



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 13, 2024 – 01:58 AM EDT

PDB ID : 3EYP
Title : Crystal structure of putative alpha-L-fucosidase from Bacteroides thetaio-
taomicron
Authors : Bonanno, J.B.; Freeman, J.; Bain, K.T.; Hu, S.; Romero, R.; Wasserman, S.;
Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for
Structural Genomics (NYSGXRC)
Deposited on : 2008-10-21
Resolution : 1.90 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
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)	:	2.36.2

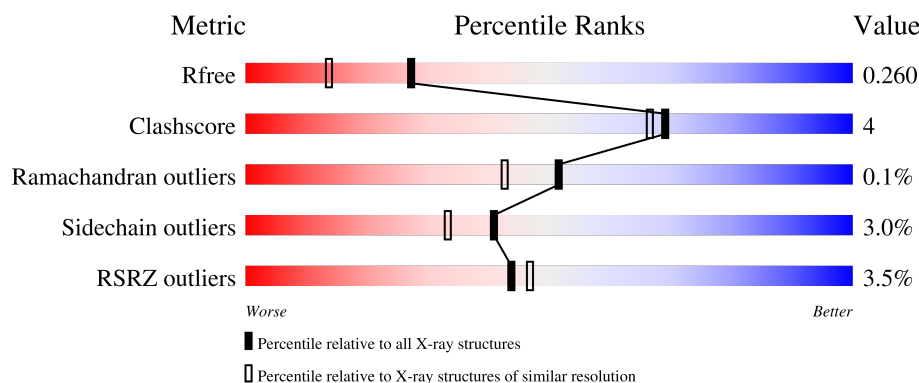
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	469	 3% 88% 9% ..
1	B	469	 4% 89% 9% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8024 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 Putative alpha-L-fucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	459	Total	C	N	O	S	0	4	0
			3681	2355	620	690	16			
1	B	459	Total	C	N	O	S	0	2	0
			3657	2338	617	686	16			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	-	expression tag	UNP Q8A5P6
A	24	SER	-	expression tag	UNP Q8A5P6
A	25	LEU	-	expression tag	UNP Q8A5P6
A	484	GLU	-	expression tag	UNP Q8A5P6
A	485	GLY	-	expression tag	UNP Q8A5P6
A	486	HIS	-	expression tag	UNP Q8A5P6
A	487	HIS	-	expression tag	UNP Q8A5P6
A	488	HIS	-	expression tag	UNP Q8A5P6
A	489	HIS	-	expression tag	UNP Q8A5P6
A	490	HIS	-	expression tag	UNP Q8A5P6
A	491	HIS	-	expression tag	UNP Q8A5P6
B	23	MET	-	expression tag	UNP Q8A5P6
B	24	SER	-	expression tag	UNP Q8A5P6
B	25	LEU	-	expression tag	UNP Q8A5P6
B	484	GLU	-	expression tag	UNP Q8A5P6
B	485	GLY	-	expression tag	UNP Q8A5P6
B	486	HIS	-	expression tag	UNP Q8A5P6
B	487	HIS	-	expression tag	UNP Q8A5P6
B	488	HIS	-	expression tag	UNP Q8A5P6
B	489	HIS	-	expression tag	UNP Q8A5P6
B	490	HIS	-	expression tag	UNP Q8A5P6
B	491	HIS	-	expression tag	UNP Q8A5P6

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

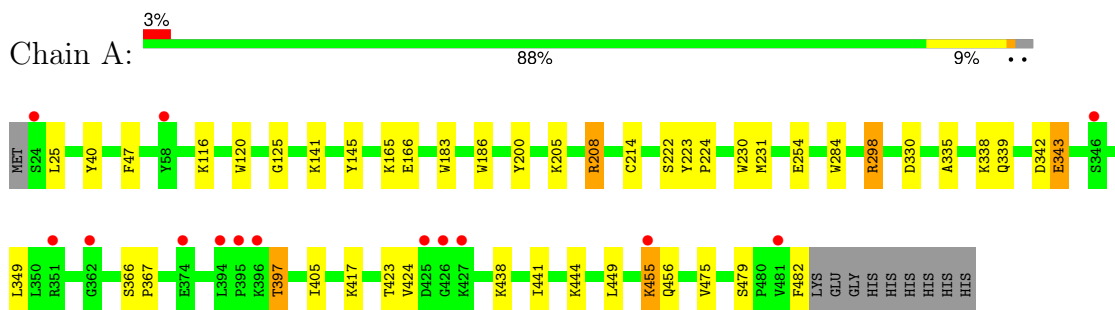
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	324	Total	O	0	0
			324	324		
3	B	314	Total	O	0	0
			314	314		

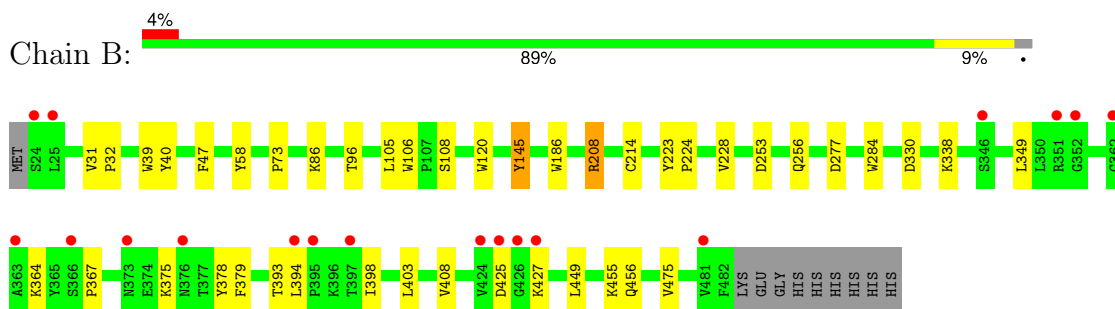
3 Residue-property plots [i](#)

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: Putative alpha-L-fucosidase



• Molecule 1: Putative alpha-L-fucosidase



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	120.71Å 125.19Å 158.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90 40.35 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-1.90) 100.0 (40.35-1.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.208 , 0.252 0.220 , 0.260	Depositor DCC
R_{free} test set	4726 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.484	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 61.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.55$, $\langle L^2 \rangle = 0.40$	Xtriage
Estimated twinning fraction	0.000 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8024	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: 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	0.94	1/3797 (0.0%)	0.88	1/5164 (0.0%)
1	B	0.91	0/3766	0.86	1/5123 (0.0%)
All	All	0.92	1/7563 (0.0%)	0.87	2/10287 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	343	GLU	CD-OE1	5.13	1.31	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	208	ARG	NE-CZ-NH2	-6.48	117.06	120.30
1	B	208	ARG	NE-CZ-NH2	-5.19	117.70	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	3681	0	3525	30	0
1	B	3657	0	3498	24	0
2	A	24	0	32	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	24	0	32	2	0
3	A	324	0	0	1	0
3	B	314	0	0	2	0
All	All	8024	0	7087	54	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 4.

All (54) 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:253:ASP:HB3	1:B:256[A]:GLN:NE2	1.99	0.78
1:A:455:LYS:O	1:A:455:LYS:HD2	1.86	0.75
1:B:284:TRP:CD1	2:B:8:GOL:H32	2.26	0.70
1:A:165:LYS:HE3	1:A:166:GLU:OE2	1.92	0.69
1:B:253:ASP:HB3	1:B:256[A]:GLN:HE21	1.56	0.69
1:A:397:THR:HG23	3:A:640:HOH:O	1.97	0.64
1:A:298:ARG:HG2	1:A:298:ARG:HH11	1.64	0.61
1:B:394:LEU:HD13	1:B:398:ILE:HD13	1.83	0.58
1:A:141:LYS:HB3	1:A:183:TRP:CD1	2.39	0.57
1:A:298:ARG:HG2	1:A:298:ARG:NH1	2.19	0.57
1:A:223:TYR:CG	1:A:224:PRO:HD3	2.40	0.57
1:A:397:THR:HG22	1:A:455:LYS:HG2	1.86	0.57
1:B:223:TYR:CG	1:B:224:PRO:HD3	2.41	0.56
1:A:423:THR:HG22	1:A:456:GLN:HB2	1.89	0.55
1:B:208:ARG:NH2	1:B:214:CYS:O	2.28	0.54
1:A:25:LEU:HD12	1:A:25:LEU:N	2.23	0.54
1:B:86:LYS:NZ	3:B:671:HOH:O	2.39	0.53
1:B:394:LEU:CD1	1:B:398:ILE:HD13	2.41	0.51
1:A:366:SER:HB2	1:A:367:PRO:HD2	1.94	0.50
1:A:223:TYR:CD2	1:A:224:PRO:HD3	2.48	0.49
1:A:284:TRP:CD1	2:A:3:GOL:H31	2.48	0.48
1:B:96:THR:HA	1:B:145:TYR:HB3	1.96	0.47
1:A:298:ARG:NH1	1:A:482:PHE:O	2.48	0.47
1:B:425:ASP:OD1	1:B:427:LYS:CB	2.63	0.47
1:A:343:GLU:HB2	1:A:479:SER:HB2	1.97	0.46
1:B:208:ARG:NE	1:B:208:ARG:HA	2.31	0.46
1:B:393:THR:HG23	1:B:456:GLN:HG2	1.98	0.46
1:B:403:LEU:C	1:B:403:LEU:HD12	2.36	0.46
1:B:39:TRP:CE2	1:B:228:VAL:HG11	2.51	0.45
1:B:284:TRP:CG	2:B:8:GOL:H32	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:TRP:CE2	2:A:2:GOL:H32	2.52	0.45
1:A:230:TRP:NE1	2:A:2:GOL:H32	2.32	0.44
1:A:335:ALA:O	1:A:339[B]:GLN:HG3	2.17	0.44
1:A:205:LYS:HB3	1:A:205:LYS:HE3	1.84	0.44
1:A:417:LYS:HA	1:A:438:LYS:O	2.18	0.43
1:B:277:ASP:HB3	3:B:545:HOH:O	2.18	0.43
1:B:367:PRO:HA	1:B:379:PHE:HB2	1.99	0.43
1:A:298:ARG:HH11	1:A:298:ARG:CG	2.31	0.43
1:A:349:LEU:HB2	1:A:475:VAL:HG23	2.00	0.43
1:B:105:LEU:HD23	1:B:105:LEU:HA	1.81	0.42
1:B:378:TYR:CG	1:B:408:VAL:HG21	2.55	0.42
1:B:73:PRO:HG2	1:B:120:TRP:CD1	2.55	0.42
1:B:223:TYR:CD2	1:B:224:PRO:HD3	2.55	0.41
1:A:208:ARG:NE	1:A:208:ARG:HA	2.35	0.41
1:A:223:TYR:CD1	1:A:224:PRO:HD3	2.55	0.41
1:A:444:LYS:HB2	1:A:444:LYS:HE2	1.89	0.41
1:A:200:TYR:CD1	1:A:222:SER:HB3	2.56	0.41
1:B:349:LEU:HB2	1:B:475:VAL:HG23	2.02	0.41
1:A:338:LYS:HE3	1:A:342:ASP:OD2	2.22	0.40
1:B:31:VAL:HB	1:B:32:PRO:HD2	2.03	0.40
1:A:208:ARG:NH2	1:A:214:CYS:O	2.30	0.40
1:A:120:TRP:CZ3	1:A:125:GLY:HA3	2.57	0.40
1:A:405:ILE:HG23	1:A:441:ILE:HG23	2.02	0.40
1:B:106:TRP:CE2	1:B:108:SER:HB2	2.56	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	461/469 (98%)	447 (97%)	13 (3%)	1 (0%)	47 38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	459/469 (98%)	445 (97%)	14 (3%)	0	100	100
All	All	920/938 (98%)	892 (97%)	27 (3%)	1 (0%)	51	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	254	GLU

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	382/393 (97%)	370 (97%)	12 (3%)	40	32
1	B	379/393 (96%)	368 (97%)	11 (3%)	42	35
All	All	761/786 (97%)	738 (97%)	23 (3%)	41	33

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	TYR
1	A	47	PHE
1	A	116	LYS
1	A	145	TYR
1	A	186	TRP
1	A	231	MET
1	A	298	ARG
1	A	330	ASP
1	A	397	THR
1	A	424	VAL
1	A	449	LEU
1	A	455	LYS
1	B	40	TYR
1	B	47	PHE
1	B	58	TYR

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Mol	Chain	Res	Type
1	B	145	TYR
1	B	186	TRP
1	B	330	ASP
1	B	338	LYS
1	B	364	LYS
1	B	375	LYS
1	B	449	LEU
1	B	455	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	373	ASN
1	A	376	ASN
1	B	122	ASN
1	B	339	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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
2	GOL	B	5	-	5,5,5	0.44	0	5,5,5	0.46	0
2	GOL	A	4	-	5,5,5	0.44	0	5,5,5	0.65	0
2	GOL	B	1	-	5,5,5	0.54	0	5,5,5	0.72	0
2	GOL	B	8	-	5,5,5	0.92	0	5,5,5	0.70	0
2	GOL	A	2	-	5,5,5	0.28	0	5,5,5	0.48	0
2	GOL	A	6	-	5,5,5	0.22	0	5,5,5	0.87	0
2	GOL	A	3	-	5,5,5	0.42	0	5,5,5	0.80	0
2	GOL	B	7	-	5,5,5	0.61	0	5,5,5	0.67	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
2	GOL	B	5	-	-	2/4/4/4	-
2	GOL	A	4	-	-	2/4/4/4	-
2	GOL	B	1	-	-	0/4/4/4	-
2	GOL	B	8	-	-	4/4/4/4	-
2	GOL	A	2	-	-	2/4/4/4	-
2	GOL	A	6	-	-	1/4/4/4	-
2	GOL	A	3	-	-	4/4/4/4	-
2	GOL	B	7	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2	GOL	C1-C2-C3-O3
2	A	3	GOL	C1-C2-C3-O3
2	A	4	GOL	C1-C2-C3-O3
2	B	5	GOL	C1-C2-C3-O3
2	B	7	GOL	O1-C1-C2-O2
2	B	7	GOL	O1-C1-C2-C3
2	B	7	GOL	C1-C2-C3-O3
2	B	7	GOL	O2-C2-C3-O3
2	B	8	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	B	8	GOL	O1-C1-C2-C3
2	B	8	GOL	C1-C2-C3-O3
2	B	5	GOL	O2-C2-C3-O3
2	A	3	GOL	O1-C1-C2-C3
2	A	2	GOL	O2-C2-C3-O3
2	A	4	GOL	O2-C2-C3-O3
2	B	8	GOL	O2-C2-C3-O3
2	A	3	GOL	O2-C2-C3-O3
2	A	6	GOL	O1-C1-C2-C3
2	A	3	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	8	GOL	2	0
2	A	2	GOL	2	0
2	A	3	GOL	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	459/469 (97%)	0.24	14 (3%) 49 51	36, 43, 56, 70	0
1	B	459/469 (97%)	0.32	18 (3%) 39 42	36, 44, 58, 76	0
All	All	918/938 (97%)	0.28	32 (3%) 44 47	36, 44, 58, 76	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	424	VAL	4.9
1	A	395	PRO	4.8
1	B	24	SER	4.7
1	B	395	PRO	4.0
1	B	426	GLY	3.5
1	A	24	SER	3.4
1	B	425	ASP	3.4
1	B	363	ALA	3.3
1	A	427	LYS	3.3
1	B	376	ASN	3.3
1	A	425	ASP	3.1
1	B	362	GLY	3.1
1	B	427	LYS	3.0
1	A	481	VAL	2.9
1	A	455	LYS	2.7
1	A	58[A]	TYR	2.7
1	B	351	ARG	2.7
1	B	481	VAL	2.6
1	A	396	LYS	2.6
1	B	366	SER	2.6
1	B	373	ASN	2.5
1	A	426	GLY	2.4
1	B	25	LEU	2.4
1	B	394	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	397	THR	2.3
1	A	394	LEU	2.3
1	A	346	SER	2.2
1	A	362	GLY	2.1
1	A	351	ARG	2.1
1	B	346	SER	2.1
1	B	352	GLY	2.1
1	A	374	GLU	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 monosaccharides 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	GOL	B	8	6/6	0.79	0.26	56,58,60,60	0
2	GOL	A	2	6/6	0.83	0.32	56,59,62,62	0
2	GOL	A	6	6/6	0.85	0.21	63,71,74,76	0
2	GOL	B	1	6/6	0.88	0.23	53,55,56,57	0
2	GOL	B	7	6/6	0.92	0.24	65,69,71,74	0
2	GOL	A	4	6/6	0.92	0.15	51,62,69,69	0
2	GOL	B	5	6/6	0.93	0.20	58,65,69,72	0
2	GOL	A	3	6/6	0.96	0.22	50,54,54,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.