



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 08:43 am BST

PDB ID : 5CVC
Title : Structure of maize serine racemase
Authors : Song, Y.; Zou, L.; Fan, J.
Deposited on : 2015-07-26
Resolution : 2.09 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
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) : 2.11

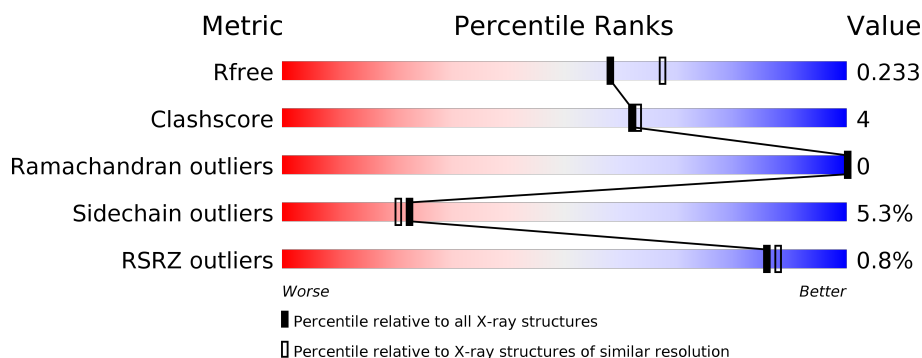
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 $\leq 5\%$. 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	346	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 88% 6% • 5% </div> </div>
1	B	346	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 81% 9% • 8% </div> </div>
1	C	346	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 82% 9% • 8% </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7265 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 racemase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	1	0
			2403	1527	410	457	9			
1	B	318	Total	C	N	O	S	0	1	0
			2299	1467	386	437	9			
1	C	318	Total	C	N	O	S	0	1	0
			2319	1479	393	438	9			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	339	LEU	-	expression tag	UNP F5CAQ9
A	340	GLU	-	expression tag	UNP F5CAQ9
A	341	HIS	-	expression tag	UNP F5CAQ9
A	342	HIS	-	expression tag	UNP F5CAQ9
A	343	HIS	-	expression tag	UNP F5CAQ9
A	344	HIS	-	expression tag	UNP F5CAQ9
A	345	HIS	-	expression tag	UNP F5CAQ9
A	346	HIS	-	expression tag	UNP F5CAQ9
B	339	LEU	-	expression tag	UNP F5CAQ9
B	340	GLU	-	expression tag	UNP F5CAQ9
B	341	HIS	-	expression tag	UNP F5CAQ9
B	342	HIS	-	expression tag	UNP F5CAQ9
B	343	HIS	-	expression tag	UNP F5CAQ9
B	344	HIS	-	expression tag	UNP F5CAQ9
B	345	HIS	-	expression tag	UNP F5CAQ9
B	346	HIS	-	expression tag	UNP F5CAQ9
C	339	LEU	-	expression tag	UNP F5CAQ9
C	340	GLU	-	expression tag	UNP F5CAQ9
C	341	HIS	-	expression tag	UNP F5CAQ9
C	342	HIS	-	expression tag	UNP F5CAQ9
C	343	HIS	-	expression tag	UNP F5CAQ9
C	344	HIS	-	expression tag	UNP F5CAQ9
C	345	HIS	-	expression tag	UNP F5CAQ9

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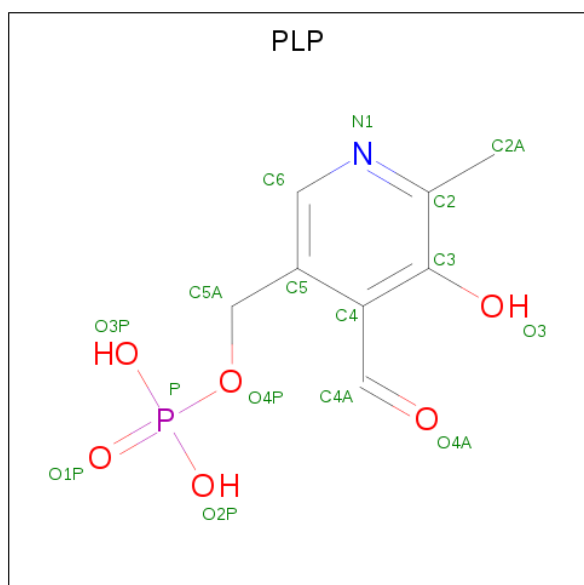
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Chain	Residue	Modelled	Actual	Comment	Reference
C	346	HIS	-	expression tag	UNP F5CAQ9

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



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

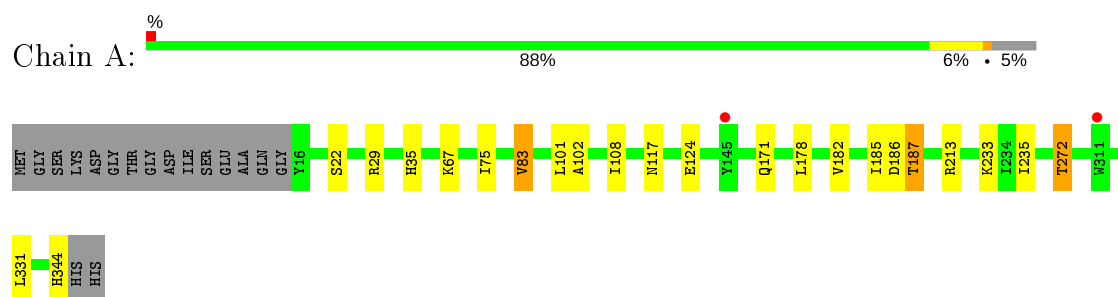
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	70	Total 70	O 70	0	0
4	B	61	Total 61	O 61	0	0
4	C	65	Total 65	O 65	0	0

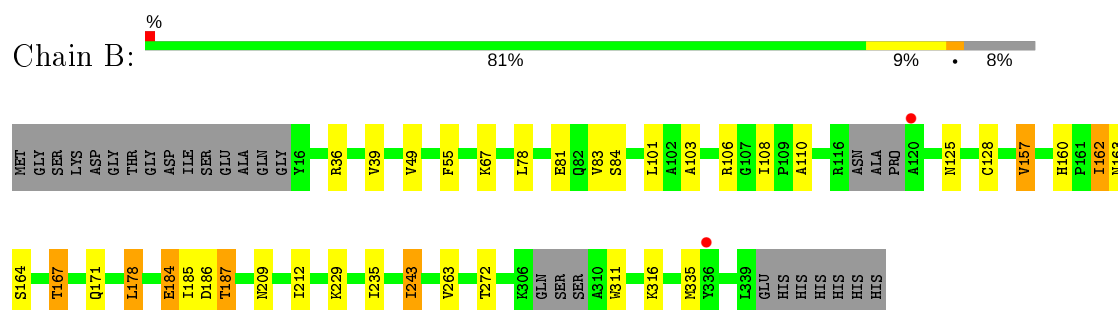
3 Residue-property plots [i](#)

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.

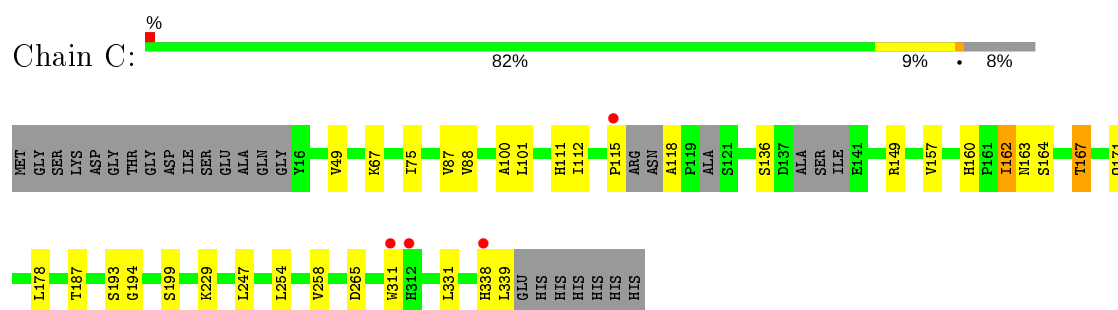
- Molecule 1: Serine racemase



- Molecule 1: Serine racemase



- Molecule 1: Serine racemase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	50.99Å 204.02Å 211.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.38 – 2.09 41.35 – 2.09	Depositor EDS
% Data completeness (in resolution range)	98.6 (41.38-2.09) 98.7 (41.35-2.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.33 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.198 , 0.229 0.206 , 0.233	Depositor DCC
R_{free} test set	3283 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.518	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7265	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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: MG, 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	0/2448	0.65	1/3336 (0.0%)
1	B	0.42	0/2337	0.62	0/3181
1	C	0.41	0/2359	0.64	1/3208 (0.0%)
All	All	0.43	0/7144	0.63	2/9725 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	29	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	C	149	ARG	NE-CZ-NH2	-5.12	117.74	120.30

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	2403	0	2362	12	0
1	B	2299	0	2280	26	0
1	C	2319	0	2302	21	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	15	0	6	3	0
3	B	15	0	6	3	0
3	C	15	0	6	2	0
4	A	70	0	0	2	1
4	B	61	0	0	0	2
4	C	65	0	0	0	1
All	All	7265	0	6962	58	2

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 4.

All (58) 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:67:LYS:HZ3	3:B:402:PLP:C4A	1.53	1.19
1:C:67:LYS:HZ3	3:C:402:PLP:C4A	1.67	1.03
1:A:67:LYS:HZ1	3:A:402:PLP:C4A	1.63	0.97
1:A:67:LYS:HZ3	3:A:402:PLP:C4A	1.66	0.97
1:B:67:LYS:HZ1	3:B:402:PLP:C4A	1.85	0.88
1:C:67:LYS:HZ1	3:C:402:PLP:C4A	1.90	0.83
1:B:83:VAL:HG22	1:B:108:ILE:HD13	1.70	0.73
1:C:163:ASN:HD21	1:C:254:LEU:H	1.35	0.73
1:B:160[B]:HIS:O	1:B:167:THR:HG21	1.89	0.72
1:B:162:ILE:HG12	1:B:163:ASN:N	2.03	0.72
1:B:160[A]:HIS:O	1:B:167:THR:HG21	1.93	0.69
1:B:164:SER:OG	1:B:167:THR:HG23	1.95	0.67
1:A:235:ILE:HG23	1:B:235:ILE:HG23	1.76	0.66
1:B:83:VAL:CG2	1:B:108:ILE:HD13	2.24	0.66
1:C:164:SER:OG	1:C:167:THR:HG23	1.97	0.64
1:A:83:VAL:HG22	1:A:108:ILE:HD13	1.79	0.64
1:C:160[B]:HIS:O	1:C:167:THR:HG21	2.01	0.60
1:C:49:VAL:O	1:C:311:TRP:CH2	2.55	0.59
1:C:162:ILE:HG12	1:C:163:ASN:N	2.17	0.59
1:C:160[A]:HIS:O	1:C:167:THR:HG21	2.02	0.57
1:A:83:VAL:CG2	1:A:108:ILE:HD13	2.36	0.55
1:C:115:PRO:HB3	1:C:118:ALA:HB2	1.89	0.53
1:B:160[B]:HIS:CD2	1:B:163:ASN:H	2.26	0.53
1:C:100:ALA:HB2	1:C:112:ILE:CD1	2.39	0.52
1:C:199:SER:HA	1:C:258:VAL:HG11	1.91	0.52
1:B:103:ALA:CB	1:B:110:ALA:HB2	2.41	0.51
1:A:272:THR:HG22	4:A:562:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:162:ILE:HD12	1:C:194:GLY:HA3	1.92	0.50
1:B:49:VAL:O	1:B:311:TRP:HH2	1.95	0.49
1:B:243:ILE:O	1:B:243:ILE:HD13	2.12	0.49
1:A:35:HIS:HD2	4:A:566:HOH:O	1.95	0.49
1:C:338:HIS:O	1:C:339:LEU:HB2	2.13	0.49
1:B:67:LYS:NZ	3:B:402:PLP:C4	2.75	0.47
1:C:49:VAL:O	1:C:311:TRP:CZ3	2.68	0.47
1:A:187:THR:HB	1:A:213:ARG:HB2	1.97	0.45
1:A:186:ASP:OD1	1:A:187:THR:HG22	2.16	0.45
1:B:160[B]:HIS:HD2	1:B:162:ILE:H	1.64	0.45
1:B:229:LYS:NZ	1:B:263:VAL:O	2.43	0.44
1:B:125:ASN:O	1:B:128:CYS:HB3	2.18	0.43
1:B:162:ILE:HG12	1:B:163:ASN:HB2	2.00	0.43
1:A:182:VAL:HB	1:A:185:ILE:HD12	2.00	0.43
1:C:162:ILE:HG12	1:C:163:ASN:HB2	2.00	0.43
1:B:186:ASP:OD1	1:B:187:THR:CG2	2.68	0.42
1:A:67:LYS:HZ1	3:A:402:PLP:C4	2.23	0.42
1:B:81:GLU:O	1:B:84:SER:OG	2.33	0.42
1:B:185:ILE:O	1:B:212:ILE:HD12	2.20	0.42
1:C:88:VAL:HA	1:C:111:HIS:O	2.19	0.42
1:C:49:VAL:O	1:C:311:TRP:HH2	2.02	0.42
1:A:75:ILE:HD12	1:A:102:ALA:HB3	2.01	0.41
1:C:100:ALA:HB2	1:C:112:ILE:HD11	2.02	0.41
1:C:75:ILE:HG12	1:C:87:VAL:HG11	2.02	0.41
1:B:78:LEU:HD11	1:B:157:VAL:HG21	2.02	0.41
1:B:178:LEU:HD22	1:B:185:ILE:HD12	2.03	0.41
1:B:39:VAL:HA	1:B:55:PHE:O	2.21	0.41
1:C:193:SER:HB2	1:C:247:LEU:HA	2.03	0.41
1:B:184:GLU:HB2	1:B:316:LYS:HD2	2.02	0.40
1:C:338:HIS:O	1:C:339:LEU:CB	2.68	0.40
1:B:186:ASP:OD1	1:B:187:THR:HG22	2.22	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:505:HOH:O	4:C:553:HOH:O[4_655]	1.73	0.47
4:A:526:HOH:O	4:B:504:HOH:O[1_455]	2.14	0.06

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	328/346 (95%)	320 (98%)	8 (2%)	0	100	100
1	B	313/346 (90%)	303 (97%)	10 (3%)	0	100	100
1	C	311/346 (90%)	302 (97%)	9 (3%)	0	100	100
All	All	952/1038 (92%)	925 (97%)	27 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	241/275 (88%)	229 (95%)	12 (5%)	24	23
1	B	231/275 (84%)	217 (94%)	14 (6%)	18	16
1	C	235/275 (86%)	224 (95%)	11 (5%)	26	25
All	All	707/825 (86%)	670 (95%)	37 (5%)	22	21

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	SER
1	A	83	VAL
1	A	101	LEU
1	A	117	ASN
1	A	124	GLU

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Mol	Chain	Res	Type
1	A	171	GLN
1	A	178	LEU
1	A	187	THR
1	A	233	LYS
1	A	272	THR
1	A	331	LEU
1	A	344	HIS
1	B	36	ARG
1	B	101	LEU
1	B	106	ARG
1	B	157	VAL
1	B	162	ILE
1	B	167	THR
1	B	171	GLN
1	B	178	LEU
1	B	184	GLU
1	B	187	THR
1	B	209	ASN
1	B	243	ILE
1	B	272	THR
1	B	335	MET
1	C	101	LEU
1	C	136	SER
1	C	157	VAL
1	C	162	ILE
1	C	167	THR
1	C	171	GLN
1	C	178	LEU
1	C	187	THR
1	C	229	LYS
1	C	265	ASP
1	C	331	LEU

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	35	HIS
1	A	90	HIS
1	A	117	ASN
1	A	132	HIS
1	A	181	GLN
1	A	209	ASN

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Mol	Chain	Res	Type
1	B	181	GLN
1	B	241	ASN
1	C	90	HIS
1	C	132	HIS
1	C	163	ASN
1	C	181	GLN
1	C	227	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](#)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	PLP	C	402	1	15,15,16	2.68	3 (20%)	20,22,23	1.52	4 (20%)
3	PLP	B	402	1	15,15,16	2.75	4 (26%)	20,22,23	1.47	3 (15%)
3	PLP	A	402	1	15,15,16	2.71	3 (20%)	20,22,23	1.41	2 (10%)

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	PLP	C	402	1	-	1/6/6/8	0/1/1/1
3	PLP	B	402	1	-	0/6/6/8	0/1/1/1
3	PLP	A	402	1	-	0/6/6/8	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	PLP	C3-C2	7.80	1.48	1.40
3	B	402	PLP	C3-C2	7.08	1.48	1.40
3	C	402	PLP	C3-C2	6.88	1.47	1.40
3	C	402	PLP	C5-C4	6.05	1.47	1.40
3	B	402	PLP	C5-C4	5.42	1.46	1.40
3	A	402	PLP	C5-C4	5.17	1.46	1.40
3	B	402	PLP	C3-C4	4.68	1.49	1.40
3	C	402	PLP	C3-C4	4.12	1.48	1.40
3	A	402	PLP	C3-C4	3.98	1.48	1.40
3	B	402	PLP	C4A-C4	-2.31	1.46	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	PLP	O4P-C5A-C5	2.99	115.06	109.35
3	A	402	PLP	O4P-C5A-C5	2.87	114.81	109.35
3	B	402	PLP	C6-N1-C2	2.83	124.41	119.17
3	C	402	PLP	C6-N1-C2	2.59	123.97	119.17
3	A	402	PLP	C6-N1-C2	2.42	123.65	119.17
3	C	402	PLP	O3P-P-O4P	-2.41	100.32	106.73
3	C	402	PLP	O4P-C5A-C5	2.35	113.83	109.35
3	C	402	PLP	C2A-C2-N1	2.09	121.75	117.67
3	B	402	PLP	C4A-C4-C5	-2.07	118.81	120.94

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	402	PLP	C5A-O4P-P-O2P

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	402	PLP	2	0
3	B	402	PLP	3	0
3	A	402	PLP	3	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, 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	329/346 (95%)	-0.34	2 (0%) 89 91	20, 29, 47, 53	0
1	B	318/346 (91%)	-0.24	2 (0%) 89 91	23, 31, 57, 73	0
1	C	318/346 (91%)	-0.23	4 (1%) 77 80	23, 32, 50, 66	0
All	All	965/1038 (92%)	-0.27	8 (0%) 86 88	20, 31, 52, 73	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	311	TRP	3.5
1	C	115	PRO	3.3
1	C	312	HIS	3.2
1	B	120	ALA	2.7
1	A	145	TYR	2.5
1	B	336	TYR	2.4
1	C	338	HIS	2.3
1	A	311	TRP	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. 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
2	MG	A	401	1/1	0.97	0.08	28,28,28,28	0
2	MG	C	401	1/1	0.98	0.05	29,29,29,29	0
3	PLP	B	402	15/16	0.98	0.15	23,25,28,28	0
2	MG	B	401	1/1	0.98	0.04	31,31,31,31	0
3	PLP	A	402	15/16	0.98	0.09	21,23,25,26	0
3	PLP	C	402	15/16	0.99	0.09	23,28,30,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.