



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2022 – 12:32 PM JST

PDB ID : 7F81
Title : Structure of the bacterial cellulose synthase subunit Z from Enterobacter sp. CJE-002
Authors : Fujiwara, T.; Fujishima, A.; Yao, M.
Deposited on : 2021-06-30
Resolution : 1.93 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.27
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)	:	2.27

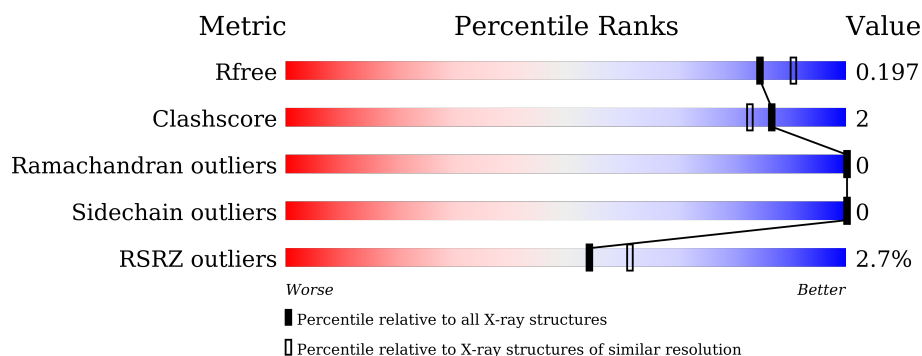
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 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	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

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	351	<div> <div>4%</div> <div>94%</div> <div>6%</div> <div>.</div> <div>.</div> </div>
1	B	351	<div> <div>4%</div> <div>90%</div> <div>6%</div> <div>.</div> </div>
1	C	351	<div> <div>%</div> <div>93%</div> <div>.</div> <div>.</div> </div>
1	D	351	<div> <div>2%</div> <div>92%</div> <div>5%</div> <div>.</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SRT	B	407	X	-	-	-
3	SRT	D	408	X	-	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11979 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 Glucanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	2	0
			2727	1752	474	495	6			
1	B	338	Total	C	N	O	S	0	1	0
			2711	1743	467	495	6			
1	C	338	Total	C	N	O	S	0	2	0
			2727	1752	474	495	6			
1	D	338	Total	C	N	O	S	0	1	0
			2711	1743	467	495	6			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	GLY	-	expression tag	UNP K0IUV6
A	18	SER	-	expression tag	UNP K0IUV6
A	19	HIS	-	expression tag	UNP K0IUV6
A	20	MET	-	expression tag	UNP K0IUV6
B	17	GLY	-	expression tag	UNP K0IUV6
B	18	SER	-	expression tag	UNP K0IUV6
B	19	HIS	-	expression tag	UNP K0IUV6
B	20	MET	-	expression tag	UNP K0IUV6
C	17	GLY	-	expression tag	UNP K0IUV6
C	18	SER	-	expression tag	UNP K0IUV6
C	19	HIS	-	expression tag	UNP K0IUV6
C	20	MET	-	expression tag	UNP K0IUV6
D	17	GLY	-	expression tag	UNP K0IUV6
D	18	SER	-	expression tag	UNP K0IUV6
D	19	HIS	-	expression tag	UNP K0IUV6
D	20	MET	-	expression tag	UNP K0IUV6

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



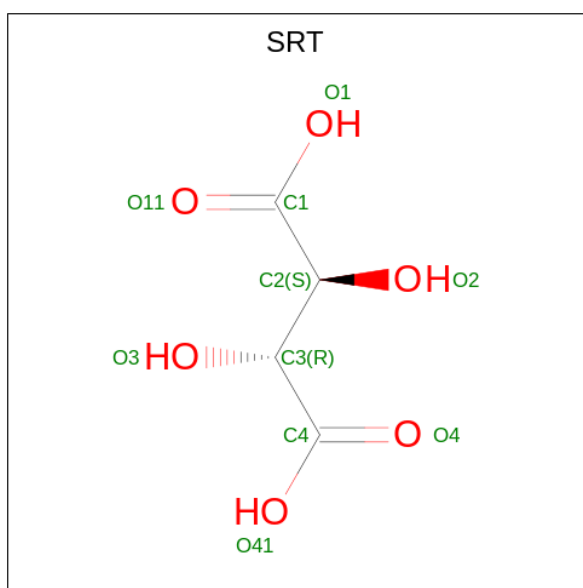
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total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

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

- Molecule 3 is S,R MESO-TARTARIC ACID (three-letter code: SRT) (formula: C₄H₆O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			10	4	6		
3	D	1	Total	C	O	0	0
			10	4	6		

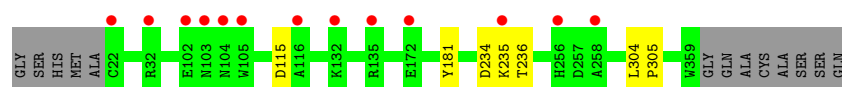
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	197	Total 197	O 197	0	0
4	B	207	Total 207	O 207	0	0
4	C	257	Total 257	O 257	0	0
4	D	296	Total 296	O 296	0	0

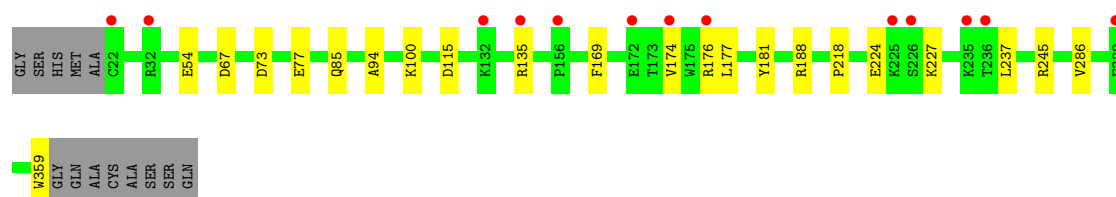
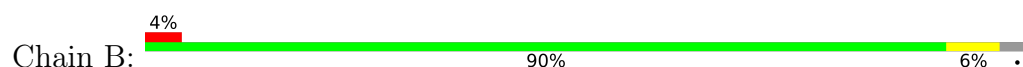
3 Residue-property plots [i](#)

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: Glucanase



- Molecule 1: Glucanase



- Molecule 1: Glucanase



- Molecule 1: Glucanase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.27Å 91.45Å 89.93Å 90.00° 98.15° 90.00°	Depositor
Resolution (Å)	38.34 – 1.93 40.02 – 1.93	Depositor EDS
% Data completeness (in resolution range)	99.7 (38.34-1.93) 99.8 (40.02-1.93)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 1.94Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.162 , 0.197 0.162 , 0.197	Depositor DCC
R_{free} test set	5343 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	25.3	Xtriage
Anisotropy	0.294	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 53.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11979	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.41% 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, SRT

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	0.41	0/2814	0.57	0/3836
1	B	0.42	0/2798	0.58	0/3816
1	C	0.51	0/2814	0.63	0/3836
1	D	0.52	0/2798	0.62	1/3816 (0.0%)
All	All	0.47	0/11224	0.60	1/15304 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	352	ARG	NE-CZ-NH2	-5.75	117.42	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	2727	0	2648	5	0
1	B	2711	0	2628	21	0
1	C	2727	0	2648	8	0
1	D	2711	0	2628	11	0
2	A	12	0	16	0	0
2	B	36	0	47	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	36	0	48	1	0
2	D	42	0	55	2	0
3	B	10	0	4	0	0
3	D	10	0	4	0	0
4	A	197	0	0	0	1
4	B	207	0	0	3	1
4	C	257	0	0	3	0
4	D	296	0	0	2	0
All	All	11979	0	10726	46	1

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 2.

All (46) 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:245:ARG:HH22	2:B:403:GOL:H31	1.24	1.01
1:B:174:VAL:HG12	1:B:224:GLU:HG2	1.60	0.84
1:B:224:GLU:HB3	1:B:227:LYS:HE3	1.71	0.71
1:B:85:GLN:NE2	2:B:404:GOL:O2	2.33	0.62
1:B:245:ARG:NH2	2:B:403:GOL:H31	2.06	0.62
1:B:135:ARG:NH2	4:B:502:HOH:O	2.36	0.59
1:C:356:GLN:NE2	4:C:507:HOH:O	2.34	0.57
1:B:94:ALA:HB2	2:B:404:GOL:H32	1.86	0.56
1:D:356:GLN:NE2	4:D:504:HOH:O	2.38	0.56
1:B:115:ASP:HB3	1:B:181:TYR:CG	2.40	0.56
1:A:115:ASP:HB3	1:A:181:TYR:CG	2.42	0.55
1:D:115:ASP:OD2	4:D:501:HOH:O	2.19	0.54
1:B:115:ASP:HB3	1:B:181:TYR:CD1	2.45	0.52
1:B:237:LEU:HG	1:B:286:VAL:HB	1.91	0.52
1:A:234:ASP:OD1	1:A:236:THR:HG22	2.10	0.51
1:D:174:VAL:HG13	2:D:407:GOL:H11	1.94	0.50
1:D:99:LYS:HE3	1:D:105:TRP:CZ2	2.47	0.50
1:D:224:GLU:OE2	1:D:227:LYS:HE3	2.12	0.50
1:B:188:ARG:O	2:B:405:GOL:H12	2.11	0.49
1:B:169:PHE:HD1	1:B:176:ARG:HD2	1.78	0.48
1:D:115:ASP:HB3	1:D:181:TYR:CG	2.50	0.47
1:C:258:ALA:O	4:C:501:HOH:O	2.20	0.47
1:C:188:ARG:NH2	1:C:259:GLN:OE1	2.48	0.47
1:B:218:PRO:HB3	1:B:286:VAL:HG21	1.97	0.46
1:B:174:VAL:CG1	1:B:224:GLU:HG2	2.39	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:ASP:HB3	1:C:181:TYR:CG	2.50	0.46
1:C:115:ASP:HB3	1:C:181:TYR:CD1	2.52	0.45
1:D:95:TRP:CE2	2:D:402:GOL:H2	2.51	0.45
1:A:115:ASP:HB3	1:A:181:TYR:CD1	2.51	0.45
1:B:67:ASP:HB2	1:B:359:TRP:CD1	2.51	0.45
2:B:402:GOL:H2	4:B:523:HOH:O	2.17	0.44
1:C:88:PRO:HA	1:C:91:HIS:O	2.17	0.44
1:B:169:PHE:CD1	1:B:176:ARG:HD2	2.53	0.43
1:B:176:ARG:C	1:B:177:LEU:HG	2.39	0.43
2:C:402:GOL:H12	4:C:514:HOH:O	2.17	0.43
1:C:224:GLU:OE1	1:C:227:LYS:HD2	2.18	0.43
1:D:176:ARG:C	1:D:177:LEU:HG	2.38	0.43
1:D:88:PRO:HA	1:D:91:HIS:O	2.19	0.42
1:B:54:GLU:OE2	4:B:501:HOH:O	2.22	0.42
1:B:100:LYS:HE2	1:D:105:TRP:CE2	2.55	0.42
1:B:169:PHE:HA	1:B:176:ARG:NH2	2.35	0.41
1:A:304:LEU:HB2	1:A:305:PRO:HD3	2.03	0.41
1:C:73:ASP:OD1	1:C:136:TYR:OH	2.33	0.41
1:B:73:ASP:O	1:B:77:GLU:HG2	2.21	0.41
1:D:67:ASP:HB2	1:D:359:TRP:CD1	2.56	0.40
1:A:234:ASP:OD1	1:A:235:LYS:N	2.55	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:615:HOH:O	4:B:648:HOH:O[1_455]	2.13	0.07

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	338/351 (96%)	335 (99%)	3 (1%)	0	100	100
1	B	337/351 (96%)	330 (98%)	7 (2%)	0	100	100
1	C	338/351 (96%)	334 (99%)	4 (1%)	0	100	100
1	D	337/351 (96%)	331 (98%)	6 (2%)	0	100	100
All	All	1350/1404 (96%)	1330 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	280/286 (98%)	280 (100%)	0	100	100
1	B	279/286 (98%)	279 (100%)	0	100	100
1	C	280/286 (98%)	280 (100%)	0	100	100
1	D	279/286 (98%)	279 (100%)	0	100	100
All	All	1118/1144 (98%)	1118 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	85	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

23 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	403	-	5,5,5	0.78	0	5,5,5	1.09	1 (20%)
2	GOL	C	405	-	5,5,5	0.98	0	5,5,5	0.91	0
2	GOL	D	403	-	5,5,5	1.15	0	5,5,5	0.77	0
2	GOL	A	401	-	5,5,5	1.08	0	5,5,5	0.95	0
2	GOL	D	402	-	5,5,5	0.91	0	5,5,5	1.17	1 (20%)
2	GOL	D	404	-	5,5,5	1.17	0	5,5,5	0.96	0
2	GOL	C	401	-	5,5,5	0.97	0	5,5,5	0.87	0
2	GOL	D	407	-	5,5,5	1.04	0	5,5,5	1.12	0
3	SRT	D	408	-	3,9,9	0.54	0	6,12,12	1.60	2 (33%)
2	GOL	B	401	-	5,5,5	0.80	0	5,5,5	0.98	0
2	GOL	C	403	-	5,5,5	0.99	0	5,5,5	0.85	0
2	GOL	D	406	-	5,5,5	1.35	1 (20%)	5,5,5	1.22	1 (20%)
2	GOL	C	402	-	5,5,5	1.01	0	5,5,5	0.91	0
2	GOL	C	406	-	5,5,5	0.64	0	5,5,5	0.97	0
2	GOL	B	402	-	5,5,5	0.86	0	5,5,5	1.05	0
2	GOL	B	404	-	5,5,5	0.91	0	5,5,5	0.89	0
2	GOL	A	402	-	5,5,5	1.06	0	5,5,5	1.09	0
2	GOL	D	405	-	5,5,5	1.18	0	5,5,5	1.35	1 (20%)
2	GOL	B	405	-	5,5,5	1.11	0	5,5,5	1.17	0
3	SRT	B	407	-	3,9,9	0.34	0	6,12,12	1.63	1 (16%)
2	GOL	C	404	-	5,5,5	0.64	0	5,5,5	0.98	0
2	GOL	B	406	-	5,5,5	1.02	0	5,5,5	0.77	0
2	GOL	D	401	-	5,5,5	0.59	0	5,5,5	0.93	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	403	-	-	4/4/4/4	-
2	GOL	C	405	-	-	0/4/4/4	-
2	GOL	D	403	-	-	1/4/4/4	-
2	GOL	A	401	-	-	0/4/4/4	-
2	GOL	D	402	-	-	0/4/4/4	-
2	GOL	D	404	-	-	2/4/4/4	-
2	GOL	C	401	-	-	0/4/4/4	-
2	GOL	D	407	-	-	3/4/4/4	-
3	SRT	D	408	-	1/1/4/4	0/4/12/12	-
2	GOL	B	401	-	-	0/4/4/4	-
2	GOL	C	403	-	-	0/4/4/4	-
2	GOL	D	406	-	-	0/4/4/4	-
2	GOL	C	402	-	-	2/4/4/4	-
2	GOL	C	406	-	-	0/4/4/4	-
2	GOL	B	402	-	-	2/4/4/4	-
2	GOL	B	404	-	-	0/4/4/4	-
2	GOL	A	402	-	-	2/4/4/4	-
2	GOL	D	405	-	-	0/4/4/4	-
2	GOL	B	405	-	-	2/4/4/4	-
3	SRT	B	407	-	1/1/4/4	0/4/12/12	-
2	GOL	C	404	-	-	2/4/4/4	-
2	GOL	B	406	-	-	2/4/4/4	-
2	GOL	D	401	-	-	0/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	406	GOL	O3-C3	-2.24	1.32	1.42

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	407	SRT	C1-C2-C3	-3.45	105.69	113.11
3	D	408	SRT	C1-C2-C3	-3.08	106.47	113.11
2	D	405	GOL	C3-C2-C1	-2.57	101.70	111.70
2	D	406	GOL	C3-C2-C1	-2.26	102.90	111.70
2	D	402	GOL	C3-C2-C1	-2.14	103.38	111.70
3	D	408	SRT	C4-C3-C2	-2.12	108.53	113.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	403	GOL	C3-C2-C1	-2.06	103.70	111.70

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	407	SRT	C2
3	D	408	SRT	C2

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	402	GOL	O1-C1-C2-O2
2	A	402	GOL	O1-C1-C2-C3
2	B	402	GOL	O2-C2-C3-O3
2	B	403	GOL	C1-C2-C3-O3
2	B	405	GOL	C1-C2-C3-O3
2	B	405	GOL	O2-C2-C3-O3
2	C	402	GOL	O1-C1-C2-C3
2	C	404	GOL	C1-C2-C3-O3
2	D	404	GOL	C1-C2-C3-O3
2	D	407	GOL	O2-C2-C3-O3
2	C	404	GOL	O2-C2-C3-O3
2	D	404	GOL	O2-C2-C3-O3
2	B	402	GOL	C1-C2-C3-O3
2	B	403	GOL	O1-C1-C2-C3
2	B	406	GOL	C1-C2-C3-O3
2	D	407	GOL	C1-C2-C3-O3
2	B	403	GOL	O1-C1-C2-O2
2	C	402	GOL	O1-C1-C2-O2
2	B	403	GOL	O2-C2-C3-O3
2	D	403	GOL	C1-C2-C3-O3
2	B	406	GOL	O2-C2-C3-O3
2	D	407	GOL	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	403	GOL	2	0
2	D	402	GOL	1	0
2	D	407	GOL	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	402	GOL	1	0
2	B	402	GOL	1	0
2	B	404	GOL	2	0
2	B	405	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	338/351 (96%)	0.03	13 (3%) 40 48	20, 32, 51, 75	0
1	B	338/351 (96%)	0.13	13 (3%) 40 48	23, 35, 58, 82	0
1	C	338/351 (96%)	-0.23	3 (0%) 84 87	14, 24, 40, 62	0
1	D	338/351 (96%)	-0.23	7 (2%) 63 70	15, 24, 42, 71	0
All	All	1352/1404 (96%)	-0.08	36 (2%) 54 61	14, 28, 49, 82	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	47	ALA	4.5
1	A	235	LYS	4.5
1	A	22	CYS	3.6
1	B	235	LYS	3.5
1	A	102	GLU	3.5
1	B	172	GLU	3.3
1	B	174	VAL	3.2
1	C	22	CYS	3.1
1	D	102	GLU	3.0
1	B	226	SER	2.9
1	B	156	PRO	2.9
1	B	225	LYS	2.9
1	B	32	ARG	2.8
1	A	258	ALA	2.8
1	C	103	ASN	2.8
1	A	103	ASN	2.7
1	A	105	TRP	2.7
1	C	258	ALA	2.7
1	A	104	ASN	2.7
1	D	45[A]	SER	2.7
1	A	256	HIS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	32	ARG	2.5
1	A	135	ARG	2.4
1	B	236	THR	2.4
1	B	338	PHE	2.4
1	A	132	LYS	2.4
1	D	105	TRP	2.3
1	B	22	CYS	2.3
1	B	132	LYS	2.3
1	A	172	GLU	2.2
1	B	176	ARG	2.2
1	B	135	ARG	2.1
1	D	226	SER	2.1
1	D	46	ASP	2.0
1	D	101	ASP	2.0
1	A	116	ALA	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	404	6/6	0.69	0.24	54,63,63,66	0
2	GOL	B	403	6/6	0.73	0.18	36,39,45,50	0
2	GOL	B	402	6/6	0.77	0.18	52,59,63,65	0
2	GOL	C	405	6/6	0.81	0.17	38,47,49,52	0
2	GOL	B	405	6/6	0.82	0.17	30,41,49,60	0
2	GOL	C	406	6/6	0.82	0.24	38,41,42,43	0
2	GOL	D	407	6/6	0.84	0.16	56,57,58,60	0
2	GOL	D	403	6/6	0.86	0.14	24,40,42,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GOL	D	405	6/6	0.89	0.15	31,34,39,39	0
2	GOL	A	402	6/6	0.91	0.19	51,52,53,54	0
2	GOL	A	401	6/6	0.93	0.22	27,30,32,32	0
2	GOL	D	404	6/6	0.94	0.10	26,37,37,37	0
2	GOL	D	402	6/6	0.94	0.13	28,37,39,43	0
2	GOL	C	404	6/6	0.94	0.12	20,25,32,37	0
3	SRT	B	407	10/10	0.94	0.09	24,27,31,33	0
2	GOL	D	406	6/6	0.95	0.13	26,27,33,40	0
2	GOL	C	402	6/6	0.95	0.12	36,39,40,43	0
2	GOL	B	406	6/6	0.95	0.11	33,37,39,39	0
2	GOL	B	401	6/6	0.96	0.12	34,36,36,37	0
2	GOL	C	403	6/6	0.96	0.09	20,23,23,24	0
2	GOL	C	401	6/6	0.97	0.14	19,20,21,21	0
3	SRT	D	408	10/10	0.97	0.07	22,24,27,29	0
2	GOL	D	401	6/6	0.99	0.06	19,21,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.