)	$35,\!66,\!66$	1.94	<mark>5 (14%)</mark>
3	PO4	А	602	-	4,4,4	0.95	0	6, 6, 6	0.43	0
3	PO4	А	603	-	4,4,4	0.93	0	$6,\!6,\!6$	0.43	0
4	MPD	А	605	-	7,7,7	0.10	0	$9,\!10,\!10$	0.29	0
4	MPD	А	604	-	$7,\!7,\!7$	0.12	0	$9,\!10,\!10$	0.31	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.



Mol	Type	Chain	\mathbf{Res}	Link	Chirals	Torsions	Rings
2	VL5	А	601	1	-	$\frac{5/30}{57/57}$	0/3/3/3
4	MPD	А	604	-	-	0/5/5/5	-
4	MPD	А	605	-	-	0/5/5/5	-

'-' means no outliers of that kind were identified.

All (3) bond length outliers are listed below:

Mol	Chain	\mathbf{Res}	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	А	601	VL5	O22-N23	-5.66	1.26	1.41
2	А	601	VL5	C27-N28	3.67	1.46	1.35
2	А	601	VL5	C30-S29	2.79	1.75	1.70

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
2	А	601	VL5	O22-N23-C24	7.69	120.55	111.75
2	А	601	VL5	C24-C31-N33	4.39	121.68	114.38
2	А	601	VL5	C10-C8-N7	4.13	120.94	114.61
2	А	601	VL5	O32-C31-C24	-3.10	116.78	120.35
2	А	601	VL5	O41-S42-O44	2.07	109.86	103.29

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	\mathbf{Res}	Type	Atoms
2	А	601	VL5	C39-C34-C35-C36
2	А	601	VL5	C39-C34-C35-C37
2	А	601	VL5	N33-C34-C35-C36
2	А	601	VL5	N33-C34-C35-C37
2	А	601	VL5	C25-C24-C31-O32

There are no ring outliers.

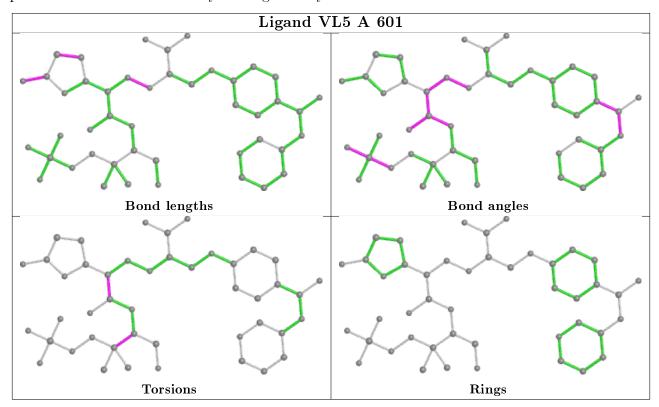
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	601	VL5	2	0
3	А	603	PO4	1	0

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 any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. 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 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.

