



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 12:32 PM GMT

PDB ID : 1DTU
Title : BACILLUS CIRCULANS STRAIN 251 CYCLODEXTRIN GLYCOSYL-TRANSFERASE: A MUTANT Y89D/S146P COMPLEXED TO AN HEX-ASACCHARIDE INHIBITOR
Authors : Uitdehaag, J.C.M.; Kalk, K.H.; Dijkstra, B.W.
Deposited on : 2000-01-13
Resolution : 2.40 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>

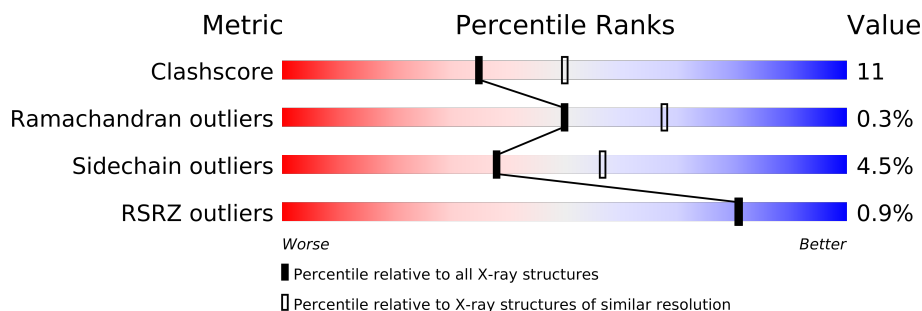
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.40 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	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5722 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 PROTEIN (CYCLODEXTRIN GLYCOSYLTRANSFERASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	686	Total	C	N	O	S	0	1	0
			5268	3322	903	1027	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	89	ASP	TYR	ENGINEERED	UNP P43379
A	146	PRO	SER	ENGINEERED	UNP P43379

- Molecule 2 is a polymer of unknown type called SUGAR (GLC-GLC).

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

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	89	ASP	TYR	ENGINEERED	UNP P43379
A	146	PRO	SER	ENGINEERED	UNP P43379

- Molecule 3 is a polymer of unknown type called SUGAR (GLD-GLC).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	2	Total	C	O	0	0
			21	12	9		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	89	ASP	TYR	ENGINEERED	UNP P43379

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Chain	Residue	Modelled	Actual	Comment	Reference
A	146	PRO	SER	ENGINEERED	UNP P43379

- Molecule 4 is a polymer of unknown type called SUGAR (GLC-GLC-GLC-ACI).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	3	Total	C	O	0	0
			34	18	16		
4	A	3	Total	C	O	0	0
			34	18	16		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	89	ASP	TYR	ENGINEERED	UNP P43379
A	146	PRO	SER	ENGINEERED	UNP P43379
A	89	ASP	TYR	ENGINEERED	UNP P43379
A	146	PRO	SER	ENGINEERED	UNP P43379

- Molecule 5 is a polymer of unknown type called SUGAR (GLC-GLC-GLC-GLC-GLC).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	5	Total	C	O	0	0
			56	30	26		

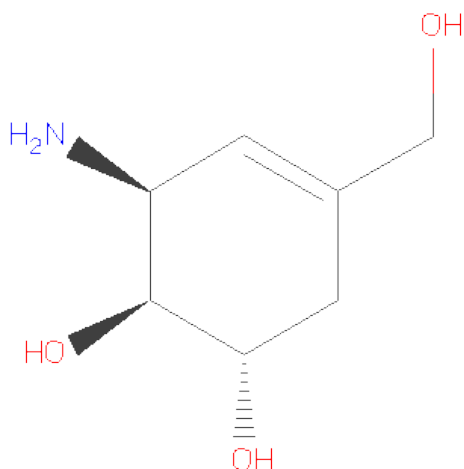
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	89	ASP	TYR	ENGINEERED	UNP P43379
A	146	PRO	SER	ENGINEERED	UNP P43379

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 7 is 1-AMINO-2,3-DIHYDROXY-5-HYDROXYMETHYLCYCLOHEX-5-ENE (three-letter code: ADH) (formula: C₇H₁₃NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			11	7	1	3		

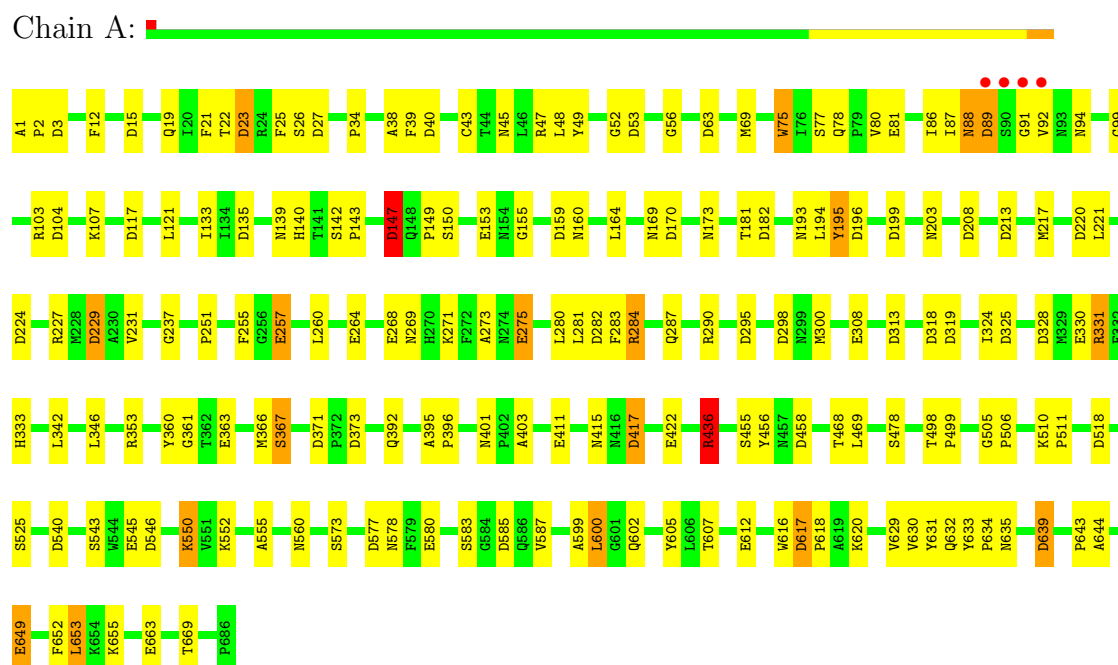
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	273	Total	O	0	0
			273	273		

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: PROTEIN (CYCLODEXTRIN GLYCOSYLTRANSFERASE)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	120.44Å 111.16Å 65.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.40 28.86 – 2.40	Depositor EDS
% Data completeness (in resolution range)	92.7 (8.00-2.40) 88.5 (28.86-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.51 (at 2.39Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.206 , 0.248 0.174 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	2 of 32700 reflections (0.006%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5722	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: GLC, CA, BGC, ADH, G6D

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	0.60	16/5402 (0.3%)	0.94	76/7363 (1.0%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	308	GLU	CD-OE2	5.41	1.31	1.25
1	A	275	GLU	CD-OE2	5.40	1.31	1.25
1	A	153	GLU	CD-OE2	5.36	1.31	1.25
1	A	580	GLU	CD-OE2	5.32	1.31	1.25
1	A	330	GLU	CD-OE2	5.32	1.31	1.25
1	A	545	GLU	CD-OE2	5.31	1.31	1.25
1	A	268	GLU	CD-OE2	5.28	1.31	1.25
1	A	264	GLU	CD-OE2	5.26	1.31	1.25
1	A	422	GLU	CD-OE2	5.25	1.31	1.25
1	A	612	GLU	CD-OE2	5.18	1.31	1.25
1	A	257	GLU	CD-OE2	5.16	1.31	1.25
1	A	363	GLU	CD-OE2	5.16	1.31	1.25
1	A	649	GLU	CD-OE2	5.15	1.31	1.25
1	A	411	GLU	CD-OE2	5.07	1.31	1.25
1	A	81	GLU	CD-OE2	5.05	1.31	1.25
1	A	663	GLU	CD-OE2	5.05	1.31	1.25

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	328	ASP	CB-CG-OD2	-7.27	111.76	118.30
1	A	371	ASP	CB-CG-OD2	-6.75	112.22	118.30
1	A	147	ASP	CB-CG-OD2	-6.75	112.23	118.30
1	A	104	ASP	CB-CG-OD2	-6.64	112.32	118.30
1	A	328	ASP	CB-CG-OD1	6.61	124.25	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	63	ASP	CB-CG-OD2	-6.58	112.38	118.30
1	A	585	ASP	CB-CG-OD2	-6.54	112.42	118.30
1	A	318	ASP	CB-CG-OD2	-6.52	112.43	118.30
1	A	104	ASP	CB-CG-OD1	6.52	124.17	118.30
1	A	458	ASP	CB-CG-OD2	-6.51	112.44	118.30
1	A	213	ASP	CB-CG-OD2	-6.47	112.47	118.30
1	A	298	ASP	CB-CG-OD2	-6.33	112.60	118.30
1	A	224	ASP	CB-CG-OD2	-6.27	112.66	118.30
1	A	135	ASP	CB-CG-OD2	-6.23	112.70	118.30
1	A	135	ASP	CB-CG-OD1	6.17	123.86	118.30
1	A	617	ASP	CB-CG-OD2	-6.17	112.75	118.30
1	A	373	ASP	CB-CG-OD2	-6.15	112.76	118.30
1	A	117	ASP	CB-CG-OD2	-6.14	112.77	118.30
1	A	229	ASP	CB-CG-OD2	-6.13	112.78	118.30
1	A	170	ASP	CB-CG-OD2	-6.11	112.80	118.30
1	A	159	ASP	CB-CG-OD2	-6.09	112.82	118.30
1	A	199	ASP	CB-CG-OD2	-6.06	112.84	118.30
1	A	546	ASP	CB-CG-OD2	-6.05	112.85	118.30
1	A	518	ASP	CB-CG-OD2	-6.05	112.86	118.30
1	A	639	ASP	CB-CG-OD2	-6.04	112.87	118.30
1	A	298	ASP	CB-CG-OD1	6.03	123.72	118.30
1	A	325	ASP	CB-CG-OD2	-5.99	112.91	118.30
1	A	89	ASP	CB-CG-OD2	-5.97	112.92	118.30
1	A	319	ASP	CB-CG-OD2	-5.94	112.95	118.30
1	A	27	ASP	CB-CG-OD2	-5.92	112.97	118.30
1	A	282	ASP	CB-CG-OD2	-5.91	112.98	118.30
1	A	23	ASP	CB-CG-OD2	-5.90	112.99	118.30
1	A	585	ASP	CB-CG-OD1	5.90	123.61	118.30
1	A	63	ASP	CB-CG-OD1	5.89	123.60	118.30
1	A	313	ASP	CB-CG-OD2	-5.88	113.01	118.30
1	A	213	ASP	CB-CG-OD1	5.84	123.56	118.30
1	A	318	ASP	CB-CG-OD1	5.82	123.54	118.30
1	A	3	ASP	CB-CG-OD1	5.78	123.50	118.30
1	A	3	ASP	CB-CG-OD2	-5.78	113.10	118.30
1	A	199	ASP	CB-CG-OD1	5.75	123.48	118.30
1	A	639	ASP	CB-CG-OD1	5.71	123.44	118.30
1	A	325	ASP	CB-CG-OD1	5.70	123.43	118.30
1	A	208	ASP	CB-CG-OD2	-5.69	113.18	118.30
1	A	220	ASP	CB-CG-OD2	-5.68	113.19	118.30
1	A	159	ASP	CB-CG-OD1	5.66	123.39	118.30
1	A	27	ASP	CB-CG-OD1	5.63	123.36	118.30
1	A	313	ASP	CB-CG-OD1	5.62	123.36	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	182	ASP	CB-CG-OD2	-5.56	113.30	118.30
1	A	295	ASP	CB-CG-OD2	-5.54	113.32	118.30
1	A	617	ASP	CB-CG-OD1	5.53	123.28	118.30
1	A	540	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	A	540	ASP	CB-CG-OD1	5.52	123.27	118.30
1	A	40	ASP	CB-CG-OD1	5.51	123.26	118.30
1	A	147	ASP	CB-CG-OD1	5.51	123.25	118.30
1	A	208	ASP	CB-CG-OD1	5.48	123.23	118.30
1	A	371	ASP	CB-CG-OD1	5.48	123.23	118.30
1	A	458	ASP	CB-CG-OD1	5.48	123.23	118.30
1	A	53	ASP	CB-CG-OD2	-5.48	113.37	118.30
1	A	117	ASP	CB-CG-OD1	5.48	123.23	118.30
1	A	40	ASP	CB-CG-OD2	-5.46	113.39	118.30
1	A	436	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	A	295	ASP	CB-CG-OD1	5.45	123.20	118.30
1	A	417	ASP	CB-CG-OD2	-5.44	113.41	118.30
1	A	319	ASP	CB-CG-OD1	5.42	123.18	118.30
1	A	229	ASP	CB-CG-OD1	5.40	123.16	118.30
1	A	15	ASP	CB-CG-OD2	-5.34	113.50	118.30
1	A	282	ASP	CB-CG-OD1	5.31	123.08	118.30
1	A	518	ASP	CB-CG-OD1	5.30	123.07	118.30
1	A	89	ASP	CB-CG-OD1	5.28	123.06	118.30
1	A	577	ASP	CB-CG-OD2	-5.27	113.56	118.30
1	A	224	ASP	CB-CG-OD1	5.26	123.04	118.30
1	A	546	ASP	CB-CG-OD1	5.25	123.03	118.30
1	A	373	ASP	CB-CG-OD1	5.25	123.02	118.30
1	A	23	ASP	CB-CG-OD1	5.24	123.02	118.30
1	A	170	ASP	CB-CG-OD1	5.17	122.95	118.30
1	A	417	ASP	CB-CG-OD1	5.02	122.82	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	5268	0	5031	100	0
2	A	23	0	21	0	0
3	A	21	0	19	5	0
4	A	68	0	59	5	0
5	A	56	0	48	5	0
6	A	2	0	0	0	0
7	A	11	0	12	6	0
8	A	273	0	0	4	0
All	All	5722	0	5190	113	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 11.

All (113) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:692:G6D:C4A	7:A:693:ADH:HN22	1.17	1.50
3:A:692:G6D:C3A	7:A:693:ADH:HN22	1.79	0.94
1:A:617:ASP:HB3	1:A:620:LYS:HE2	1.52	0.91
1:A:75:TRP:CZ2	1:A:227:ARG:HD3	2.07	0.90
1:A:139:ASN:HD22	1:A:140:HIS:HD2	1.20	0.89
1:A:227:ARG:HG3	1:A:255:PHE:CE2	2.08	0.88
1:A:395:ALA:HB3	1:A:396:PRO:HD3	1.64	0.79
1:A:89:ASP:O	1:A:92:VAL:HG22	1.84	0.76
4:A:695:GLC:H61	4:A:696:GLC:H5	1.72	0.72
1:A:607:THR:HG22	1:A:653:LEU:HD21	1.73	0.71
1:A:147:ASP:O	1:A:149:PRO:HD3	1.92	0.69
4:A:695:GLC:C6	4:A:696:GLC:H5	2.21	0.69
1:A:560:ASN:HD21	1:A:578:ASN:HA	1.57	0.68
1:A:333:HIS:O	1:A:367:SER:HB3	1.94	0.68
1:A:88:ASN:HD21	1:A:91:GLY:H	1.39	0.67
1:A:630:VAL:HG12	1:A:631:TYR:CD1	2.31	0.66
1:A:498:THR:CG2	1:A:499:PRO:HD2	2.25	0.66
5:A:701:GLC:H62	5:A:702:GLC:H1	1.77	0.66
1:A:498:THR:HG22	1:A:499:PRO:HD2	1.77	0.65
1:A:633:TYR:HA	1:A:635:ASN:N	2.11	0.65
4:A:695:GLC:H61	4:A:696:GLC:C5	2.27	0.64
1:A:632:GLN:O	1:A:635:ASN:HB2	1.98	0.64
1:A:260:LEU:HB2	1:A:283:PHE:HB3	1.80	0.63
1:A:38:ALA:HB2	1:A:86:ILE:HD11	1.80	0.63
1:A:271:LYS:O	1:A:275:GLU:HG3	1.99	0.63
1:A:194:LEU:O	1:A:195:TYR:HB2	1.99	0.62
1:A:251:PRO:HB3	1:A:506:PRO:HG3	1.82	0.61
1:A:607:THR:CG2	1:A:653:LEU:HD21	2.31	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:617:ASP:CB	1:A:620:LYS:HE2	2.31	0.58
1:A:227:ARG:HD2	1:A:281:LEU:HD11	1.85	0.58
1:A:47[A]:ARG:HD3	1:A:94:ASN:OD1	2.04	0.58
1:A:34:PRO:HG2	1:A:49:TYR:CG	2.39	0.58
3:A:692:G6D:C4A	7:A:693:ADH:C7	2.82	0.57
1:A:88:ASN:ND2	1:A:91:GLY:H	2.03	0.57
3:A:692:G6D:C4A	7:A:693:ADH:C1	2.78	0.56
1:A:142:SER:HB2	1:A:143:PRO:HD2	1.86	0.56
1:A:633:TYR:CD2	1:A:634:PRO:HA	2.40	0.56
1:A:139:ASN:HD22	1:A:140:HIS:CD2	2.12	0.56
1:A:633:TYR:CG	1:A:634:PRO:HA	2.41	0.56
1:A:34:PRO:HG2	1:A:49:TYR:CB	2.37	0.55
1:A:633:TYR:HA	1:A:635:ASN:H	1.71	0.55
1:A:587:VAL:HG13	1:A:644:ALA:HB2	1.88	0.55
1:A:139:ASN:ND2	1:A:140:HIS:HD2	1.98	0.55
1:A:543:SER:HB3	1:A:550:LYS:HB2	1.89	0.55
1:A:231:VAL:HG22	1:A:257:GLU:O	2.06	0.54
1:A:45:ASN:HD22	1:A:48:LEU:HD11	1.73	0.53
1:A:1:ALA:HB1	1:A:2:PRO:HD2	1.91	0.53
1:A:633:TYR:CA	1:A:635:ASN:H	2.22	0.53
1:A:630:VAL:HG12	1:A:631:TYR:CE1	2.44	0.52
1:A:121:LEU:HD12	1:A:121:LEU:O	2.11	0.51
1:A:499:PRO:HB2	1:A:573:SER:HB3	1.92	0.51
1:A:45:ASN:ND2	1:A:48:LEU:HD11	2.26	0.51
1:A:616:TRP:O	1:A:618:PRO:HD3	2.11	0.51
1:A:284:ARG:HD3	8:A:837:HOH:O	2.11	0.50
1:A:92:VAL:HG23	1:A:94:ASN:HD21	1.76	0.50
1:A:456:TYR:O	1:A:468:THR:HG23	2.11	0.50
1:A:26:SER:O	1:A:56:GLY:HA3	2.11	0.50
1:A:652:PHE:C	1:A:653:LEU:HD23	2.32	0.50
1:A:505:GLY:HA2	1:A:506:PRO:C	2.30	0.50
1:A:617:ASP:HB3	1:A:620:LYS:CE	2.35	0.49
1:A:80:VAL:HA	1:A:107:LYS:O	2.12	0.49
1:A:12:PHE:CE2	1:A:133:ILE:HD11	2.47	0.49
1:A:510:LYS:HB2	1:A:511:PRO:HD2	1.95	0.49
1:A:649:GLU:HG2	1:A:669:THR:HG22	1.93	0.49
1:A:290:ARG:HD2	1:A:324:ILE:O	2.13	0.48
1:A:583:SER:HB2	1:A:643:PRO:HG3	1.95	0.48
1:A:395:ALA:HB3	1:A:396:PRO:CD	2.37	0.48
1:A:417:ASP:OD1	1:A:436:ARG:HD2	2.13	0.48
1:A:34:PRO:O	1:A:39:PHE:HB2	2.14	0.48
1:A:19:GLN:O	1:A:360:TYR:HB3	2.14	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:300:MET:HB2	1:A:415:ASN:O	2.14	0.47
1:A:599:ALA:O	1:A:602:GLN:HB3	2.14	0.47
5:A:701:GLC:H62	5:A:702:GLC:C1	2.42	0.47
1:A:284:ARG:HB2	8:A:720:HOH:O	2.14	0.47
1:A:78:GLN:HB3	1:A:99:GLY:O	2.14	0.47
1:A:361:GLY:HA3	1:A:366:MET:SD	2.54	0.47
1:A:229:ASP:OD1	7:A:693:ADH:H1	2.14	0.47
1:A:417:ASP:O	1:A:436:ARG:HG3	2.15	0.47
1:A:237:GLY:HA3	1:A:639:ASP:O	2.14	0.47
3:A:692:G6D:C5A	7:A:693:ADH:N1	2.68	0.47
1:A:605:TYR:CE1	1:A:655:LYS:HB2	2.50	0.47
1:A:23:ASP:HA	1:A:52:GLY:HA3	1.97	0.46
1:A:401:ASN:OD1	1:A:403:ALA:HB3	2.16	0.46
1:A:195:TYR:HE2	4:A:696:GLC:HO4	1.62	0.46
1:A:333:HIS:HB2	1:A:342:LEU:HD13	1.98	0.46
1:A:203:ASN:HA	8:A:942:HOH:O	2.16	0.45
1:A:25:PHE:HB3	8:A:742:HOH:O	2.16	0.45
5:A:698:BGC:H6C1	5:A:699:GLC:O5	2.16	0.45
1:A:142:SER:HB2	1:A:143:PRO:CD	2.46	0.45
1:A:69:MET:O	1:A:392:GLN:HA	2.17	0.45
1:A:260:LEU:HB2	1:A:283:PHE:CB	2.44	0.44
1:A:468:THR:HG22	1:A:469:LEU:N	2.32	0.44
4:A:695:GLC:H62	4:A:696:GLC:H5	1.99	0.44
1:A:92:VAL:O	1:A:92:VAL:HG23	2.18	0.43
1:A:395:ALA:N	1:A:396:PRO:HD2	2.33	0.43
1:A:87:ILE:CD1	1:A:143:PRO:HG2	2.47	0.43
1:A:600:LEU:HD22	5:A:700:GLC:H3	2.00	0.43
1:A:511:PRO:HB3	1:A:555:ALA:HA	2.01	0.43
1:A:143:PRO:HB3	1:A:196:ASP:OD1	2.19	0.43
1:A:273:ALA:HB2	1:A:280:LEU:HD12	2.00	0.42
1:A:164:LEU:HD23	1:A:164:LEU:HA	1.85	0.42
1:A:22:THR:O	1:A:52:GLY:HA3	2.20	0.42
1:A:181:THR:HG22	1:A:193:ASN:O	2.20	0.42
1:A:498:THR:HG22	1:A:499:PRO:CD	2.47	0.41
1:A:231:VAL:HG22	1:A:257:GLU:C	2.41	0.41
1:A:560:ASN:ND2	1:A:578:ASN:HA	2.30	0.41
1:A:78:GLN:HG2	1:A:80:VAL:HG22	2.03	0.41
5:A:698:BGC:C6	5:A:699:GLC:C1	2.98	0.41
1:A:221:LEU:HA	1:A:221:LEU:HD23	1.99	0.40
1:A:283:PHE:O	1:A:287:GLN:HG2	2.22	0.40
1:A:331:ARG:HB3	1:A:331:ARG:HE	1.72	0.40
1:A:142:SER:OG	1:A:155:GLY:HA2	2.22	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:346:LEU:HA	1:A:346:LEU:HD23	1.71	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	685/686 (100%)	652 (95%)	31 (4%)	2 (0%)	50 68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	629	VAL
1	A	195	TYR

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	557/556 (100%)	532 (96%)	25 (4%)	38 57

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

Mol	Chain	Res	Type
1	A	21	PHE
1	A	43	CYS
1	A	75	TRP
1	A	77	SER
1	A	88	ASN

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Mol	Chain	Res	Type
1	A	103	ARG
1	A	147	ASP
1	A	150	SER
1	A	160	ASN
1	A	169	ASN
1	A	173	ASN
1	A	217	MET
1	A	269	ASN
1	A	284	ARG
1	A	331	ARG
1	A	353	ARG
1	A	367	SER
1	A	436	ARG
1	A	455	SER
1	A	478	SER
1	A	525	SER
1	A	550	LYS
1	A	552	LYS
1	A	600	LEU
1	A	653	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	62	ASN
1	A	88	ASN
1	A	120	ASN
1	A	140	HIS
1	A	263	ASN
1	A	269	ASN
1	A	392	GLN
1	A	410	GLN
1	A	416	ASN
1	A	453	GLN
1	A	479	ASN
1	A	560	ASN
1	A	685	GLN

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 ⓘ

15 carbohydrates 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	GLC	A	689	2	11,11,12	0.54	0	11,15,17	1.03	1 (9%)
2	GLC	A	690	2	12,12,12	0.37	0	17,17,17	0.53	0
3	GLC	A	691	3	11,11,12	0.43	0	11,15,17	0.92	1 (9%)
3	G6D	A	692	3,7	10,10,11	0.53	0	10,14,16	1.50	2 (20%)
4	GLC	A	694	4,7	11,11,12	0.43	0	11,15,17	0.66	0
4	GLC	A	695	4	11,11,12	0.42	0	11,15,17	0.94	1 (9%)
4	GLC	A	696	4	12,12,12	0.37	0	17,17,17	0.85	0
5	BGC	A	698	5	11,11,12	0.43	0	11,15,17	0.82	0
5	GLC	A	699	5	11,11,12	0.44	0	11,15,17	1.27	2 (18%)
5	GLC	A	700	5	11,11,12	0.44	0	11,15,17	1.02	1 (9%)
5	GLC	A	701	5	11,11,12	0.45	0	11,15,17	0.79	0
5	GLC	A	702	5	12,12,12	0.37	0	17,17,17	1.06	1 (5%)
4	GLC	A	705	4	11,11,12	0.45	0	11,15,17	0.67	0
4	GLC	A	706	4	11,11,12	0.43	0	11,15,17	0.84	0
4	GLC	A	707	4	12,12,12	0.37	0	17,17,17	1.15	2 (11%)

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	GLC	A	689	2	-	0/2/18/22	0/1/1/1
2	GLC	A	690	2	-	0/2/22/22	0/1/1/1
3	GLC	A	691	3	-	0/2/18/22	0/1/1/1
3	G6D	A	692	3,7	-	0/0/16/20	0/1/1/1
4	GLC	A	694	4,7	-	0/2/18/22	0/1/1/1
4	GLC	A	695	4	-	0/2/18/22	0/1/1/1
4	GLC	A	696	4	-	0/2/22/22	0/1/1/1
5	BGC	A	698	5	-	0/2/18/22	0/1/1/1
5	GLC	A	699	5	-	0/2/18/22	1/1/1/1
5	GLC	A	700	5	-	0/2/18/22	0/1/1/1
5	GLC	A	701	5	-	0/2/18/22	0/1/1/1
5	GLC	A	702	5	-	0/2/22/22	0/1/1/1
4	GLC	A	705	4	-	0/2/18/22	0/1/1/1
4	GLC	A	706	4	-	0/2/18/22	0/1/1/1
4	GLC	A	707	4	-	0/2/22/22	0/1/1/1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	692	G6D	O5-C1-C2	3.40	115.13	109.86
2	A	689	GLC	O6-C6-C5	-2.66	108.48	112.42
4	A	695	GLC	C1-O5-C5	2.64	118.07	112.41
5	A	700	GLC	C1-O5-C5	2.60	117.98	112.41
5	A	699	GLC	C1-O5-C5	2.53	117.83	112.41
3	A	692	G6D	C1-O5-C5A	2.46	117.68	112.41
5	A	699	GLC	O5-C1-C2	2.37	113.54	109.86
4	A	707	GLC	O5-C5-C4	2.37	114.15	109.76
4	A	707	GLC	C3-C4-C5	2.20	114.13	110.20
3	A	691	GLC	O6-C6-C5	-2.15	109.23	112.42
5	A	702	GLC	O5-C5-C4	2.10	113.65	109.76

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	699	GLC	C1-C2-C3-C4-C5-O5

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	$\# Z > 2$
7	ADH	A	693	3,4	11,11,11	1.52	1 (9%)	15,15,15	1.92	1 (6%)

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
7	ADH	A	693	3,4	-	0/2/18/18	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	693	ADH	C4-C5	-4.43	1.40	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	693	ADH	C6-C5-C7	-6.17	110.28	123.13

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%)	-0.57	4 (0%) 86 86	11, 20, 37, 70	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	89	ASP	3.9
1	A	90	SER	3.0
1	A	92	VAL	2.8
1	A	91	GLY	2.1

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 ⓘ

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
5	GLC	A	702	12/12	0.25	6.22	46,49,58,66	0
4	GLC	A	707	12/12	0.24	5.55	36,38,48,90	0
4	GLC	A	694	11/12	0.18	2.10	30,32,36,39	0
4	GLC	A	695	11/12	0.18	1.60	39,44,49,52	0
2	GLC	A	690	12/12	0.11	0.51	19,21,25,30	0
5	GLC	A	701	11/12	0.15	0.31	41,42,47,49	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	GLC	A	691	11/12	0.15	0.24	26,27,32,36	0
4	GLC	A	706	11/12	0.09	0.06	30,31,36,39	0
2	GLC	A	689	11/12	0.10	0.02	16,18,23,29	0
5	GLC	A	700	11/12	0.13	-0.29	44,46,51,52	0
4	GLC	A	705	11/12	0.09	-0.50	31,31,36,39	0
3	G6D	A	692	10/11	0.10	-0.61	22,23,26,28	0
5	GLC	A	699	11/12	0.33	-	55,57,63,65	0
4	GLC	A	696	12/12	0.47	-	51,54,63,90	0
5	BGC	A	698	11/12	0.41	-	65,66,71,74	0

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(\AA^2)	Q<0.9
7	ADH	A	693	11/11	0.19	1.90	25,26,30,35	0
6	CA	A	688	1/1	0.07	-2.54	15,15,15,15	0
6	CA	A	687	1/1	0.02	-4.07	24,24,24,24	0

6.5 Other polymers

There are no such residues in this entry.