



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:31 PM UTC

PDB ID : 3FVZ / pdb_00003fvz
Title : Structure of Peptidyl-alpha-hydroxyglycine alpha-Amidating Lyase (PAL)
Authors : Chufan, E.E.; De, M.; Eipper, B.A.; Mains, R.E.; Amzel, L.M.
Deposited on : 2009-01-16
Resolution : 2.35 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

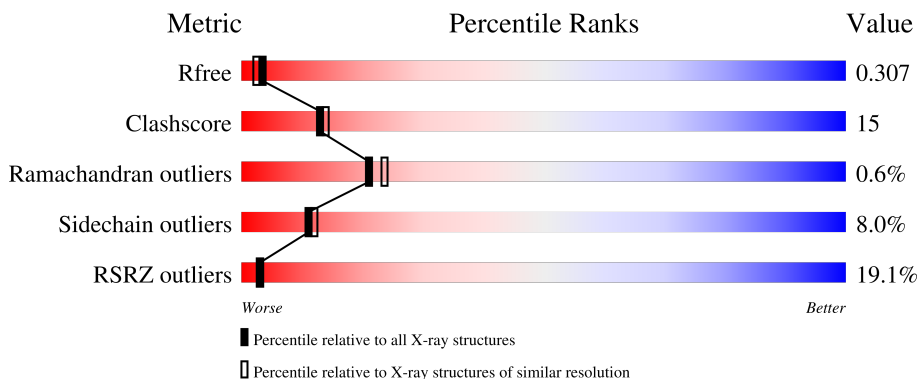
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	329	19% 76% 20% .

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 2891 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 Peptidyl-glycine alpha-amidating monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	13	0
			2724	1728	477	511	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	492	HIS	-	expression tag	UNP P14925
A	493	HIS	-	expression tag	UNP P14925
A	494	HIS	-	expression tag	UNP P14925
A	495	HIS	-	expression tag	UNP P14925
A	496	HIS	-	expression tag	UNP P14925
A	497	HIS	-	expression tag	UNP P14925

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

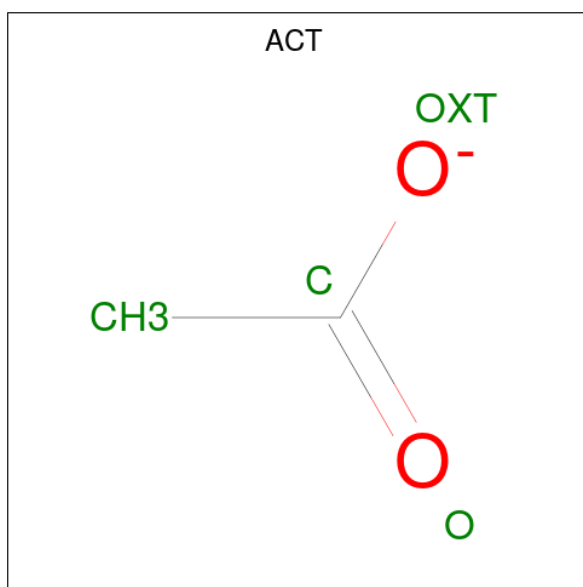
- Molecule 3 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Fe	0	0
			1	1		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

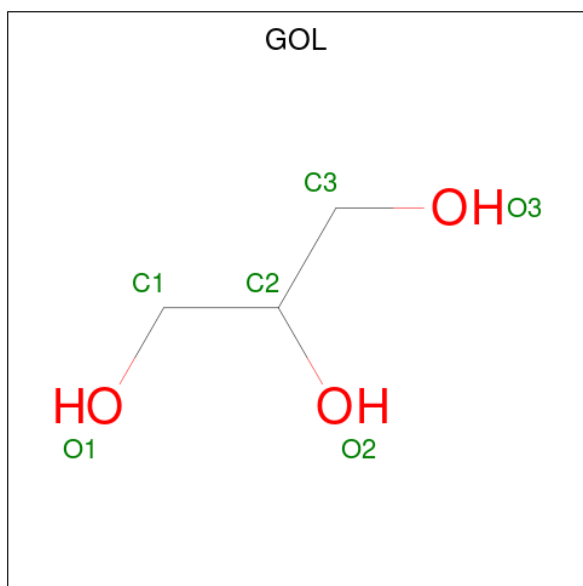
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

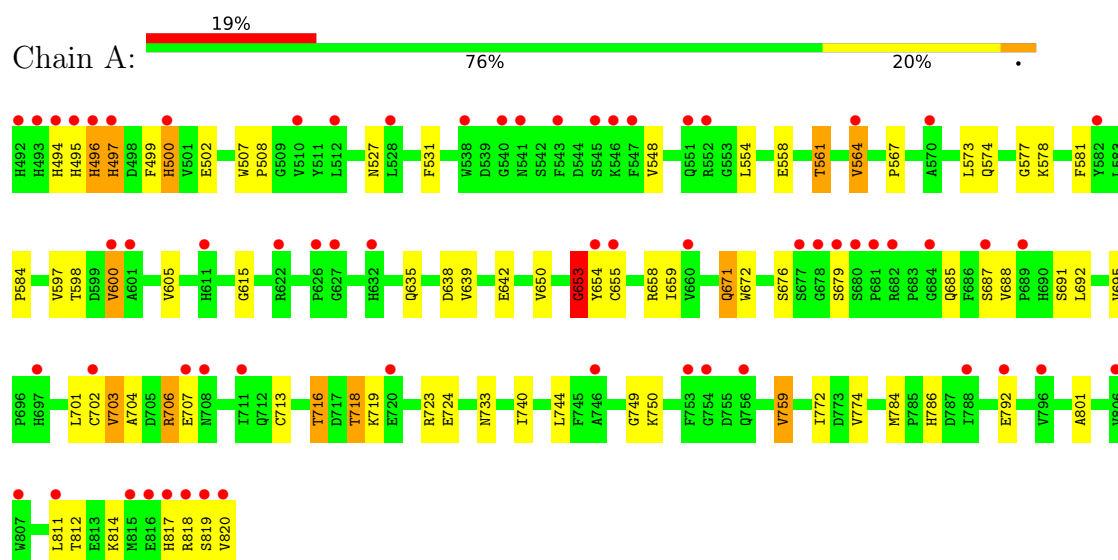
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	148	Total 148	O 148	0	0

3 Residue-property plots [i](#)

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: Peptidyl-glycine alpha-amidating monooxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.18Å 75.06Å 97.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.22 – 2.35 39.22 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.6 (39.22-2.35) 99.6 (39.22-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.82 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.203 , 0.263 0.268 , 0.307	Depositor DCC
R_{free} test set	847 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.439	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2891	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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: FE, ACT, ZN, CA, 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	0.50	0/2807	0.81	5/3814 (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	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	695	VAL	CA-C-N	6.26	125.75	119.24
1	A	695	VAL	C-N-CA	6.26	125.75	119.24
1	A	615	GLY	CA-C-N	6.10	126.12	119.90
1	A	615	GLY	C-N-CA	6.10	126.12	119.90
1	A	740	ILE	N-CA-C	-5.53	104.66	109.19

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	653	GLY	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	2724	0	2556	80	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	4	0	3	0	0
6	A	12	0	16	2	0
7	A	148	0	0	0	0
All	All	2891	0	2575	80	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 15.

All (80) 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:502[B]:GLU:OE1	1:A:818[B]:ARG:HD2	1.37	1.25
1:A:819[B]:SER:O	1:A:820[B]:VAL:HG22	1.43	1.12
1:A:784:MET:HE2	1:A:801:ALA:HB3	1.31	1.12
1:A:500[B]:HIS:HB2	1:A:819[B]:SER:HB3	1.29	1.08
1:A:500[B]:HIS:HB2	1:A:819[B]:SER:CB	1.88	1.04
1:A:496[B]:HIS:NE2	1:A:817[B]:HIS:CD2	2.32	0.96
1:A:819[B]:SER:O	1:A:820[B]:VAL:CG2	2.17	0.93
1:A:564:VAL:HG13	1:A:574:GLN:HG2	1.53	0.90
1:A:819[B]:SER:O	1:A:820[B]:VAL:HG13	1.78	0.83
1:A:500[B]:HIS:CG	1:A:819[B]:SER:HA	2.15	0.82
1:A:819[B]:SER:C	1:A:820[B]:VAL:HG22	2.06	0.79
1:A:500[B]:HIS:CB	1:A:819[B]:SER:HB3	2.13	0.78
1:A:638:ASP:CG	6:A:824:GOL:H12	2.09	0.77
1:A:500[B]:HIS:CD2	1:A:819[B]:SER:HA	2.20	0.77
1:A:502[B]:GLU:OE1	1:A:818[B]:ARG:CD	2.28	0.76
1:A:500[B]:HIS:CB	1:A:819[B]:SER:CB	2.66	0.73
1:A:558:GLU:O	1:A:578:LYS:HG3	1.93	0.69
1:A:819[B]:SER:O	1:A:820[B]:VAL:CG1	2.40	0.69
1:A:496[B]:HIS:CE1	1:A:817[B]:HIS:CD2	2.81	0.68
1:A:750:LYS:HA	1:A:759:VAL:HG22	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:527:ASN:HD22	1:A:567:PRO:HD3	1.60	0.67
1:A:819[B]:SER:O	1:A:820[B]:VAL:CB	2.44	0.65
1:A:564:VAL:CG1	1:A:574:GLN:HG2	2.24	0.65
1:A:561:THR:HG22	1:A:577:GLY:HA3	1.78	0.64
1:A:564:VAL:HG13	1:A:574:GLN:CG	2.25	0.63
1:A:784:MET:HE2	1:A:801:ALA:CB	2.19	0.62
1:A:496[B]:HIS:CG	1:A:497[B]:HIS:H	2.15	0.62
1:A:497[B]:HIS:CE1	1:A:772:ILE:HD12	2.35	0.61
1:A:500[B]:HIS:HB2	1:A:819[B]:SER:OG	2.00	0.59
1:A:502[B]:GLU:OE1	1:A:818[B]:ARG:NH1	2.35	0.59
1:A:812:THR:HG21	1:A:814:LYS:HE3	1.85	0.58
1:A:496[B]:HIS:NE2	1:A:817[B]:HIS:CG	2.72	0.57
1:A:497[B]:HIS:CE1	1:A:499:PHE:HD1	2.25	0.55
1:A:500[B]:HIS:CD2	1:A:502[B]:GLU:HG3	2.42	0.55
1:A:497[B]:HIS:CD2	1:A:772:ILE:O	2.60	0.54
1:A:496[B]:HIS:CG	1:A:497[B]:HIS:N	2.73	0.54
1:A:659:ILE:HD11	1:A:703:VAL:HG21	1.89	0.54
1:A:702:CYS:HG	1:A:713:CYS:HG	1.56	0.54
1:A:638:ASP:OD1	6:A:824:GOL:H12	2.08	0.53
1:A:564:VAL:HG22	1:A:573:LEU:HB2	1.91	0.52
1:A:642:GLU:OE2	1:A:716:THR:HG21	2.09	0.52
1:A:497[B]:HIS:CE1	1:A:772:ILE:HB	2.44	0.52
1:A:817[B]:HIS:HB3	1:A:820[B]:VAL:H	1.75	0.52
1:A:497[B]:HIS:O	1:A:817[B]:HIS:CD2	2.64	0.51
1:A:671:GLN:HG2	1:A:672:TRP:N	2.24	0.51
1:A:792[A]:GLU:H	1:A:792[A]:GLU:CD	2.19	0.51
1:A:497[B]:HIS:HE1	1:A:499:PHE:HD1	1.60	0.49
1:A:561:THR:HG21	1:A:581:PHE:CD1	2.48	0.48
1:A:692:LEU:HD22	1:A:701:LEU:HD11	1.95	0.48
1:A:733:ASN:HD22	1:A:749:GLY:HA3	1.79	0.48
1:A:676:SER:HA	1:A:685:GLN:HB3	1.95	0.47
1:A:723:ARG:HG2	1:A:724:GLU:N	2.29	0.47
1:A:500[B]:HIS:CB	1:A:819[B]:SER:OG	2.61	0.46
1:A:500[B]:HIS:CG	1:A:819[B]:SER:CA	2.93	0.46
1:A:819[B]:SER:C	1:A:820[B]:VAL:CG2	2.76	0.46
1:A:496[A]:HIS:HB3	1:A:497[A]:HIS:H	1.27	0.46
1:A:718:THR:O	1:A:719:LYS:HB2	2.15	0.45
1:A:784:MET:CE	1:A:801:ALA:HB3	2.22	0.45
1:A:639:VAL:HG22	1:A:650:VAL:HG22	1.98	0.45
1:A:502[B]:GLU:CD	1:A:818[B]:ARG:HD2	2.29	0.45
1:A:598:THR:HG22	1:A:605:VAL:HG22	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:496[B]:HIS:CD2	1:A:497[B]:HIS:H	2.35	0.44
1:A:819[A]:SER:O	1:A:820[A]:VAL:C	2.62	0.43
1:A:500[B]:HIS:CG	1:A:819[B]:SER:HB3	2.53	0.43
1:A:653:GLY:O	1:A:688[A]:VAL:HA	2.19	0.43
1:A:691:SER:HB3	1:A:704:ALA:HB3	1.99	0.43
1:A:507:TRP:HA	1:A:508:PRO:HD3	1.86	0.43
1:A:784:MET:HE3	1:A:786:HIS:CD2	2.54	0.42
1:A:531:PHE:CE2	1:A:584:PRO:HB3	2.55	0.41
1:A:687[A]:SER:C	1:A:688[A]:VAL:HG23	2.45	0.41
1:A:561:THR:HG21	1:A:581:PHE:HD1	1.85	0.41
1:A:600:VAL:HG13	1:A:635:GLN:HG2	2.02	0.41
1:A:500[A]:HIS:CE1	1:A:814:LYS:HB2	2.55	0.41
1:A:706:ARG:O	1:A:733:ASN:HA	2.20	0.41
1:A:531:PHE:CE1	1:A:561:THR:HG23	2.56	0.41
1:A:784:MET:HE3	1:A:786:HIS:HD2	1.86	0.40
1:A:494:HIS:HB2	1:A:817[B]:HIS:CE1	2.56	0.40
1:A:497[B]:HIS:NE2	1:A:499:PHE:HB2	2.37	0.40
1:A:500[B]:HIS:CG	1:A:819[B]:SER:CB	3.05	0.40
1:A:687[A]:SER:O	1:A:688[A]:VAL:HB	2.21	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	339/329 (103%)	320 (94%)	17 (5%)	2 (1%)	21 24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	495	HIS
1	A	653	GLY

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	302/289 (104%)	276 (91%)	26 (9%)	10	10

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

Mol	Chain	Res	Type
1	A	496[A]	HIS
1	A	496[B]	HIS
1	A	497[A]	HIS
1	A	497[B]	HIS
1	A	500[A]	HIS
1	A	500[B]	HIS
1	A	548	VAL
1	A	554	LEU
1	A	561	THR
1	A	564	VAL
1	A	597	VAL
1	A	600	VAL
1	A	654	TYR
1	A	655	CYS
1	A	658	ARG
1	A	671	GLN
1	A	679	SER
1	A	703	VAL
1	A	706	ARG
1	A	707	GLU
1	A	716	THR
1	A	718	THR
1	A	744	LEU
1	A	759	VAL
1	A	774	VAL
1	A	811	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	495	HIS
1	A	526	ASN
1	A	527	ASN
1	A	550	GLN
1	A	551	GLN
1	A	579	ASN
1	A	708	ASN
1	A	733	ASN
1	A	781	HIS

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, 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	$\# Z > 2$
5	ACT	A	1	2	3,3,3	0.82	0	3,3,3	1.21	0
6	GOL	A	824	4	5,5,5	0.49	0	5,5,5	0.67	0
6	GOL	A	2	-	5,5,5	0.47	0	5,5,5	0.24	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
6	GOL	A	2	-	-	4/4/4/4	-
6	GOL	A	824	4	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	824	GOL	O1-C1-C2-C3
6	A	2	GOL	O1-C1-C2-C3
6	A	2	GOL	C1-C2-C3-O3
6	A	824	GOL	O1-C1-C2-O2
6	A	2	GOL	O1-C1-C2-O2
6	A	2	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	824	GOL	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	329/329 (100%)	1.42	63 (19%) 3 3	27, 54, 61, 69	13 (3%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	820[A]	VAL	9.1
1	A	819[A]	SER	6.8
1	A	496[A]	HIS	5.7
1	A	818[A]	ARG	5.5
1	A	678	GLY	5.3
1	A	600	VAL	4.6
1	A	679	SER	4.2
1	A	492	HIS	4.0
1	A	495	HIS	3.9
1	A	547	PHE	3.8
1	A	817[A]	HIS	3.8
1	A	680	SER	3.6
1	A	654	TYR	3.5
1	A	753	PHE	3.5
1	A	552[A]	ARG	3.4
1	A	627	GLY	3.4
1	A	497[A]	HIS	3.3
1	A	601	ALA	3.1
1	A	510	VAL	3.1
1	A	681	PRO	3.1
1	A	707	GLU	3.1
1	A	754	GLY	3.0
1	A	494	HIS	3.0
1	A	788	ILE	2.8
1	A	687[A]	SER	2.8
1	A	682	ARG	2.7
1	A	720	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	622	ARG	2.6
1	A	655	CYS	2.6
1	A	697	HIS	2.6
1	A	543	PHE	2.5
1	A	545	SER	2.5
1	A	677	SER	2.5
1	A	806	VAL	2.5
1	A	570	ALA	2.5
1	A	816	GLU	2.4
1	A	546	LYS	2.4
1	A	626	PRO	2.4
1	A	538	TRP	2.4
1	A	807	TRP	2.4
1	A	500[A]	HIS	2.4
1	A	551	GLN	2.4
1	A	493	HIS	2.3
1	A	796	VAL	2.3
1	A	689	PRO	2.3
1	A	708	ASN	2.3
1	A	540	GLY	2.2
1	A	541	ASN	2.2
1	A	582	TYR	2.2
1	A	512	LEU	2.2
1	A	564	VAL	2.2
1	A	702	CYS	2.2
1	A	756	GLN	2.1
1	A	611	HIS	2.1
1	A	711	ILE	2.1
1	A	811	LEU	2.1
1	A	746	ALA	2.1
1	A	528	LEU	2.1
1	A	660	VAL	2.1
1	A	684	GLY	2.1
1	A	815	MET	2.0
1	A	792[A]	GLU	2.0
1	A	632	HIS	2.0

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

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
6	GOL	A	2	6/6	0.75	0.23	62,62,63,63	0
6	GOL	A	824	6/6	0.85	0.26	50,53,53,55	0
5	ACT	A	1	4/4	0.90	0.14	59,59,59,59	0
3	FE	A	822	1/1	0.98	0.03	48,48,48,48	1
2	ZN	A	821	1/1	0.98	0.11	61,61,61,61	0
4	CA	A	823	1/1	0.99	0.09	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.