

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

Dec 16, 2023 – 03:40 PM EST

PDB ID : 4PFF

> Title : Crystal structure of Plasmodium vivax SHMT with PLP Schiff base

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2014-04-29 Deposited on

2.30 Å(reported) Resolution

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

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

Xtriage (Phenix) 1.13

EDS 2.36

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

> Refmac 5.8.0158

CCP4 7.0.044 (Gargrove)

Ideal geometry (proteins) Engh & Huber (2001) Ideal geometry (DNA, RNA) Parkinson et al. (1996)

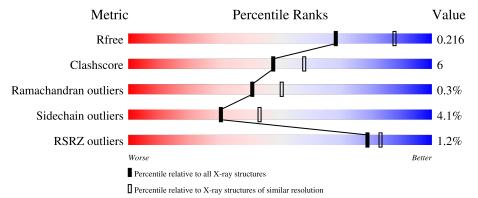
Validation Pipeline (wwPDB-VP) 2.36

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	$(\#  ext{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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% 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	442	84%	15%	•
1	В	442	86%	14%	•
1	С	442	83%	16%	•



# 2 Entry composition (i)

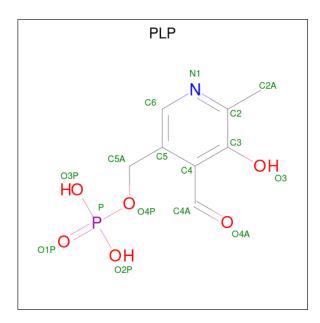
There are 3 unique types of molecules in this entry. The entry contains 10911 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Serine hydroxymethyltransferase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	٨	442	Total	С	N	О	S	0	1	0
1	A	442	3464	2189	601	656	18	0	1	0
1	В	442	Total	С	N	О	S	0	1	0
1	Б	D 442	3464	2189	601	656	18	0	1	
1	С	442	Total	С	N	О	S	0	1	0
	442	3464	2189	601	656	18		1		

• Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C<sub>8</sub>H<sub>10</sub>NO<sub>6</sub>P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
2	Δ	1	Total	С	N	О	Р	0	0	
	2 A	1	15	8	1	5	1	U	0	
9	R	1	Total	$\mathbf{C}$	Ν	Ο	Р	0	0	
2	Ъ	1	15	8	1	5	1	0		
2	С	1	Total	С	N	О	Р	0	0	
2		C		15	8	1	5	1	0	



### • Molecule 3 is water.

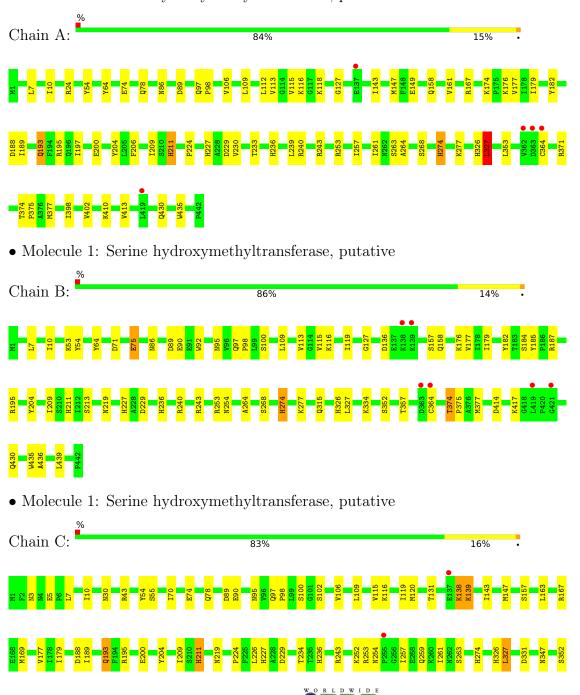
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	155	Total O 155 155	0	0
3	В	170	Total O 170 170	0	0
3	С	149	Total O 149 149	0	0



# 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 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: Serine hydroxymethyltransferase, putative







# 4 Data and refinement statistics (i)

Property	Value	Source	
Space group	C 1 2 1	Depositor	
Cell constants	101.05Å 58.33Å 237.14Å	Depositor	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.01^{\circ}$ $90.00^{\circ}$	-	
Resolution (Å)	30.00 - 2.30	Depositor	
resolution (A)	29.64 - 2.30	EDS	
% Data completeness	94.4 (30.00-2.30)	Depositor	
(in resolution range)	94.4 (29.64-2.30)	EDS	
$R_{merge}$	0.03	Depositor	
$\frac{R_{sym}}{\langle I/\sigma(I)\rangle^{-1}}$	(Not available)	Depositor	
$< I/\sigma(I) > 1$	10.57  (at  2.29Å)	Xtriage	
Refinement program	REFMAC 5.6.0117	Depositor	
P.P.	0.221 , $0.240$	Depositor	
$R, R_{free}$	0.217 , $0.216$	DCC	
$R_{free}$ test set	5989 reflections $(10.18\%)$	wwPDB-VP	
Wilson B-factor (Å <sup>2</sup> )	31.3	Xtriage	
Anisotropy	0.815	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.33, 18.2	EDS	
L-test for twinning <sup>2</sup>	$< L >=0.50, < L^2>=0.33$	Xtriage	
	0.034  for  -1/2 *h- 3/2 *k, -1/2 *h+ 1/2 *k, -1		
	0.034  for  -1/2 * h + 3/2 * k, 1/2 * h + 1/2 * k, - l		
Estimated twinning fraction	0.480  for  1/2*h-3/2*k,-1/2*h-1/2*k,-1	Xtriage	
	0.487  for  1/2*h+3/2*k,1/2*h-1/2*k,-l		
	0.029  for  -h,-k,l		
$F_o, F_c$ correlation	0.94	EDS	
Total number of atoms	10911	wwPDB-VP	
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.26% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<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: 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	Bo	nd lengths	Bond angles		
IVIOI		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.45	1/3527~(0.0%)	0.59	$1/4762 \ (0.0\%)$	
1	В	0.46	2/3527~(0.1%)	0.59	0/4762	
1	С	0.45	1/3527~(0.0%)	0.58	0/4762	
All	All	0.45	4/10581 (0.0%)	0.59	1/14286 (0.0%)	

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(A)
1	С	435	TRP	CD2-CE2	5.54	1.48	1.41
1	В	435	TRP	CD2-CE2	5.14	1.47	1.41
1	A	435	TRP	CD2-CE2	5.03	1.47	1.41
1	В	92	TRP	CD2-CE2	5.01	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(^{o})$	$\operatorname{Ideal}({}^o)$
1	A	327	LEU	CA-CB-CG	5.29	127.47	115.30

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	3464	0	3477	48	0
1	В	3464	0	3477	38	0
1	С	3464	0	3477	45	0
2	A	15	0	6	0	0
2	В	15	0	6	1	0
2	С	15	0	6	1	0
3	A	155	0	0	8	0
3	В	170	0	0	6	0
3	С	149	0	0	7	0
All	All	10911	0	10449	127	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 6.

The worst 5 of 127 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:B:236:HIS:CE1	1:B:243:ARG:HD2	2.00	0.96
1:A:236:HIS:CE1	1:A:243:ARG:HD2	2.18	0.78
2:B:501:PLP:C4A	3:B:761:HOH:O	2.35	0.74
1:C:436:ALA:HA	1:C:439:LEU:HD12	1.70	0.73
1:B:53:LYS:NZ	3:B:770:HOH:O	2.24	0.71

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	Percentiles
1	A	441/442 (100%)	427 (97%)	13 (3%)	1 (0%)	47 58
1	В	441/442 (100%)	423 (96%)	16 (4%)	2 (0%)	29 35
1	$\mathbf{C}$	441/442 (100%)	421 (96%)	19 (4%)	1 (0%)	47 58

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1323/1326 (100%)	1271 (96%)	48 (4%)	4 (0%)	41 50	

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	С	89	ASP
1	В	89	ASP
1	A	89	ASP
1	В	374	THR

#### 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	Perce	ntiles
1	A	382/381 (100%)	370 (97%)	12 (3%)	40	55
1	В	382/381 (100%)	364 (95%)	18 (5%)	26	37
1	С	382/381 (100%)	365 (96%)	17 (4%)	27	39
All	All	1146/1143 (100%)	1099 (96%)	47 (4%)	30	43

5 of 47 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	364	CYS
1	С	157	SER
1	В	430	GLN
1	С	100	SER
1	С	211	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
1	В	274	HIS
1	С	86	ASN

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Mol	Chain	Res	Type
1	С	426	GLN
1	С	211	HIS
1	С	227	HIS

#### 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)

3 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	Dec	Res Link Bond lengths			Bond angles			
WIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	PLP	A	501	1	15,15,16	0.89	1 (6%)	20,22,23	1.65	4 (20%)
2	PLP	С	501	1	15,15,16	0.81	0	20,22,23	1.93	5 (25%)
2	PLP	В	501	1	15,15,16	0.82	0	20,22,23	1.67	3 (15%)

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	A	501	1	-	4/6/6/8	0/1/1/1
2	PLP	С	501	1	-	5/6/6/8	0/1/1/1
2	PLP	В	501	1	-	2/6/6/8	0/1/1/1

#### All (1) bond length outliers are listed below:

N	/Iol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(\text{\AA})$	Ideal(A)
	2	A	501	PLP	C3-C2	-2.42	1.38	1.40

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}(^{o})$
2	С	501	PLP	C4A-C4-C5	4.32	125.39	120.94
2	С	501	PLP	C5A-C5-C6	-4.03	112.74	119.37
2	В	501	PLP	C4A-C4-C5	3.96	125.01	120.94
2	A	501	PLP	C4A-C4-C5	3.85	124.90	120.94
2	С	501	PLP	O4P-C5A-C5	3.74	116.47	109.35

There are no chirality outliers.

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	PLP	C5A-O4P-P-O2P
2	В	501	PLP	C5A-O4P-P-O2P
2	С	501	PLP	C4-C5-C5A-O4P
2	С	501	PLP	C6-C5-C5A-O4P
2	С	501	PLP	C5A-O4P-P-O1P

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	С	501	PLP	1	0
2	В	501	PLP	1	0

# 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 $>$	$\#\mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q<0.9
1	A	442/442 (100%)	-0.10	5 (1%) 80 85	22, 33, 51, 81	0
1	В	442/442 (100%)	-0.05	6 (1%) 75 80	22, 34, 51, 72	0
1	С	442/442 (100%)	-0.03	5 (1%) 80 85	21, 33, 51, 77	0
All	All	1326/1326 (100%)	-0.06	16 (1%) 79 83	21, 33, 51, 81	0

The worst 5 of 16 RSRZ outliers are listed below:

Mol	Chain	$\operatorname{Res}$	Type	RSRZ
1	В	364	CYS	4.4
1	A	363	ASP	4.2
1	A	364	CYS	4.1
1	С	363	ASP	3.7
1	С	137	GLU	3.5

## 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 monosaccharides 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	PLP	A	501	15/16	0.97	0.13	28,33,35,37	0
2	PLP	В	501	15/16	0.97	0.14	27,31,34,35	0
2	PLP	С	501	15/16	0.97	0.12	27,30,32,33	0

# 6.5 Other polymers (i)

There are no such residues in this entry.

