



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 03:34 PM GMT

PDB ID : 1CDG
Title : NUCLEOTIDE SEQUENCE AND X-RAY STRUCTURE OF CYCLODEX-
TRIN GLYCOSYLTRANSFERASE FROM BACILLUS CIRCULANS
STRAIN 251 IN A MALTOSE-DEPENDENT CRYSTAL FORM
Authors : Lawson, C.L.; Van Montfort, R.; Strokopytov, B.V.; Kalk, K.H.; Rozeboom,
H.J.; Dijkstra, B.W.
Deposited on : 1993-08-02
Resolution : 2.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

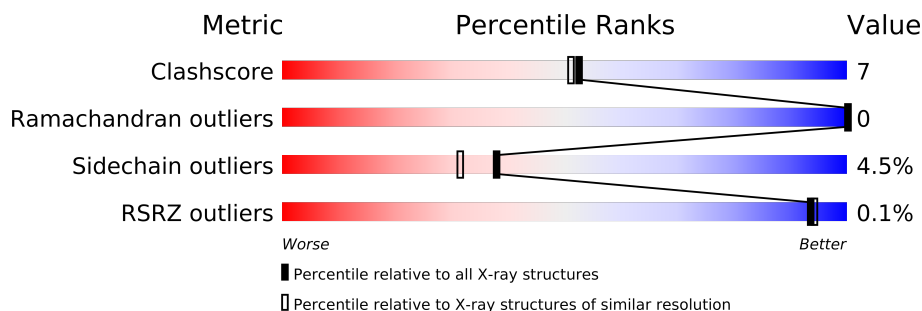
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance


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))
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	686	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	MAL	A	688	-	X
2	MAL	A	689	-	X
2	MAL	A	690	-	X

2 Entry composition i

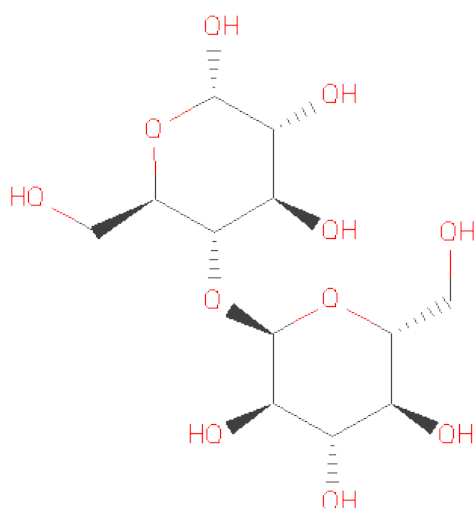
There are 4 unique types of molecules in this entry. The entry contains 5806 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CYCLODEXTRIN GLYCOSYL-TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	686	5264	3321	900	1027	16	0	0	0

- Molecule 2 is SUGAR (MALTOSE) (three-letter code: MAL) (formula: C₁₂H₂₂O₁₁).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	23	12	11	0	0
2	A	1	23	12	11	0	0
2	A	1	23	12	11	0	0

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total 2	Ca 2	0	0

- Molecule 4 is water.

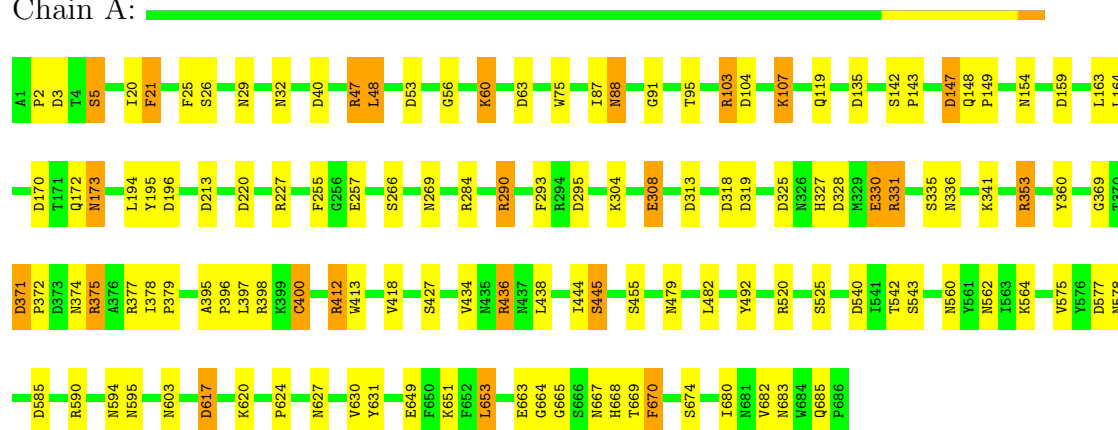
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	471	Total 471	O 471	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: CYCLODEXTRIN GLYCOSYL-TRANSFERASE

Chain A:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	120.36Å 110.94Å 66.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.00 15.03 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00) 80.5 (15.03-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.05 (at 2.00Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.164 , (Not available) 0.157 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	21.0	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 98.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 48567 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5806	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, MAL

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	0.87	0/5394	1.39	64/7352 (0.9%)

There are no bond length outliers.

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	412	ARG	NE-CZ-NH1	13.48	127.04	120.30
1	A	375	ARG	NE-CZ-NH1	12.84	126.72	120.30
1	A	331	ARG	NE-CZ-NH1	11.98	126.29	120.30
1	A	331	ARG	NE-CZ-NH2	-11.18	114.71	120.30
1	A	577	ASP	CB-CG-OD1	10.76	127.98	118.30
1	A	375	ARG	NE-CZ-NH2	-10.54	115.03	120.30
1	A	213	ASP	CB-CG-OD2	-10.07	109.24	118.30
1	A	328	ASP	CB-CG-OD2	-9.61	109.65	118.30
1	A	284	ARG	NE-CZ-NH2	-9.18	115.71	120.30
1	A	590	ARG	NE-CZ-NH1	8.86	124.73	120.30
1	A	398	ARG	NE-CZ-NH1	8.83	124.72	120.30
1	A	104	ASP	CB-CG-OD2	-8.41	110.73	118.30
1	A	353	ARG	NE-CZ-NH1	8.33	124.46	120.30
1	A	147	ASP	CB-CG-OD2	-8.15	110.97	118.30
1	A	577	ASP	CB-CG-OD2	-8.07	111.03	118.30
1	A	585	ASP	CB-CG-OD1	8.02	125.51	118.30
1	A	412	ARG	NE-CZ-NH2	-7.98	116.31	120.30
1	A	213	ASP	CB-CG-OD1	7.81	125.33	118.30
1	A	53	ASP	CB-CG-OD2	-7.81	111.27	118.30
1	A	159	ASP	CB-CG-OD2	-7.61	111.45	118.30
1	A	585	ASP	CB-CG-OD2	-7.59	111.47	118.30
1	A	436	ARG	NE-CZ-NH2	-7.58	116.51	120.30
1	A	353	ARG	NE-CZ-NH2	-7.46	116.57	120.30
1	A	135	ASP	CB-CG-OD1	7.15	124.74	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	159	ASP	CB-CG-OD1	7.15	124.74	118.30
1	A	543	SER	N-CA-CB	7.12	121.17	110.50
1	A	147	ASP	CB-CG-OD1	6.99	124.59	118.30
1	A	325	ASP	CB-CG-OD2	-6.92	112.08	118.30
1	A	540	ASP	CB-CG-OD2	-6.77	112.20	118.30
1	A	47	ARG	NE-CZ-NH2	-6.76	116.92	120.30
1	A	436	ARG	NE-CZ-NH1	6.72	123.66	120.30
1	A	540	ASP	CB-CG-OD1	6.71	124.33	118.30
1	A	328	ASP	CB-CG-OD1	6.67	124.31	118.30
1	A	103	ARG	NE-CZ-NH1	6.37	123.48	120.30
1	A	377	ARG	NE-CZ-NH2	-6.35	117.12	120.30
1	A	617	ASP	CB-CG-OD2	-6.33	112.60	118.30
1	A	104	ASP	CB-CG-OD1	6.32	123.99	118.30
1	A	170	ASP	CB-CG-OD2	-6.24	112.69	118.30
1	A	196	ASP	CB-CG-OD1	6.14	123.83	118.30
1	A	135	ASP	CB-CG-OD2	-6.13	112.78	118.30
1	A	319	ASP	CB-CG-OD2	-6.12	112.79	118.30
1	A	325	ASP	CB-CG-OD1	6.05	123.75	118.30
1	A	5	SER	N-CA-CB	-5.94	101.59	110.50
1	A	400	CYS	CA-CB-SG	-5.89	103.39	114.00
1	A	40	ASP	CB-CG-OD2	-5.86	113.02	118.30
1	A	525	SER	N-CA-CB	-5.85	101.73	110.50
1	A	3	ASP	CB-CG-OD1	5.80	123.52	118.30
1	A	40	ASP	CB-CG-OD1	5.66	123.40	118.30
1	A	520	ARG	NE-CZ-NH2	-5.59	117.50	120.30
1	A	266	SER	N-CA-CB	-5.57	102.15	110.50
1	A	492	TYR	CB-CG-CD1	-5.57	117.66	121.00
1	A	313	ASP	CB-CG-OD1	5.53	123.28	118.30
1	A	53	ASP	CB-CG-OD1	5.49	123.24	118.30
1	A	398	ARG	NE-CZ-NH2	-5.45	117.58	120.30
1	A	371	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	A	295	ASP	CB-CG-OD1	5.39	123.15	118.30
1	A	3	ASP	CB-CG-OD2	-5.34	113.49	118.30
1	A	220	ASP	CB-CG-OD1	5.32	123.09	118.30
1	A	445	SER	N-CA-CB	5.30	118.45	110.50
1	A	377	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	A	375	ARG	CD-NE-CZ	5.21	130.89	123.60
1	A	220	ASP	CB-CG-OD2	-5.08	113.73	118.30
1	A	617	ASP	CB-CG-OD1	5.02	122.82	118.30
1	A	318	ASP	CB-CG-OD2	-5.00	113.80	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5264	0	5025	71	0
2	A	69	0	66	5	0
3	A	2	0	0	0	0
4	A	471	0	0	8	0
All	All	5806	0	5091	74	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (74) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:88:ASN:HD21	1:A:91:GLY:H	1.10	0.91
1:A:560:ASN:HD21	1:A:578:ASN:HA	1.37	0.87
1:A:620:LYS:HE3	4:A:937:HOH:O	1.84	0.77
1:A:88:ASN:HD21	1:A:91:GLY:N	1.83	0.75
1:A:336:ASN:OD1	1:A:336:ASN:N	2.25	0.69
1:A:562:ASN:HB3	1:A:575:VAL:HG13	1.76	0.67
1:A:617:ASP:OD2	1:A:620:LYS:HE2	1.95	0.66
1:A:47:ARG:HD2	4:A:908:HOH:O	1.95	0.65
1:A:653:LEU:N	1:A:653:LEU:HD23	2.13	0.64
1:A:293:PHE:O	1:A:341:LYS:HD2	1.98	0.64
1:A:142:SER:HB2	1:A:143:PRO:HD2	1.81	0.63
2:A:690:MAL:H6'2	2:A:690:MAL:C5	2.29	0.63
1:A:142:SER:HB2	1:A:143:PRO:CD	2.30	0.60
2:A:690:MAL:H6'2	2:A:690:MAL:O5	2.01	0.60
1:A:664:GLY:N	1:A:685:GLN:O	2.32	0.59
1:A:395:ALA:HB3	1:A:396:PRO:HD3	1.83	0.59
1:A:374:ASN:N	1:A:374:ASN:OD1	2.35	0.59
1:A:445:SER:HB3	1:A:479:ASN:ND2	2.18	0.59
1:A:670:PHE:CE2	1:A:680:ILE:HD11	2.39	0.57
1:A:669:THR:HG22	1:A:670:PHE:N	2.20	0.57
1:A:674:SER:HB2	4:A:1126:HOH:O	2.05	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:26:SER:O	1:A:56:GLY:HA3	2.07	0.55
1:A:227:ARG:HG2	1:A:255:PHE:CE2	2.44	0.53
1:A:107:LYS:NZ	4:A:776:HOH:O	2.41	0.53
1:A:227:ARG:HH11	1:A:257:GLU:HB2	1.74	0.52
1:A:562:ASN:HB3	1:A:575:VAL:CG1	2.40	0.51
1:A:87:ILE:HD12	1:A:143:PRO:HG2	1.92	0.51
1:A:649:GLU:OE1	1:A:667:ASN:ND2	2.28	0.51
1:A:444:ILE:HD12	1:A:482:LEU:HB2	1.93	0.50
1:A:479:ASN:ND2	4:A:856:HOH:O	2.44	0.50
1:A:327:HIS:ND1	1:A:327:HIS:N	2.57	0.50
1:A:148:GLN:N	1:A:149:PRO:HD3	2.26	0.50
1:A:48:LEU:HD12	1:A:95:THR:HG23	1.93	0.50
1:A:2:PRO:HD2	1:A:5:SER:HB2	1.95	0.49
1:A:651:LYS:NZ	1:A:665:GLY:O	2.46	0.49
1:A:20:ILE:HG23	1:A:25:PHE:HB2	1.96	0.48
1:A:594:ASN:HB2	1:A:683:ASN:OD1	2.14	0.48
1:A:444:ILE:HD12	1:A:482:LEU:CB	2.44	0.48
1:A:603:ASN:HB3	1:A:624:PRO:HB3	1.96	0.48
1:A:542:THR:HG21	4:A:892:HOH:O	2.14	0.47
1:A:669:THR:CG2	1:A:670:PHE:N	2.78	0.47
1:A:397:LEU:HA	1:A:400:CYS:SG	2.54	0.47
2:A:690:MAL:H6'2	2:A:690:MAL:H5	1.95	0.47
1:A:304:LYS:O	1:A:308:GLU:HG2	2.15	0.47
1:A:60:LYS:NZ	1:A:63:ASP:OD2	2.45	0.47
1:A:594:ASN:ND2	1:A:682:VAL:O	2.48	0.46
1:A:330:GLU:HB3	1:A:369:GLY:HA2	1.96	0.46
1:A:627:ASN:HD21	2:A:690:MAL:H61	1.80	0.46
1:A:29:ASN:ND2	1:A:32:ASN:ND2	2.64	0.46
1:A:564:LYS:HE3	1:A:564:LYS:HB3	1.82	0.46
1:A:680:ILE:HG21	1:A:680:ILE:HD13	1.74	0.45
1:A:630:VAL:HG12	1:A:631:TYR:CE1	2.51	0.45
1:A:331:ARG:NH2	1:A:360:TYR:O	2.35	0.45
1:A:119:GLN:HG2	4:A:883:HOH:O	2.16	0.45
1:A:445:SER:CB	1:A:479:ASN:ND2	2.80	0.44
1:A:664:GLY:O	1:A:685:GLN:HB2	2.17	0.44
1:A:371:ASP:OD1	1:A:372:PRO:HA	2.18	0.44
1:A:149:PRO:HA	1:A:154:ASN:ND2	2.33	0.44
1:A:119:GLN:HB2	1:A:119:GLN:HE21	1.54	0.44
1:A:413:TRP:CE2	2:A:688:MAL:H2'	2.52	0.44
1:A:87:ILE:CD1	1:A:143:PRO:HG2	2.47	0.43
1:A:395:ALA:N	1:A:396:PRO:CD	2.80	0.43
1:A:378:ILE:HA	1:A:379:PRO:HD3	1.89	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:194:LEU:O	1:A:195:TYR:HB2	2.19	0.43
1:A:670:PHE:CZ	1:A:680:ILE:HD11	2.54	0.42
1:A:227:ARG:C	1:A:227:ARG:HD2	2.40	0.42
1:A:663:GLU:CD	1:A:668:HIS:HE2	2.21	0.41
1:A:21:PHE:CE1	1:A:327:HIS:HB3	2.55	0.41
1:A:418:VAL:HA	1:A:434:VAL:O	2.21	0.41
1:A:172:GLN:O	1:A:173:ASN:HB2	2.21	0.41
1:A:331:ARG:HG2	4:A:809:HOH:O	2.20	0.41
1:A:378:ILE:HG13	1:A:379:PRO:HD2	2.03	0.41
1:A:164:LEU:HD23	1:A:164:LEU:HA	1.88	0.40
1:A:290:ARG:NH2	1:A:330:GLU:O	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	684/686 (100%)	655 (96%)	29 (4%)	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	556/556 (100%)	531 (96%)	25 (4%)	38	31

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	PHE
1	A	48	LEU
1	A	60	LYS
1	A	75	TRP
1	A	88	ASN
1	A	103	ARG
1	A	107	LYS
1	A	147	ASP
1	A	163	LEU
1	A	173	ASN
1	A	269	ASN
1	A	290	ARG
1	A	308	GLU
1	A	330	GLU
1	A	335	SER
1	A	353	ARG
1	A	375	ARG
1	A	412	ARG
1	A	427	SER
1	A	436	ARG
1	A	438	LEU
1	A	455	SER
1	A	595	ASN
1	A	653	LEU
1	A	670	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	88	ASN
1	A	119	GLN
1	A	120	ASN
1	A	269	ASN
1	A	296	ASN
1	A	410	GLN
1	A	416	ASN
1	A	479	ASN
1	A	560	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	$\# Z > 2$
2	MAL	A	688	-	24,24,24	0.37	0	35,35,35	0.95	1 (2%)
2	MAL	A	689	-	24,24,24	0.28	0	35,35,35	1.21	2 (5%)
2	MAL	A	690	-	24,24,24	0.39	0	35,35,35	1.74	5 (14%)

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	MAL	A	688	-	-	0/8/48/48	0/2/2/2
2	MAL	A	689	-	-	0/8/48/48	0/2/2/2
2	MAL	A	690	-	-	0/8/48/48	0/2/2/2

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	690	MAL	O5-C1-C2	4.59	119.71	110.31
2	A	689	MAL	O5'-C1'-C2'	3.91	115.91	109.86
2	A	690	MAL	O5'-C1'-C2'	3.70	115.60	109.86
2	A	690	MAL	O1-C1-C2	-3.64	99.40	108.12
2	A	690	MAL	C1-O5-C5	3.61	120.74	113.73
2	A	690	MAL	O1-C4'-C3'	-3.39	98.46	107.16
2	A	689	MAL	C1'-C2'-C3'	2.68	114.79	110.53
2	A	688	MAL	O2-C2-C1	2.17	114.77	110.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	686/686 (100%)	-0.83	0 100 100	12, 23, 40, 76	0

There are no RSRZ outliers to report.

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MAL	A	689	23/23	0.24	8.63	60,65,76,79	0
2	MAL	A	690	23/23	0.29	5.40	77,81,87,90	0
2	MAL	A	688	23/23	0.10	2.45	25,36,38,40	0
3	CA	A	692	1/1	0.04	-2.04	21,21,21,21	0
3	CA	A	691	1/1	0.02	-4.04	26,26,26,26	0

6.5 Other polymers ⓘ

There are no such residues in this entry.