



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:31 AM UTC

PDB ID : 4PX7 / pdb_00004px7
Title : Crystal structure of lipid phosphatase E. coli PgpB
Authors : Fan, J.; Jiang, D.; Zhao, Y.; Zhang, X.C.
Deposited on : 2014-03-22
Resolution : 3.20 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

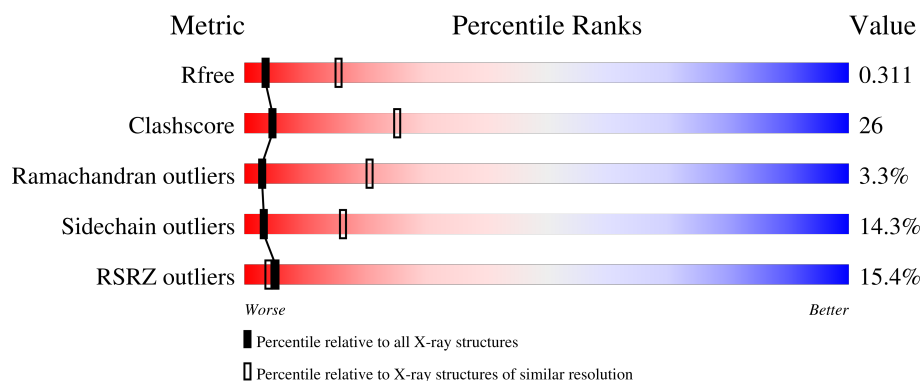
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	262	15% 60% 26% 9% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1943 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.

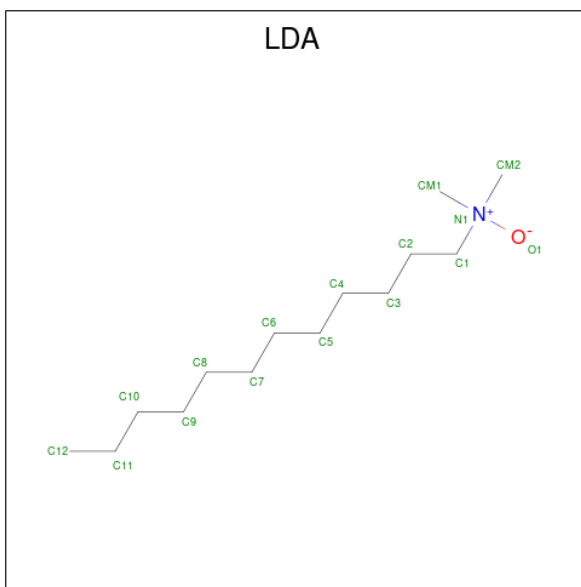
- Molecule 1 is a protein called Phosphatidylglycerophosphatase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	251	1921	1282	328	304	2	5	6	0	0

There are 11 discrepancies between the modelled and reference sequences:

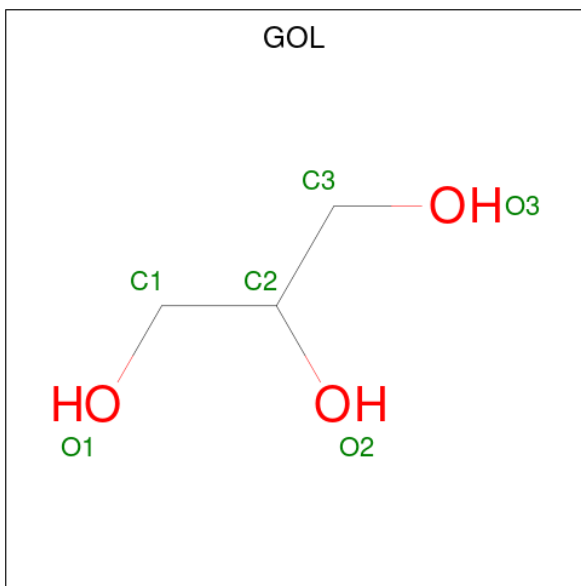
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP C6EFV6
A	116	MSE	ILE	engineered mutation	UNP C6EFV6
A	120	LYS	GLU	engineered mutation	UNP C6EFV6
A	255	LEU	-	expression tag	UNP C6EFV6
A	256	GLU	-	expression tag	UNP C6EFV6
A	257	HIS	-	expression tag	UNP C6EFV6
A	258	HIS	-	expression tag	UNP C6EFV6
A	259	HIS	-	expression tag	UNP C6EFV6
A	260	HIS	-	expression tag	UNP C6EFV6
A	261	HIS	-	expression tag	UNP C6EFV6
A	262	HIS	-	expression tag	UNP C6EFV6

- Molecule 2 is LAURYL DIMETHYLAMINE-N-OXIDE (CCD ID: LDA) (formula: C₁₄H₃₁NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			16	14	1	1		

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

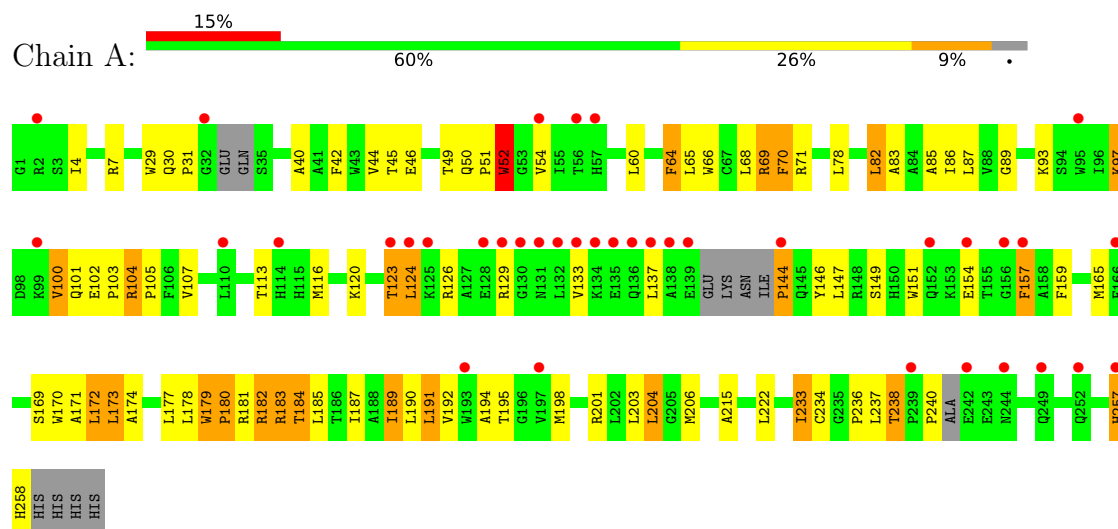


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

3 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: Phosphatidylglycerophosphatase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	36.72Å 73.67Å 131.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.86 – 3.20 32.86 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (32.86-3.20) 98.9 (32.86-3.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.08 (at 3.18Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.270 , 0.302 0.301 , 0.311	Depositor DCC
R_{free} test set	290 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	93.1	Xtriage
Anisotropy	0.674	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 62.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	1943	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.72% 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: LDA, GOL

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	0.44	0/1980	0.95	4/2711 (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

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	144	PRO	CA-N-CD	-8.87	99.59	112.00
1	A	182	ARG	N-CA-C	-8.75	95.72	109.72
1	A	238	THR	CA-C-N	-6.74	113.44	120.38
1	A	238	THR	C-N-CA	-6.74	113.44	120.38

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	180	PRO	Peptide
1	A	181	ARG	Peptide
1	A	50	GLN	Peptide

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	1921	0	1883	100	0
2	A	16	0	31	2	0
3	A	6	0	8	0	0
All	All	1943	0	1922	100	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 26.

All (100) 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:A:144:PRO:HG2	1:A:146:TYR:CE1	1.29	1.65
1:A:144:PRO:CG	1:A:146:TYR:CE1	1.90	1.50
1:A:144:PRO:HG2	1:A:146:TYR:CD1	1.59	1.37
1:A:69:ARG:HD2	1:A:70:PHE:CE1	1.62	1.32
1:A:144:PRO:CG	1:A:146:TYR:HE1	1.34	1.27
1:A:29:TRP:O	1:A:30:GLN:HG3	1.38	1.20
1:A:144:PRO:HG3	1:A:146:TYR:HE1	1.17	1.05
1:A:144:PRO:CD	1:A:146:TYR:CE1	2.52	0.92
1:A:66:TRP:HE1	1:A:69:ARG:NH2	1.70	0.89
1:A:69:ARG:CD	1:A:70:PHE:CE1	2.55	0.87
1:A:144:PRO:CD	1:A:147:LEU:HB2	2.08	0.83
1:A:69:ARG:HG3	1:A:70:PHE:CD1	2.15	0.82
1:A:51:PRO:HB2	1:A:52:TRP:NE1	1.99	0.77
1:A:66:TRP:HE1	1:A:69:ARG:HH21	1.32	0.76
1:A:144:PRO:HD3	1:A:147:LEU:HB2	1.69	0.74
1:A:66:TRP:NE1	1:A:69:ARG:NH2	2.36	0.73
1:A:69:ARG:HD2	1:A:70:PHE:CD1	2.22	0.73
1:A:29:TRP:O	1:A:30:GLN:CG	2.29	0.73
1:A:144:PRO:HD2	1:A:147:LEU:CB	2.18	0.73
1:A:144:PRO:CD	1:A:147:LEU:CB	2.68	0.72
1:A:51:PRO:HB2	1:A:52:TRP:CD1	2.24	0.71
1:A:68:LEU:HD21	1:A:177:LEU:HD13	1.74	0.69
1:A:4:ILE:HD11	1:A:237:LEU:O	1.93	0.67
1:A:179:TRP:O	1:A:179:TRP:HE3	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:PRO:CG	1:A:146:TYR:CD1	2.50	0.66
1:A:7:ARG:HD2	1:A:179:TRP:CD1	2.31	0.66
1:A:144:PRO:HD2	1:A:147:LEU:HB2	1.76	0.65
1:A:30:GLN:HB3	1:A:31:PRO:HD2	1.79	0.65
1:A:183:ARG:HH11	1:A:185:LEU:HD23	1.62	0.64
1:A:69:ARG:HG3	1:A:70:PHE:HD1	1.62	0.63
1:A:29:TRP:C	1:A:30:GLN:HG3	2.21	0.63
1:A:257:HIS:O	1:A:258:HIS:C	2.43	0.62
1:A:82:LEU:O	1:A:86:ILE:HG13	2.00	0.62
1:A:69:ARG:CD	1:A:70:PHE:CD1	2.82	0.62
1:A:165:MSE:HG3	1:A:215:ALA:HB1	1.81	0.61
1:A:144:PRO:HD2	1:A:147:LEU:H	1.66	0.60
1:A:204:LEU:HB3	1:A:206:MSE:HG2	1.83	0.60
1:A:69:ARG:O	1:A:71:ARG:N	2.36	0.59
1:A:69:ARG:CG	1:A:70:PHE:CD1	2.85	0.59
1:A:144:PRO:HG3	1:A:146:TYR:CE1	2.00	0.58
1:A:70:PHE:HB3	1:A:240:PRO:HB3	1.85	0.58
1:A:51:PRO:C	1:A:52:TRP:CD1	2.83	0.55
1:A:7:ARG:HD2	1:A:179:TRP:HD1	1.71	0.55
1:A:101:GLN:H	1:A:126:ARG:HH12	1.55	0.54
1:A:42:PHE:HA	1:A:45:THR:HG22	1.89	0.54
1:A:60:LEU:HD21	1:A:190:LEU:HD11	1.90	0.54
1:A:7:ARG:CD	1:A:179:TRP:HD1	2.22	0.53
1:A:137:LEU:HD21	1:A:151:TRP:HE3	1.74	0.53
1:A:7:ARG:NE	1:A:179:TRP:HB2	2.25	0.52
1:A:68:LEU:O	1:A:69:ARG:O	2.28	0.52
1:A:93:LYS:HD2	2:A:501:LDA:H123	1.92	0.52
1:A:194:ALA:O	1:A:198:MSE:HG3	2.10	0.52
1:A:71:ARG:HB2	1:A:237:LEU:HD12	1.92	0.52
1:A:173:LEU:O	1:A:177:LEU:HG	2.11	0.50
1:A:104:ARG:O	1:A:105:PRO:C	2.54	0.50
1:A:236:PRO:HB2	1:A:238:THR:HG23	1.93	0.50
1:A:97:LYS:HG3	1:A:159:PHE:CZ	2.48	0.49
1:A:144:PRO:HG2	1:A:146:TYR:HD1	1.57	0.48
1:A:179:TRP:O	1:A:179:TRP:CE3	2.63	0.48
1:A:144:PRO:CD	1:A:147:LEU:HB3	2.41	0.48
1:A:97:LYS:HG3	1:A:159:PHE:HZ	1.78	0.48
1:A:184:THR:HG22	1:A:187:ILE:HD12	1.95	0.48
1:A:85:ALA:HB1	1:A:222:LEU:HD11	1.96	0.47
1:A:45:THR:O	1:A:49:THR:HG23	2.15	0.46
1:A:69:ARG:HD2	1:A:70:PHE:CZ	2.37	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:ARG:CD	1:A:179:TRP:HB2	2.46	0.46
1:A:179:TRP:N	1:A:180:PRO:CD	2.78	0.46
1:A:7:ARG:CD	1:A:179:TRP:CD1	2.97	0.46
1:A:85:ALA:O	1:A:89:GLY:N	2.40	0.46
1:A:69:ARG:HD2	1:A:70:PHE:HE1	1.58	0.46
1:A:102:GLU:HG2	1:A:103:PRO:HD2	1.97	0.46
1:A:144:PRO:HD2	1:A:147:LEU:HB3	1.95	0.46
1:A:101:GLN:H	1:A:126:ARG:NH1	2.14	0.45
1:A:171:ALA:HB1	1:A:191:LEU:HD12	1.99	0.45
1:A:64:PHE:O	1:A:68:LEU:N	2.49	0.45
1:A:120:LYS:O	1:A:123:THR:HG22	2.16	0.45
1:A:129:ARG:O	1:A:133:VAL:HG23	2.17	0.44
1:A:237:LEU:HA	1:A:238:THR:OG1	2.18	0.44
1:A:174:ALA:HB2	1:A:190:LEU:HD22	1.99	0.43
1:A:83:ALA:O	1:A:86:ILE:HB	2.19	0.43
1:A:65:LEU:HD23	1:A:65:LEU:HA	1.75	0.43
1:A:233:ILE:HA	1:A:234:CYS:HA	1.71	0.42
1:A:120:LYS:O	1:A:124:LEU:HD23	2.19	0.42
1:A:204:LEU:HD12	1:A:204:LEU:HA	1.84	0.42
1:A:182:ARG:O	1:A:183:ARG:C	2.62	0.42
1:A:183:ARG:HD2	1:A:185:LEU:HB3	2.02	0.42
1:A:171:ALA:HA	1:A:190:LEU:HB3	2.02	0.42
1:A:68:LEU:HD23	1:A:68:LEU:HA	1.69	0.42
1:A:189:ILE:HD13	1:A:189:ILE:HA	1.80	0.41
1:A:102:GLU:O	1:A:157:PHE:HA	2.20	0.41
1:A:173:LEU:HG	1:A:177:LEU:HD21	2.01	0.41
1:A:104:ARG:O	1:A:107:VAL:N	2.50	0.41
1:A:146:TYR:CG	1:A:147:LEU:N	2.89	0.41
1:A:170:TRP:C	1:A:190:LEU:HD23	2.46	0.41
1:A:146:TYR:O	1:A:149:SER:HB2	2.21	0.41
1:A:169:SER:HB2	2:A:501:LDA:H12	2.02	0.40
1:A:201:ARG:HA	1:A:201:ARG:HD2	1.81	0.40
1:A:172:LEU:HD12	1:A:172:LEU:HA	1.84	0.40
1:A:123:THR:O	1:A:123:THR:HG23	2.21	0.40
1:A:40:ALA:O	1:A:44:VAL:HG23	2.21	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	243/262 (93%)	209 (86%)	26 (11%)	8 (3%)	3 21

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	54	VAL
1	A	69	ARG
1	A	70	PHE
1	A	183	ARG
1	A	52	TRP
1	A	257	HIS
1	A	154	GLU
1	A	100	VAL

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	182/212 (86%)	156 (86%)	26 (14%)	3 16

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	GLU
1	A	52	TRP
1	A	64	PHE

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Mol	Chain	Res	Type
1	A	78	LEU
1	A	82	LEU
1	A	87	LEU
1	A	97	LYS
1	A	100	VAL
1	A	104	ARG
1	A	113	THR
1	A	116	MSE
1	A	123	THR
1	A	124	LEU
1	A	157	PHE
1	A	172	LEU
1	A	173	LEU
1	A	178	LEU
1	A	179	TRP
1	A	184	THR
1	A	189	ILE
1	A	191	LEU
1	A	192	VAL
1	A	195	THR
1	A	203	LEU
1	A	204	LEU
1	A	233	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	207	HIS
1	A	231	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

2 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$
3	GOL	A	502	-	5,5,5	0.38	0	5,5,5	0.23	0
2	LDA	A	501	-	13,15,15	2.26	2 (15%)	14,17,17	0.47	0

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	GOL	A	502	-	-	2/4/4/4	-
2	LDA	A	501	-	-	5/13/13/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	LDA	O1-N1	-6.68	1.25	1.42
2	A	501	LDA	C1-N1	-4.58	1.46	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

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

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Mol	Chain	Res	Type	Atoms
3	A	502	GOL	O1-C1-C2-O2
2	A	501	LDA	C3-C4-C5-C6
2	A	501	LDA	C1-C2-C3-C4
2	A	501	LDA	C7-C8-C9-C10
2	A	501	LDA	C9-C10-C11-C12
2	A	501	LDA	C4-C5-C6-C7

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	LDA	2	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 ⓘ

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	246/262 (93%)	1.04	38 (15%) 5 4	38, 75, 109, 149	1 (0%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	138	ALA	16.1
1	A	137	LEU	13.3
1	A	139	GLU	11.9
1	A	128	GLU	8.4
1	A	242	GLU	7.6
1	A	133	VAL	6.5
1	A	56	THR	5.8
1	A	152	GLN	5.8
1	A	135	GLU	5.4
1	A	130	GLY	4.7
1	A	125	LYS	4.7
1	A	166	PHE	4.5
1	A	244	ASN	4.5
1	A	144	PRO	4.4
1	A	136	GLN	4.4
1	A	124	LEU	3.9
1	A	57	HIS	3.6
1	A	129	ARG	3.5
1	A	54	VAL	3.4
1	A	132	LEU	3.2
1	A	110	LEU	3.1
1	A	239	PRO	3.0
1	A	257	HIS	2.9
1	A	131	ASN	2.7
1	A	156	GLY	2.7
1	A	95	TRP	2.7
1	A	99	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	157	PHE	2.4
1	A	32	GLY	2.4
1	A	114	HIS	2.3
1	A	197	VAL	2.3
1	A	154	GLU	2.2
1	A	193	TRP	2.1
1	A	123	THR	2.1
1	A	249	GLN	2.0
1	A	134	LYS	2.0
1	A	252	GLN	2.0
1	A	2	ARG	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	A	502	6/6	0.55	0.32	99,110,121,125	0
2	LDA	A	501	16/16	0.79	0.26	76,92,110,111	0

6.5 Other polymers [i](#)

There are no such residues in this entry.