

Full wwPDB X-ray Structure Validation Report (i)

Jul 26, 2023 – 01:43 AM EDT

PDB ID	:	1BLI
Title	:	BACILLUS LICHENIFORMIS ALPHA-AMYLASE
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Deposited on		
Resolution	:	1.90 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:

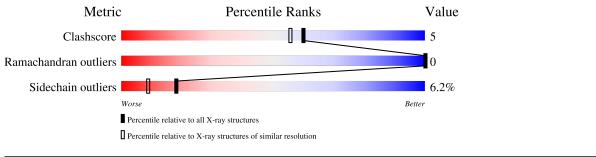
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.34

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 1.90 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	Similar resolution
Metric	$(\# {\rm Entries})$	$(\# { m Entries}, { m resolution} { m range}({ m \AA}))$
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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	483	87%	11%	•



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4130 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 ALPHA-AMYLASE.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	481	Total 3905	C 2491	N 678	0 729	S 7	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	134	LEU	ARG	conflict	UNP P06278
А	190	PHE	ASN	engineered mutation	UNP P06278
А	264	SER	GLN	engineered mutation	UNP P06278
А	265	TYR	ASN	engineered mutation	UNP P06278
А	310	GLY	SER	conflict	UNP P06278
А	320	SER	ALA	conflict	UNP P06278
А	465	ALA	GLU	conflict	UNP P06278

• Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Ca 3 3	0	0

• Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	Total Na 1 1	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	221	Total O 221 221	0	0

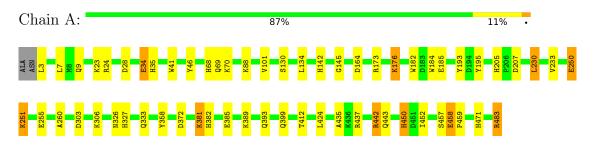


3 Residue-property plots (i)

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: ALPHA-AMYLASE





4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 61	Depositor
Cell constants	91.30Å 91.30Å 137.70Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 - 1.90	Depositor
% Data completeness	89.4 (30.00-1.90)	Depositor
(in resolution range)	05.4 (00.00-1.00)	Depositor
R_{merge}	0.08	Depositor
R _{sym}	0.08	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.154 , 0.185	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4130	wwPDB-VP
Average B, all atoms $(Å^2)$	21.0	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: NA, CA

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	0.63	0/4029	0.75	1/5472~(0.0%)	

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	7	LEU	CA-CB-CG	6.10	129.33	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	195	TYR	Sidechain

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	3905	0	3645	35	0
2	А	3	0	0	0	0
3	А	1	0	0	0	0
4	А	221	0	0	2	0
All	All	4130	0	3645	35	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 5.

All (35) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:389:LYS:HE2	1:A:393:GLN:HE22	1.29	0.96
1:A:389:LYS:HG3	1:A:393:GLN:NE2	1.87	0.89
1:A:389:LYS:HG3	1:A:393:GLN:HE21	1.38	0.88
1:A:450:HIS:HB3	1:A:483:ARG:NH1	1.96	0.80
1:A:458:GLU:H	1:A:458:GLU:CD	1.81	0.79
1:A:459:PRO:HB3	1:A:483:ARG:HH12	1.49	0.78
1:A:450:HIS:HB3	1:A:483:ARG:HH12	1.51	0.76
1:A:389:LYS:CE	1:A:393:GLN:HE22	2.03	0.71
1:A:9:GLN:HE22	1:A:326:ASN:HB2	1.58	0.67
1:A:303:ASP:HB3	1:A:306:LYS:HD2	1.77	0.66
1:A:389:LYS:HG2	1:A:452:ILE:HD11	1.85	0.59
1:A:459:PRO:HB3	1:A:483:ARG:NH1	2.16	0.58
1:A:389:LYS:HE2	1:A:393:GLN:NE2	2.10	0.58
1:A:35:HIS:CE1	1:A:385:GLU:HG2	2.40	0.56
1:A:250:GLU:HG3	1:A:251:LYS:N	2.19	0.55
1:A:381:LYS:HG2	1:A:382:HIS:N	2.21	0.54
1:A:24:ARG:NH1	4:A:2012:HOH:O	2.40	0.54
1:A:164:ASP:HA	1:A:173:ARG:O	2.09	0.53
1:A:101:VAL:HB	1:A:230:LEU:HD12	1.93	0.49
1:A:68:HIS:HD2	4:A:1028:HOH:O	1.95	0.48
1:A:184:TRP:CD2	1:A:185:GLU:HA	2.48	0.48
1:A:458:GLU:CD	1:A:458:GLU:N	2.61	0.48
1:A:424:LEU:C	1:A:424:LEU:HD12	2.33	0.48
1:A:34:GLU:HG2	1:A:35:HIS:CD2	2.48	0.47
1:A:205:HIS:CE1	1:A:207:ASP:HB2	2.50	0.46
1:A:233:VAL:HG11	1:A:260:ALA:HB1	1.98	0.45
1:A:435:ALA:HB2	1:A:471:HIS:CE1	2.52	0.44
1:A:176:LYS:HG3	1:A:182:TRP:CZ2	2.54	0.42
1:A:389:LYS:CG	1:A:393:GLN:NE2	2.71	0.42
1:A:69:GLN:HG2	1:A:145:GLY:HA3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:GLN:HA	1:A:412:THR:O	2.20	0.41
1:A:184:TRP:CG	1:A:185:GLU:HA	2.55	0.41
1:A:327:HIS:O	1:A:333:GLN:HG3	2.20	0.41
1:A:442:ARG:HH11	1:A:443:GLN:HA	1.85	0.41
1:A:70:LYS:HA	1:A:70:LYS:HD2	1.77	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	Percenti	les
1	А	479/483~(99%)	466~(97%)	13 (3%)	0	100 10)0

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	401/402~(100%)	376~(94%)	25~(6%)	18 9

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	3	LEU
	<i>a i</i> :	7	,

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	\mathbf{Res}	Type
А	23	LYS
А	28	ASP
А	34	GLU
А	41	TRP
	46	TYR
	88	LYS
А	130	SER
	134	LEU
	142	HIS
	176	LYS
А	193	TYR
А	230	LEU
А	250	GLU
А	251	LYS
А	255	GLU
А	358	TYR
	372	ASP
	381	LYS
	437	ARG
А	442	ARG
	450	HIS
	457	SER
	458	GLU
А	483	ARG
	A A	A 34 A 41 A 46 A 88 A 130 A 134 A 134 A 134 A 134 A 142 A 142 A 142 A 142 A 1230 A 230 A 250 A 251 A 255 A 358 A 372 A 381 A 437 A 442 A 450 A 457 A 458

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Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such side chains are listed below:

Mol	Chain	Res	Type
1	А	9	GLN
1	А	27	ASN
1	А	35	HIS
1	А	68	HIS
1	А	96	ASN
1	А	393	GLN
1	А	421	ASN
1	А	482	GLN

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 4 ligands modelled in this entry, 4 are monoatomic - 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.

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.

