



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:24 PM UTC

PDB ID : 4PFN / pdb_00004pfn
Title : Crystal structure of Plasmodium vivax SHMT with L-serine Schiff base
Authors : Chitnumsub, P.; Jaruwat, A.; Leartsakulpanich, U.
Deposited on : 2014-04-30
Resolution : 2.50 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

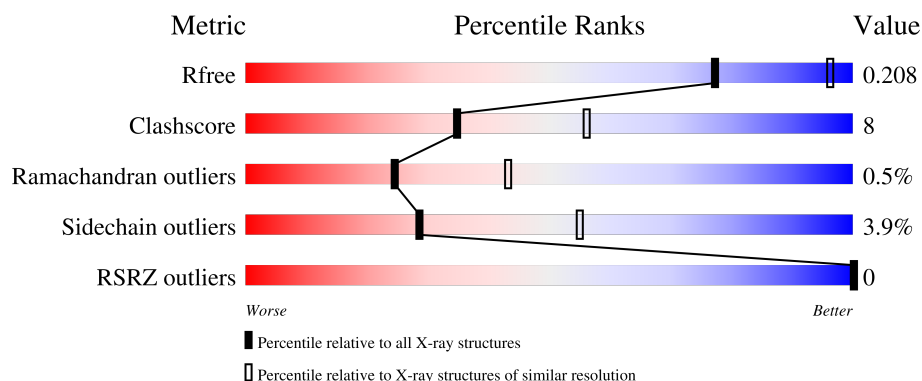
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	 78% 20% .
1	B	442	 81% 19% .
1	C	442	 79% 19% .

2 Entry composition [i](#)

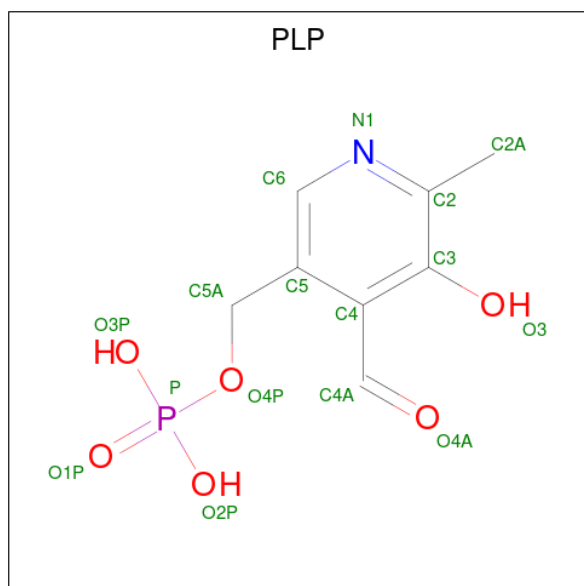
There are 5 unique types of molecules in this entry. The entry contains 10823 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

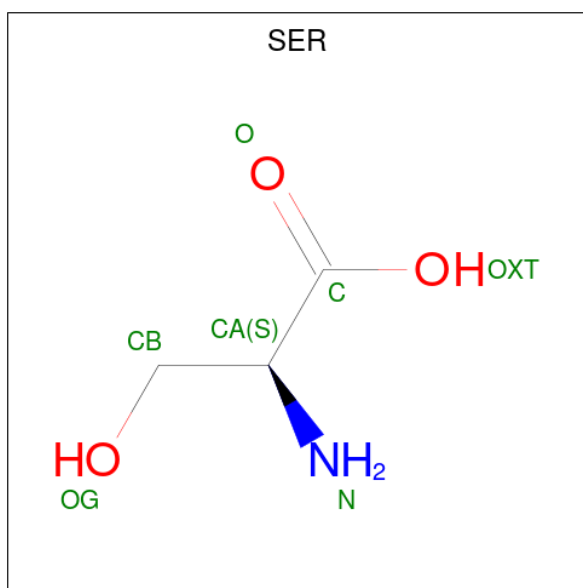
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	0	0
			3458	2186	600	655	17			
1	B	442	Total	C	N	O	S	0	0	0
			3458	2186	600	655	17			
1	C	442	Total	C	N	O	S	0	0	0
			3458	2186	600	655	17			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: 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 SERINE (CCD ID: SER) (formula: $C_3H_7NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			7	3	1	3		
3	B	1	Total	C	N	O	0	0
			7	3	1	3		
3	C	1	Total	C	N	O	0	0
			7	3	1	3		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	122	Total	O	0	0
			122	122		
5	B	141	Total	O	0	0
			141	141		
5	C	118	Total	O	0	0
			118	118		

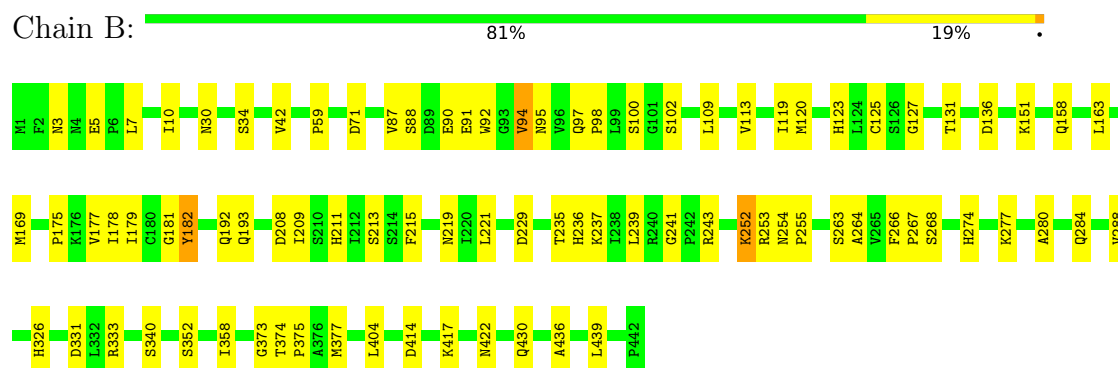
3 Residue-property plots

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.

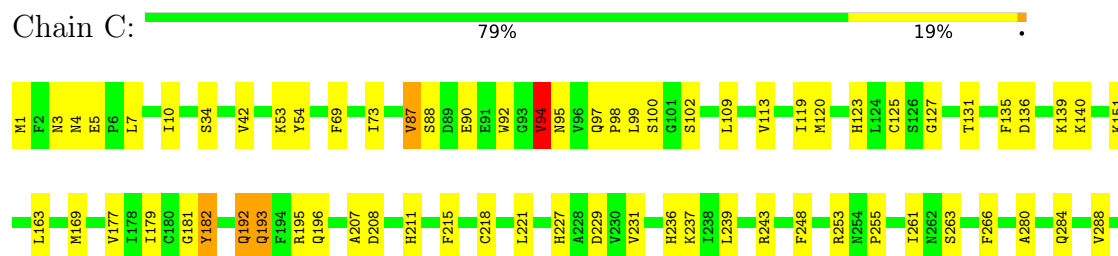
- Molecule 1: Serine hydroxymethyltransferase, putative

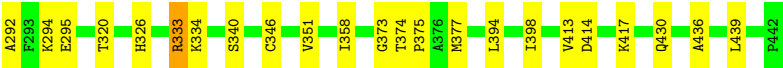


- Molecule 1: Serine hydroxymethyltransferase, putative



- 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, α , β , γ	100.40Å 57.94Å 235.92Å 90.00° 90.01° 90.00°	Depositor
Resolution (Å)	30.00 – 2.50 30.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	93.9 (30.00-2.50) 94.0 (30.00-2.50)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.83 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.207 , 0.263 0.207 , 0.208	Depositor DCC
R_{free} test set	4498 reflections (9.47%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	1.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 10.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.025 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.026 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.477 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.477 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.022 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10823	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: CL, 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	0.62	0/3521	0.87	5/4754 (0.1%)
1	B	0.62	0/3521	0.86	4/4754 (0.1%)
1	C	0.62	0/3521	0.85	2/4754 (0.0%)
All	All	0.62	0/10563	0.86	11/14262 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	254	ASN	CA-C-N	6.03	127.38	119.84
1	B	254	ASN	C-N-CA	6.03	127.38	119.84
1	A	254	ASN	CA-C-N	5.92	127.23	119.84
1	A	254	ASN	C-N-CA	5.92	127.23	119.84
1	C	182	TYR	N-CA-C	5.53	117.38	109.14
1	C	94	VAL	CB-CA-C	-5.40	102.51	110.82
1	A	182	TYR	N-CA-C	5.37	117.14	109.14
1	A	366	SER	CA-C-N	5.36	125.30	119.78
1	A	366	SER	C-N-CA	5.36	125.30	119.78
1	B	182	TYR	N-CA-C	5.33	117.08	109.14
1	B	94	VAL	CB-CA-C	-5.07	103.02	110.82

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	3458	0	3477	59	0
1	B	3458	0	3477	53	0
1	C	3458	0	3477	50	0
2	A	15	0	6	0	0
2	B	15	0	6	0	0
2	C	15	0	6	0	0
3	A	7	0	4	0	0
3	B	7	0	4	0	0
3	C	7	0	4	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	122	0	0	1	0
5	B	141	0	0	0	0
5	C	118	0	0	1	0
All	All	10823	0	10461	157	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 8.

All (157) 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:C:120:MET:HE3	1:C:169:MET:HE2	1.53	0.91
1:C:377:MET:HE2	1:C:377:MET:HA	1.58	0.83
1:A:209:ILE:HD12	1:A:213:SER:HA	1.73	0.70
1:A:119:ILE:HG22	1:A:177:VAL:CG1	2.22	0.69
1:C:215:PHE:HB3	1:C:221:LEU:HG	1.77	0.67
1:A:120:MET:HE3	1:A:169:MET:HE2	1.77	0.67
1:B:123:HIS:HD2	1:B:125:CYS:H	1.43	0.67
1:B:333:ARG:HD2	1:B:358:ILE:CD1	2.26	0.66
1:B:252:LYS:O	1:B:252:LYS:HG3	1.95	0.64
1:A:151:LYS:HB3	1:A:169:MET:HE3	1.80	0.63
1:C:436:ALA:HA	1:C:439:LEU:HD12	1.81	0.62
1:A:3:ASN:OD1	1:A:5:GLU:HG2	1.99	0.62
1:A:123:HIS:HD2	1:A:125:CYS:H	1.48	0.62
1:B:163:LEU:HD13	1:B:193:GLN:HG2	1.81	0.62
1:A:333:ARG:HD2	1:A:358:ILE:CD1	2.30	0.61
1:A:377:MET:HA	1:A:377:MET:HE2	1.83	0.61
1:B:91:GLU:HG3	1:B:252:LYS:HB3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:333:ARG:HD2	1:C:358:ILE:HD11	1.84	0.60
1:C:119:ILE:HG22	1:C:177:VAL:HG12	1.85	0.59
1:C:374:THR:N	1:C:375:PRO:CD	2.66	0.58
1:A:374:THR:N	1:A:375:PRO:CD	2.67	0.58
1:C:109:LEU:O	1:C:113:VAL:HG22	2.04	0.58
1:C:123:HIS:HD2	1:C:125:CYS:H	1.52	0.58
1:C:229:ASP:OD1	1:C:253:ARG:HD2	2.03	0.57
1:A:215:PHE:HB3	1:A:221:LEU:HG	1.87	0.57
1:B:109:LEU:HD21	1:B:179:ILE:HD11	1.86	0.57
1:C:94:VAL:HG13	1:C:248:PHE:CD1	2.40	0.57
1:C:333:ARG:HD2	1:C:358:ILE:CD1	2.35	0.57
1:A:95:ASN:OD1	1:A:98:PRO:HD3	2.05	0.56
1:C:109:LEU:HD21	1:C:179:ILE:HD11	1.87	0.56
1:C:119:ILE:HG22	1:C:177:VAL:CG1	2.36	0.56
1:B:326:HIS:H	1:B:326:HIS:CD2	2.23	0.55
1:C:135:PHE:HD2	1:C:136:ASP:O	1.90	0.55
1:A:109:LEU:HD21	1:A:179:ILE:HD11	1.88	0.55
1:B:414:ASP:HA	1:B:417:LYS:HD2	1.87	0.55
1:C:374:THR:H	1:C:375:PRO:HD3	1.72	0.55
1:A:436:ALA:HA	1:A:439:LEU:HD12	1.89	0.55
1:B:119:ILE:HG22	1:B:177:VAL:CG1	2.37	0.54
1:C:218:CYS:HB3	1:C:294:LYS:HG3	1.89	0.54
1:A:326:HIS:CD2	1:A:326:HIS:H	2.24	0.54
1:B:7:LEU:HA	1:B:10:ILE:HG22	1.90	0.54
1:B:211:HIS:ND1	1:B:326:HIS:HE1	2.06	0.53
1:A:91:GLU:HG3	1:A:252:LYS:HB3	1.91	0.53
1:A:274:HIS:HD2	1:B:277:LYS:HZ1	1.54	0.53
1:B:87:VAL:HG22	1:B:92:TRP:CD1	2.43	0.53
1:A:218:CYS:HB3	1:A:294:LYS:HG3	1.90	0.52
1:B:120:MET:HE2	1:B:178:ILE:HG12	1.91	0.52
1:B:263:SER:HA	1:B:266:PHE:O	2.09	0.52
1:C:374:THR:N	1:C:375:PRO:HD3	2.24	0.52
1:C:326:HIS:H	1:C:326:HIS:CD2	2.27	0.52
1:C:97:GLN:N	1:C:98:PRO:CD	2.73	0.51
1:B:374:THR:N	1:B:375:PRO:CD	2.74	0.51
1:C:99:LEU:HD11	5:C:633:HOH:O	2.10	0.51
1:B:97:GLN:N	1:B:98:PRO:CD	2.74	0.51
1:C:236:HIS:CD2	1:C:243:ARG:HA	2.45	0.50
1:B:404:LEU:HD21	1:B:422:ASN:ND2	2.26	0.50
1:A:181:GLY:HA3	1:A:208:ASP:O	2.12	0.49
1:A:211:HIS:ND1	1:A:326:HIS:HE1	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:377:MET:HE2	1:C:377:MET:CA	2.37	0.49
1:A:236:HIS:CD2	1:A:243:ARG:HA	2.48	0.49
1:A:277:LYS:NZ	1:B:274:HIS:CD2	2.81	0.49
1:C:192:GLN:HG3	1:C:196:GLN:HE21	1.78	0.49
1:C:195:ARG:HB2	1:C:227:HIS:HB3	1.94	0.49
1:A:195:ARG:HB2	1:A:227:HIS:HB3	1.94	0.49
1:B:120:MET:HE2	1:B:175:PRO:HG3	1.95	0.48
1:B:377:MET:HA	1:B:377:MET:HE2	1.95	0.48
1:A:135:PHE:HD2	1:A:136:ASP:O	1.96	0.48
1:B:87:VAL:CG2	1:B:92:TRP:CD1	2.96	0.48
1:B:209:ILE:HD12	1:B:213:SER:HA	1.96	0.48
1:C:263:SER:HA	1:C:266:PHE:O	2.13	0.48
1:A:242:PRO:HG3	1:A:280:ALA:HB3	1.95	0.47
1:C:151:LYS:HB3	1:C:169:MET:HE3	1.95	0.47
1:A:274:HIS:HD2	1:B:277:LYS:NZ	2.12	0.47
1:B:436:ALA:HA	1:B:439:LEU:HD12	1.95	0.47
1:A:151:LYS:HB3	1:A:169:MET:CE	2.44	0.47
1:C:88:SER:C	1:C:90:GLU:H	2.22	0.47
1:A:326:HIS:O	1:A:373:GLY:HA2	2.15	0.47
1:B:236:HIS:CD2	1:B:243:ARG:HA	2.49	0.46
1:C:102:SER:HA	1:C:131:THR:HG21	1.98	0.46
1:A:119:ILE:HG22	1:A:177:VAL:HG12	1.94	0.46
1:A:263:SER:HA	1:A:266:PHE:O	2.15	0.46
1:C:326:HIS:O	1:C:373:GLY:HA2	2.15	0.46
1:A:374:THR:H	1:A:375:PRO:HD3	1.81	0.46
1:C:284:GLN:O	1:C:288:VAL:HG23	2.16	0.46
1:A:284:GLN:O	1:A:288:VAL:HG23	2.16	0.46
1:C:346:CYS:HB3	1:C:351:VAL:HB	1.98	0.45
1:A:122:MET:HE3	1:A:153:TYR:CD2	2.51	0.45
1:A:97:GLN:N	1:A:98:PRO:CD	2.80	0.45
1:B:102:SER:HA	1:B:131:THR:HG21	1.99	0.45
1:A:274:HIS:CD2	1:B:277:LYS:NZ	2.85	0.45
1:B:30:ASN:H	1:B:377:MET:HE3	1.81	0.45
1:A:7:LEU:HA	1:A:10:ILE:HG22	1.99	0.45
1:A:374:THR:N	1:A:375:PRO:HD3	2.30	0.45
1:B:374:THR:N	1:B:375:PRO:HD3	2.31	0.45
1:B:34:SER:HB2	1:B:237:LYS:HE3	1.98	0.45
1:B:59:PRO:HA	1:B:71:ASP:OD1	2.17	0.45
1:B:215:PHE:HB3	1:B:221:LEU:HG	1.99	0.45
1:B:151:LYS:HB3	1:B:169:MET:HE3	1.99	0.44
1:B:252:LYS:O	1:B:252:LYS:CG	2.64	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:PRO:HA	1:A:71:ASP:OD1	2.17	0.44
1:B:229:ASP:OD1	1:B:253:ARG:HD2	2.17	0.44
1:A:207:ALA:HB3	1:A:231:VAL:HG22	1.99	0.44
1:C:181:GLY:HA3	1:C:208:ASP:O	2.18	0.44
1:A:255:PRO:HB2	1:A:256:GLY:H	1.59	0.44
1:C:94:VAL:HG13	1:C:248:PHE:CE1	2.52	0.44
1:C:394:LEU:O	1:C:398:ILE:HG13	2.17	0.44
1:A:358:ILE:CD1	1:A:368:SER:HB2	2.48	0.44
1:A:229:ASP:OD1	1:A:253:ARG:HD2	2.18	0.44
1:B:333:ARG:HD2	1:B:358:ILE:HD12	1.99	0.44
1:A:277:LYS:NZ	1:B:274:HIS:HD2	2.16	0.43
1:C:377:MET:HA	1:C:377:MET:CE	2.35	0.43
1:C:7:LEU:HA	1:C:10:ILE:HG22	1.99	0.43
1:C:120:MET:CE	1:C:169:MET:HE2	2.38	0.43
1:B:264:ALA:O	1:B:268:SER:HB2	2.18	0.43
1:A:333:ARG:HD2	1:A:358:ILE:HD11	2.00	0.43
1:B:374:THR:H	1:B:375:PRO:HD3	1.84	0.43
1:C:87:VAL:HG22	1:C:92:TRP:CD1	2.53	0.43
1:C:42:VAL:HG13	1:C:280:ALA:HB1	2.00	0.43
1:A:113:VAL:HG11	1:A:177:VAL:HG12	2.01	0.43
1:B:95:ASN:OD1	1:B:98:PRO:HD3	2.18	0.42
1:A:109:LEU:O	1:A:113:VAL:HG22	2.19	0.42
1:C:207:ALA:HB3	1:C:231:VAL:HG22	2.00	0.42
1:B:331:ASP:OD1	1:B:331:ASP:C	2.62	0.42
1:A:118:LYS:HE2	1:A:149:GLU:OE1	2.20	0.42
1:C:292:ALA:O	1:C:295:GLU:HB2	2.18	0.42
1:C:95:ASN:OD1	1:C:98:PRO:HD3	2.19	0.42
1:A:251:LYS:C	1:A:253:ARG:H	2.28	0.42
1:B:88:SER:C	1:B:90:GLU:H	2.27	0.42
1:A:87:VAL:CG2	1:A:92:TRP:CD1	3.02	0.42
1:B:127:GLY:O	1:B:182:TYR:HB3	2.19	0.42
1:C:127:GLY:O	1:C:182:TYR:HB3	2.20	0.42
1:B:284:GLN:O	1:B:288:VAL:HG23	2.20	0.42
1:C:3:ASN:OD1	1:C:5:GLU:HG2	2.19	0.42
1:C:69:PHE:O	1:C:73:ILE:HG13	2.19	0.42
1:A:216:VAL:CG1	1:A:224:PRO:HD2	2.50	0.42
1:A:349:ILE:HA	1:A:429:LYS:HB2	2.01	0.42
1:B:109:LEU:O	1:B:113:VAL:HG22	2.20	0.42
1:C:34:SER:HB3	1:C:211:HIS:CE1	2.55	0.42
1:A:209:ILE:HG21	1:A:224:PRO:HG3	2.02	0.41
1:A:127:GLY:O	1:A:182:TYR:HB3	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:ASN:OD1	1:B:5:GLU:HG2	2.20	0.41
1:B:42:VAL:HG13	1:B:280:ALA:HB1	2.01	0.41
1:A:115:VAL:O	1:A:116:LYS:HB2	2.20	0.41
5:A:678:HOH:O	1:B:267:PRO:HD2	2.19	0.41
1:B:404:LEU:HD21	1:B:422:ASN:HD22	1.84	0.41
1:A:91:GLU:HG3	1:A:252:LYS:CB	2.49	0.41
1:B:235:THR:O	1:B:241:GLY:N	2.50	0.41
1:A:333:ARG:HD2	1:A:358:ILE:HD12	2.01	0.41
1:A:34:SER:HB2	1:A:237:LYS:HE3	2.03	0.41
1:A:366:SER:HA	1:A:367:PRO:HD2	1.91	0.41
1:C:163:LEU:HD13	1:C:193:GLN:HG2	2.02	0.41
1:C:414:ASP:HA	1:C:417:LYS:HD2	2.02	0.41
1:A:176:LYS:HA	1:A:176:LYS:HD3	1.95	0.40
1:B:326:HIS:O	1:B:373:GLY:HA2	2.22	0.40
1:B:181:GLY:HA3	1:B:208:ASP:O	2.21	0.40
1:C:34:SER:HB2	1:C:237:LYS:HE3	2.03	0.40
1:A:239:LEU:HD22	1:A:288:VAL:HG21	2.04	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	440/442 (100%)	416 (94%)	22 (5%)	2 (0%)	24	43
1	B	440/442 (100%)	419 (95%)	19 (4%)	2 (0%)	24	43
1	C	440/442 (100%)	420 (96%)	18 (4%)	2 (0%)	24	43
All	All	1320/1326 (100%)	1255 (95%)	59 (4%)	6 (0%)	24	43

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	255	PRO
1	B	255	PRO
1	A	334	LYS
1	C	334	LYS
1	B	252	LYS
1	C	255	PRO

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	381/381 (100%)	364 (96%)	17 (4%)	24	49
1	B	381/381 (100%)	371 (97%)	10 (3%)	40	68
1	C	381/381 (100%)	363 (95%)	18 (5%)	23	47
All	All	1143/1143 (100%)	1098 (96%)	45 (4%)	28	55

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ASN
1	A	54	TYR
1	A	87	VAL
1	A	94	VAL
1	A	124	LEU
1	A	136	ASP
1	A	137	GLU
1	A	158	GLN
1	A	192	GLN
1	A	193	GLN
1	A	239	LEU
1	A	261	ILE
1	A	333	ARG
1	A	340	SER
1	A	352	SER
1	A	364	CYS
1	A	430	GLN

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Mol	Chain	Res	Type
1	B	94	VAL
1	B	100	SER
1	B	136	ASP
1	B	158	GLN
1	B	192	GLN
1	B	219	ASN
1	B	239	LEU
1	B	340	SER
1	B	352	SER
1	B	430	GLN
1	C	1	MET
1	C	4	ASN
1	C	53	LYS
1	C	54	TYR
1	C	87	VAL
1	C	94	VAL
1	C	100	SER
1	C	139	LYS
1	C	140	LYS
1	C	192	GLN
1	C	193	GLN
1	C	239	LEU
1	C	261	ILE
1	C	320	THR
1	C	333	ARG
1	C	340	SER
1	C	413	VAL
1	C	430	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	ASN
1	A	9	GLN
1	A	52	ASN
1	A	123	HIS
1	A	236	HIS
1	A	274	HIS
1	A	289	HIS
1	A	326	HIS
1	A	356	ASN
1	B	4	ASN

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Mol	Chain	Res	Type
1	B	9	GLN
1	B	52	ASN
1	B	123	HIS
1	B	158	GLN
1	B	196	GLN
1	B	236	HIS
1	B	250	ASN
1	B	274	HIS
1	B	315	GLN
1	B	326	HIS
1	B	356	ASN
1	C	4	ASN
1	C	52	ASN
1	C	123	HIS
1	C	158	GLN
1	C	196	GLN
1	C	236	HIS
1	C	259	GLN
1	C	326	HIS
1	C	356	ASN
1	C	424	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 for Mogul analysis.

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
3	SER	B	502	2	4,6,6	0.92	0	2,7,7	0.77	0
3	SER	C	502	2	4,6,6	1.28	1 (25%)	2,7,7	0.81	0
3	SER	A	502	2	4,6,6	1.25	0	2,7,7	2.16	1 (50%)
2	PLP	C	501	3	15,15,16	3.11	3 (20%)	21,22,23	1.78	5 (23%)
2	PLP	B	501	3	15,15,16	3.11	3 (20%)	21,22,23	1.61	3 (14%)
2	PLP	A	501	3	15,15,16	3.10	3 (20%)	21,22,23	1.81	6 (28%)

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
3	SER	B	502	2	-	0/6/6/6	-
3	SER	C	502	2	-	0/6/6/6	-
3	SER	A	502	2	-	2/6/6/6	-
2	PLP	C	501	3	-	3/6/6/8	0/1/1/1
2	PLP	B	501	3	-	3/6/6/8	0/1/1/1
2	PLP	A	501	3	-	1/6/6/8	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	PLP	C5-C4	8.76	1.50	1.40
2	B	501	PLP	C5-C4	8.39	1.50	1.40
2	A	501	PLP	C5-C4	8.17	1.49	1.40
2	A	501	PLP	C3-C2	7.21	1.48	1.41
2	B	501	PLP	C3-C2	6.82	1.48	1.41
2	C	501	PLP	C3-C2	6.51	1.47	1.41
2	B	501	PLP	C3-C4	4.70	1.49	1.40
2	C	501	PLP	C3-C4	4.48	1.48	1.40
2	A	501	PLP	C3-C4	4.39	1.48	1.40
3	C	502	SER	OXT-C	-2.14	1.23	1.30

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	PLP	C4A-C4-C5	4.41	125.48	120.94
2	A	501	PLP	C4A-C4-C5	3.63	124.68	120.94
2	B	501	PLP	C4A-C4-C5	3.40	124.44	120.94
2	A	501	PLP	O3P-P-O4P	-3.39	97.83	106.67
2	B	501	PLP	C6-N1-C2	3.07	124.76	119.20
3	A	502	SER	OXT-C-O	-3.05	117.17	124.08
2	A	501	PLP	C5A-C5-C6	-3.02	114.44	119.36
2	C	501	PLP	C6-N1-C2	2.84	124.34	119.20
2	B	501	PLP	O4P-C5A-C5	2.84	114.67	109.36
2	C	501	PLP	O4P-C5A-C5	2.72	114.46	109.36
2	A	501	PLP	C3-C4-C5	-2.59	115.48	118.59
2	A	501	PLP	C6-C5-C4	2.36	120.03	118.10
2	C	501	PLP	O2P-P-O4P	-2.35	100.53	106.67
2	A	501	PLP	C6-N1-C2	2.22	123.23	119.20
2	C	501	PLP	O3P-P-O2P	2.09	115.63	107.80

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	PLP	C5A-O4P-P-O2P
2	B	501	PLP	C5A-O4P-P-O3P
2	C	501	PLP	C5A-O4P-P-O2P
2	C	501	PLP	C5A-O4P-P-O3P
2	B	501	PLP	C5A-O4P-P-O1P
2	A	501	PLP	C5A-O4P-P-O3P
2	C	501	PLP	C5A-O4P-P-O1P
3	A	502	SER	OXT-C-CA-N
3	A	502	SER	C-CA-CB-OG

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, 95th 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 ²)	Q<0.9
1	A	442/442 (100%)	-1.49	0 100 100	14, 27, 49, 71	0
1	B	442/442 (100%)	-1.48	0 100 100	13, 27, 50, 68	0
1	C	442/442 (100%)	-1.48	0 100 100	14, 27, 50, 66	0
All	All	1326/1326 (100%)	-1.49	0 100 100	13, 27, 50, 71	0

There are no RSRZ outliers to report.

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 oligosaccharides 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, 95th 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	B-factors(Å ²)	Q<0.9
3	SER	A	502	7/7	0.99	0.02	13,13,14,14	0
3	SER	B	502	7/7	0.99	0.03	17,17,18,19	0
3	SER	C	502	7/7	0.99	0.03	16,16,16,17	0
4	CL	B	503	1/1	0.99	0.02	23,23,23,23	0
4	CL	C	503	1/1	0.99	0.04	28,28,28,28	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PLP	C	501	15/16	1.00	0.03	16,17,17,17	0
2	PLP	A	501	15/16	1.00	0.03	12,13,13,13	0
2	PLP	B	501	15/16	1.00	0.02	16,17,17,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.