

Full wwPDB X-ray Structure Validation Report (i)

May 16, 2020 – 09:31 am BST

PDB ID : 5YYC

Title : Crystal structure of alanine racemase from Bacillus pseudofirmus (OF4) Authors

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1.80 Å(reported) Resolution

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:

MolProbity 4.02b-467

> 1.8.5 (274361), CSD as541be (2020) Mogul

Xtriage (Phenix) 1.13 EDS 2.11

20191225.v01 (using entries in the PDB archive December 25th 2019) Percentile statistics

> Refmac 5.8.0158

7.0.044 (Gargrove) CCP4 Engh & Huber (2001)

Ideal geometry (proteins) Ideal geometry (DNA, RNA) Parkinson et al. (1996)

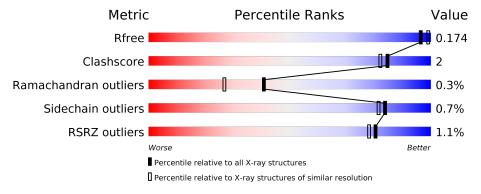
Validation Pipeline (wwPDB-VP) 2.11

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 1.80 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain	
1	A	369	93%	6%
1	С	369	95%	• •



2 Entry composition (i)

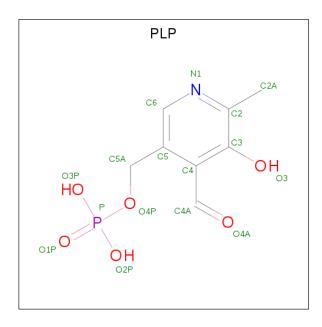
There are 3 unique types of molecules in this entry. The entry contains 6565 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Alanine racemase.

\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Δ	368	Total	С	N	О	S	0	0	0
1	Λ	300	2821	1778	481	546	16	U	U	
1	C	368	Total	С	N	О	S	0	0	0
T		300	2821	1778	481	546	16			

• Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf		
2	٨	1	Total	С	N	О	Р	0	0
	2 A	1	15	8	1	5	1	0	0
2	C	1	Total	С	N	О	Р	0	0
		1	15	8	1	5	1	0	U

• Molecule 3 is water.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	451	Total O 451 451	0	0
3	С	442	Total O 442 442	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). 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: Alanine racemase

Chain A:

93%

6%

Molecule 1: Alanine racemase

Chain C:

95%

• Molecule 1: Alanine racemase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	97.30Å 97.31Å 112.57Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.77 - 1.80	Depositor
rtesolution (A)	21.77 - 1.80	EDS
% Data completeness	99.7 (21.77-1.80)	Depositor
(in resolution range)	99.7 (21.77-1.80)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.42 (at 1.80Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
D D	0.149 , 0.174	Depositor
R, R_{free}	0.151 , 0.174	DCC
R_{free} test set	5003 reflections $(5.06%)$	wwPDB-VP
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.44 , 59.6	EDS
L-test for twinning ²	$< L >=0.51, < L^2>=0.34$	Xtriage
Estimated twinning fraction	0.488 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6565	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.84% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $< L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities 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: PLP

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		
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.40	0/2872	0.53	0/3890	
1	С	0.42	0/2872	0.55	1/3890 (0.0%)	
All	All	0.41	0/5744	0.54	1/7780 (0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
1	С	123	ALA	N-CA-C	5.84	126.78	111.00

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	A	2821	0	2812	17	0
1	С	2821	0	2812	13	0
2	A	15	0	6	0	0
2	С	15	0	6	0	0
3	A	451	0	0	2	0
3	С	442	0	0	4	0
All	All	6565	0	5636	26	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 2.

All (26) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
		$\operatorname{distance}\left(\mathrm{\AA}\right)$	overlap (Å)
1:A:16:GLN:N	1:A:368:GLU:OE2	2.09	0.85
1:C:12:GLN:NE2	3:C:502:HOH:O	2.20	0.75
1:C:345:GLU:OE2	3:C:501:HOH:O	2.18	0.55
1:A:145:LYS:NZ	3:A:504:HOH:O	2.40	0.53
1:A:95:LEU:HD13	1:A:116:LEU:HD13	1.90	0.53
1:A:271:ARG:HD3	1:C:180:TYR:CE2	2.43	0.52
1:C:266:SER:OG	1:C:271:ARG:HD2	2.10	0.52
1:A:180:TYR:CE2	1:C:271:ARG:HD3	2.46	0.50
1:A:266:SER:OG	1:A:271:ARG:HD2	2.13	0.49
1:A:15:LEU:HB2	1:A:368:GLU:HG3	1.94	0.49
1:A:260:LYS:HG2	1:A:280:VAL:HG22	1.94	0.47
1:C:145:LYS:NZ	3:C:505:HOH:O	2.48	0.47
1:A:12:GLN:HG3	1:A:367:GLU:HG3	1.96	0.47
1:A:15:LEU:N	1:A:368:GLU:OE2	2.48	0.47
1:A:345:GLU:HG2	3:A:913:HOH:O	2.14	0.46
1:C:95:LEU:HD13	1:C:116:LEU:HD13	1.97	0.46
1:C:120:ALA:C	1:C:122:GLU:N	2.70	0.45
1:A:289:ALA:HB2	1:A:317:MET:HG3	1.99	0.44
1:A:68:ILE:HG21	1:C:7:ARG:HD3	2.00	0.43
1:C:367:GLU:OE2	3:C:503:HOH:O	2.21	0.42
1:A:193:GLU:O	1:A:197:LYS:HG3	2.19	0.42
1:A:96:ARG:O	1:A:99:ILE:HG13	2.20	0.41
1:C:7:ARG:HD2	1:C:285:PRO:HB2	2.01	0.41
1:A:7:ARG:HD2	1:A:285:PRO:HB2	2.01	0.41
1:C:289:ALA:HB2	1:C:317:MET:HG3	2.02	0.41
1:A:271:ARG:HD3	1:C:180:TYR:CZ	2.56	0.40

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	\mathbf{ntiles}
1	A	366/369 (99%)	350 (96%)	16 (4%)	0	100	100
1	С	366/369~(99%)	349 (95%)	15 (4%)	2 (0%)	29	15
All	All	732/738 (99%)	699 (96%)	31 (4%)	2 (0%)	41	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	С	123	ALA
1	С	121	SER

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	${f Analysed}$	Rotameric	Outliers	Percentiles		
1	A	$307/308 \; (100\%)$	305 (99%)	2 (1%)	84 81		
1	С	307/308 (100%)	305 (99%)	2 (1%)	84 81		
All	All	614/616 (100%)	610 (99%)	4 (1%)	84 81		

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	TYR
1	A	368	GLU
1	С	63	TYR
1	С	121	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

2 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 Type	Chain	Chain	Chain	Chain	asin Des	Res Link	Bo	Bond lengths			Bond angles		
MIGI	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	$\mid \# Z > 2$			
2	PLP	С	401	1	15,15,16	2.53	7 (46%)	20,22,23	1.63	5 (25%)			
2	PLP	A	401	1	15,15,16	2.56	7 (46%)	20,22,23	1.53	6 (30%)			

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	PLP	С	401	1	-	0/6/6/8	0/1/1/1
2	PLP	A	401	1	_	0/6/6/8	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	${ m Observed}({ m \AA})$	$Ideal(\AA)$
2	С	401	PLP	C5-C4	5.26	1.46	1.40
2	A	401	PLP	C5-C4	5.06	1.46	1.40
2	A	401	PLP	C3-C2	4.55	1.45	1.40
2	С	401	PLP	C3-C2	4.23	1.45	1.40
2	A	401	PLP	P-O2P	-3.63	1.40	1.54
2	A	401	PLP	P-O3P	-3.62	1.40	1.54

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Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(\mathbf{\mathring{A}})$	$Ideal(\AA)$
2	С	401	PLP	P-O3P	-3.40	1.41	1.54
2	С	401	PLP	P-O2P	-3.37	1.41	1.54
2	A	401	PLP	P-O1P	-2.83	1.41	1.50
2	С	401	PLP	P-O1P	-2.82	1.41	1.50
2	С	401	PLP	C3-C4	2.74	1.45	1.40
2	A	401	PLP	C3-C4	2.61	1.45	1.40
2	С	401	PLP	O3-C3	-2.27	1.31	1.37
2	A	401	PLP	O3-C3	-2.17	1.31	1.37

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
2	С	401	PLP	C2A-C2-C3	-3.16	116.98	120.89
2	A	401	PLP	C2A-C2-C3	-3.04	117.13	120.89
2	С	401	PLP	C2A-C2-N1	2.98	123.50	117.67
2	С	401	PLP	C6-N1-C2	2.76	124.27	119.17
2	A	401	PLP	C2A-C2-N1	2.75	123.03	117.67
2	A	401	PLP	C5A-C5-C6	2.63	123.70	119.37
2	A	401	PLP	C6-N1-C2	2.37	123.56	119.17
2	С	401	PLP	C4A-C4-C5	2.24	123.24	120.94
2	С	401	PLP	C5A-C5-C6	2.22	123.02	119.37
2	A	401	PLP	O3-C3-C2	2.13	122.13	117.49
2	A	401	PLP	C4A-C4-C5	2.05	123.05	120.94

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)

In the following table, the column labelled '#RSRZ>2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ $>$	$\#\mathrm{RSRZ}{>}2$		$\mathbf{OWAB}(\mathbf{\AA}^2)$	Q < 0.9
1	A	368/369 (99%)	-0.52	5 (1%) 75	72	16, 22, 39, 87	0
1	С	368/369 (99%)	-0.51	3 (0%) 86	84	15, 22, 39, 105	0
All	All	736/738 (99%)	-0.52	8 (1%) 80	78	15, 22, 39, 105	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	С	121	SER	3.8
1	A	124	GLU	3.0
1	A	368	GLU	2.5
1	A	122	GLU	2.4
1	A	367	GLU	2.4
1	С	368	GLU	2.4
1	С	367	GLU	2.2
1	A	123	ALA	2.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,



median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\operatorname{\textbf{B-factors}}(\AA^2)$	Q < 0.9
2	PLP	С	401	15/16	0.98	0.09	16,18,23,35	0
2	PLP	A	401	15/16	0.99	0.08	16,18,23,36	0

6.5 Other polymers (i)

There are no such residues in this entry.

