



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:41 AM UTC

PDB ID : 4OWJ / pdb_00004owj
Title : Crystal Structure of the Vibrio vulnificus Hemolysin/Cytolysin Beta-Trefoil Lectin
Authors : Kaus, K.; Olson, R.
Deposited on : 2014-02-02
Resolution : 2.00 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

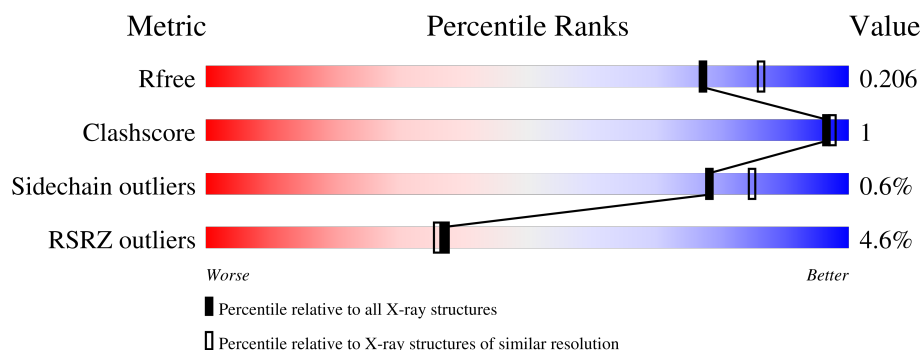
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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.



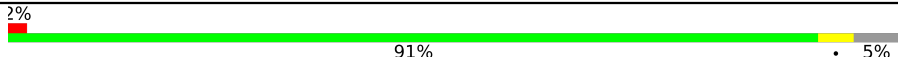
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	138	6% 89% 6% 5%
1	B	138	4% 93% 5%
1	C	138	2% 92% 5%
1	D	138	7% 95% 5%
1	E	138	7% 91% 5%
1	F	138	3% 93% 5%

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Mol	Chain	Length	Quality of chain
1	G	138	 A horizontal bar chart showing the quality of chain G. The bar is green, indicating a high quality score. The bar is labeled with '2%' at the start and '91%' at the end. A small yellow segment at the end of the bar is labeled '5%'.

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14653 atoms, of which 6631 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 Cytolysin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	131	Total	C	H	N	O	S	0	0	0
			1920	613	922	183	197	5			
1	B	131	Total	C	H	N	O	S	0	0	0
			1944	621	938	186	194	5			
1	C	131	Total	C	H	N	O	S	0	0	0
			1939	617	935	184	198	5			
1	D	131	Total	C	H	N	O	S	0	0	0
			1926	619	927	182	193	5			
1	E	131	Total	C	H	N	O	S	0	0	0
			1943	623	936	183	196	5			
1	F	132	Total	C	H	N	O	S	0	0	0
			1942	619	937	186	195	5			
1	G	131	Total	C	H	N	O	S	0	0	0
			1928	620	924	183	196	5			

There are 28 discrepancies between the modelled and reference sequences:

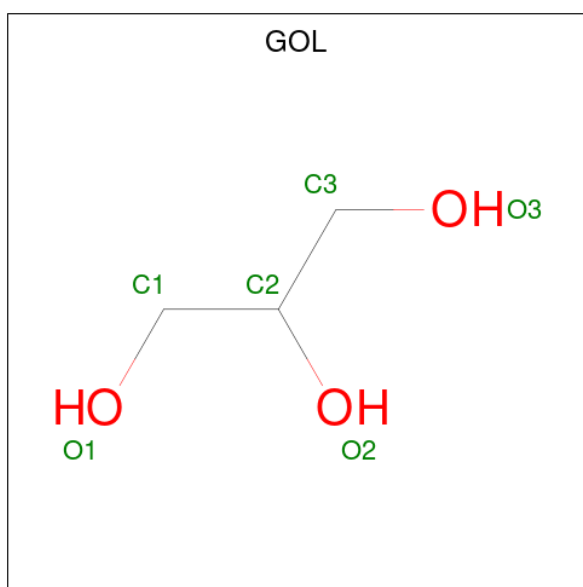
Chain	Residue	Modelled	Actual	Comment	Reference
A	334	GLY	-	expression tag	UNP P19247
A	335	SER	-	expression tag	UNP P19247
A	336	ALA	-	expression tag	UNP P19247
A	337	MET	-	expression tag	UNP P19247
B	334	GLY	-	expression tag	UNP P19247
B	335	SER	-	expression tag	UNP P19247
B	336	ALA	-	expression tag	UNP P19247
B	337	MET	-	expression tag	UNP P19247
C	334	GLY	-	expression tag	UNP P19247
C	335	SER	-	expression tag	UNP P19247
C	336	ALA	-	expression tag	UNP P19247
C	337	MET	-	expression tag	UNP P19247
D	334	GLY	-	expression tag	UNP P19247
D	335	SER	-	expression tag	UNP P19247
D	336	ALA	-	expression tag	UNP P19247

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Chain	Residue	Modelled	Actual	Comment	Reference
D	337	MET	-	expression tag	UNP P19247
E	334	GLY	-	expression tag	UNP P19247
E	335	SER	-	expression tag	UNP P19247
E	336	ALA	-	expression tag	UNP P19247
E	337	MET	-	expression tag	UNP P19247
F	334	GLY	-	expression tag	UNP P19247
F	335	SER	-	expression tag	UNP P19247
F	336	ALA	-	expression tag	UNP P19247
F	337	MET	-	expression tag	UNP P19247
G	334	GLY	-	expression tag	UNP P19247
G	335	SER	-	expression tag	UNP P19247
G	336	ALA	-	expression tag	UNP P19247
G	337	MET	-	expression tag	UNP P19247

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			14	3	8	3		
2	A	1	Total	C	H	O	0	0
			14	3	8	3		
2	B	1	Total	C	H	O	0	0
			14	3	8	3		
2	B	1	Total	C	H	O	0	0
			14	3	8	3		
2	C	1	Total	C	H	O	0	0
			14	3	8	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	H	O	0	0
			14	3	8	3		
2	D	1	Total	C	H	O	0	0
			14	3	8	3		
2	D	1	Total	C	H	O	0	0
			14	3	8	3		
2	E	1	Total	C	H	O	0	0
			14	3	8	3		
2	E	1	Total	C	H	O	0	0
			14	3	8	3		
2	F	1	Total	C	H	O	0	0
			14	3	8	3		
2	F	1	Total	C	H	O	0	0
			14	3	8	3		
2	G	1	Total	C	H	O	0	0
			14	3	8	3		
2	G	1	Total	C	H	O	0	0
			14	3	8	3		

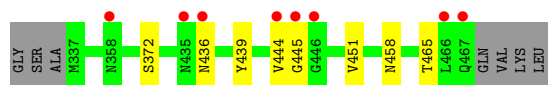
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	123	Total	O	0	0
			123	123		
3	B	147	Total	O	0	0
			147	147		
3	C	118	Total	O	0	0
			118	118		
3	D	99	Total	O	0	0
			99	99		
3	E	123	Total	O	0	0
			123	123		
3	F	152	Total	O	0	0
			152	152		
3	G	153	Total	O	0	0
			153	153		

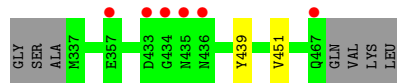
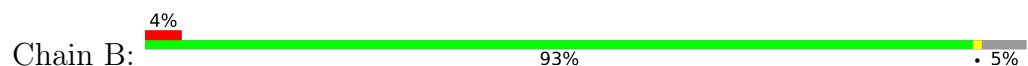
3 Residue-property plots [i](#)

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: Cytolysin



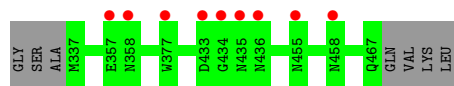
- Molecule 1: Cytolysin



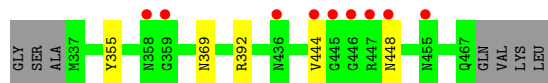
- Molecule 1: Cytolysin



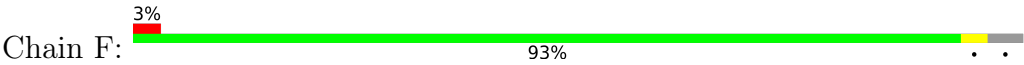
- Molecule 1: Cytolysin



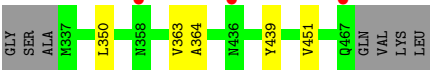
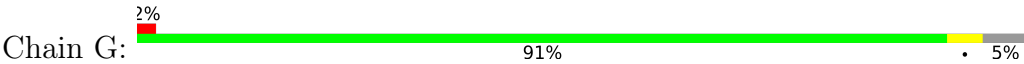
- Molecule 1: Cytolysin



- Molecule 1: Cytolysin



● Molecule 1: Cytolysin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.83Å 118.21Å 149.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.07 – 2.00 38.07 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.07-2.00) 99.9 (38.07-2.00)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.8.4-1496	Depositor
R, R_{free}	0.166 , 0.200 0.176 , 0.206	Depositor DCC
R_{free} test set	4568 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	26.3	Xtriage
Anisotropy	0.881	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14653	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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: 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	0.64	0/1016	0.76	0/1383
1	B	0.67	0/1026	0.76	0/1397
1	C	0.57	0/1022	0.72	0/1391
1	D	0.63	0/1019	0.73	0/1388
1	E	0.65	0/1027	0.76	0/1399
1	F	0.60	0/1023	0.74	0/1392
1	G	0.64	0/1024	0.77	0/1395
All	All	0.63	0/7157	0.75	0/9745

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	998	922	924	3	1
1	B	1006	938	935	1	0
1	C	1004	935	937	2	0
1	D	999	927	924	0	0
1	E	1007	936	937	2	1
1	F	1005	937	938	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	1004	924	928	3	0
2	A	12	16	16	0	0
2	B	12	16	16	0	0
2	C	12	16	16	0	0
2	D	12	16	16	0	0
2	E	12	16	16	0	0
2	F	12	16	16	0	0
2	G	12	16	16	0	0
3	A	123	0	0	1	0
3	B	147	0	0	0	0
3	C	118	0	0	1	0
3	D	99	0	0	0	0
3	E	123	0	0	0	0
3	F	152	0	0	1	0
3	G	153	0	0	0	0
All	All	8022	6631	6635	13	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 1.

All (13) 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:C:456:GLU:OE2	3:C:601:HOH:O	2.10	0.69
1:A:436:ASN:N	3:A:711:HOH:O	2.27	0.67
1:F:459:GLN:NE2	3:F:739:HOH:O	2.45	0.48
1:G:363:VAL:HG22	1:G:364:ALA:O	2.13	0.48
1:E:444:VAL:HG22	1:E:448:ASN:O	2.15	0.47
1:C:439:TYR:HB3	1:C:451:VAL:HB	1.98	0.46
1:A:444:VAL:HG23	1:A:445:GLY:N	2.32	0.45
1:B:439:TYR:HB3	1:B:451:VAL:HB	2.01	0.42
1:G:350:LEU:HD23	1:G:350:LEU:C	2.45	0.42
1:G:439:TYR:HB3	1:G:451:VAL:HB	2.02	0.41
1:F:439:TYR:HB3	1:F:451:VAL:HB	2.01	0.41
1:E:355:TYR:CZ	1:E:369:ASN:HB3	2.55	0.41
1:A:439:TYR:HB3	1:A:451:VAL:HB	2.03	0.41

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)
1:A:458:ASN:OD1	1:E:392:ARG:HH11[4_477]	1.57	0.03

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	103/116 (89%)	101 (98%)	2 (2%)	50	56
1	B	103/116 (89%)	103 (100%)	0	100	100
1	C	105/116 (90%)	104 (99%)	1 (1%)	68	75
1	D	101/116 (87%)	101 (100%)	0	100	100
1	E	104/116 (90%)	104 (100%)	0	100	100
1	F	103/116 (89%)	102 (99%)	1 (1%)	68	75
1	G	103/116 (89%)	103 (100%)	0	100	100
All	All	722/812 (89%)	718 (99%)	4 (1%)	78	85

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	372	SER
1	A	465	THR
1	C	466	LEU
1	F	392	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	435	ASN
1	C	402	ASN
1	E	459	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

14 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	G	502	-	5,5,5	0.46	0	5,5,5	0.41	0
2	GOL	D	502	-	5,5,5	0.40	0	5,5,5	0.32	0
2	GOL	F	501	-	5,5,5	0.45	0	5,5,5	0.95	0
2	GOL	C	501	-	5,5,5	0.47	0	5,5,5	0.69	0
2	GOL	D	501	-	5,5,5	0.42	0	5,5,5	0.27	0
2	GOL	C	502	-	5,5,5	0.37	0	5,5,5	0.33	0
2	GOL	B	502	-	5,5,5	0.38	0	5,5,5	0.31	0
2	GOL	E	501	-	5,5,5	0.44	0	5,5,5	0.51	0
2	GOL	B	501	-	5,5,5	0.54	0	5,5,5	1.04	0
2	GOL	E	502	-	5,5,5	0.38	0	5,5,5	0.44	0
2	GOL	G	501	-	5,5,5	0.97	1 (20%)	5,5,5	1.47	1 (20%)
2	GOL	F	502	-	5,5,5	0.35	0	5,5,5	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	A	502	-	5,5,5	0.34	0	5,5,5	0.43	0
2	GOL	A	501	-	5,5,5	0.63	0	5,5,5	1.26	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	G	502	-	-	1/4/4/4	-
2	GOL	D	502	-	-	2/4/4/4	-
2	GOL	F	501	-	-	3/4/4/4	-
2	GOL	C	501	-	-	2/4/4/4	-
2	GOL	D	501	-	-	1/4/4/4	-
2	GOL	C	502	-	-	2/4/4/4	-
2	GOL	B	502	-	-	2/4/4/4	-
2	GOL	E	501	-	-	0/4/4/4	-
2	GOL	B	501	-	-	2/4/4/4	-
2	GOL	E	502	-	-	2/4/4/4	-
2	GOL	G	501	-	-	2/4/4/4	-
2	GOL	F	502	-	-	2/4/4/4	-
2	GOL	A	502	-	-	2/4/4/4	-
2	GOL	A	501	-	-	3/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	501	GOL	O2-C2	-2.10	1.37	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	501	GOL	O1-C1-C2	-2.25	100.26	110.38

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
2	A	501	GOL	C1-C2-C3-O3
2	B	501	GOL	O1-C1-C2-O2
2	B	501	GOL	O1-C1-C2-C3
2	B	502	GOL	O1-C1-C2-C3
2	C	501	GOL	O1-C1-C2-C3
2	C	502	GOL	C1-C2-C3-O3
2	C	502	GOL	O2-C2-C3-O3
2	F	501	GOL	O1-C1-C2-C3
2	G	501	GOL	O1-C1-C2-C3
2	D	501	GOL	C1-C2-C3-O3
2	D	502	GOL	C1-C2-C3-O3
2	E	502	GOL	O1-C1-C2-C3
2	F	502	GOL	C1-C2-C3-O3
2	B	502	GOL	O1-C1-C2-O2
2	D	502	GOL	O2-C2-C3-O3
2	G	501	GOL	O1-C1-C2-O2
2	E	502	GOL	O1-C1-C2-O2
2	F	501	GOL	O2-C2-C3-O3
2	A	501	GOL	O1-C1-C2-O2
2	A	502	GOL	C1-C2-C3-O3
2	F	501	GOL	O1-C1-C2-O2
2	F	502	GOL	O2-C2-C3-O3
2	G	502	GOL	C1-C2-C3-O3
2	A	502	GOL	O2-C2-C3-O3
2	C	501	GOL	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	131/138 (94%)	-0.14	8 (6%)	27 26	23, 36, 65, 84	0
1	B	131/138 (94%)	-0.25	6 (4%)	37 36	27, 35, 61, 90	0
1	C	131/138 (94%)	-0.21	3 (2%)	61 60	28, 41, 67, 90	0
1	D	131/138 (94%)	0.01	9 (6%)	23 21	26, 40, 70, 90	0
1	E	131/138 (94%)	-0.14	9 (6%)	23 21	26, 35, 64, 90	0
1	F	132/138 (95%)	-0.31	4 (3%)	52 51	27, 35, 54, 76	0
1	G	131/138 (94%)	-0.34	3 (2%)	61 60	22, 33, 58, 79	0
All	All	918/966 (95%)	-0.20	42 (4%)	37 36	22, 37, 65, 90	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	435	ASN	5.0
1	D	455	ASN	4.8
1	A	446	GLY	4.8
1	D	434	GLY	4.6
1	E	358	ASN	4.4
1	E	446	GLY	4.3
1	D	433	ASP	4.2
1	B	467	GLN	4.2
1	E	436	ASN	4.1
1	B	434	GLY	4.1
1	A	358	ASN	3.4
1	A	467	GLN	3.4
1	A	445	GLY	3.4
1	D	436	ASN	3.4
1	G	358	ASN	3.3
1	D	357	GLU	3.3
1	B	433	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	436	ASN	3.3
1	B	435	ASN	3.2
1	B	436	ASN	3.2
1	G	467	GLN	3.1
1	E	444	VAL	3.1
1	E	447	ARG	3.0
1	A	466	LEU	2.8
1	A	444	VAL	2.8
1	C	358	ASN	2.8
1	D	358	ASN	2.8
1	F	436	ASN	2.7
1	E	445	GLY	2.7
1	B	357	GLU	2.6
1	F	455	ASN	2.6
1	F	358	ASN	2.5
1	D	377	TRP	2.5
1	D	458	ASN	2.5
1	E	448	ASN	2.5
1	C	435	ASN	2.3
1	C	467	GLN	2.3
1	F	468	GLN	2.3
1	E	359	GLY	2.2
1	G	436	ASN	2.2
1	E	455	ASN	2.1
1	A	435	ASN	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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
2	GOL	D	502	6/6	0.82	0.14	63,75,78,78	0
2	GOL	C	502	6/6	0.83	0.14	61,74,77,77	0
2	GOL	A	502	6/6	0.85	0.15	66,80,85,85	0
2	GOL	B	502	6/6	0.86	0.15	69,83,85,87	0
2	GOL	G	502	6/6	0.86	0.14	57,68,71,75	0
2	GOL	F	502	6/6	0.89	0.10	63,76,77,78	0
2	GOL	A	501	6/6	0.89	0.14	37,47,57,62	0
2	GOL	D	501	6/6	0.90	0.11	43,56,62,68	0
2	GOL	E	502	6/6	0.90	0.11	59,71,75,75	0
2	GOL	F	501	6/6	0.91	0.11	41,50,57,60	0
2	GOL	E	501	6/6	0.91	0.12	31,38,49,51	0
2	GOL	B	501	6/6	0.91	0.10	37,47,54,56	0
2	GOL	G	501	6/6	0.94	0.08	31,37,42,42	0
2	GOL	C	501	6/6	0.94	0.09	43,52,59,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.