



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:28 PM UTC

PDB ID : 7AGZ / pdb_00007agz
Title : BsrV no-histagged
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Deposited on : 2020-09-23
Resolution : 1.52 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

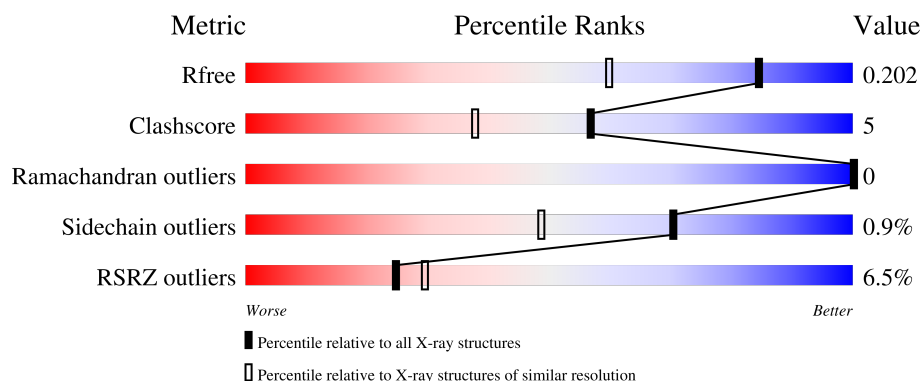
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.52 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	5890 (1.54-1.50)
Clashscore	190562	6116 (1.54-1.50)
Ramachandran outliers	187476	6002 (1.54-1.50)
Sidechain outliers	187428	5999 (1.54-1.50)
RSRZ outliers	180081	5891 (1.54-1.50)

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	390	5% 86% 12% .
1	B	390	8% 88% 10% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6712 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.

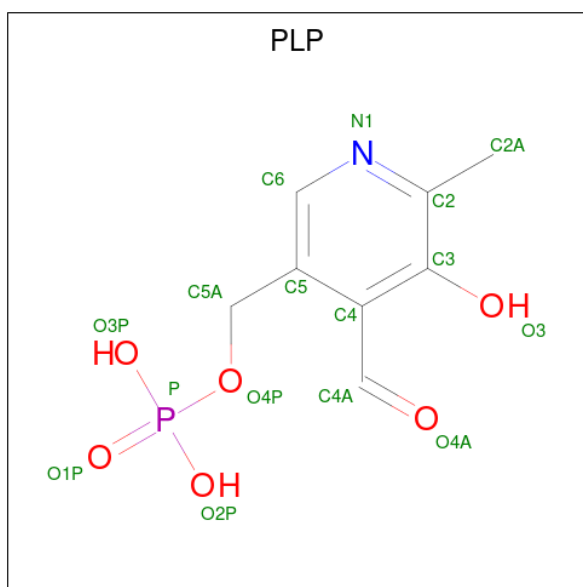
- Molecule 1 is a protein called Broad specificity amino-acid racemase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	386	Total	C	N	O	S	0	4	0
			2975	1862	518	580	15			
1	B	388	Total	C	N	O	S	0	4	0
			2990	1874	517	584	15			

There are 12 discrepancies between the modelled and reference sequences:

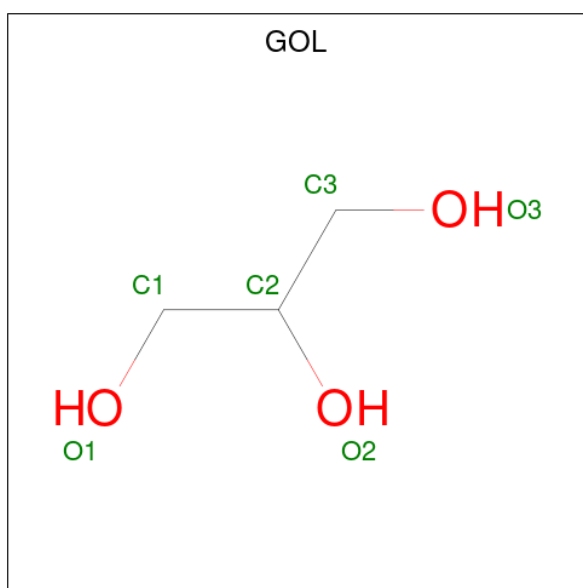
Chain	Residue	Modelled	Actual	Comment	Reference
A	408	GLU	-	expression tag	UNP Q9KSE5
A	409	ASN	-	expression tag	UNP Q9KSE5
A	410	LEU	-	expression tag	UNP Q9KSE5
A	411	TYR	-	expression tag	UNP Q9KSE5
A	412	PHE	-	expression tag	UNP Q9KSE5
A	413	GLN	-	expression tag	UNP Q9KSE5
B	408	GLU	-	expression tag	UNP Q9KSE5
B	409	ASN	-	expression tag	UNP Q9KSE5
B	410	LEU	-	expression tag	UNP Q9KSE5
B	411	TYR	-	expression tag	UNP Q9KSE5
B	412	PHE	-	expression tag	UNP Q9KSE5
B	413	GLN	-	expression tag	UNP Q9KSE5

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: C₈H₁₀NO₆P).



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		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

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

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

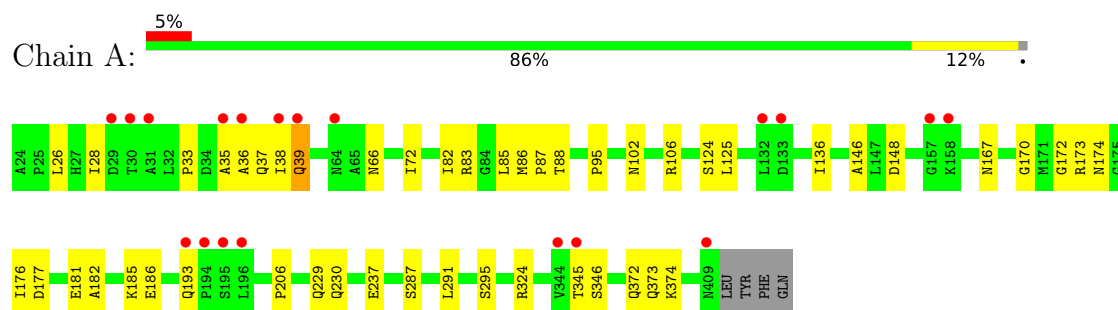
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	339	Total 339	O 339	0	0
5	B	364	Total 364	O 364	0	0

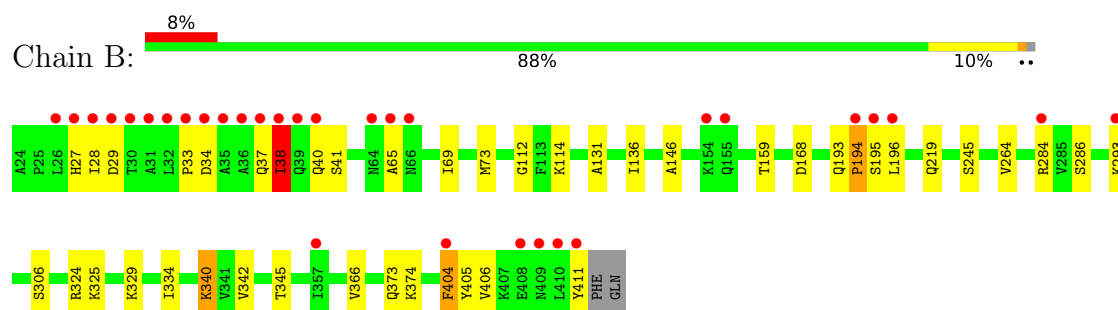
3 Residue-property plots

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: Broad specificity amino-acid racemase



- Molecule 1: Broad specificity amino-acid racemase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.03Å 82.09Å 160.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.20 – 1.52 51.20 – 1.52	Depositor EDS
% Data completeness (in resolution range)	99.8 (51.20-1.52) 95.5 (51.20-1.52)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.54 (at 1.52Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874, PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.185 , 0.200 0.186 , 0.202	Depositor DCC
R_{free} test set	7982 reflections (7.18%)	wwPDB-VP
Wilson B-factor (Å ²)	13.5	Xtriage
Anisotropy	0.756	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6712	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, 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	1.08	26/3025 (0.9%)	0.99	1/4099 (0.0%)
1	B	0.81	9/3041 (0.3%)	1.00	7/4122 (0.2%)
All	All	0.95	35/6066 (0.6%)	1.00	8/8221 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	6
All	All	0	9

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	373	GLN	C-O	-8.20	1.17	1.24
1	A	86	MET	C-O	-8.18	1.17	1.24
1	A	186	GLU	C-O	-7.49	1.15	1.24
1	A	373	GLN	C-O	-7.41	1.17	1.24
1	A	82	ILE	C-O	-7.25	1.16	1.24
1	A	85	LEU	C-O	-7.16	1.14	1.24
1	A	185	LYS	C-O	-7.05	1.15	1.24
1	A	167	ASN	C-O	-6.89	1.15	1.24
1	B	340	LYS	C-O	-6.85	1.15	1.24
1	A	372	GLN	C-O	-6.79	1.15	1.24
1	B	406	VAL	C-O	-6.70	1.16	1.24
1	A	87	PRO	C-O	-6.66	1.15	1.24
1	A	177	ASP	C-O	-6.55	1.16	1.24
1	A	181	GLU	C-O	-6.45	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	237	GLU	C-N	-6.38	1.25	1.33
1	A	173	ARG	C-O	-6.37	1.16	1.24
1	A	88	THR	C-O	-6.19	1.16	1.24
1	A	229	GLN	C-N	-6.17	1.24	1.33
1	A	174	ASN	C-O	-6.10	1.15	1.23
1	A	95	PRO	C-O	-6.07	1.17	1.24
1	B	73[A]	MET	C-N	-6.07	1.25	1.33
1	B	73[B]	MET	C-N	-6.07	1.25	1.33
1	A	72	ILE	C-O	-6.05	1.17	1.24
1	B	69	ILE	C-O	-5.80	1.17	1.24
1	B	405	TYR	C-O	-5.47	1.17	1.24
1	B	286	SER	C-N	-5.44	1.26	1.33
1	A	287	SER	C-O	-5.43	1.17	1.23
1	A	172	GLY	C-O	-5.42	1.18	1.24
1	A	176	ILE	C-O	-5.41	1.18	1.24
1	A	83	ARG	C-O	-5.38	1.17	1.24
1	B	404	PHE	C-O	-5.30	1.17	1.24
1	A	230	GLN	C-N	-5.29	1.26	1.33
1	A	206	PRO	C-O	-5.17	1.17	1.24
1	A	170	GLY	C-O	-5.11	1.16	1.24
1	A	182	ALA	C-O	-5.06	1.18	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	65	ALA	CA-C-N	-9.13	108.15	122.67
1	B	65	ALA	C-N-CA	-9.13	108.15	122.67
1	B	159	THR	N-CA-C	-6.38	98.13	108.52
1	A	39	GLN	CB-CG-CD	-5.57	103.12	112.60
1	B	34	ASP	CA-C-N	5.42	128.95	120.82
1	B	34	ASP	C-N-CA	5.42	128.95	120.82
1	B	194	PRO	CA-C-O	-5.28	115.23	121.67
1	B	38	ILE	CA-C-O	-5.14	115.35	121.05

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	28[A]	ILE	Mainchain
1	A	28[B]	ILE	Mainchain
1	A	346[A]	SER	Mainchain
1	B	194	PRO	Mainchain

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Mol	Chain	Res	Type	Group
1	B	195[A]	SER	Mainchain
1	B	195[B]	SER	Mainchain
1	B	27	HIS	Mainchain

5.2 Too-close contacts

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	2975	0	2971	21	0
1	B	2990	0	2983	42	0
2	A	15	0	6	0	0
2	B	15	0	6	0	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	339	0	0	4	0
5	B	364	0	0	4	0
All	All	6712	0	5982	55	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 5.

All (55) 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:38:ILE:HG13	1:B:404:PHE:CE2	1.64	1.33
1:B:38:ILE:HG13	1:B:404:PHE:CZ	1.81	1.13
1:B:340:LYS:HE3	5:B:812:HOH:O	1.49	1.12
1:B:38:ILE:CG1	1:B:404:PHE:CE2	2.34	1.11
1:B:38:ILE:HG13	1:B:404:PHE:CD2	1.91	1.05
1:A:106[A]:ARG:HH22	1:B:40:GLN:HE22	1.01	0.95
1:B:38:ILE:CG1	1:B:404:PHE:CD2	2.52	0.91
1:A:102:ASN:HD22	1:B:40:GLN:HE21	1.17	0.91
1:A:106[A]:ARG:HH22	1:B:40:GLN:NE2	1.71	0.87
1:B:340:LYS:HE2	1:B:342:VAL:HG12	1.65	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106[A]:ARG:NH2	1:B:40:GLN:HE22	1.79	0.77
1:A:26:LEU:HD21	1:B:37:GLN:HB3	1.65	0.77
1:A:146:ALA:HB1	1:A:193:GLN:HG3	1.66	0.75
1:B:146:ALA:HB1	1:B:193:GLN:HG3	1.70	0.73
1:B:38:ILE:HG13	1:B:404:PHE:CE1	2.24	0.73
1:A:124:SER:HB2	5:B:815:HOH:O	1.88	0.73
1:B:38:ILE:HG12	1:B:404:PHE:CE2	2.24	0.72
1:A:106[B]:ARG:NH1	5:A:601:HOH:O	2.27	0.67
1:A:35:ALA:O	1:A:38:ILE:HG22	1.96	0.65
1:B:38:ILE:CG1	1:B:404:PHE:CZ	2.67	0.65
1:B:38:ILE:HG13	1:B:404:PHE:CG	2.34	0.63
1:B:33:PRO:HB2	1:B:404:PHE:CE1	2.37	0.59
1:B:38:ILE:CD1	1:B:404:PHE:CD2	2.86	0.59
1:B:38:ILE:HD11	1:B:404:PHE:CD2	2.40	0.56
1:B:245:SER:HB3	1:B:264:VAL:HB	1.90	0.54
1:B:293:LYS:HG3	1:B:306:SER:C	2.35	0.52
1:B:168:ASP:OD1	1:B:219:GLN:HG3	2.11	0.51
1:A:102:ASN:HD22	1:B:40:GLN:NE2	1.99	0.51
1:B:38:ILE:HG13	1:B:404:PHE:CD1	2.46	0.50
1:B:40:GLN:O	1:B:40:GLN:HG3	2.11	0.50
1:A:136:ILE:HG22	5:A:708:HOH:O	2.12	0.50
1:A:36:ALA:O	1:A:39:GLN:HG2	2.12	0.49
1:B:334:ILE:HG12	1:B:366:VAL:HG22	1.97	0.47
1:A:106[B]:ARG:HD3	5:A:628:HOH:O	2.16	0.46
1:B:131:ALA:HB1	1:B:136:ILE:HD12	1.96	0.46
1:B:374:LYS:HG2	1:B:374:LYS:O	2.16	0.45
1:B:324:ARG:HA	1:B:345:THR:HB	1.99	0.45
1:A:324:ARG:HA	1:A:345:THR:HB	1.99	0.45
1:B:411:TYR:OH	5:B:601:HOH:O	2.17	0.45
1:A:102:ASN:ND2	1:B:40:GLN:HE21	1.99	0.44
1:A:125:LEU:HD13	1:A:148:ASP:HB3	1.99	0.44
1:A:106[B]:ARG:HH22	1:B:40:GLN:NE2	2.15	0.44
1:B:340:LYS:CE	5:B:812:HOH:O	2.31	0.43
1:B:325:LYS:O	1:B:329:LYS:HE3	2.17	0.43
1:A:66:ASN:ND2	5:A:605:HOH:O	2.36	0.43
1:B:38:ILE:CD1	1:B:404:PHE:CG	3.02	0.42
1:B:33:PRO:CB	1:B:404:PHE:CE1	3.02	0.42
1:B:112:GLY:O	1:B:114:LYS:HD3	2.20	0.41
1:A:291:LEU:HG	1:A:295:SER:CB	2.51	0.41
1:B:196:LEU:HD23	1:B:196:LEU:HA	1.88	0.41
1:B:340:LYS:HE2	1:B:342:VAL:CG1	2.42	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:GLN:O	1:B:40:GLN:CG	2.69	0.41
1:A:33:PRO:HA	1:A:37:GLN:OE1	2.21	0.40
1:A:26:LEU:HD12	1:A:26:LEU:HA	1.89	0.40
1:B:193:GLN:HE21	1:B:193:GLN:HB3	1.76	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	Percentiles	
1	A	388/390 (100%)	378 (97%)	10 (3%)	0	100	100
1	B	390/390 (100%)	380 (97%)	10 (3%)	0	100	100
All	All	778/780 (100%)	758 (97%)	20 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	Percentiles	
1	A	327/327 (100%)	326 (100%)	1 (0%)	86	75
1	B	329/327 (101%)	323 (98%)	6 (2%)	51	23
All	All	656/654 (100%)	649 (99%)	7 (1%)	70	40

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	374	LYS
1	B	28[A]	ILE
1	B	28[B]	ILE
1	B	29	ASP
1	B	38	ILE
1	B	41	SER
1	B	284	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	219	GLN
1	B	40	GLN
1	B	51	GLN
1	B	193	GLN
1	B	348	ASN
1	B	371	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 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z > 2$
2	PLP	A	501	-	15,15,16	0.94	0	21,22,23	1.61	1 (4%)
3	GOL	B	502	-	5,5,5	0.35	0	5,5,5	0.58	0
3	GOL	A	502	-	5,5,5	0.36	0	5,5,5	0.25	0
2	PLP	B	501	-	15,15,16	0.96	0	21,22,23	1.63	1 (4%)

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	-	-	0/6/6/8	0/1/1/1
3	GOL	B	502	-	-	2/4/4/4	-
3	GOL	A	502	-	-	4/4/4/4	-
2	PLP	B	501	-	-	0/6/6/8	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	PLP	C4A-C4-C5	-6.26	114.48	120.94
2	A	501	PLP	C4A-C4-C5	-6.25	114.50	120.94

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	502	GOL	C1-C2-C3-O3
3	B	502	GOL	O2-C2-C3-O3
3	A	502	GOL	C1-C2-C3-O3
3	A	502	GOL	O2-C2-C3-O3
3	A	502	GOL	O1-C1-C2-C3
3	A	502	GOL	O1-C1-C2-O2

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

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, 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	386/390 (98%)	0.25	19 (4%)	35 40	7, 17, 35, 66	4 (1%)
1	B	388/390 (99%)	0.27	31 (7%)	18 22	6, 16, 34, 64	4 (1%)
All	All	774/780 (99%)	0.26	50 (6%)	25 29	6, 17, 35, 66	8 (1%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	411	TYR	5.9
1	A	409	ASN	5.5
1	B	38	ILE	5.4
1	B	29	ASP	4.8
1	B	28[A]	ILE	4.7
1	B	31	ALA	4.1
1	B	33	PRO	4.0
1	B	410	LEU	3.8
1	B	194	PRO	3.8
1	B	32	LEU	3.8
1	B	34	ASP	3.7
1	A	195	SER	3.6
1	B	66	ASN	3.6
1	A	157	GLY	3.6
1	B	409	ASN	3.4
1	B	30	THR	3.3
1	A	36	ALA	3.3
1	A	38	ILE	3.2
1	A	196	LEU	3.2
1	B	35	ALA	3.2
1	B	36	ALA	3.1
1	A	31	ALA	3.0
1	A	158	LYS	2.9
1	A	132	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	35	ALA	2.8
1	B	65	ALA	2.8
1	B	284	ARG	2.8
1	B	155	GLN	2.7
1	B	39	GLN	2.6
1	A	29	ASP	2.5
1	B	37	GLN	2.4
1	B	26	LEU	2.4
1	A	345	THR	2.4
1	A	344	VAL	2.4
1	B	357	ILE	2.3
1	A	193	GLN	2.3
1	A	39	GLN	2.3
1	B	154	LYS	2.3
1	A	30	THR	2.3
1	B	196	LEU	2.2
1	B	195[A]	SER	2.2
1	B	404	PHE	2.2
1	B	27	HIS	2.2
1	A	133	ASP	2.1
1	A	194	PRO	2.1
1	B	408	GLU	2.1
1	B	40	GLN	2.1
1	A	64	ASN	2.0
1	B	293	LYS	2.0
1	B	64	ASN	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 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($\text{\AA}^2$)	Q<0.9
3	GOL	B	502	6/6	0.68	0.18	35,36,37,40	0
3	GOL	A	502	6/6	0.75	0.15	30,31,34,38	0
2	PLP	B	501	15/16	0.90	0.11	12,17,22,24	0
2	PLP	A	501	15/16	0.91	0.10	13,17,26,27	0
4	CL	A	503	1/1	0.99	0.17	2,2,2,2	0
4	CL	B	503	1/1	1.00	0.15	2,2,2,2	0

6.5 Other polymers [i](#)

There are no such residues in this entry.