



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:25 PM UTC

PDB ID : 1FQD / pdb_00001fqd
Title : CRYSTAL STRUCTURE OF MALTOTETRAITOL BOUND TO CLOSED-FORM MALTODEXTRIN BINDING PROTEIN
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Deposited on : 2000-09-04
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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

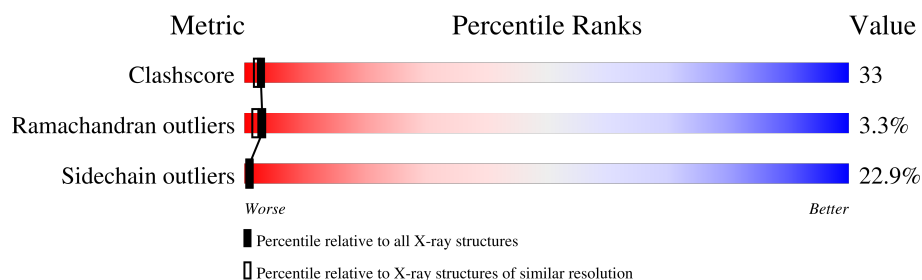
1 Overall quality at a glance

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))
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	370	38% 46% 15% .
2	B	4	50% 50%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLO	B	1	-	-	X	-
2	GLC	B	2	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3012 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 MALTODEXTRIN-BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	370	Total	C	N	O	S	0	0	0
			2862	1843	468	545	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	ALA	LYS	cloning artifact	UNP P02928

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-D-glucose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	4	Total	C	O	0	0	0
			45	24	21			

- Molecule 3 is water.

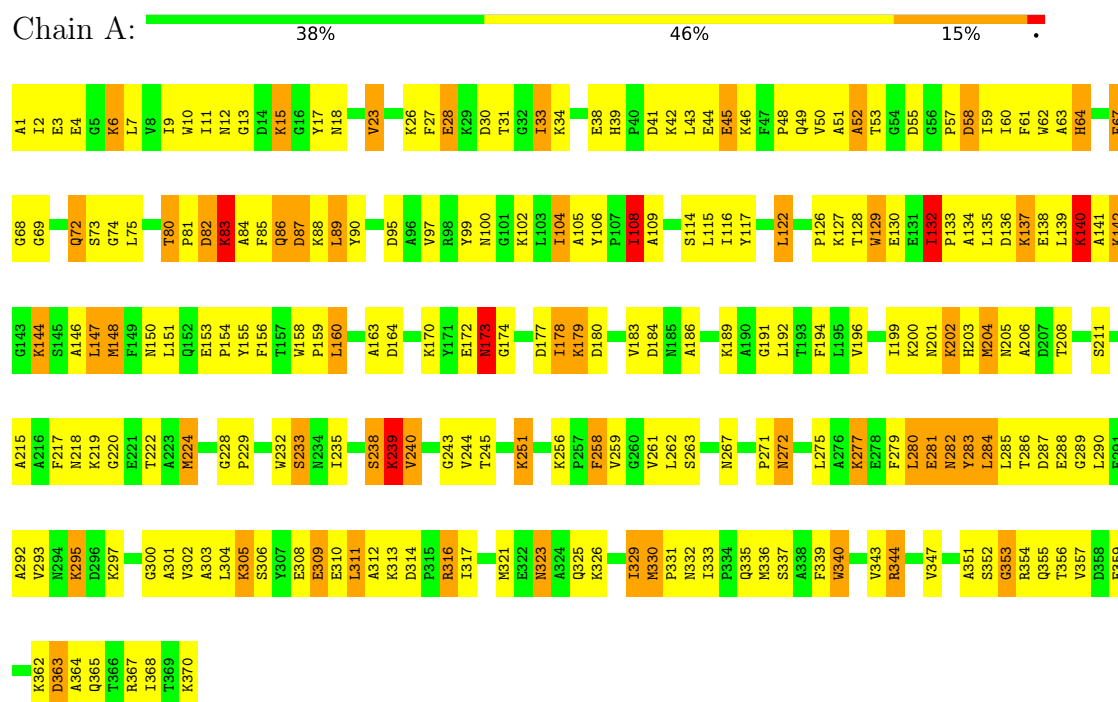
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	105	Total	O	0	0
			105	105		

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. 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. 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.

Note EDS was not executed.

• Molecule 1: MALTODEXTRIN-BINDING PROTEIN



• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-D-glucose



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	104.30Å 69.40Å 56.30Å 90.00° 112.40° 90.00°	Depositor
Resolution (Å)	10.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.30)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.191 , 0.244	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3012	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

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, GLO

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	0.51	1/2931 (0.0%)	1.03	23/3979 (0.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	132	ILE	CA-CB	5.21	1.57	1.54

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	58	ASP	N-CA-C	-7.32	103.38	111.36
1	A	353	GLY	N-CA-C	-6.35	105.86	115.00
1	A	132	ILE	N-CA-C	6.28	119.89	112.35
1	A	67	PHE	N-CA-C	6.02	119.88	112.54
1	A	174	GLY	N-CA-C	5.90	122.91	115.42
1	A	300	GLY	N-CA-C	-5.90	104.19	111.45
1	A	13	GLY	N-CA-C	5.87	121.07	113.27
1	A	228	GLY	CA-C-N	5.82	125.68	119.28
1	A	228	GLY	C-N-CA	5.82	125.68	119.28
1	A	95	ASP	N-CA-C	-5.77	104.99	111.28
1	A	52	ALA	N-CA-C	-5.60	105.09	111.14
1	A	129	TRP	N-CA-C	-5.48	105.41	111.71
1	A	238	SER	N-CA-C	-5.44	106.48	113.23
1	A	64	HIS	N-CA-C	5.37	118.63	111.75
1	A	301	ALA	N-CA-C	-5.28	102.70	110.52
1	A	362	LYS	N-CA-C	-5.27	106.69	113.23
1	A	153	GLU	CA-C-N	5.25	125.31	119.32
1	A	153	GLU	C-N-CA	5.25	125.31	119.32
1	A	323	ASN	N-CA-C	-5.19	105.80	111.82
1	A	108	ILE	N-CA-C	5.10	116.20	111.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	144	LYS	N-CA-C	5.07	118.37	110.32
1	A	28	GLU	N-CA-C	5.06	116.48	111.07
1	A	23	VAL	N-CA-C	-5.00	105.52	110.62

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	2862	0	2831	186	7
2	B	45	0	39	14	0
3	A	105	0	0	3	3
All	All	3012	0	2870	193	8

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 33.

All (193) 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:B:1:GLO:H62	2:B:2:GLC:H5	1.17	1.13
2:B:1:GLO:H62	2:B:2:GLC:C5	1.95	0.96
2:B:1:GLO:C6	2:B:2:GLC:H5	1.94	0.96
1:A:31:THR:HG23	1:A:33:ILE:H	1.29	0.94
1:A:1:ALA:HB1	1:A:271:PRO:HG3	1.61	0.82
1:A:43:LEU:CD1	1:A:60:ILE:HD11	2.14	0.77
1:A:6:LYS:HA	1:A:33:ILE:HG23	1.68	0.74
1:A:200:LYS:HE3	1:A:201:ASN:HD21	1.53	0.73
1:A:43:LEU:HD13	1:A:60:ILE:HD11	1.69	0.73
2:B:1:GLO:H5	2:B:2:GLC:H3	1.69	0.73
1:A:232:TRP:CH2	1:A:316:ARG:HB3	2.25	0.71
1:A:343:VAL:O	1:A:347:VAL:HG23	1.91	0.70
1:A:364:ALA:O	1:A:368:ILE:HD12	1.91	0.70
1:A:277:LYS:HD2	1:A:281:GLU:OE1	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:ILE:HD13	1:A:178:ILE:H	1.57	0.69
1:A:108:ILE:HG23	1:A:262:LEU:O	1.92	0.69
1:A:128:THR:HB	1:A:130:GLU:OE1	1.95	0.67
1:A:64:HIS:CD2	1:A:261:VAL:H	2.14	0.66
1:A:146:ALA:O	1:A:224:MET:HG3	1.96	0.65
1:A:272:ASN:HB2	1:A:275:LEU:HD12	1.78	0.65
1:A:258:PHE:HB3	1:A:330:MET:HE3	1.79	0.65
1:A:85:PHE:CE2	1:A:281:GLU:HG3	2.33	0.64
1:A:11:ILE:CG2	1:A:17:TYR:HB3	2.29	0.63
1:A:97:VAL:HG21	1:A:105:ALA:O	1.99	0.63
1:A:27:PHE:O	1:A:31:THR:HG22	2.00	0.61
1:A:333:ILE:HD12	1:A:335:GLN:HE21	1.65	0.61
2:B:1:GLO:H3	2:B:2:GLC:O2	2.01	0.61
1:A:215:ALA:O	1:A:219:LYS:HB2	2.00	0.61
1:A:12:ASN:ND2	2:B:1:GLO:H4	2.16	0.60
1:A:122:LEU:HD11	1:A:135:LEU:HD21	1.83	0.60
1:A:331:PRO:HD2	1:A:336:MET:SD	2.41	0.60
1:A:64:HIS:HE1	1:A:330:MET:O	1.85	0.59
1:A:139:LEU:HD23	1:A:142:LYS:HD3	1.85	0.59
1:A:12:ASN:HB3	2:B:1:GLO:O3	2.03	0.59
1:A:90:TYR:CE1	1:A:305:LYS:HD3	2.38	0.59
1:A:148:MET:HE2	1:A:208:THR:OG1	2.03	0.59
1:A:173:ASN:O	1:A:173:ASN:CG	2.45	0.59
2:B:1:GLO:C5	2:B:2:GLC:H5	2.33	0.59
1:A:23:VAL:HG12	1:A:279:PHE:HE1	1.68	0.59
1:A:116:ILE:HA	1:A:243:GLY:O	2.03	0.59
1:A:330:MET:CE	1:A:340:TRP:HZ2	2.16	0.58
1:A:363:ASP:O	1:A:367:ARG:HG3	2.03	0.58
1:A:11:ILE:HG21	1:A:17:TYR:HB3	1.87	0.57
1:A:147:LEU:HG	1:A:148:MET:N	2.19	0.57
1:A:136:ASP:HA	1:A:146:ALA:HB2	1.85	0.57
1:A:81:PRO:HB2	1:A:86:GLN:HG2	1.87	0.57
1:A:129:TRP:HA	1:A:132:ILE:HD12	1.86	0.57
1:A:340:TRP:HA	1:A:340:TRP:CE3	2.40	0.56
1:A:48:PRO:HA	1:A:75:LEU:HD13	1.86	0.56
1:A:139:LEU:C	1:A:141:ALA:H	2.13	0.56
1:A:178:ILE:HD13	1:A:178:ILE:N	2.21	0.56
1:A:339:PHE:HA	1:A:368:ILE:HG12	1.88	0.56
1:A:199:ILE:O	1:A:202:LYS:HE2	2.06	0.55
1:A:229:PRO:HA	1:A:232:TRP:CE2	2.40	0.55
1:A:69:GLY:O	1:A:72:GLN:HB3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:ILE:HD12	1:A:335:GLN:HB2	1.89	0.55
1:A:89:LEU:N	1:A:89:LEU:HD23	2.22	0.55
1:A:31:THR:HG23	1:A:33:ILE:N	2.11	0.54
1:A:46:LYS:O	1:A:49:GLN:HB2	2.07	0.54
1:A:154:PRO:HD3	1:A:344:ARG:HG3	1.89	0.54
1:A:292:ALA:O	1:A:295:LYS:HG3	2.08	0.54
1:A:58:ASP:CG	1:A:272:ASN:HD21	2.15	0.54
1:A:330:MET:HE1	1:A:340:TRP:CZ2	2.42	0.54
1:A:356:THR:OG1	1:A:359:GLU:HB2	2.06	0.54
1:A:139:LEU:C	1:A:141:ALA:N	2.66	0.53
1:A:55:ASP:O	1:A:267:ASN:ND2	2.40	0.53
1:A:82:ASP:C	1:A:84:ALA:N	2.64	0.53
1:A:178:ILE:H	1:A:178:ILE:CD1	2.21	0.53
1:A:192:LEU:HB3	1:A:357:VAL:HG13	1.91	0.53
1:A:99:TYR:O	1:A:100:ASN:HB2	2.08	0.53
1:A:203:HIS:O	1:A:204:MET:HB3	2.08	0.53
1:A:280:LEU:HD12	1:A:284:LEU:HD12	1.90	0.53
1:A:97:VAL:O	1:A:104:ILE:HG13	2.09	0.52
1:A:129:TRP:O	1:A:132:ILE:HB	2.09	0.52
1:A:62:TRP:CH2	2:B:1:GLO:O5	2.61	0.52
1:A:82:ASP:O	1:A:84:ALA:N	2.43	0.52
1:A:330:MET:HE1	1:A:340:TRP:HZ2	1.75	0.52
1:A:67:PHE:HB3	1:A:104:ILE:HD12	1.91	0.52
2:B:1:GLO:H5	2:B:2:GLC:H5	1.91	0.52
1:A:68:GLY:HA3	1:A:332:ASN:O	2.10	0.51
1:A:64:HIS:CE1	1:A:330:MET:O	2.63	0.51
1:A:50:VAL:HG23	1:A:51:ALA:N	2.25	0.51
1:A:82:ASP:N	1:A:82:ASP:OD1	2.43	0.51
1:A:314:ASP:HB3	1:A:317:ILE:HG12	1.91	0.51
1:A:367:ARG:O	1:A:370:LYS:HB2	2.11	0.51
1:A:85:PHE:O	1:A:87:ASP:N	2.44	0.51
1:A:311:LEU:C	1:A:313:LYS:H	2.19	0.51
1:A:164:ASP:OD2	1:A:251:LYS:HG2	2.11	0.50
1:A:330:MET:CE	1:A:340:TRP:CZ2	2.94	0.50
1:A:28:GLU:C	1:A:30:ASP:H	2.19	0.50
1:A:244:VAL:HG21	1:A:316:ARG:HG2	1.93	0.50
1:A:177:ASP:OD1	1:A:179:LYS:N	2.42	0.50
1:A:311:LEU:O	1:A:313:LYS:N	2.45	0.50
1:A:286:THR:O	1:A:287:ASP:C	2.55	0.49
1:A:117:TYR:CE1	1:A:243:GLY:HA3	2.47	0.49
1:A:136:ASP:HB3	1:A:203:HIS:HD2	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:LYS:O	1:A:240:VAL:C	2.55	0.49
1:A:80:THR:O	1:A:80:THR:OG1	2.21	0.49
1:A:1:ALA:HB1	1:A:271:PRO:CG	2.39	0.49
1:A:43:LEU:HD12	1:A:60:ILE:HD11	1.94	0.49
1:A:139:LEU:O	1:A:141:ALA:N	2.46	0.49
1:A:72:GLN:C	1:A:74:GLY:H	2.21	0.49
1:A:259:VAL:HB	1:A:329:ILE:HA	1.95	0.49
1:A:356:THR:HG1	1:A:359:GLU:HB2	1.78	0.49
1:A:57:PRO:HB2	1:A:59:ILE:O	2.13	0.49
1:A:155:TYR:HB2	2:B:3:GLC:O6	2.13	0.49
1:A:9:ILE:HD11	1:A:279:PHE:CE2	2.48	0.48
1:A:331:PRO:HD2	1:A:336:MET:CE	2.44	0.48
1:A:48:PRO:HA	1:A:75:LEU:CD1	2.43	0.48
1:A:147:LEU:HB2	1:A:224:MET:HE3	1.95	0.48
1:A:154:PRO:CD	1:A:344:ARG:HG3	2.43	0.48
1:A:88:LYS:C	1:A:89:LEU:HD23	2.38	0.48
1:A:292:ALA:O	1:A:295:LYS:HE3	2.14	0.47
1:A:43:LEU:HD13	1:A:60:ILE:CD1	2.42	0.47
1:A:192:LEU:O	1:A:196:VAL:HG23	2.13	0.47
1:A:64:HIS:HD2	1:A:261:VAL:H	1.57	0.47
1:A:277:LYS:O	1:A:281:GLU:HB2	2.14	0.47
1:A:199:ILE:O	1:A:202:LYS:N	2.36	0.47
1:A:62:TRP:HD1	1:A:63:ALA:O	1.97	0.47
1:A:10:TRP:CD1	1:A:57:PRO:HB3	2.50	0.47
1:A:11:ILE:O	1:A:39:HIS:HA	2.14	0.47
1:A:60:ILE:O	1:A:60:ILE:HG23	2.14	0.47
1:A:45:GLU:O	1:A:49:GLN:HG3	2.15	0.46
1:A:259:VAL:O	1:A:330:MET:HG3	2.15	0.46
1:A:308:GLU:OE2	1:A:321:MET:HE2	2.15	0.46
1:A:311:LEU:C	1:A:313:LYS:N	2.72	0.46
1:A:12:ASN:HD22	2:B:1:GLO:H4	1.80	0.46
1:A:239:LYS:O	1:A:240:VAL:O	2.33	0.46
1:A:108:ILE:HG12	1:A:262:LEU:HD23	1.97	0.46
1:A:144:LYS:HE3	1:A:220:GLY:O	2.15	0.46
1:A:279:PHE:O	1:A:283:TYR:HB2	2.15	0.46
1:A:82:ASP:C	1:A:84:ALA:H	2.24	0.46
1:A:282:ASN:O	1:A:283:TYR:CD2	2.69	0.46
1:A:217:PHE:HA	1:A:222:THR:HG22	1.99	0.45
1:A:85:PHE:C	1:A:87:ASP:H	2.24	0.45
1:A:136:ASP:CA	1:A:146:ALA:HB2	2.46	0.45
1:A:184:ASP:HB2	1:A:365:GLN:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:TYR:O	1:A:289:GLY:HA3	2.17	0.45
1:A:311:LEU:HD21	3:A:625:HOH:O	2.17	0.45
1:A:353:GLY:C	1:A:355:GLN:H	2.24	0.45
1:A:217:PHE:HD1	1:A:218:ASN:HD22	1.65	0.45
1:A:10:TRP:CD2	1:A:57:PRO:HG3	2.52	0.45
1:A:44:GLU:HB3	1:A:62:TRP:CZ3	2.52	0.45
1:A:85:PHE:C	1:A:87:ASP:N	2.74	0.45
1:A:81:PRO:HG2	1:A:86:GLN:NE2	2.31	0.45
1:A:41:ASP:O	1:A:42:LYS:C	2.60	0.45
1:A:122:LEU:HD21	1:A:135:LEU:HD21	1.99	0.45
1:A:160:LEU:HD22	1:A:191:GLY:HA2	1.98	0.44
1:A:356:THR:O	1:A:357:VAL:C	2.60	0.44
1:A:109:ALA:HA	1:A:302:VAL:HA	2.00	0.44
1:A:262:LEU:HD13	2:B:1:GLO:O1	2.18	0.44
1:A:158:TRP:N	1:A:159:PRO:CD	2.81	0.43
1:A:233:SER:HB2	1:A:297:LYS:HD3	2.00	0.43
1:A:26:LYS:HG2	1:A:283:TYR:HE1	1.82	0.43
1:A:102:LYS:HA	3:A:559:HOH:O	2.18	0.43
1:A:134:ALA:O	1:A:137:LYS:HG2	2.18	0.43
1:A:106:TYR:O	1:A:263:SER:HA	2.19	0.43
1:A:133:PRO:HA	1:A:203:HIS:CD2	2.53	0.43
1:A:218:ASN:HD21	1:A:235:ILE:HG12	1.83	0.43
1:A:206:ALA:HB1	1:A:351:ALA:HB1	2.01	0.43
1:A:82:ASP:O	1:A:83:LYS:C	2.62	0.43
1:A:31:THR:HG23	1:A:33:ILE:HB	2.01	0.43
1:A:305:LYS:HB2	3:A:476:HOH:O	2.18	0.42
1:A:83:LYS:O	1:A:87:ASP:OD1	2.37	0.42
1:A:304:LEU:HD12	1:A:304:LEU:HA	1.84	0.42
1:A:122:LEU:HD21	1:A:126:PRO:HD3	2.02	0.42
1:A:10:TRP:CG	1:A:57:PRO:HG3	2.54	0.42
1:A:62:TRP:CD1	1:A:63:ALA:N	2.88	0.42
1:A:172:GLU:O	1:A:173:ASN:C	2.62	0.42
1:A:272:ASN:CB	1:A:275:LEU:HD12	2.48	0.42
1:A:109:ALA:O	1:A:262:LEU:N	2.44	0.42
1:A:280:LEU:O	1:A:285:LEU:HB2	2.20	0.42
1:A:114:SER:OG	1:A:323:ASN:ND2	2.53	0.41
1:A:354:ARG:HE	1:A:354:ARG:HB2	1.66	0.41
1:A:130:GLU:HA	1:A:194:PHE:HZ	1.84	0.41
1:A:150:ASN:HB3	1:A:156:PHE:CD2	2.55	0.41
1:A:333:ILE:HB	1:A:335:GLN:NE2	2.35	0.41
1:A:50:VAL:C	1:A:52:ALA:N	2.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:PHE:HD1	1:A:258:PHE:HA	1.77	0.41
1:A:15:LYS:HA	1:A:15:LYS:HD2	1.74	0.41
1:A:151:LEU:HD12	1:A:206:ALA:HA	2.03	0.41
1:A:163:ALA:HA	1:A:256:LYS:HG3	2.02	0.41
1:A:130:GLU:O	1:A:133:PRO:HD2	2.21	0.41
1:A:23:VAL:HG12	1:A:279:PHE:CE1	2.53	0.40
1:A:89:LEU:HA	1:A:303:ALA:O	2.20	0.40
1:A:179:LYS:HA	1:A:179:LYS:HD3	1.65	0.40
1:A:27:PHE:HD1	1:A:283:TYR:CE2	2.39	0.40
1:A:108:ILE:HD12	1:A:290:LEU:CD2	2.51	0.40
1:A:117:TYR:O	1:A:243:GLY:N	2.53	0.40
1:A:340:TRP:CD1	2:B:3:GLC:H4	2.56	0.40
1:A:136:ASP:O	1:A:140:LYS:HB2	2.21	0.40
1:A:238:SER:O	1:A:239:LYS:CB	2.70	0.40
1:A:61:PHE:HA	1:A:263:SER:O	2.22	0.40

All (8) 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:180:ASP:OD2	1:A:186:ALA:CB[2_657]	1.42	0.78
3:A:504:HOH:O	3:A:504:HOH:O[2_657]	1.43	0.77
1:A:72:GLN:OE1	1:A:309:GLU:CD[4_557]	1.53	0.67
1:A:189:LYS:NZ	3:A:447:HOH:O[2_657]	1.53	0.67
1:A:354:ARG:NH2	3:A:510:HOH:O[4_556]	1.78	0.42
1:A:72:GLN:OE1	1:A:309:GLU:CG[4_557]	1.80	0.40
1:A:72:GLN:OE1	1:A:309:GLU:OE2[4_557]	1.84	0.36
1:A:99:TYR:OH	1:A:309:GLU:OE2[4_557]	2.17	0.03

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	368/370 (100%)	323 (88%)	33 (9%)	12 (3%)	3 2

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	240	VAL
1	A	3	GLU
1	A	4	GLU
1	A	86	GLN
1	A	173	ASN
1	A	239	LYS
1	A	283	TYR
1	A	312	ALA
1	A	83	LYS
1	A	140	LYS
1	A	204	MET
1	A	2	ILE

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	292/296 (99%)	225 (77%)	67 (23%)	1 1

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	7	LEU
1	A	15	LYS
1	A	18	ASN
1	A	33	ILE
1	A	34	LYS
1	A	38	GLU
1	A	45	GLU
1	A	53	THR

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Mol	Chain	Res	Type
1	A	72	GLN
1	A	73	SER
1	A	80	THR
1	A	82	ASP
1	A	83	LYS
1	A	87	ASP
1	A	89	LEU
1	A	104	ILE
1	A	108	ILE
1	A	115	LEU
1	A	122	LEU
1	A	127	LYS
1	A	132	ILE
1	A	137	LYS
1	A	138	GLU
1	A	140	LYS
1	A	142	LYS
1	A	147	LEU
1	A	148	MET
1	A	160	LEU
1	A	170	LYS
1	A	173	ASN
1	A	178	ILE
1	A	179	LYS
1	A	183	VAL
1	A	202	LYS
1	A	205	ASN
1	A	211	SER
1	A	224	MET
1	A	233	SER
1	A	239	LYS
1	A	245	THR
1	A	251	LYS
1	A	258	PHE
1	A	272	ASN
1	A	277	LYS
1	A	280	LEU
1	A	281	GLU
1	A	282	ASN
1	A	284	LEU
1	A	288	GLU
1	A	293	VAL

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Mol	Chain	Res	Type
1	A	295	LYS
1	A	305	LYS
1	A	306	SER
1	A	309	GLU
1	A	310	GLU
1	A	311	LEU
1	A	316	ARG
1	A	325	GLN
1	A	326	LYS
1	A	329	ILE
1	A	330	MET
1	A	337	SER
1	A	340	TRP
1	A	344	ARG
1	A	352	SER
1	A	363	ASP

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

Mol	Chain	Res	Type
1	A	12	ASN
1	A	18	ASN
1	A	49	GLN
1	A	64	HIS
1	A	100	ASN
1	A	173	ASN
1	A	201	ASN
1	A	218	ASN
1	A	272	ASN
1	A	335	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

4 monosaccharides 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	GLO	B	1	2	10,11,11	2.33	5 (50%)	13,14,14	2.61	3 (23%)
2	GLC	B	2	2	11,11,12	0.60	0	15,15,17	1.15	2 (13%)
2	GLC	B	3	2	11,11,12	0.52	0	15,15,17	0.56	0
2	GLC	B	4	2	11,11,12	0.64	0	15,15,17	0.71	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
2	GLO	B	1	2	-	7/15/16/16	-
2	GLC	B	2	2	-	0/2/19/22	0/1/1/1
2	GLC	B	3	2	-	0/2/19/22	0/1/1/1
2	GLC	B	4	2	-	2/2/19/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	GLO	C3-C2	4.38	1.59	1.53
2	B	1	GLO	C4-C3	3.43	1.59	1.53
2	B	1	GLO	O5-C5	-2.51	1.38	1.43
2	B	1	GLO	O2-C2	-2.43	1.39	1.43
2	B	1	GLO	C6-C5	2.34	1.58	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	GLO	C5-C4-C3	8.05	124.84	112.48
2	B	1	GLO	O6-C6-C5	2.94	117.32	111.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	4	GLC	C1-O5-C5	2.32	115.29	112.19
2	B	1	GLO	C4-C3-C2	2.28	117.48	113.49
2	B	2	GLC	C1-O5-C5	2.20	115.14	112.19
2	B	2	GLC	C6-C5-C4	2.12	118.22	113.02

There are no chirality outliers.

All (9) torsion outliers are listed below:

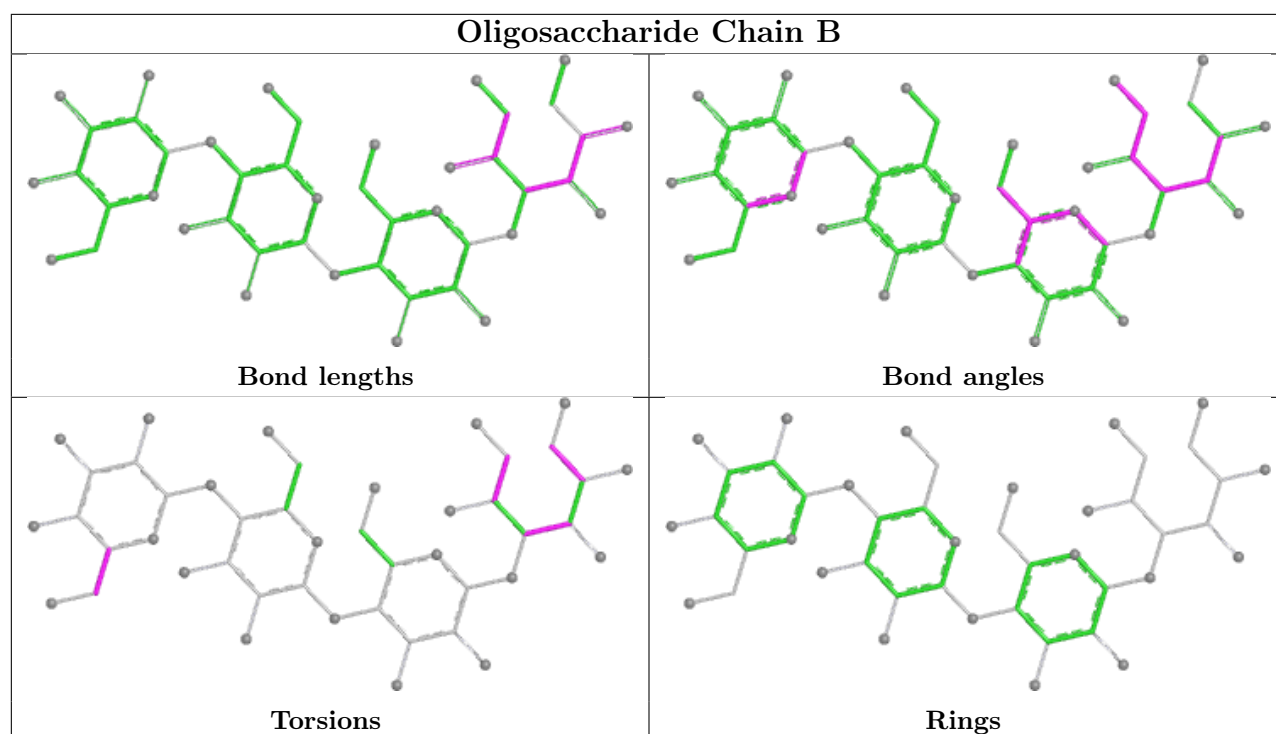
Mol	Chain	Res	Type	Atoms
2	B	1	GLO	O1-C1-C2-C3
2	B	1	GLO	C2-C3-C4-C5
2	B	1	GLO	C2-C3-C4-O4
2	B	1	GLO	O3-C3-C4-C5
2	B	1	GLO	O3-C3-C4-O4
2	B	1	GLO	O5-C5-C6-O6
2	B	4	GLC	C4-C5-C6-O6
2	B	4	GLC	O5-C5-C6-O6
2	B	1	GLO	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	GLC	7	0
2	B	3	GLC	2	0
2	B	1	GLO	12	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.