



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:13 AM UTC

PDB ID : 8QRE / pdb_00008qre
Title : Cholera holotoxin (wildtype)
Authors : Mojica, N.; Cordara, G.; Krengel, U.
Deposited on : 2023-10-06
Resolution : 2.30 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

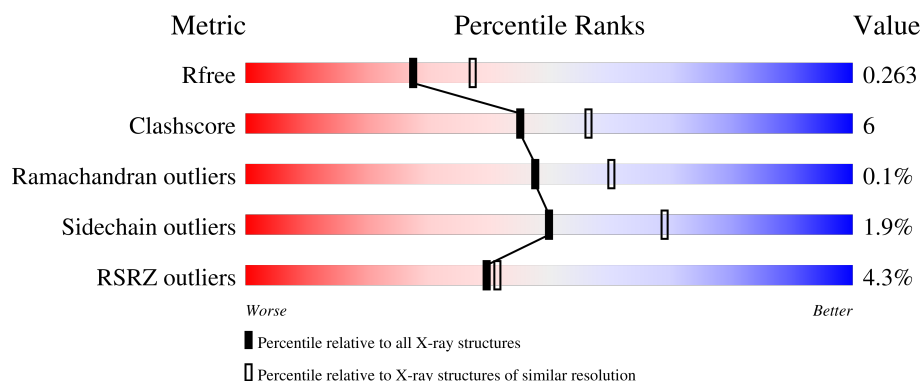
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	240	
1	B	240	
2	C	103	
2	D	103	
2	E	103	

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Mol	Chain	Length	Quality of chain
2	F	103	% 85% 15%
2	G	103	2% 90% 7%
2	H	103	% 85% 14%
2	I	103	8% 91% 8%
2	J	103	91% 9%
2	K	103	91% 7%
2	L	103	95% 5%

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 13002 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 Cholera enterotoxin subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	10	0
			1989	1236	369	378	6			
1	B	231	Total	C	N	O	S	0	11	0
			1947	1212	355	374	6			

- Molecule 2 is a protein called Cholera enterotoxin subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	103	Total	C	N	O	S	0	5	0
			853	532	150	164	7			
2	E	103	Total	C	N	O	S	0	5	0
			851	533	149	162	7			
2	F	103	Total	C	N	O	S	0	6	0
			859	537	150	165	7			
2	G	103	Total	C	N	O	S	0	3	0
			837	523	148	159	7			
2	H	103	Total	C	N	O	S	0	6	0
			859	539	150	163	7			
2	C	103	Total	C	N	O	S	0	4	0
			844	529	148	160	7			
2	I	103	Total	C	N	O	S	0	7	0
			877	550	156	164	7			
2	J	103	Total	C	N	O	S	0	2	0
			826	517	144	158	7			
2	K	103	Total	C	N	O	S	0	3	0
			834	522	145	159	8			
2	L	103	Total	C	N	O	S	0	3	0
			833	521	145	160	7			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	18	HIS	TYR	engineered mutation	UNP P01556

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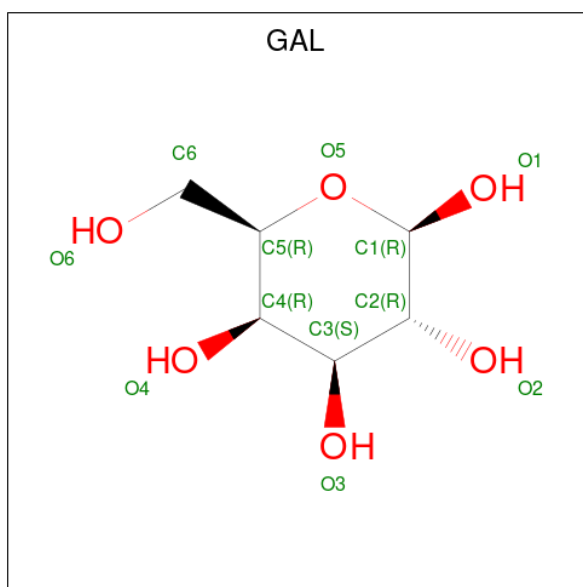
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Chain	Residue	Modelled	Actual	Comment	Reference
D	47	THR	ILE	engineered mutation	UNP P01556
E	18	HIS	TYR	engineered mutation	UNP P01556
E	47	THR	ILE	engineered mutation	UNP P01556
F	18	HIS	TYR	engineered mutation	UNP P01556
F	47	THR	ILE	engineered mutation	UNP P01556
G	18	HIS	TYR	engineered mutation	UNP P01556
G	47	THR	ILE	engineered mutation	UNP P01556
H	18	HIS	TYR	engineered mutation	UNP P01556
H	47	THR	ILE	engineered mutation	UNP P01556
C	18	HIS	TYR	engineered mutation	UNP P01556
C	47	THR	ILE	engineered mutation	UNP P01556
I	18	HIS	TYR	engineered mutation	UNP P01556
I	47	THR	ILE	engineered mutation	UNP P01556
J	18	HIS	TYR	engineered mutation	UNP P01556
J	47	THR	ILE	engineered mutation	UNP P01556
K	18	HIS	TYR	engineered mutation	UNP P01556
K	47	THR	ILE	engineered mutation	UNP P01556
L	18	HIS	TYR	engineered mutation	UNP P01556
L	47	THR	ILE	engineered mutation	UNP P01556

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Na 1 1	0	0
3	D	1	Total Na 1 1	0	0
3	B	2	Total Na 2 2	0	0

- Molecule 4 is beta-D-galactopyranose (CCD ID: GAL) (formula: C₆H₁₂O₆).



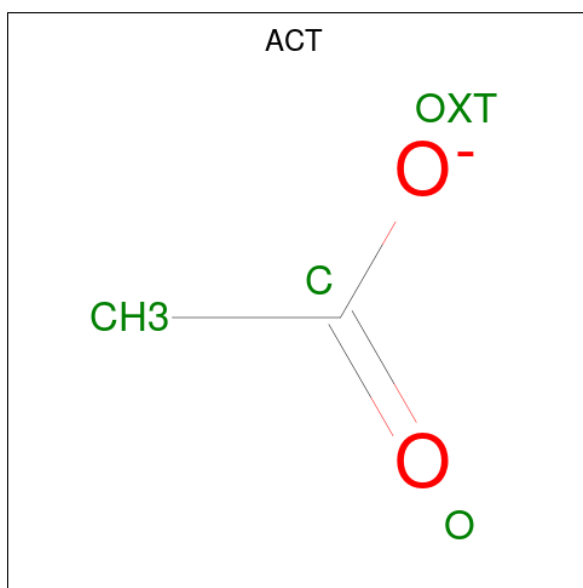
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			12	6	6		
4	D	1	Total	C	O	0	0
			12	6	6		
4	E	1	Total	C	O	0	0
			12	6	6		
4	F	1	Total	C	O	0	0
			12	6	6		
4	G	1	Total	C	O	0	0
			12	6	6		
4	H	1	Total	C	O	0	0
			12	6	6		
4	I	1	Total	C	O	0	0
			12	6	6		
4	J	1	Total	C	O	0	0
			12	6	6		
4	K	1	Total	C	O	0	0
			12	6	6		
4	L	1	Total	C	O	0	0
			12	6	6		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



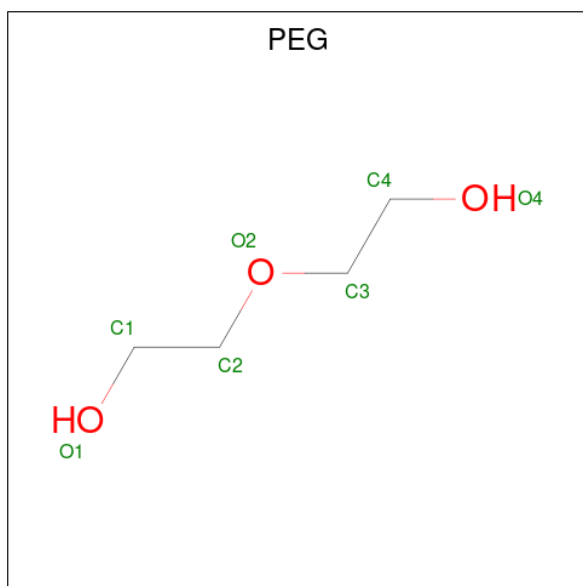
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		
5	G	1	Total	C	O	0	0
			6	3	3		
5	H	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	I	1	Total	C	O	0	0
			6	3	3		
5	I	1	Total	C	O	0	0
			6	3	3		
5	J	1	Total	C	O	0	0
			6	3	3		
5	L	1	Total	C	O	0	0
			6	3	3		
5	L	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is ACETATE ION (CCD ID: ACT) (formula: $\text{C}_2\text{H}_3\text{O}_2$).



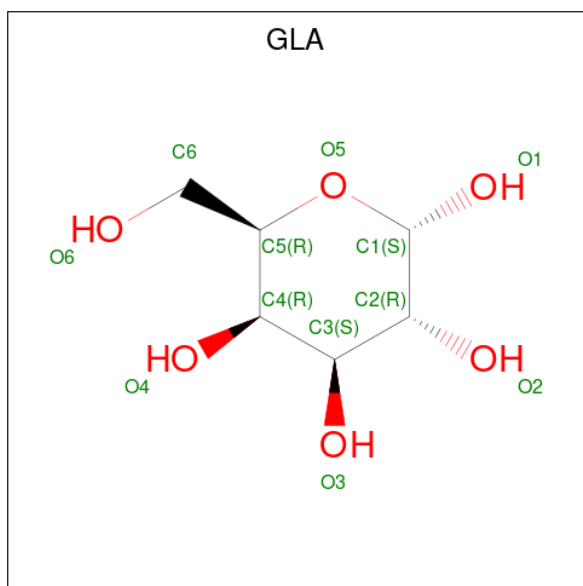
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	F	1	Total	C	O	0	0
			4	2	2		
6	G	1	Total	C	O	0	0
			4	2	2		
6	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $\text{C}_4\text{H}_{10}\text{O}_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	G	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is alpha-D-galactopyranose (CCD ID: GLA) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	C	O	0	0
			12	6	6		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	46	Total	O	0	0
			46	46		
9	D	35	Total	O	0	0
			35	35		
9	E	30	Total	O	0	0
			30	30		
9	F	31	Total	O	0	0
			31	31		
9	G	36	Total	O	0	0
			36	36		
9	H	28	Total	O	0	0
			28	28		
9	B	37	Total	O	0	0
			37	37		
9	C	28	Total	O	0	0
			28	28		

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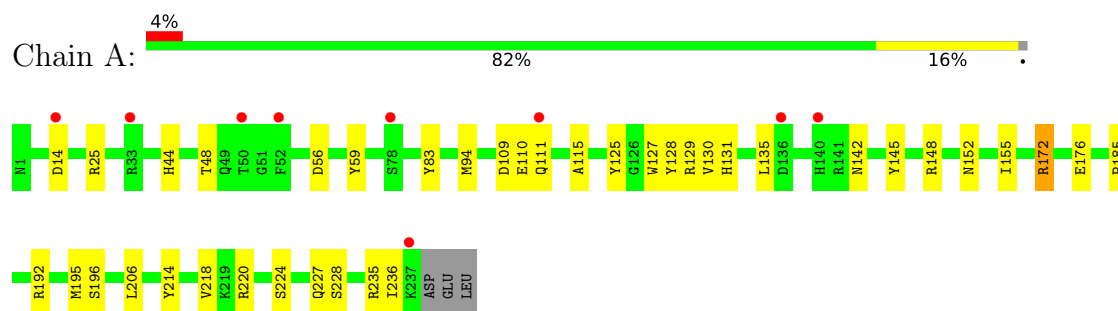
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	I	16	Total 16	O 16	0	0
9	J	23	Total 23	O 23	0	0
9	K	16	Total 16	O 16	0	0
9	L	28	Total 28	O 28	0	0

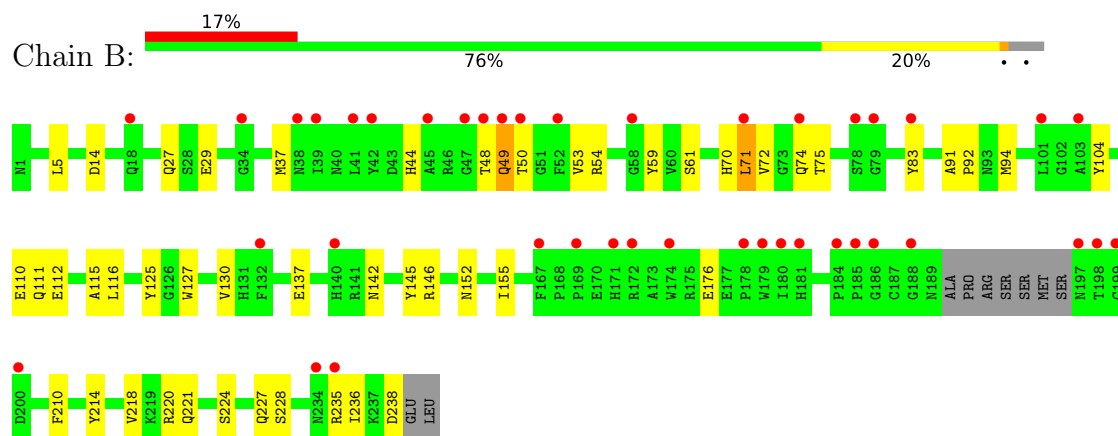
3 Residue-property plots

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.

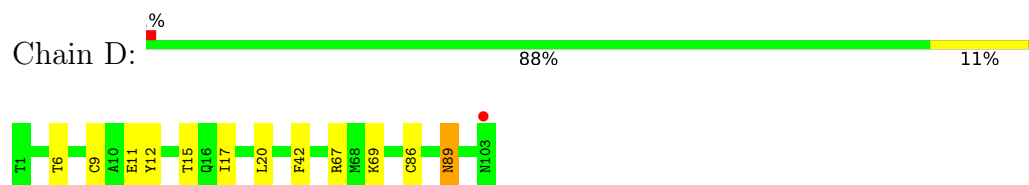
• Molecule 1: Cholera enterotoxin subunit A



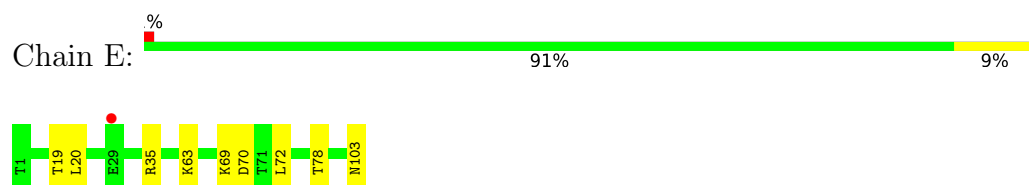
• Molecule 1: Cholera enterotoxin subunit A



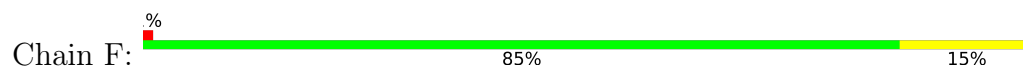
• Molecule 2: Cholera enterotoxin subunit B



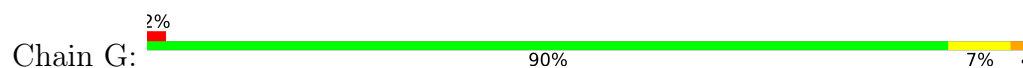
• Molecule 2: Cholera enterotoxin subunit B



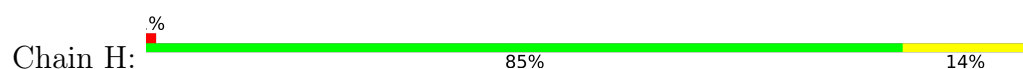
● Molecule 2: Cholera enterotoxin subunit B



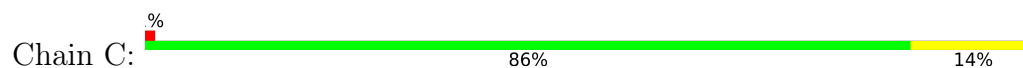
● Molecule 2: Cholera enterotoxin subunit B



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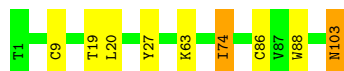
● Molecule 2: Cholera enterotoxin subunit B



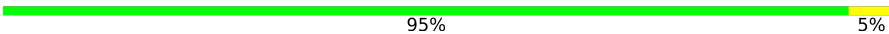
● Molecule 2: Cholera enterotoxin subunit B



● Molecule 2: Cholera enterotoxin subunit B



- Molecule 2: Cholera enterotoxin subunit B

Chain L:  95% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.66Å 108.13Å 124.77Å 90.00° 95.96° 90.00°	Depositor
Resolution (Å)	124.09 – 2.30 124.09 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.6 (124.09-2.30) 98.6 (124.09-2.30)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.98 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, R_{free}	0.216 , 0.263 0.220 , 0.263	Depositor DCC
R_{free} test set	7803 reflections (3.04%)	wwPDB-VP
Wilson B-factor (Å ²)	16.7	Xtriage
Anisotropy	0.185	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 41.6	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	0.94	EDS
Total number of atoms	13002	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.58% 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: ACT, NA, PEG, GAL, GOL, GLA

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	0.59	0/2044	0.82	0/2768
1	B	0.55	0/1999	0.82	0/2709
2	C	0.56	0/858	0.77	0/1156
2	D	0.56	0/867	0.78	0/1170
2	E	0.57	0/865	0.78	0/1166
2	F	0.54	0/873	0.80	0/1177
2	G	0.56	0/851	0.79	0/1148
2	H	0.56	0/873	0.77	0/1177
2	I	0.53	0/891	0.78	0/1199
2	J	0.55	0/840	0.79	0/1134
2	K	0.54	0/848	0.76	0/1144
2	L	0.54	0/847	0.76	0/1144
All	All	0.56	0/12656	0.79	0/17092

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	1989	0	1861	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1947	0	1809	42	0
2	C	844	0	849	9	0
2	D	853	0	848	16	0
2	E	851	0	853	11	0
2	F	859	0	856	17	0
2	G	837	0	837	10	0
2	H	859	0	866	13	0
2	I	877	0	890	14	0
2	J	826	0	825	9	0
2	K	834	0	833	10	0
2	L	833	0	831	7	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	D	1	0	0	0	0
4	A	12	0	12	1	0
4	D	12	0	12	0	0
4	E	12	0	12	0	0
4	F	12	0	12	0	0
4	G	12	0	11	0	0
4	H	12	0	12	0	0
4	I	12	0	12	0	0
4	J	12	0	12	0	0
4	K	12	0	12	1	0
4	L	12	0	12	0	0
5	A	18	0	24	0	0
5	B	6	0	8	0	0
5	C	6	0	8	0	0
5	D	6	0	8	0	0
5	E	6	0	8	0	0
5	G	6	0	8	0	0
5	H	6	0	8	1	0
5	I	12	0	16	0	0
5	J	6	0	8	0	0
5	L	12	0	16	0	0
6	C	4	0	3	0	0
6	F	4	0	3	0	0
6	G	4	0	3	0	0
7	G	7	0	10	0	0
8	C	12	0	12	0	0
9	A	46	0	0	2	0
9	B	37	0	0	3	0
9	C	28	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	D	35	0	0	1	0
9	E	30	0	0	0	0
9	F	31	0	0	1	0
9	G	36	0	0	0	0
9	H	28	0	0	1	0
9	I	16	0	0	1	0
9	J	23	0	0	1	0
9	K	16	0	0	2	0
9	L	28	0	0	1	0
All	All	13002	0	12420	149	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 6.

All (149) 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)
2:D:9[B]:CYS:SG	2:D:15[B]:THR:CG2	2.22	1.28
2:F:9[B]:CYS:SG	2:F:15[B]:THR:CG2	2.26	1.24
1:B:146[B]:ARG:HH11	1:B:146[B]:ARG:HG2	1.01	1.17
2:F:9[B]:CYS:SG	2:F:15[B]:THR:HG23	1.85	1.12
1:B:227:GLN:OE1	2:I:74[A]:ILE:HG22	1.53	1.08
2:F:51:GLU:OE2	2:F:91:LYS:HE2	1.56	1.04
2:D:9[B]:CYS:SG	2:D:15[B]:THR:HG22	1.95	1.03
2:F:9[B]:CYS:SG	2:F:15[B]:THR:HG21	2.03	0.98
1:B:146[B]:ARG:HG2	1:B:146[B]:ARG:NH1	1.78	0.92
2:D:6:THR:HA	2:D:17:ILE:HD11	1.56	0.86
2:D:9[B]:CYS:SG	2:D:15[B]:THR:HG23	2.17	0.83
2:D:9[B]:CYS:SG	2:D:15[B]:THR:HG21	2.19	0.82
2:K:74:ILE:HD13	2:L:77:LEU:HD13	1.62	0.81
1:A:94:MET:HE3	1:A:115:ALA:HB2	1.64	0.79
2:G:74:ILE:HD13	2:H:77:LEU:HD13	1.64	0.78
2:I:81[B]:LYS:HD2	2:I:103:ASN:ND2	2.00	0.76
1:A:235:ARG:NH2	2:E:70:ASP:OD2	2.17	0.76
2:H:103:ASN:ND2	9:H:301:HOH:O	2.17	0.75
2:F:65:ILE:O	2:F:69[B]:LYS:HD3	1.88	0.74
2:D:6:THR:HA	2:D:17:ILE:CD1	2.17	0.74
1:B:49[B]:GLN:C	1:B:50:THR:HG1	1.97	0.72
1:B:224:SER:O	2:J:74:ILE:HD11	1.89	0.72
1:B:49[B]:GLN:O	1:B:50:THR:OG1	2.05	0.72
2:I:103:ASN:O	9:I:301:HOH:O	2.10	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:GLN:CD	2:I:74[A]:ILE:HG22	2.19	0.67
1:B:228:SER:CB	2:J:74:ILE:HG12	2.24	0.67
1:A:224:SER:O	2:F:74:ILE:HD11	1.94	0.67
2:D:11[B]:GLU:OE1	2:E:35:ARG:CZ	2.42	0.67
1:B:146[B]:ARG:HH11	1:B:146[B]:ARG:CG	1.90	0.66
2:L:22:ASP:OD2	9:L:301:HOH:O	2.14	0.66
1:B:94[A]:MET:SD	1:B:115:ALA:HB2	2.36	0.65
2:I:74[A]:ILE:HD12	2:I:75:ALA:N	2.11	0.65
2:C:2:PRO:HG3	2:C:8:LEU:HD23	1.80	0.64
1:B:228:SER:HB3	2:J:74:ILE:HG12	1.80	0.63
2:K:27:TYR:OH	9:K:301:HOH:O	2.14	0.63
2:J:13:HIS:CD2	2:J:14:ASN:OD1	2.52	0.63
2:F:20:LEU:HB3	2:F:22[B]:ASP:OD1	1.99	0.62
2:G:103:ASN:O	2:G:103:ASN:ND2	2.33	0.61
2:E:103:ASN:HD22	2:F:25:PHE:HA	1.65	0.60
1:B:48[A]:THR:O	1:B:49[A]:GLN:HB2	2.02	0.60
2:D:11[B]:GLU:OE1	2:E:35:ARG:NE	2.35	0.59
1:B:49[A]:GLN:O	1:B:50:THR:OG1	2.13	0.59
1:B:236:ILE:HD13	2:I:63:LYS:HB2	1.84	0.59
2:J:63:LYS:HG2	9:J:306:HOH:O	2.02	0.58
1:B:48[B]:THR:HG22	1:B:48[B]:THR:O	2.01	0.58
1:B:70:HIS:O	1:B:74:GLN:HG2	2.03	0.58
1:B:137:GLU:HG3	1:B:137:GLU:O	2.03	0.58
1:B:220:ARG:HA	2:I:78:THR:HB	1.85	0.58
1:A:185:PRO:HB3	2:I:45:GLY:HA3	1.85	0.57
1:A:235:ARG:HD3	4:A:302:GAL:H62	1.84	0.57
1:B:146[B]:ARG:NH1	1:B:221:GLN:OE1	2.37	0.57
1:A:111[A]:GLN:O	1:A:111[A]:GLN:HG3	2.05	0.56
2:I:81[B]:LYS:HD2	2:I:103:ASN:HD21	1.69	0.56
2:I:74[A]:ILE:HD13	2:J:77:LEU:CD1	2.36	0.55
1:A:125:TYR:HD1	1:A:145:TYR:CE2	2.25	0.55
1:A:220:ARG:HA	2:E:78:THR:HB	1.89	0.55
2:K:9[B]:CYS:HB2	2:K:86[B]:CYS:SG	2.47	0.55
2:H:47:THR:H	5:H:202:GOL:H12	1.72	0.54
1:A:228:SER:CB	2:F:74:ILE:HG12	2.37	0.54
1:B:104:TYR:OH	1:B:176:GLU:OE2	2.16	0.53
2:F:51:GLU:OE2	2:F:91:LYS:CE	2.45	0.53
2:G:9[B]:CYS:HB2	2:G:86[B]:CYS:SG	2.49	0.53
2:C:67:ARG:NH1	9:C:302:HOH:O	2.32	0.52
2:H:20:LEU:HD13	2:H:42:PHE:CZ	2.45	0.52
1:A:228:SER:HB3	2:F:74:ILE:HG12	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:74[A]:ILE:HD13	2:J:77:LEU:HD11	1.93	0.51
2:C:20:LEU:HD13	2:C:42:PHE:CZ	2.46	0.51
1:A:228:SER:OG	2:G:74:ILE:CG1	2.59	0.50
1:A:155:ILE:O	9:A:401:HOH:O	2.19	0.50
2:L:20:LEU:HD13	2:L:42:PHE:CZ	2.46	0.50
2:K:63:LYS:HG2	9:K:303:HOH:O	2.11	0.50
2:D:9[B]:CYS:HB2	2:D:86[B]:CYS:SG	2.52	0.49
2:H:9[A]:CYS:HB2	2:H:86[A]:CYS:SG	2.52	0.49
2:C:15:THR:OG1	9:C:301:HOH:O	2.14	0.49
2:H:20:LEU:HD22	2:H:42:PHE:CG	2.47	0.49
1:B:125:TYR:HD1	1:B:145:TYR:CE2	2.31	0.49
2:H:22:ASP:O	2:H:82:VAL:HG13	2.13	0.49
2:K:74:ILE:HD13	2:L:77:LEU:CD1	2.39	0.49
2:D:20:LEU:HD13	2:D:42:PHE:CZ	2.48	0.49
1:B:155:ILE:O	9:B:402:HOH:O	2.20	0.49
2:G:67[B]:ARG:NH2	2:H:69:LYS:O	2.46	0.49
1:A:236:ILE:HD13	2:E:63[B]:LYS:HB2	1.95	0.48
2:F:67:ARG:NH1	9:F:304:HOH:O	2.34	0.48
2:J:9[A]:CYS:HB2	2:J:86[A]:CYS:SG	2.53	0.48
2:D:12:TYR:OH	2:E:35:ARG:HD2	2.13	0.48
1:B:72:VAL:O	1:B:75:THR:OG1	2.28	0.48
2:G:67[B]:ARG:HH21	2:H:70:ASP:HA	1.79	0.48
2:C:20:LEU:HD22	2:C:42:PHE:CG	2.49	0.47
2:D:20:LEU:HD22	2:D:42:PHE:CG	2.50	0.47
1:B:110:GLU:O	1:B:111[B]:GLN:C	2.58	0.47
2:C:9[B]:CYS:HB2	2:C:86[B]:CYS:SG	2.55	0.47
1:B:83:TYR:CZ	1:B:130:VAL:HG11	2.50	0.47
1:A:236:ILE:HD13	2:E:63[A]:LYS:HB2	1.97	0.47
2:K:19:THR:C	2:K:20:LEU:HD22	2.40	0.47
1:B:146[B]:ARG:NH1	1:B:146[B]:ARG:CG	2.59	0.47
2:F:50:VAL:HG11	2:F:69[B]:LYS:HD2	1.98	0.46
2:J:19:THR:C	2:J:20:LEU:HD22	2.40	0.46
2:L:20:LEU:HD22	2:L:42:PHE:CG	2.50	0.46
1:A:148:ARG:HD2	2:F:103:ASN:OXT	2.16	0.46
2:I:74[A]:ILE:HD12	2:I:74[A]:ILE:C	2.41	0.46
2:G:74:ILE:HD13	2:H:77:LEU:CD1	2.41	0.46
2:G:19:THR:C	2:G:20:LEU:HD22	2.40	0.46
2:E:19:THR:C	2:E:20:LEU:HD22	2.41	0.46
2:D:89:ASN:HD22	2:D:89:ASN:C	2.24	0.45
2:F:15[B]:THR:HG23	2:F:16:GLN:N	2.32	0.45
2:F:19:THR:C	2:F:20:LEU:HD22	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:GLN:HG3	1:B:37:MET:SD	2.57	0.45
2:H:8[B]:LEU:HD12	2:H:8[B]:LEU:O	2.17	0.45
2:D:67[B]:ARG:NH2	2:E:69:LYS:O	2.49	0.45
2:I:19:THR:C	2:I:20:LEU:HD22	2.40	0.45
1:B:235[C]:ARG:NH2	9:B:410:HOH:O	2.50	0.45
1:B:48[B]:THR:O	1:B:48[B]:THR:CG2	2.64	0.45
2:K:88:TRP:CH2	4:K:201:GAL:H62	2.52	0.45
2:D:69:LYS:NZ	9:D:305:HOH:O	2.48	0.44
2:G:22:ASP:O	2:G:82:VAL:HG13	2.17	0.44
1:A:227:GLN:NE2	9:A:402:HOH:O	2.27	0.44
1:B:146[B]:ARG:NH2	9:B:404:HOH:O	2.37	0.44
2:K:74:ILE:CD1	2:L:77:LEU:HD13	2.42	0.43
1:B:29:GLU:OE2	2:C:23:LYS:NZ	2.52	0.43
1:B:71[A]:LEU:HD13	1:B:71[A]:LEU:HA	1.87	0.43
1:B:127:TRP:CZ3	1:B:142:ASN:HB2	2.52	0.43
1:A:192:ARG:O	1:A:196:SER:HB2	2.17	0.43
1:A:129:ARG:NH1	1:A:131:HIS:HB2	2.33	0.43
1:A:25[A]:ARG:HD2	1:A:25[A]:ARG:C	2.44	0.43
1:B:214:TYR:O	1:B:218:VAL:HG23	2.18	0.43
1:A:214:TYR:O	1:A:218:VAL:HG23	2.18	0.43
1:A:83:TYR:CZ	1:A:130:VAL:HG11	2.54	0.42
1:B:53:VAL:O	1:B:54:ARG:C	2.62	0.42
1:A:110:GLU:O	1:A:111[A]:GLN:C	2.62	0.42
2:C:25:PHE:CD2	2:C:43[A]:LYS:HG3	2.54	0.42
1:A:127:TRP:CZ3	1:A:142:ASN:HB2	2.54	0.42
2:F:69[B]:LYS:HZ2	2:F:69[B]:LYS:HB2	1.84	0.42
1:B:94[A]:MET:CE	1:B:115:ALA:HB2	2.50	0.42
1:A:128:TYR:CB	1:A:135:LEU:HD11	2.50	0.41
1:B:49[A]:GLN:C	1:B:50:THR:HG1	2.00	0.41
2:D:67[B]:ARG:CZ	2:E:69:LYS:HB3	2.51	0.41
2:H:33:GLY:O	2:H:34[B]:LYS:HB2	2.21	0.41
1:A:172:ARG:HD2	1:A:176:GLU:OE1	2.21	0.41
2:G:67[B]:ARG:NH1	2:H:29:GLU:OE1	2.53	0.40
1:B:5:LEU:HD12	1:B:94[B]:MET:CE	2.51	0.40
1:A:110:GLU:O	1:A:111[B]:GLN:C	2.62	0.40
1:B:228:SER:OG	2:K:74:ILE:CG1	2.69	0.40
1:A:44:HIS:CG	1:A:59:TYR:HB2	2.57	0.40
1:A:185:PRO:HB3	2:I:45:GLY:CA	2.51	0.40
1:B:116:LEU:HD13	1:B:210:PHE:CD1	2.56	0.40
1:B:44:HIS:CG	1:B:59:TYR:HB2	2.56	0.40
1:B:91:ALA:HB1	1:B:92:PRO:HD2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:92:THR:HA	2:C:93:PRO:HA	1.95	0.40
2:K:103:ASN:HB3	2:L:25:PHE:CE1	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	245/240 (102%)	235 (96%)	10 (4%)	0	100	100
1	B	238/240 (99%)	227 (95%)	9 (4%)	2 (1%)	16	20
2	C	105/103 (102%)	103 (98%)	2 (2%)	0	100	100
2	D	106/103 (103%)	104 (98%)	2 (2%)	0	100	100
2	E	106/103 (103%)	103 (97%)	3 (3%)	0	100	100
2	F	107/103 (104%)	104 (97%)	3 (3%)	0	100	100
2	G	104/103 (101%)	101 (97%)	3 (3%)	0	100	100
2	H	107/103 (104%)	104 (97%)	3 (3%)	0	100	100
2	I	109/103 (106%)	105 (96%)	4 (4%)	0	100	100
2	J	103/103 (100%)	101 (98%)	2 (2%)	0	100	100
2	K	104/103 (101%)	101 (97%)	3 (3%)	0	100	100
2	L	104/103 (101%)	101 (97%)	3 (3%)	0	100	100
All	All	1538/1510 (102%)	1489 (97%)	47 (3%)	2 (0%)	48	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	49[A]	GLN
1	B	49[B]	GLN

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	211/204 (103%)	203 (96%)	8 (4%)	29	44
1	B	206/204 (101%)	199 (97%)	7 (3%)	32	49
2	C	93/89 (104%)	92 (99%)	1 (1%)	65	81
2	D	94/89 (106%)	93 (99%)	1 (1%)	65	81
2	E	94/89 (106%)	93 (99%)	1 (1%)	65	81
2	F	95/89 (107%)	95 (100%)	0	100	100
2	G	92/89 (103%)	88 (96%)	4 (4%)	26	39
2	H	95/89 (107%)	94 (99%)	1 (1%)	65	81
2	I	96/89 (108%)	95 (99%)	1 (1%)	68	82
2	J	91/89 (102%)	91 (100%)	0	100	100
2	K	92/89 (103%)	90 (98%)	2 (2%)	45	65
2	L	92/89 (103%)	92 (100%)	0	100	100
All	All	1351/1298 (104%)	1325 (98%)	26 (2%)	50	69

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

Mol	Chain	Res	Type
1	A	14	ASP
1	A	48	THR
1	A	56	ASP
1	A	109	ASP
1	A	152	ASN
1	A	172	ARG
1	A	195	MET
1	A	206	LEU
2	D	89	ASN
2	E	72	LEU
2	G	43	LYS
2	G	74	ILE
2	G	82	VAL
2	G	103	ASN

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Mol	Chain	Res	Type
2	H	82	VAL
1	B	14	ASP
1	B	61	SER
1	B	71[A]	LEU
1	B	71[B]	LEU
1	B	112	GLU
1	B	152	ASN
1	B	238	ASP
2	C	103	ASN
2	I	103	ASN
2	K	74	ILE
2	K	103	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	221	GLN
2	D	16	GLN
2	D	89	ASN
2	E	103	ASN
2	F	49	GLN
2	G	49	GLN
2	H	49	GLN
2	H	103	ASN
1	B	107	HIS
1	B	233	HIS
2	C	3	GLN
2	C	21	ASN
2	C	103	ASN
2	I	103	ASN
2	J	49	GLN
2	K	13	HIS
2	K	49	GLN
2	L	49	GLN

5.3.3 RNA ⓘ

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 33 ligands modelled in this entry, 4 are monoatomic - leaving 29 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	$\# Z > 2$
4	GAL	F	202	-	12,12,12	0.62	0	17,17,17	0.32	0
4	GAL	J	201	-	12,12,12	0.67	0	17,17,17	0.53	0
4	GAL	K	201	-	12,12,12	0.20	0	17,17,17	0.28	0
5	GOL	G	204	-	5,5,5	0.12	0	5,5,5	0.25	0
4	GAL	D	202	-	12,12,12	0.68	0	17,17,17	0.53	0
5	GOL	I	202	-	5,5,5	0.12	0	5,5,5	0.25	0
4	GAL	E	201	-	12,12,12	0.55	0	17,17,17	0.38	0
6	ACT	C	201	-	3,3,3	1.09	0	3,3,3	0.80	0
5	GOL	J	202	-	5,5,5	0.10	0	5,5,5	0.26	0
5	GOL	A	304	-	5,5,5	0.10	0	5,5,5	0.21	0
5	GOL	D	203	-	5,5,5	0.11	0	5,5,5	0.24	0
5	GOL	B	303	-	5,5,5	0.11	0	5,5,5	0.26	0
5	GOL	H	202	-	5,5,5	0.11	0	5,5,5	0.25	0
5	GOL	A	305	-	5,5,5	0.11	0	5,5,5	0.24	0
7	PEG	G	201	-	6,6,6	0.19	0	5,5,5	0.09	0
5	GOL	C	203	-	5,5,5	0.16	0	5,5,5	0.27	0
4	GAL	L	201	-	12,12,12	0.45	0	17,17,17	0.42	0
6	ACT	F	201	-	3,3,3	1.06	0	3,3,3	0.81	0
4	GAL	G	203	-	12,12,12	0.57	0	17,17,17	0.30	0
4	GAL	I	201	-	12,12,12	0.49	0	17,17,17	0.26	0
5	GOL	L	203	-	5,5,5	0.10	0	5,5,5	0.25	0
6	ACT	G	202	-	3,3,3	1.23	0	3,3,3	0.78	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GOL	I	203	-	5,5,5	0.11	0	5,5,5	0.23	0
5	GOL	E	202	-	5,5,5	0.11	0	5,5,5	0.26	0
5	GOL	L	202	-	5,5,5	0.14	0	5,5,5	0.27	0
5	GOL	A	303	-	5,5,5	0.08	0	5,5,5	0.22	0
4	GAL	H	201	-	12,12,12	0.53	0	17,17,17	0.42	0
8	GLA	C	202	-	12,12,12	0.79	0	17,17,17	0.47	0
4	GAL	A	302	-	12,12,12	0.87	1 (8%)	17,17,17	0.46	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
4	GAL	F	202	-	-	0/2/22/22	0/1/1/1
4	GAL	J	201	-	-	0/2/22/22	0/1/1/1
4	GAL	K	201	-	-	2/2/22/22	0/1/1/1
5	GOL	G	204	-	-	0/4/4/4	-
4	GAL	D	202	-	-	0/2/22/22	0/1/1/1
5	GOL	I	202	-	-	4/4/4/4	-
4	GAL	E	201	-	-	0/2/22/22	0/1/1/1
5	GOL	J	202	-	-	2/4/4/4	-
5	GOL	A	304	-	-	2/4/4/4	-
5	GOL	D	203	-	-	4/4/4/4	-
5	GOL	B	303	-	-	4/4/4/4	-
5	GOL	H	202	-	-	4/4/4/4	-
5	GOL	A	305	-	-	0/4/4/4	-
7	PEG	G	201	-	-	1/4/4/4	-
5	GOL	C	203	-	-	2/4/4/4	-
4	GAL	L	201	-	-	0/2/22/22	0/1/1/1
4	GAL	G	203	-	-	0/2/22/22	0/1/1/1
4	GAL	I	201	-	-	0/2/22/22	0/1/1/1
5	GOL	L	203	-	-	2/4/4/4	-
5	GOL	I	203	-	-	2/4/4/4	-
5	GOL	E	202	-	-	2/4/4/4	-
5	GOL	L	202	-	-	4/4/4/4	-
5	GOL	A	303	-	-	2/4/4/4	-
4	GAL	H	201	-	-	1/2/22/22	0/1/1/1
8	GLA	C	202	-	-	1/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GAL	A	302	-	-	2/2/22/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	302	GAL	C4-C5	2.69	1.58	1.53

There are no bond angle outliers.

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	303	GOL	O2-C2-C3-O3
5	A	304	GOL	C1-C2-C3-O3
5	H	202	GOL	C1-C2-C3-O3
5	B	303	GOL	O1-C1-C2-C3
5	B	303	GOL	C1-C2-C3-O3
5	C	203	GOL	O1-C1-C2-C3
5	I	202	GOL	O1-C1-C2-C3
5	J	202	GOL	C1-C2-C3-O3
5	L	202	GOL	O1-C1-C2-C3
4	K	201	GAL	O5-C5-C6-O6
5	A	303	GOL	C1-C2-C3-O3
5	D	203	GOL	O1-C1-C2-C3
5	D	203	GOL	C1-C2-C3-O3
5	E	202	GOL	C1-C2-C3-O3
5	H	202	GOL	O1-C1-C2-C3
5	I	202	GOL	C1-C2-C3-O3
5	I	203	GOL	O1-C1-C2-C3
5	L	202	GOL	C1-C2-C3-O3
5	L	203	GOL	C1-C2-C3-O3
5	A	304	GOL	O2-C2-C3-O3
5	D	203	GOL	O1-C1-C2-O2
5	D	203	GOL	O2-C2-C3-O3
5	H	202	GOL	O2-C2-C3-O3
5	B	303	GOL	O1-C1-C2-O2
5	I	202	GOL	O2-C2-C3-O3
5	I	203	GOL	O1-C1-C2-O2
5	L	202	GOL	O1-C1-C2-O2
5	H	202	GOL	O1-C1-C2-O2
5	B	303	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
5	C	203	GOL	O1-C1-C2-O2
5	J	202	GOL	O2-C2-C3-O3
5	L	202	GOL	O2-C2-C3-O3
4	A	302	GAL	C4-C5-C6-O6
5	I	202	GOL	O1-C1-C2-O2
4	H	201	GAL	O5-C5-C6-O6
5	E	202	GOL	O2-C2-C3-O3
4	A	302	GAL	O5-C5-C6-O6
4	K	201	GAL	C4-C5-C6-O6
7	G	201	PEG	C4-C3-O2-C2
5	L	203	GOL	O2-C2-C3-O3
8	C	202	GLA	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	K	201	GAL	1	0
5	H	202	GOL	1	0
4	A	302	GAL	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	237/240 (98%)	0.25	9 (3%) 44 46	12, 37, 72, 90	10 (4%)
1	B	231/240 (96%)	1.14	41 (17%) 4 4	17, 58, 89, 106	11 (4%)
2	C	103/103 (100%)	0.38	1 (0%) 79 80	20, 42, 58, 84	4 (3%)
2	D	103/103 (100%)	0.10	1 (0%) 79 80	14, 35, 49, 78	5 (4%)
2	E	103/103 (100%)	0.15	1 (0%) 79 80	15, 38, 50, 61	5 (4%)
2	F	103/103 (100%)	0.18	1 (0%) 79 80	15, 35, 48, 79	6 (5%)
2	G	103/103 (100%)	0.06	2 (1%) 66 68	12, 34, 51, 64	3 (2%)
2	H	103/103 (100%)	0.19	1 (0%) 79 80	16, 34, 51, 87	6 (5%)
2	I	103/103 (100%)	0.76	8 (7%) 19 21	16, 49, 67, 80	7 (6%)
2	J	103/103 (100%)	0.38	0 100 100	20, 45, 63, 88	2 (1%)
2	K	103/103 (100%)	0.28	0 100 100	21, 41, 59, 78	3 (2%)
2	L	103/103 (100%)	0.27	0 100 100	17, 40, 56, 97	3 (2%)
All	All	1498/1510 (99%)	0.40	65 (4%) 40 41	12, 40, 73, 106	65 (4%)

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	48[A]	THR	10.9
2	I	10[A]	ALA	4.1
1	B	52	PHE	4.0
1	B	235[A]	ARG	3.9
2	I	102	ALA	3.8
1	B	198	THR	3.8
1	B	71[A]	LEU	3.4
1	B	47[A]	GLY	3.4
1	B	199	CYS	3.3
1	A	237	LYS	3.2
1	B	38[A]	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	185	PRO	3.1
2	D	103	ASN	3.1
2	I	9[A]	CYS	3.1
1	B	132	PHE	3.0
1	B	101	LEU	3.0
1	B	174	TRP	3.0
1	B	186	GLY	3.0
1	A	78	SER	2.9
1	B	167	PHE	2.9
1	B	180	ILE	2.9
1	A	52	PHE	2.8
1	A	140[A]	HIS	2.7
2	H	103	ASN	2.7
2	I	81[A]	LYS	2.7
1	B	45	ALA	2.6
1	B	34	GLY	2.6
2	F	103	ASN	2.6
2	I	43[A]	LYS	2.6
1	A	33[A]	ARG	2.5
1	B	169	PRO	2.5
1	B	179	TRP	2.5
2	G	103	ASN	2.5
1	B	188	GLY	2.5
2	E	29	GLU	2.4
1	B	79	GLY	2.4
1	B	184	PRO	2.4
1	B	172	ARG	2.4
1	A	50	THR	2.3
2	I	13	HIS	2.3
1	B	49[A]	GLN	2.3
1	A	136	ASP	2.3
1	B	140	HIS	2.3
1	B	39	ILE	2.3
2	C	103	ASN	2.3
1	B	58	GLY	2.3
1	B	42	TYR	2.3
1	B	103	ALA	2.3
2	G	1	THR	2.2
1	B	200	ASP	2.2
1	A	111[A]	GLN	2.2
1	B	197	ASN	2.2
1	B	41	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	171	HIS	2.2
1	B	78	SER	2.2
1	B	234	ASN	2.1
1	B	18[A]	GLN	2.1
2	I	54	GLY	2.1
1	A	14	ASP	2.1
1	B	74	GLN	2.1
2	I	20	LEU	2.1
1	B	178	PRO	2.0
1	B	83	TYR	2.0
1	B	50	THR	2.0
1	B	181	HIS	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(Å ²)	Q<0.9
6	ACT	G	202	4/4	0.65	0.27	65,69,75,78	0
3	NA	B	302	1/1	0.70	0.13	75,75,75,75	0
5	GOL	L	203	6/6	0.76	0.16	74,76,82,83	0
6	ACT	C	201	4/4	0.76	0.18	61,65,65,66	0
5	GOL	I	203	6/6	0.80	0.14	60,70,71,73	0
5	GOL	H	202	6/6	0.81	0.16	68,72,73,73	0
5	GOL	C	203	6/6	0.82	0.14	61,69,71,72	0
4	GAL	A	302	12/12	0.83	0.14	57,88,94,99	0
5	GOL	A	303	6/6	0.84	0.15	54,63,65,68	0
5	GOL	A	305	6/6	0.84	0.15	63,67,71,74	0
5	GOL	D	203	6/6	0.85	0.11	51,59,60,63	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	GOL	G	204	6/6	0.85	0.16	60,65,66,70	0
6	ACT	F	201	4/4	0.86	0.19	59,59,62,63	0
5	GOL	A	304	6/6	0.86	0.12	55,58,61,61	0
5	GOL	I	202	6/6	0.86	0.12	66,76,81,84	0
5	GOL	L	202	6/6	0.87	0.16	59,62,65,68	0
5	GOL	J	202	6/6	0.88	0.13	60,64,68,70	0
5	GOL	E	202	6/6	0.88	0.13	56,57,67,77	0
7	PEG	G	201	7/7	0.88	0.15	53,64,67,67	0
5	GOL	B	303	6/6	0.89	0.13	68,74,75,76	0
3	NA	D	201	1/1	0.90	0.09	58,58,58,58	0
4	GAL	G	203	12/12	0.91	0.09	41,43,47,48	0
4	GAL	J	201	12/12	0.91	0.11	40,49,57,74	0
4	GAL	I	201	12/12	0.92	0.09	47,52,58,65	0
4	GAL	H	201	12/12	0.93	0.08	39,45,52,53	0
4	GAL	K	201	12/12	0.93	0.08	47,50,58,68	0
4	GAL	L	201	12/12	0.93	0.08	37,44,46,48	0
8	GLA	C	202	12/12	0.93	0.08	32,39,42,45	0
4	GAL	E	201	12/12	0.94	0.07	35,39,47,47	0
4	GAL	D	202	12/12	0.94	0.07	29,34,40,44	0
4	GAL	F	202	12/12	0.95	0.06	30,33,36,42	0
3	NA	A	301	1/1	0.99	0.03	17,17,17,17	0
3	NA	B	301	1/1	0.99	0.06	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.