



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:36 PM GMT

PDB ID : 4PFF  
Title : Crystal structure of Plasmodium vivax SHMT with PLP Schiff base  
Authors : Chitnumsub, P.; Jaruwat, A.; Leartsakulpanich, U.  
Deposited on : 2014-04-29  
Resolution : 2.30 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

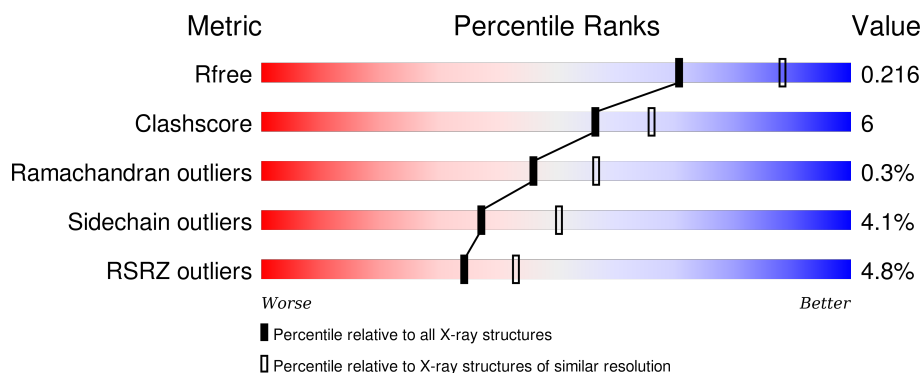
# 1 Overall quality at a glance

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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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  $\leq 5\%$ . 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	<div> <div>5%</div> <div>84%</div> <div>15%</div> </div>
1	B	442	<div> <div>5%</div> <div>86%</div> <div>14%</div> </div>
1	C	442	<div> <div>4%</div> <div>83%</div> <div>16%</div> </div>

## 2 Entry composition [i](#)

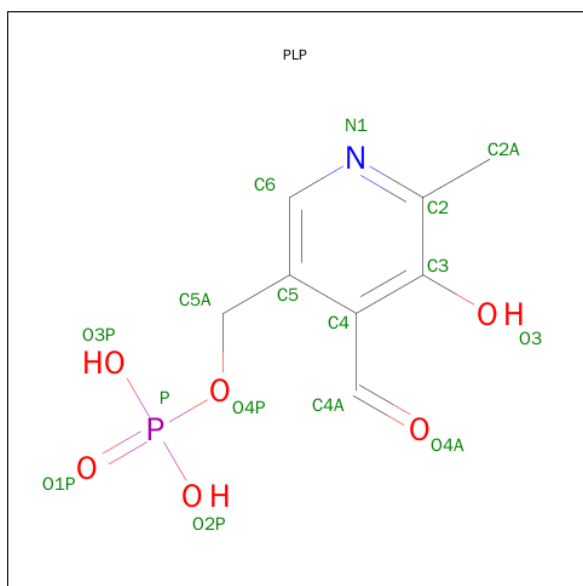
There are 3 unique types of molecules in this entry. The entry contains 10911 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	A	442	Total	C	N	O	S	0	1	0
			3464	2189	601	656	18			
1	B	442	Total	C	N	O	S	0	1	0
			3464	2189	601	656	18			
1	C	442	Total	C	N	O	S	0	1	0
			3464	2189	601	656	18			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula:  $C_8H_{10}NO_6P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	C	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

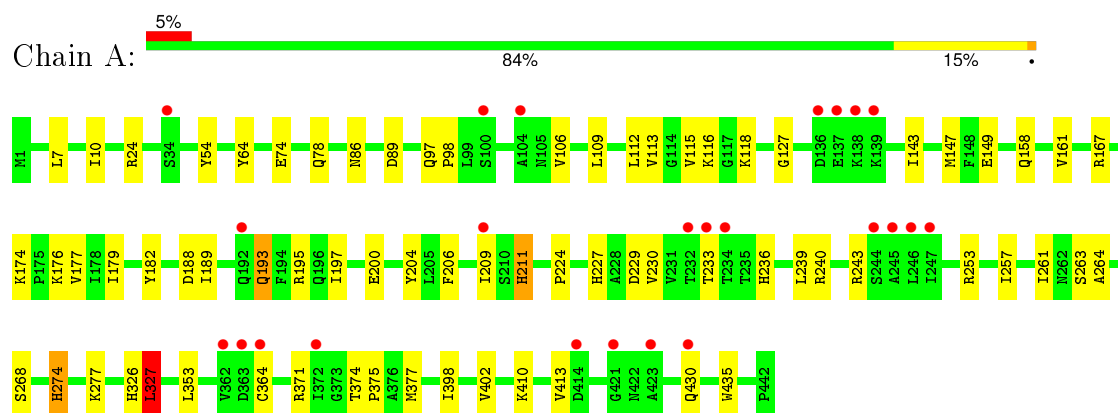
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	155	Total 155	O 155	0	0
3	B	170	Total 170	O 170	0	0
3	C	149	Total 149	O 149	0	0

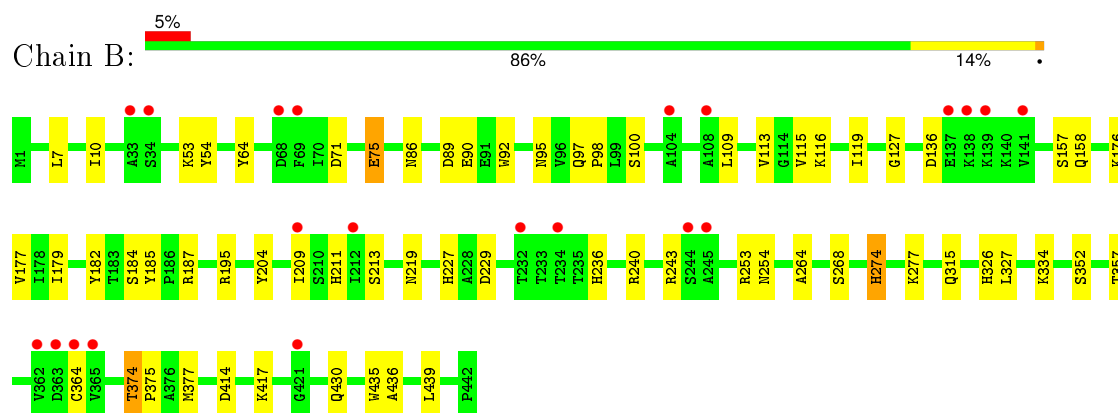
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

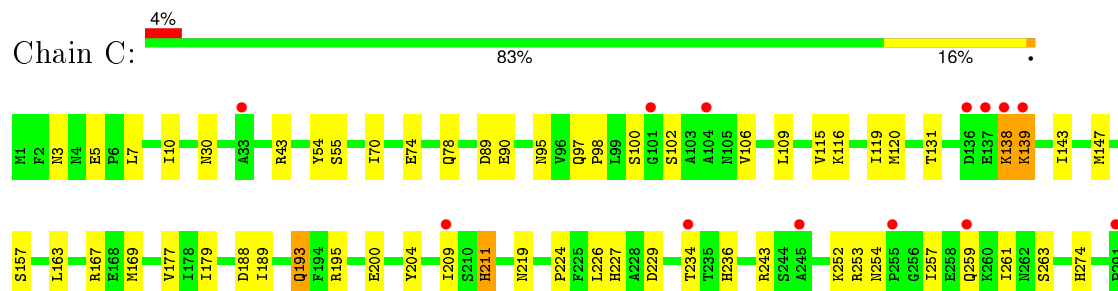
- Molecule 1: Serine hydroxymethyltransferase, putative

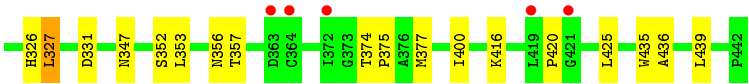


- Molecule 1: Serine hydroxymethyltransferase, putative



- Molecule 1: Serine hydroxymethyltransferase, putative





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.05Å 58.33Å 237.14Å 90.00° 90.01° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 29.64 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.4 (30.00-2.30) 94.4 (29.64-2.30)	Depositor EDS
$R_{merge}$	0.03	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	10.57 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.221 , 0.240 0.217 , 0.216	Depositor DCC
$R_{free}$ test set	5962 reflections (11.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.3	Xtriage
Anisotropy	0.815	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 35.7	EDS
Estimated twinning fraction	0.034 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.034 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.480 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.487 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.029 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 58848 reflections	Xtriage
$F_o, F_c$ 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>1</sup> Intensities estimated from amplitudes.

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

## 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	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.45	1/3527 (0.0%)	0.59	1/4762 (0.0%)
1	B	0.46	2/3527 (0.1%)	0.59	0/4762
1	C	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	Observed(Å)	Ideal(Å)
1	C	435	TRP	CD2-CE2	5.54	1.48	1.41
1	B	435	TRP	CD2-CE2	5.14	1.47	1.41
1	A	435	TRP	CD2-CE2	5.03	1.47	1.41
1	B	92	TRP	CD2-CE2	5.01	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
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	B	3464	0	3477	38	0
1	C	3464	0	3477	45	0
2	A	15	0	6	0	0
2	B	15	0	6	1	0
2	C	15	0	6	1	0
3	A	155	0	0	8	0
3	B	170	0	0	6	0
3	C	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.

All (127) 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: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
1:B:436:ALA:HA	1:B:439:LEU:HD12	1.72	0.71
1:C:236:HIS:CE1	1:C:243:ARG:HD2	2.27	0.70
1:B:109:LEU:HD21	1:B:179:ILE:HD11	1.74	0.69
1:A:158:GLN:HG3	3:A:618:HOH:O	1.91	0.68
1:C:138:LYS:HG2	1:C:139:LYS:N	2.08	0.68
1:C:163:LEU:HD13	1:C:193:GLN:HG2	1.76	0.67
1:B:240:ARG:HD3	3:B:766:HOH:O	1.94	0.67
1:A:377:MET:HE3	3:A:662:HOH:O	1.96	0.66
1:C:109:LEU:HD21	1:C:179:ILE:HD11	1.79	0.64
1:A:64:TYR:HD2	3:B:769:HOH:O	1.79	0.64
1:A:74:GLU:O	1:A:78:GLN:HG3	1.98	0.62
1:A:374:THR:N	1:A:375:PRO:CD	2.63	0.61
1:A:374:THR:N	1:A:375:PRO:HD3	2.16	0.61
1:B:374:THR:N	1:B:375:PRO:CD	2.64	0.60
1:A:240:ARG:HD3	3:A:753:HOH:O	2.02	0.60
1:C:229:ASP:OD1	1:C:253:ARG:HD2	2.00	0.60
1:A:188:ASP:OD1	1:A:189:ILE:N	2.34	0.59
1:B:64:TYR:HD2	3:B:770:HOH:O	1.84	0.59
1:C:374:THR:N	1:C:375:PRO:CD	2.65	0.58
1:C:167:ARG:NH1	1:C:200:GLU:OE1	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ILE:CG1	1:A:147:MET:HE2	2.34	0.57
1:A:277:LYS:HZ1	1:B:274:HIS:HD2	1.52	0.57
1:A:161:VAL:HG21	3:A:709:HOH:O	2.04	0.57
1:B:195:ARG:HD2	1:B:227:HIS:O	2.03	0.57
1:C:7:LEU:HA	1:C:10:ILE:HG22	1.87	0.57
1:C:211:HIS:HD2	1:C:326:HIS:NE2	2.03	0.57
1:B:7:LEU:HA	1:B:10:ILE:HG22	1.85	0.57
1:B:236:HIS:ND1	1:B:243:ARG:HD2	2.19	0.56
1:A:277:LYS:NZ	1:B:274:HIS:HD2	2.04	0.56
1:A:229:ASP:OD1	1:A:253:ARG:HD2	2.06	0.55
1:B:414:ASP:HA	1:B:417:LYS:HE2	1.88	0.55
1:A:374:THR:H	1:A:375:PRO:HD3	1.71	0.55
1:B:229:ASP:OD1	1:B:253:ARG:HD2	2.07	0.55
1:C:374:THR:N	1:C:375:PRO:HD3	2.21	0.54
2:C:501:PLP:C4A	3:C:646:HOH:O	2.54	0.54
1:A:193:GLN:O	1:A:197:ILE:HG13	2.07	0.54
1:A:398:ILE:O	1:A:402:VAL:HG23	2.08	0.54
1:B:374:THR:H	1:B:375:PRO:HD3	1.73	0.54
1:C:211:HIS:HE1	3:C:658:HOH:O	1.91	0.54
1:B:71:ASP:O	1:B:75:GLU:HB2	2.08	0.53
1:C:195:ARG:HD2	1:C:227:HIS:O	2.08	0.53
1:A:430:GLN:HG2	3:A:679:HOH:O	2.09	0.53
1:A:167:ARG:NH1	1:A:200:GLU:OE1	2.41	0.53
1:C:326:HIS:H	1:C:326:HIS:HD1	1.56	0.52
1:B:236:HIS:CE1	1:B:243:ARG:CD	2.86	0.52
1:C:138:LYS:HG2	1:C:139:LYS:H	1.74	0.52
1:C:120:MET:SD	1:C:169:MET:HE3	2.50	0.52
1:A:143:ILE:HG13	1:A:147:MET:HE2	1.90	0.52
1:B:315:GLN:HB3	1:B:334:LYS:NZ	2.25	0.52
1:A:7:LEU:HA	1:A:10:ILE:HG22	1.92	0.52
1:B:177:VAL:HG22	1:B:204:TYR:HB2	1.92	0.51
1:A:243:ARG:NH1	3:A:660:HOH:O	2.40	0.51
1:B:374:THR:N	1:B:375:PRO:HD3	2.26	0.51
1:C:243:ARG:NH1	3:C:602:HOH:O	2.43	0.50
1:C:74:GLU:O	1:C:78:GLN:HG3	2.12	0.50
1:A:327:LEU:HD22	1:A:371:ARG:HD2	1.94	0.50
1:C:115:VAL:O	1:C:116:LYS:HB2	2.12	0.50
1:C:30:ASN:H	1:C:377:MET:HE3	1.77	0.49
1:A:112:LEU:HD11	1:A:230:VAL:HG21	1.95	0.49
1:C:259:GLN:HG2	3:C:734:HOH:O	2.11	0.49
1:B:377:MET:HE3	3:B:641:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ILE:HG12	1:A:147:MET:HE2	1.95	0.49
1:B:326:HIS:H	1:B:326:HIS:HD1	1.61	0.49
1:C:400:ILE:HG21	1:C:425:LEU:HD21	1.94	0.49
1:A:209:ILE:HG21	1:A:224:PRO:HG3	1.94	0.48
1:B:115:VAL:O	1:B:116:LYS:HB2	2.14	0.48
1:A:195:ARG:HD2	1:A:227:HIS:O	2.14	0.48
1:A:177:VAL:HG22	1:A:204:TYR:HB2	1.96	0.48
1:B:109:LEU:O	1:B:113:VAL:HG22	2.14	0.47
1:C:177:VAL:HG22	1:C:204:TYR:HB2	1.96	0.47
1:C:102:SER:HA	1:C:131:THR:HG21	1.95	0.47
1:C:95:ASN:OD1	1:C:98:PRO:HD3	2.14	0.47
1:B:127:GLY:O	1:B:182:TYR:HB3	2.15	0.47
1:A:109:LEU:O	1:A:113:VAL:HG22	2.15	0.47
1:A:211:HIS:HD2	1:A:326:HIS:NE2	2.13	0.47
1:C:3:ASN:OD1	1:C:5:GLU:HG2	2.14	0.47
1:C:219:ASN:HB2	3:C:733:HOH:O	2.14	0.47
1:C:374:THR:H	1:C:375:PRO:HD3	1.79	0.47
1:C:347:ASN:ND2	3:C:603:HOH:O	2.48	0.46
1:B:195:ARG:HB2	1:B:227:HIS:HB3	1.96	0.46
1:A:127:GLY:O	1:A:182:TYR:HB3	2.14	0.46
1:A:257:ILE:HG12	1:A:261:ILE:HD13	1.98	0.46
1:A:115:VAL:HG12	1:A:116:LYS:HG3	1.98	0.46
1:B:119:ILE:HG22	1:B:177:VAL:HB	1.98	0.46
1:A:118:LYS:NZ	1:A:149:GLU:OE1	2.48	0.45
1:B:113:VAL:HG12	1:B:176:LYS:HB3	1.98	0.45
1:C:188:ASP:OD1	1:C:189:ILE:N	2.42	0.45
1:A:274:HIS:HD2	1:B:277:LYS:NZ	2.14	0.45
1:A:410:LYS:HD3	3:A:733:HOH:O	2.17	0.44
1:A:209:ILE:HG13	1:A:233:THR:HB	1.99	0.44
1:B:264:ALA:O	1:B:268:SER:HB2	2.17	0.44
1:A:353:LEU:HD12	1:A:353:LEU:C	2.38	0.44
1:A:24:ARG:NH2	1:B:53:LYS:HD3	2.33	0.44
1:C:327:LEU:C	1:C:327:LEU:HD12	2.39	0.43
1:C:331:ASP:OD1	1:C:331:ASP:C	2.57	0.43
1:C:119:ILE:HG22	1:C:177:VAL:HB	2.00	0.43
1:A:97:GLN:N	1:A:98:PRO:CD	2.82	0.43
1:A:274:HIS:HD2	1:B:277:LYS:HZ1	1.65	0.43
1:C:257:ILE:HG12	1:C:261:ILE:HD13	2.01	0.43
1:C:109:LEU:HD22	1:C:119:ILE:HG21	2.01	0.43
1:C:43:ARG:HD3	3:C:607:HOH:O	2.19	0.43
1:A:161:VAL:CG2	3:A:709:HOH:O	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:GLN:N	1:B:98:PRO:CD	2.82	0.42
1:A:264:ALA:O	1:A:268:SER:HB2	2.20	0.42
1:A:106:VAL:HG13	1:A:147:MET:HE1	2.02	0.42
1:A:109:LEU:HD21	1:A:179:ILE:HD11	2.02	0.41
1:C:55:SER:OG	1:C:70:ILE:HG21	2.21	0.41
1:B:185:TYR:CD2	1:B:187:ARG:O	2.72	0.41
1:C:226:LEU:HB2	1:C:227:HIS:CD2	2.56	0.41
1:C:97:GLN:N	1:C:98:PRO:CD	2.84	0.41
1:C:143:ILE:HG13	1:C:147:MET:HB2	2.01	0.41
1:A:239:LEU:HA	1:A:239:LEU:HD12	1.92	0.41
1:C:234:THR:HB	1:C:236:HIS:CE1	2.55	0.41
1:C:209:ILE:HG21	1:C:224:PRO:HG3	2.01	0.41
1:B:95:ASN:OD1	1:B:98:PRO:HD3	2.21	0.41
1:C:416:LYS:O	1:C:420:PRO:HD3	2.21	0.41
1:C:102:SER:O	1:C:106:VAL:HG23	2.21	0.40
1:B:209:ILE:HD12	1:B:213:SER:HA	2.03	0.40
1:B:109:LEU:HD22	1:B:119:ILE:HG21	2.03	0.40
1:A:277:LYS:NZ	1:B:274:HIS:CD2	2.87	0.40
1:A:116:LYS:HB2	1:A:116:LYS:HE3	1.93	0.40
1:C:139:LYS:HB3	1:C:139:LYS:HE3	1.89	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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%)	52	64
1	B	441/442 (100%)	423 (96%)	16 (4%)	2 (0%)	34	41
1	C	441/442 (100%)	421 (96%)	19 (4%)	1 (0%)	52	64
All	All	1323/1326 (100%)	1271 (96%)	48 (4%)	4 (0%)	46	57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	89	ASP
1	B	89	ASP
1	A	89	ASP
1	B	374	THR

### 5.3.2 Protein sidechains ⓘ

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	382/381 (100%)	370 (97%)	12 (3%)	47	64
1	B	382/381 (100%)	364 (95%)	18 (5%)	32	43
1	C	382/381 (100%)	365 (96%)	17 (4%)	34	46
All	All	1146/1143 (100%)	1099 (96%)	47 (4%)	37	50

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	TYR
1	A	86	ASN
1	A	174	LYS
1	A	176	LYS
1	A	193	GLN
1	A	206	PHE
1	A	211	HIS
1	A	263	SER
1	A	274	HIS
1	A	327	LEU
1	A	364	CYS
1	A	413	VAL
1	B	54	TYR
1	B	75	GLU
1	B	86	ASN
1	B	90	GLU
1	B	100	SER

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Mol	Chain	Res	Type
1	B	136	ASP
1	B	157	SER
1	B	158	GLN
1	B	184	SER
1	B	211	HIS
1	B	219	ASN
1	B	254	ASN
1	B	274	HIS
1	B	327	LEU
1	B	352	SER
1	B	357	THR
1	B	364	CYS
1	B	430	GLN
1	C	54	TYR
1	C	90	GLU
1	C	100	SER
1	C	138	LYS
1	C	139	LYS
1	C	157	SER
1	C	193	GLN
1	C	211	HIS
1	C	252	LYS
1	C	254	ASN
1	C	263	SER
1	C	274	HIS
1	C	327	LEU
1	C	352	SER
1	C	353	LEU
1	C	356	ASN
1	C	357	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	GLN
1	A	211	HIS
1	A	274	HIS
1	B	107	GLN
1	B	193	GLN
1	B	211	HIS
1	B	227	HIS
1	B	274	HIS

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Mol	Chain	Res	Type
1	C	86	ASN
1	C	158	GLN
1	C	211	HIS
1	C	227	HIS
1	C	426	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 carbohydrates 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	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	PLP	A	501	1	15,15,16	1.05	1 (6%)	21,22,23	1.77	4 (19%)
2	PLP	B	501	1	15,15,16	0.89	1 (6%)	21,22,23	1.78	4 (19%)
2	PLP	C	501	1	15,15,16	0.88	1 (6%)	21,22,23	2.18	5 (23%)

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	-	0/6/6/8	0/1/1/1
2	PLP	B	501	1	-	0/6/6/8	0/1/1/1
2	PLP	C	501	1	-	0/6/6/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	PLP	C3-C2	-3.31	1.38	1.40
2	B	501	PLP	C3-C2	-2.34	1.39	1.40
2	C	501	PLP	C3-C2	-2.32	1.39	1.40

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	PLP	C5A-C5-C6	-3.46	112.74	119.28
2	B	501	PLP	C5A-C5-C6	-2.24	115.03	119.28
2	A	501	PLP	C5A-C5-C6	-2.18	115.17	119.28
2	C	501	PLP	O3P-P-O4P	-2.02	100.74	106.56
2	B	501	PLP	C5A-C5-C4	2.81	125.37	121.65
2	A	501	PLP	C5A-C5-C4	3.01	125.64	121.65
2	A	501	PLP	C4A-C4-C5	3.86	124.90	120.88
2	A	501	PLP	O4P-C5A-C5	3.94	115.51	108.99
2	B	501	PLP	C4A-C4-C5	3.96	125.01	120.88
2	B	501	PLP	O4P-C5A-C5	4.11	115.78	108.99
2	C	501	PLP	C4A-C4-C5	4.32	125.39	120.88
2	C	501	PLP	O4P-C5A-C5	4.52	116.47	108.99
2	C	501	PLP	C5A-C5-C4	4.75	127.94	121.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	PLP	1	0
2	C	501	PLP	1	0

## 5.7 Other polymers ⓘ

There are no such residues in this entry.



## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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<sup>th</sup> 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>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	442/442 (100%)	0.26	24 (5%) 29 38	22, 33, 51, 81	0
1	B	442/442 (100%)	0.28	21 (4%) 34 43	22, 34, 51, 72	0
1	C	442/442 (100%)	0.25	18 (4%) 41 50	21, 33, 51, 77	0
All	All	1326/1326 (100%)	0.26	63 (4%) 34 43	21, 33, 51, 81	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	363	ASP	5.7
1	B	362	VAL	5.3
1	A	137	GLU	5.1
1	B	364	CYS	5.0
1	A	364	CYS	4.6
1	B	363	ASP	4.6
1	C	137	GLU	4.5
1	B	137	GLU	4.2
1	B	421	GLY	4.2
1	A	138	LYS	4.1
1	C	363	ASP	3.8
1	A	136	ASP	3.6
1	C	138	LYS	3.4
1	B	365	VAL	3.3
1	A	245	ALA	3.2
1	A	233	THR	2.8
1	C	364	CYS	2.8
1	A	232	THR	2.7
1	A	104	ALA	2.7
1	A	362	VAL	2.7
1	B	138	LYS	2.7
1	C	255	PRO	2.7
1	C	104	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	209	ILE	2.6
1	A	139	LYS	2.6
1	B	33	ALA	2.6
1	B	104	ALA	2.6
1	B	209	ILE	2.6
1	B	234	THR	2.6
1	A	414	ASP	2.6
1	C	136	ASP	2.4
1	C	245	ALA	2.4
1	B	34	SER	2.4
1	C	209	ILE	2.4
1	B	139	LYS	2.4
1	C	419	LEU	2.4
1	B	141	VAL	2.4
1	B	69	PHE	2.3
1	B	232	THR	2.3
1	A	421	GLY	2.3
1	C	234	THR	2.3
1	A	430	GLN	2.3
1	C	33	ALA	2.3
1	C	101	GLY	2.3
1	C	421	GLY	2.3
1	A	247	ILE	2.2
1	C	139	LYS	2.2
1	B	68	ASP	2.2
1	A	192	GLN	2.2
1	A	246	LEU	2.2
1	A	244	SER	2.2
1	A	234	THR	2.2
1	B	244	SER	2.2
1	A	34	SER	2.2
1	B	108	ALA	2.1
1	A	372	ILE	2.1
1	A	100	SER	2.1
1	C	372	ILE	2.1
1	A	423	ALA	2.1
1	C	291	PRO	2.1
1	B	245	ALA	2.1
1	C	259	GLN	2.0
1	B	212	ILE	2.0

## 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	PLP	A	501	15/16	0.98	0.22	0.74	28,33,35,37	0
2	PLP	B	501	15/16	0.98	0.19	0.13	27,31,34,35	0
2	PLP	C	501	15/16	0.98	0.19	0.07	27,30,32,33	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.