



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:10 PM UTC

PDB ID : 7AQU / pdb_00007aqu
Title : Flavine-dependent tryptophan halogenase Thal: N-terminally His-tagged form of quintuple mutant (NHis-Thal-RebH5)
Authors : Moritzer, A.C.; Prior, T.; Niemann, H.H.
Deposited on : 2020-10-23
Resolution : 1.63 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

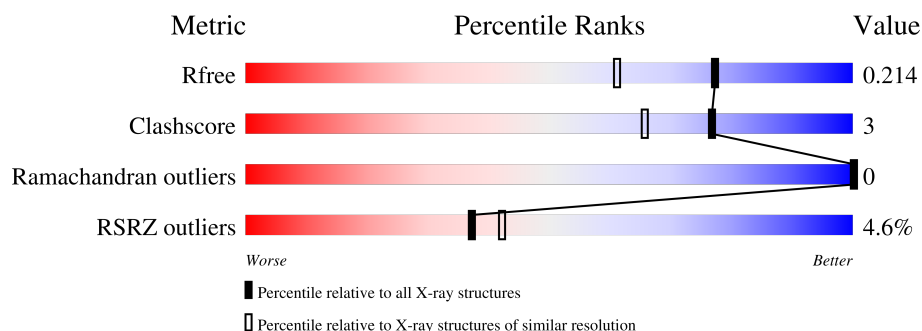
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.63 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	551	3% 88% 7% 5%
1	B	551	6% 90% 5% 5%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GLY	A	605	-	X	-	-
4	GLY	A	606	-	X	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GLY	A	607	-	X	-	-
4	GLY	B	604	-	X	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9667 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 Tryptophan 6-halogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	525	Total	C	N	O	S	0	16	0
			4319	2752	741	804	22			
1	B	523	Total	C	N	O	S	0	12	0
			4271	2731	735	784	21			

There are 50 discrepancies between the modelled and reference sequences:

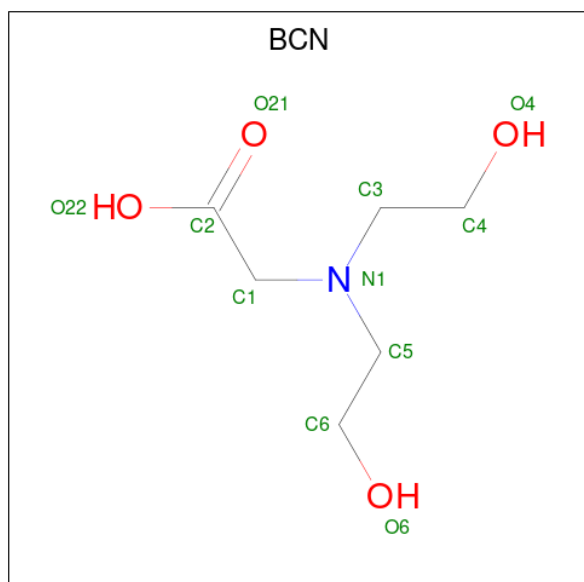
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A1E280
A	-18	GLY	-	expression tag	UNP A1E280
A	-17	SER	-	expression tag	UNP A1E280
A	-16	SER	-	expression tag	UNP A1E280
A	-15	HIS	-	expression tag	UNP A1E280
A	-14	HIS	-	expression tag	UNP A1E280
A	-13	HIS	-	expression tag	UNP A1E280
A	-12	HIS	-	expression tag	UNP A1E280
A	-11	HIS	-	expression tag	UNP A1E280
A	-10	HIS	-	expression tag	UNP A1E280
A	-9	SER	-	expression tag	UNP A1E280
A	-8	SER	-	expression tag	UNP A1E280
A	-7	GLY	-	expression tag	UNP A1E280
A	-6	LEU	-	expression tag	UNP A1E280
A	-5	VAL	-	expression tag	UNP A1E280
A	-4	PRO	-	expression tag	UNP A1E280
A	-3	ARG	-	expression tag	UNP A1E280
A	-2	GLY	-	expression tag	UNP A1E280
A	-1	SER	-	expression tag	UNP A1E280
A	0	HIS	-	expression tag	UNP A1E280
A	52	ILE	VAL	engineered mutation	UNP A1E280
A	82	ILE	VAL	engineered mutation	UNP A1E280
A	360	THR	SER	engineered mutation	UNP A1E280
A	469	SER	GLY	engineered mutation	UNP A1E280
A	470	ASN	SER	engineered mutation	UNP A1E280

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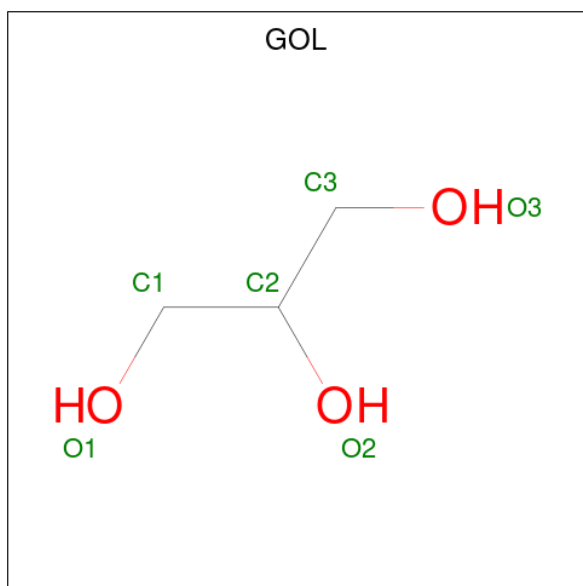
Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP A1E280
B	-18	GLY	-	expression tag	UNP A1E280
B	-17	SER	-	expression tag	UNP A1E280
B	-16	SER	-	expression tag	UNP A1E280
B	-15	HIS	-	expression tag	UNP A1E280
B	-14	HIS	-	expression tag	UNP A1E280
B	-13	HIS	-	expression tag	UNP A1E280
B	-12	HIS	-	expression tag	UNP A1E280
B	-11	HIS	-	expression tag	UNP A1E280
B	-10	HIS	-	expression tag	UNP A1E280
B	-9	SER	-	expression tag	UNP A1E280
B	-8	SER	-	expression tag	UNP A1E280
B	-7	GLY	-	expression tag	UNP A1E280
B	-6	LEU	-	expression tag	UNP A1E280
B	-5	VAL	-	expression tag	UNP A1E280
B	-4	PRO	-	expression tag	UNP A1E280
B	-3	ARG	-	expression tag	UNP A1E280
B	-2	GLY	-	expression tag	UNP A1E280
B	-1	SER	-	expression tag	UNP A1E280
B	0	HIS	-	expression tag	UNP A1E280
B	52	ILE	VAL	engineered mutation	UNP A1E280
B	82	ILE	VAL	engineered mutation	UNP A1E280
B	360	THR	SER	engineered mutation	UNP A1E280
B	469	SER	GLY	engineered mutation	UNP A1E280
B	470	ASN	SER	engineered mutation	UNP A1E280

- Molecule 2 is BICINE (CCD ID: BCN) (formula: $C_6H_{13}NO_4$).



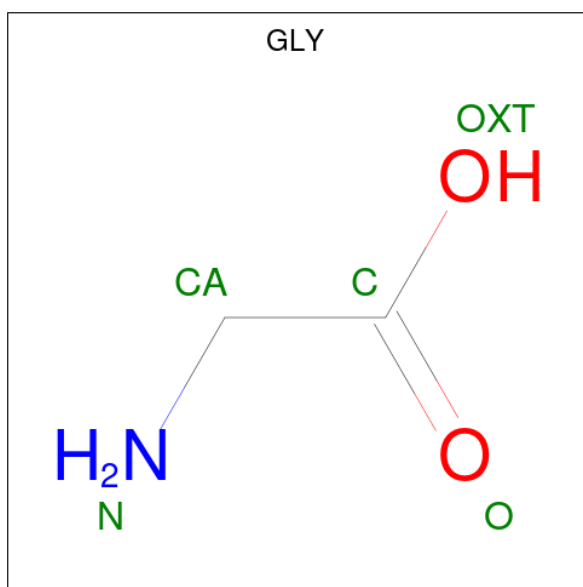
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	6	1	4		

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



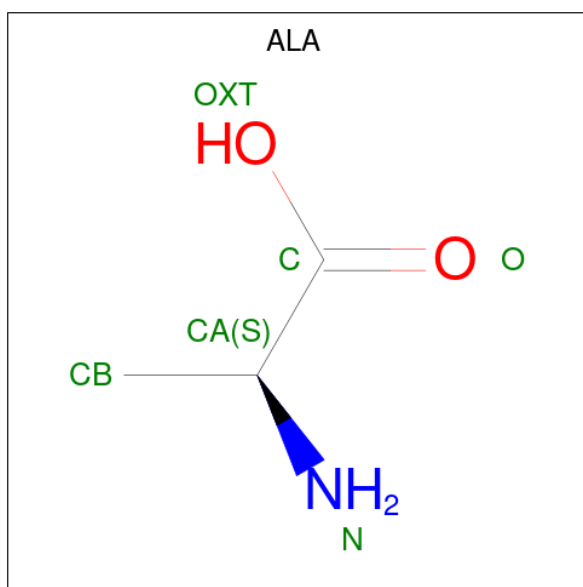
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			5	3	2		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is GLYCINE (CCD ID: GLY) (formula: $C_2H_5NO_2$).



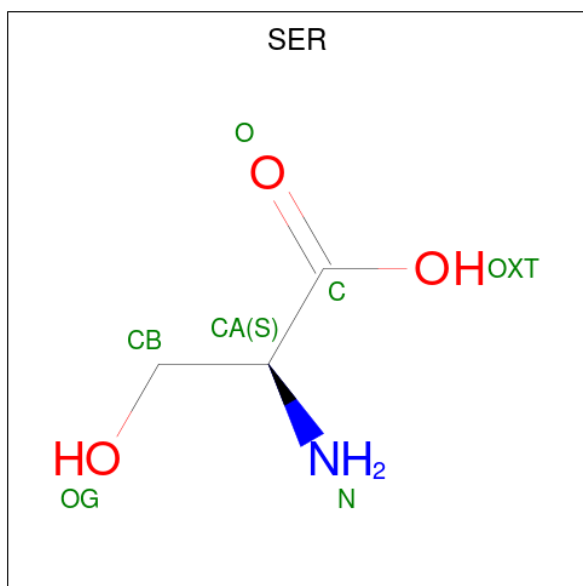
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			5	2	1	2		
4	A	1	Total	C	N	O	0	0
			5	2	1	2		
4	A	1	Total	C	N	O	0	0
			5	2	1	2		
4	B	1	Total	C	N	O	0	0
			5	2	1	2		
4	B	1	Total	C	N	O	0	0
			5	2	1	2		
4	B	1	Total	C	N	O	0	0
			5	2	1	2		

- Molecule 5 is ALANINE (CCD ID: ALA) (formula: C₃H₇NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			6	3	1	2		

- Molecule 6 is SERINE (CCD ID: SER) (formula: C₃H₇NO₃).



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

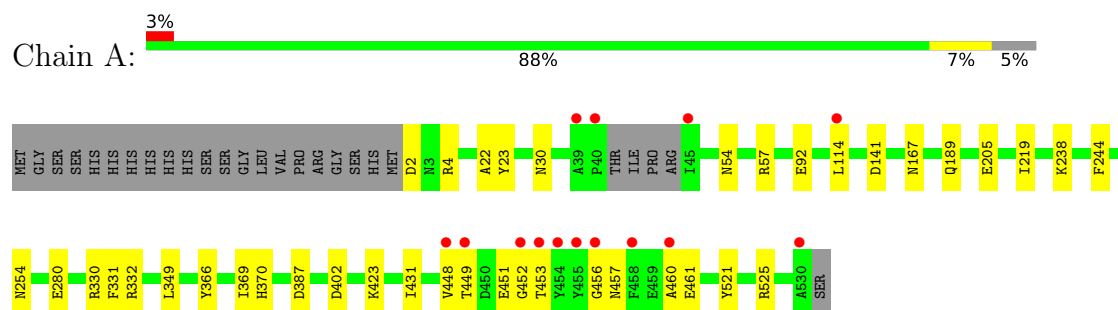
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	594	Total 604	O 604	0	10
7	B	381	Total 385	O 385	0	4

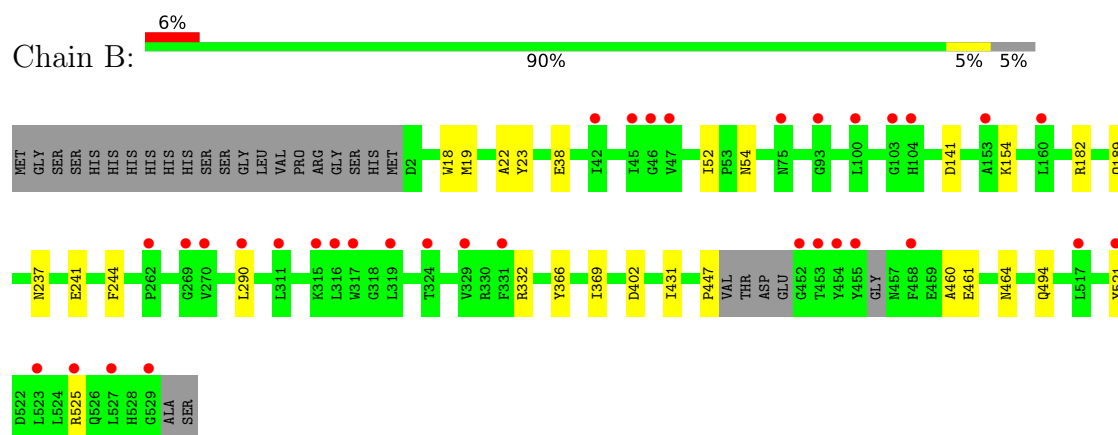
3 Residue-property plots [i](#)

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: Tryptophan 6-halogenase



• Molecule 1: Tryptophan 6-halogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.58Å 118.57Å 87.10Å 90.00° 104.04° 90.00°	Depositor
Resolution (Å)	48.58 – 1.63 48.58 – 1.63	Depositor EDS
% Data completeness (in resolution range)	96.6 (48.58-1.63) 96.6 (48.58-1.63)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 1.63Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.167 , 0.202 (Not available) , 0.214	Depositor DCC
R_{free} test set	6474 reflections (4.56%)	wwPDB-VP
Wilson B-factor (Å ²)	22.7	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9667	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 3.96% 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: BCN, 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	1.05	0/4442	1.18	4/6026 (0.1%)
1	B	0.99	0/4402	1.16	2/5969 (0.0%)
All	All	1.02	0/8844	1.17	6/11995 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	141	ASP	CA-CB-CG	7.24	119.84	112.60
1	B	141	ASP	CA-CB-CG	6.28	118.88	112.60
1	A	219	ILE	CB-CA-C	5.96	117.70	111.09
1	A	331	PHE	CA-CB-CG	5.56	119.36	113.80
1	B	52	ILE	N-CA-CB	-5.44	108.07	111.83
1	A	423	LYS	CB-CA-C	5.26	119.47	110.01

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	4319	0	4151	32	0
1	B	4271	0	4149	18	0
2	A	11	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	17	0	22	2	0
3	B	12	0	16	0	0
4	A	15	0	6	1	0
4	B	20	0	8	1	0
5	A	6	0	4	0	0
6	A	7	0	4	0	0
7	A	604	0	0	6	0
7	B	385	0	0	3	0
All	All	9667	0	8372	49	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 3.

All (49) 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:205:GLU:H	1:A:205:GLU:CD	1.99	0.70
1:A:366:TYR:HA	1:A:369[B]:ILE:HG12	1.76	0.67
1:A:448:VAL:CG2	1:B:460:ALA:HA	2.27	0.64
1:A:451:GLU:HG3	1:A:452:GLY:N	2.13	0.64
1:B:38:GLU:OE1	4:B:604:GLY:N	2.31	0.64
1:A:167:ASN:HD21	1:A:451:GLU:CD	2.10	0.60
1:A:448:VAL:HG22	1:B:460:ALA:HA	1.84	0.59
1:A:449:THR:HG21	1:A:453:THR:HG21	1.87	0.56
1:A:30:ASN:OD1	2:A:601:BCN:H32	2.06	0.56
1:B:54:ASN:ND2	1:B:461:GLU:OE1	2.40	0.55
1:B:23:TYR:HB3	1:B:369[A]:ILE:HG21	1.90	0.54
1:B:18:TRP:CZ2	1:B:182[B]:ARG:HG3	2.43	0.54
1:A:244:PHE:CZ	1:A:332:ARG:HG2	2.43	0.54
1:A:254:ASN:O	1:A:330:ARG:HD2	2.09	0.52
1:B:244:PHE:CZ	1:B:332:ARG:HG2	2.44	0.52
1:A:54:ASN:ND2	1:A:461:GLU:OE1	2.42	0.52
1:A:2:ASP:N	7:A:710:HOH:O	2.42	0.52
1:A:57:ARG:HD3	1:A:456:GLY:HA2	1.92	0.51
1:A:521:TYR:CZ	1:A:525:ARG:HD2	2.46	0.51
1:B:521:TYR:CZ	1:B:525:ARG:HD2	2.47	0.50
1:B:290:LEU:HB2	7:B:762:HOH:O	2.12	0.49
1:A:114:LEU:HD21	1:A:451:GLU:HB3	1.95	0.48
1:A:238:LYS:HE2	7:A:736:HOH:O	2.13	0.48
1:B:244:PHE:CE2	1:B:332:ARG:HG2	2.49	0.48
1:B:494:GLN:HG3	7:B:1033:HOH:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:402:ASP:HB3	1:B:431:ILE:HB	1.95	0.47
1:A:370:HIS:HE1	7:A:1195:HOH:O	1.95	0.47
1:A:2:ASP:OD2	1:A:4:ARG:NH2	2.49	0.45
1:B:447:PRO:HA	1:B:464[A]:ASN:ND2	2.31	0.45
1:A:244:PHE:CE2	1:A:332:ARG:HG2	2.52	0.45
1:B:23:TYR:CD1	1:B:369[A]:ILE:HG22	2.51	0.45
1:B:154:LYS:HE2	7:B:951:HOH:O	2.15	0.45
1:A:114:LEU:HD12	1:A:114:LEU:N	2.32	0.45
1:A:349:LEU:HD13	3:A:604:GOL:O3	2.18	0.44
1:A:448:VAL:HG13	1:A:449:THR:HG23	1.99	0.44
1:A:92:GLU:HG2	1:A:280[B]:GLU:OE2	2.17	0.44
1:B:22:ALA:HA	1:B:189:GLN:OE1	2.17	0.44
1:A:457:ASN:HD22	1:A:460:ALA:HB2	1.82	0.44
1:A:23:TYR:HB3	1:A:369[A]:ILE:HG21	2.01	0.42
1:B:19:MET:HE1	1:B:366:TYR:HD2	1.83	0.42
1:A:349:LEU:HD22	7:A:1249:HOH:O	2.20	0.42
1:B:237:ASN:O	1:B:241:GLU:HA	2.19	0.42
1:A:349:LEU:CD1	3:A:604:GOL:O3	2.68	0.42
1:A:387[B]:ASP:OD1	7:A:1138[B]:HOH:O	2.21	0.41
1:A:22:ALA:HA	1:A:189:GLN:OE1	2.21	0.41
1:A:402:ASP:HB3	1:A:431:ILE:HB	2.03	0.41
1:A:449:THR:HG21	1:A:453:THR:CG2	2.50	0.41
4:A:607:GLY:C	7:A:888:HOH:O	2.64	0.40
1:A:23:TYR:CD1	1:A:369[A]:ILE:HG22	2.57	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	537/551 (98%)	525 (98%)	12 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	529/551 (96%)	517 (98%)	12 (2%)	0	100	100
All	All	1066/1102 (97%)	1042 (98%)	24 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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](#)

15 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	604	-	5,5,5	0.21	0	5,5,5	0.46	0
6	SER	A	609	-	4,6,6	1.18	1 (25%)	2,7,7	1.34	0
4	GLY	B	605	-	4,4,4	0.93	0	3,4,4	1.46	1 (33%)
4	GLY	A	607	-	4,4,4	1.29	1 (25%)	3,4,4	1.39	1 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	A	602	-	5,5,5	0.19	0	5,5,5	0.44	0
2	BCN	A	601	-	10,10,10	0.66	0	11,11,11	0.83	0
4	GLY	A	606	-	4,4,4	1.32	1 (25%)	3,4,4	1.93	2 (66%)
3	GOL	B	601	-	5,5,5	0.33	0	5,5,5	0.65	0
3	GOL	B	602	-	5,5,5	0.21	0	5,5,5	0.48	0
4	GLY	B	604	-	4,4,4	1.03	1 (25%)	3,4,4	1.85	1 (33%)
3	GOL	A	603	-	4,4,5	0.20	0	3,3,5	0.37	0
4	GLY	B	606	-	4,4,4	1.15	1 (25%)	3,4,4	1.61	0
5	ALA	A	608	-	5,5,5	1.24	1 (20%)	6,6,6	0.55	0
4	GLY	A	605	-	4,4,4	1.22	1 (25%)	3,4,4	1.43	1 (33%)
4	GLY	B	603	-	4,4,4	1.00	0	3,4,4	2.02	2 (66%)

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	604	-	-	0/4/4/4	-
6	SER	A	609	-	-	0/6/6/6	-
4	GLY	B	605	-	-	2/2/2/2	-
4	GLY	A	607	-	-	2/2/2/2	-
3	GOL	A	602	-	-	0/4/4/4	-
2	BCN	A	601	-	-	2/10/10/10	-
4	GLY	A	606	-	-	2/2/2/2	-
3	GOL	B	601	-	-	3/4/4/4	-
3	GOL	B	602	-	-	3/4/4/4	-
4	GLY	B	604	-	-	2/2/2/2	-
3	GOL	A	603	-	-	1/2/2/4	-
4	GLY	B	606	-	-	2/2/2/2	-
5	ALA	A	608	-	-	2/4/4/4	-
4	GLY	A	605	-	-	2/2/2/2	-
4	GLY	B	603	-	-	0/2/2/2	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	606	GLY	OXT-C	-2.46	1.22	1.30
4	A	607	GLY	OXT-C	-2.45	1.22	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	609	SER	OXT-C	-2.29	1.23	1.30
4	A	605	GLY	OXT-C	-2.23	1.23	1.30
5	A	608	ALA	OXT-C	-2.16	1.23	1.30
4	B	606	GLY	OXT-C	-2.14	1.23	1.30
4	B	604	GLY	OXT-C	-2.01	1.24	1.30

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	603	GLY	OXT-C-O	-2.87	115.96	123.33
4	B	604	GLY	OXT-C-CA	2.48	123.24	113.38
4	A	606	GLY	OXT-C-O	-2.45	117.03	123.33
4	A	606	GLY	OXT-C-CA	2.24	122.28	113.38
4	A	605	GLY	OXT-C-O	-2.16	117.77	123.33
4	A	607	GLY	OXT-C-O	-2.13	117.85	123.33
4	B	603	GLY	OXT-C-CA	2.01	121.37	113.38
4	B	605	GLY	OXT-C-CA	2.00	121.35	113.38

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	GOL	O1-C1-C2-O2
3	B	602	GOL	O1-C1-C2-C3
4	A	605	GLY	O-C-CA-N
4	A	605	GLY	OXT-C-CA-N
4	A	606	GLY	O-C-CA-N
4	A	606	GLY	OXT-C-CA-N
4	B	606	GLY	O-C-CA-N
4	B	606	GLY	OXT-C-CA-N
4	A	607	GLY	O-C-CA-N
4	A	607	GLY	OXT-C-CA-N
4	B	605	GLY	OXT-C-CA-N
4	B	605	GLY	O-C-CA-N
4	B	604	GLY	O-C-CA-N
3	B	601	GOL	O1-C1-C2-C3
3	B	602	GOL	O1-C1-C2-O2
5	A	608	ALA	O-C-CA-CB
5	A	608	ALA	OXT-C-CA-CB
4	B	604	GLY	OXT-C-CA-N
3	B	601	GOL	O2-C2-C3-O3
2	A	601	BCN	C2-C1-N1-C3

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Mol	Chain	Res	Type	Atoms
2	A	601	BCN	C2-C1-N1-C5
3	B	602	GOL	C1-C2-C3-O3
3	A	603	GOL	O1-C1-C2-C3

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	604	GOL	2	0
4	A	607	GLY	1	0
2	A	601	BCN	1	0
4	B	604	GLY	1	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	525/551 (95%)	-0.16	14 (2%)	56 61	8, 20, 40, 77	16 (3%)
1	B	523/551 (94%)	0.56	34 (6%)	25 28	11, 31, 58, 77	12 (2%)
All	All	1048/1102 (95%)	0.20	48 (4%)	37 42	8, 25, 52, 77	28 (2%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	45	ILE	5.7
1	A	455	TYR	4.0
1	B	47	VAL	4.0
1	B	455	TYR	3.9
1	A	45	ILE	3.9
1	B	521	TYR	3.5
1	B	290	LEU	3.4
1	B	529	GLY	3.4
1	B	46	GLY	3.2
1	A	454	TYR	3.2
1	A	114	LEU	3.1
1	B	319	LEU	2.9
1	A	448	VAL	2.9
1	B	270	VAL	2.9
1	B	453	THR	2.8
1	A	453	THR	2.8
1	B	160	LEU	2.8
1	A	460	ALA	2.7
1	B	104	HIS	2.7
1	B	316	LEU	2.7
1	A	530	ALA	2.7
1	B	311	LEU	2.7
1	B	458	PHE	2.6
1	B	269	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	317	TRP	2.5
1	B	454	TYR	2.5
1	A	456	GLY	2.5
1	B	452	GLY	2.5
1	B	93	GLY	2.4
1	B	42	ILE	2.4
1	B	324	THR	2.4
1	B	262	PRO	2.3
1	B	523	LEU	2.3
1	B	525	ARG	2.3
1	B	75	ASN	2.3
1	A	40	PRO	2.3
1	B	153	ALA	2.3
1	B	329	VAL	2.3
1	A	452	GLY	2.3
1	B	517	LEU	2.3
1	B	527	LEU	2.2
1	B	100	LEU	2.2
1	B	331	PHE	2.1
1	A	449	THR	2.1
1	B	315	LYS	2.1
1	B	103	GLY	2.0
1	A	39	ALA	2.0
1	A	458	PHE	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	603	5/6	0.69	0.20	34,37,40,48	0
4	GLY	A	606	5/5	0.76	0.15	42,43,46,48	5
5	ALA	A	608	6/6	0.77	0.20	50,56,57,58	0
4	GLY	B	603	5/5	0.79	0.13	40,48,59,60	0
4	GLY	A	607	5/5	0.79	0.17	36,42,47,49	0
6	SER	A	609	7/7	0.80	0.16	31,45,53,67	0
4	GLY	B	605	5/5	0.81	0.15	56,56,59,59	0
4	GLY	B	606	5/5	0.81	0.15	33,39,39,47	0
3	GOL	B	601	6/6	0.84	0.14	27,32,33,33	6
4	GLY	B	604	5/5	0.84	0.14	33,38,41,45	5
4	GLY	A	605	5/5	0.84	0.17	43,53,57,65	0
3	GOL	A	604	6/6	0.88	0.13	20,23,27,30	6
3	GOL	B	602	6/6	0.90	0.10	34,37,39,41	0
2	BCN	A	601	11/11	0.90	0.11	21,33,36,37	11
3	GOL	A	602	6/6	0.92	0.08	24,26,28,28	6

6.5 Other polymers [i](#)

There are no such residues in this entry.