



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 04:14 AM UTC

PDB ID : 1CGY / pdb_00001cgy
Title : SITE DIRECTED MUTATIONS OF THE ACTIVE SITE RESIDUE TYROSINE 195 OF CYCLODEXTRIN GLYXOSYLTRANSFERASE FROM BACILLUS CIRCULANS STRAIN 251 AFFECTING ACTIVITY AND PRODUCT SPECIFICITY
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Deposited on : 1994-08-05
Resolution : 2.50 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

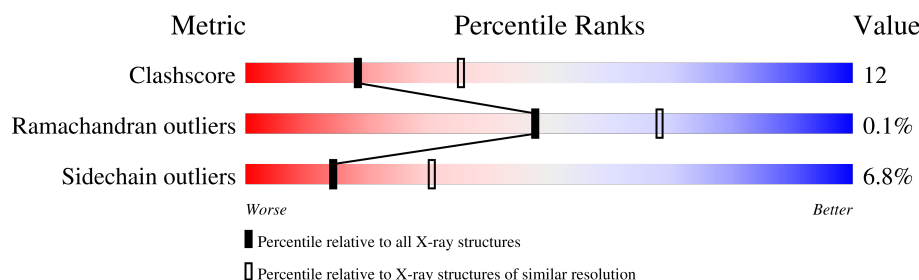
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

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	686	 68% 28% . .
2	B	2	 100%
2	C	2	 100%
2	D	2	 100%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5510 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 CYCLOMALTODEXTRIN GLUCANOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	686	Total	C	N	O	S	0	0	0
			5266	3323	901	1026	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	195	TRP	TYR	conflict	UNP P43379

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	2	Total	C	O	0	0	0
			23	12	11			
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 4 is water.

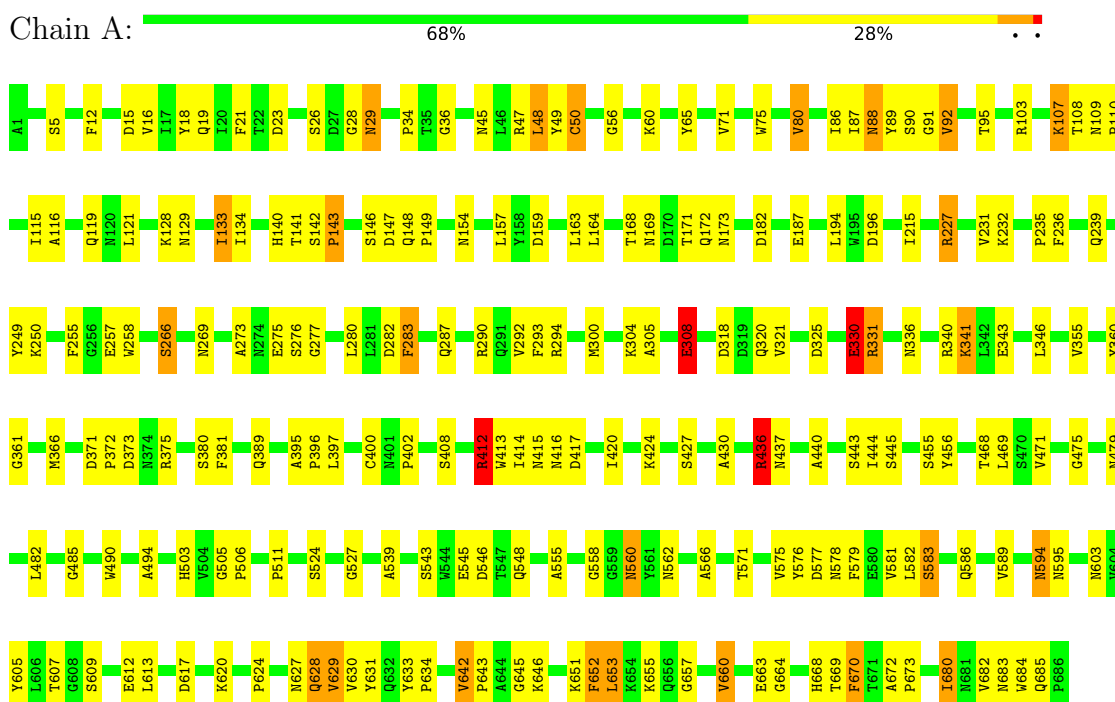
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	173	Total 173	O 173	0	0

3 Residue-property plots

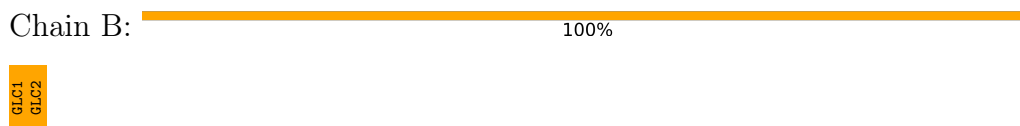
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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.

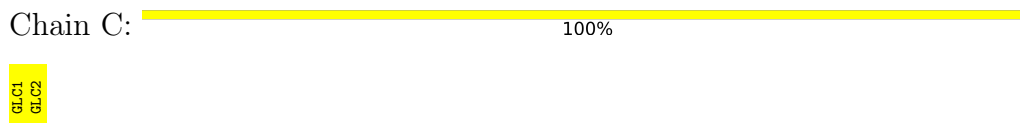
• Molecule 1: CYCLOMALTODEXTRIN GLUCANOTRANSFERASE



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



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



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

Chain D:  100%

GLC1
GLC2

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	120.18Å 110.70Å 66.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, TNT	Depositor
R, R_{free}	0.162 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5510	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	1.31	5/5397 (0.1%)	1.70	70/7357 (1.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	543	SER	CA-C	-5.50	1.45	1.52
1	A	440	ALA	CA-C	-5.30	1.48	1.52
1	A	50	CYS	C-O	-5.20	1.17	1.23
1	A	28	GLY	C-N	-5.17	1.29	1.33
1	A	305	ALA	N-CA	-5.01	1.39	1.46

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	331	ARG	NE-CZ-NH2	-9.06	111.05	119.20
1	A	412	ARG	N-CA-CB	8.15	121.44	110.59
1	A	539	ALA	N-CA-C	7.33	120.20	111.33
1	A	545	GLU	CG-CD-OE2	-7.28	101.66	118.40
1	A	430	ALA	N-CA-CB	7.05	122.66	110.81
1	A	92	VAL	N-CA-C	6.80	117.86	108.89
1	A	546	ASP	CA-CB-CG	-6.79	105.81	112.60
1	A	277	GLY	N-CA-C	-6.70	105.93	114.37
1	A	129	ASN	CA-CB-CG	-6.64	105.96	112.60
1	A	331	ARG	NE-CZ-NH1	6.63	128.13	121.50
1	A	236	PHE	N-CA-CB	6.54	119.49	110.01
1	A	629	VAL	N-CA-C	6.51	122.89	109.34
1	A	668	HIS	N-CA-C	-6.49	101.79	110.55
1	A	29	ASN	CA-C-N	6.30	126.21	119.28
1	A	29	ASN	C-N-CA	6.30	126.21	119.28
1	A	545	GLU	CG-CD-OE1	6.27	132.83	118.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	276	SER	N-CA-C	6.27	119.09	111.82
1	A	36	GLY	O-C-N	6.20	129.62	122.61
1	A	543	SER	N-CA-CB	6.13	121.84	110.99
1	A	373	ASP	N-CA-C	6.10	119.82	112.38
1	A	414	ILE	N-CA-C	6.08	116.27	107.88
1	A	660	VAL	N-CA-C	6.03	118.06	108.89
1	A	65	TYR	N-CA-C	6.01	118.33	111.11
1	A	440	ALA	O-C-N	5.95	127.09	121.38
1	A	275	GLU	N-CA-CB	-5.91	102.72	110.88
1	A	576	TYR	O-C-N	-5.89	116.44	123.27
1	A	485	GLY	CA-C-N	5.88	126.31	120.31
1	A	485	GLY	C-N-CA	5.88	126.31	120.31
1	A	109	ASN	N-CA-C	-5.88	101.07	109.24
1	A	436	ARG	N-CA-C	5.82	118.40	111.71
1	A	628	GLN	CB-CA-C	-5.78	101.70	110.83
1	A	579	PHE	CA-CB-CG	-5.77	108.03	113.80
1	A	16	VAL	N-CA-CB	5.71	119.59	111.82
1	A	308	GLU	N-CA-CB	5.68	118.57	110.16
1	A	412	ARG	NE-CZ-NH1	5.68	127.18	121.50
1	A	236	PHE	CA-C-N	5.68	126.39	120.03
1	A	236	PHE	C-N-CA	5.68	126.39	120.03
1	A	437	ASN	CA-CB-CG	-5.62	106.98	112.60
1	A	140	HIS	CA-CB-CG	-5.60	108.20	113.80
1	A	416	ASN	CB-CA-C	-5.59	101.50	110.79
1	A	330	GLU	N-CA-CB	5.56	118.18	110.17
1	A	292	VAL	N-CA-C	5.56	115.75	110.53
1	A	609	SER	N-CA-C	5.55	119.53	112.87
1	A	652	PHE	CA-CB-CG	-5.50	108.30	113.80
1	A	560	ASN	OD1-CG-ND2	5.50	128.10	122.60
1	A	503	HIS	CA-CB-CG	5.47	119.27	113.80
1	A	673	PRO	N-CA-CB	5.47	108.53	103.39
1	A	548	GLN	OE1-CD-NE2	-5.43	117.17	122.60
1	A	577	ASP	N-CA-C	5.34	118.24	110.17
1	A	50	CYS	CA-C-O	-5.30	113.30	118.97
1	A	182	ASP	N-CA-C	-5.27	106.16	112.59
1	A	28	GLY	CA-C-N	-5.26	116.42	122.52
1	A	28	GLY	C-N-CA	-5.26	116.42	122.52
1	A	402	PRO	N-CA-CB	5.26	108.74	103.48
1	A	594	ASN	OD1-CG-ND2	-5.24	117.36	122.60
1	A	266	SER	CA-C-N	5.22	124.83	119.56
1	A	266	SER	C-N-CA	5.22	124.83	119.56
1	A	336	ASN	N-CA-CB	-5.22	102.64	110.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	143	PRO	N-CA-CB	5.19	107.79	103.17
1	A	346	LEU	N-CA-CB	5.15	117.48	110.01
1	A	408	SER	N-CA-C	-5.15	104.08	110.41
1	A	657	GLY	N-CA-C	-5.15	101.39	110.86
1	A	389	GLN	CB-CG-CD	-5.08	103.97	112.60
1	A	586	GLN	N-CA-C	5.08	118.01	110.14
1	A	642	VAL	O-C-N	5.05	126.07	121.57
1	A	92	VAL	N-CA-CB	5.05	116.03	110.53
1	A	133	ILE	CA-C-N	-5.04	114.25	122.57
1	A	133	ILE	C-N-CA	-5.04	114.25	122.57
1	A	490	TRP	CA-C-O	-5.04	115.41	121.11
1	A	283	PHE	N-CA-CB	5.01	117.93	110.22

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	5266	0	5026	122	0
2	B	23	0	21	3	0
2	C	23	0	21	0	0
2	D	23	0	21	5	0
3	A	2	0	0	0	0
4	A	173	0	0	4	0
All	All	5510	0	5089	125	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 12.

All (125) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:560:ASN:HD21	1:A:578:ASN:HA	1.24	1.01
1:A:88:ASN:HD21	1:A:91:GLY:H	1.04	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:SER:HB2	1:A:143:PRO:HD2	1.61	0.82
1:A:88:ASN:HD21	1:A:91:GLY:N	1.81	0.76
1:A:159:ASP:HB2	1:A:164:LEU:HD11	1.73	0.70
1:A:603:ASN:HB3	1:A:624:PRO:HB3	1.75	0.69
1:A:583:SER:HB2	1:A:612:GLU:OE2	1.95	0.66
1:A:231:VAL:HG22	1:A:257:GLU:O	1.96	0.65
1:A:45:ASN:ND2	1:A:48:LEU:HD22	2.13	0.63
1:A:48:LEU:HD12	1:A:95:THR:HG23	1.82	0.61
1:A:395:ALA:HB3	1:A:396:PRO:HD3	1.83	0.61
1:A:121:LEU:HD12	1:A:121:LEU:O	2.01	0.60
1:A:227:ARG:HG2	1:A:255:PHE:CE2	2.37	0.60
1:A:560:ASN:ND2	1:A:578:ASN:HA	2.07	0.59
1:A:331:ARG:HA	4:A:1164:HOH:O	2.02	0.59
2:D:1:GLC:H62	2:D:2:GLC:C5	2.33	0.59
1:A:645:GLY:HA2	1:A:672:ALA:O	2.03	0.58
1:A:142:SER:HB2	1:A:143:PRO:CD	2.33	0.58
1:A:417:ASP:OD1	1:A:436:ARG:HD2	2.03	0.58
1:A:427:SER:HB2	1:A:494:ALA:O	2.04	0.57
1:A:60:LYS:HE2	1:A:381:PHE:CD2	2.40	0.57
1:A:321:VAL:HG22	1:A:355:VAL:HB	1.87	0.57
1:A:18:TYR:HB2	1:A:71:VAL:HG11	1.87	0.56
1:A:18:TYR:HB2	1:A:71:VAL:CG1	2.35	0.56
1:A:669:THR:HG22	1:A:670:PHE:N	2.21	0.56
1:A:227:ARG:HD2	1:A:227:ARG:C	2.31	0.56
1:A:87:ILE:CD1	1:A:143:PRO:HG2	2.36	0.55
1:A:653:LEU:HD23	1:A:653:LEU:N	2.22	0.55
1:A:583:SER:HB2	1:A:612:GLU:CD	2.32	0.55
1:A:60:LYS:HE2	1:A:381:PHE:CE2	2.42	0.54
1:A:26:SER:O	1:A:56:GLY:HA3	2.07	0.54
1:A:235:PRO:O	1:A:239:GLN:HG3	2.08	0.54
1:A:227:ARG:HH11	1:A:257:GLU:HB2	1.75	0.52
1:A:304:LYS:O	1:A:308:GLU:HG2	2.10	0.52
1:A:589:VAL:HG11	1:A:680:ILE:HD11	1.91	0.52
1:A:630:VAL:HG12	1:A:631:TYR:CE1	2.45	0.52
1:A:88:ASN:ND2	1:A:91:GLY:H	1.88	0.52
1:A:627:ASN:HD21	2:D:2:GLC:H61	1.74	0.52
1:A:562:ASN:HB3	1:A:575:VAL:HG13	1.92	0.51
1:A:282:ASP:HB2	1:A:320:GLN:HB3	1.93	0.51
1:A:607:THR:HG22	1:A:653:LEU:CD2	2.42	0.49
1:A:232:LYS:O	1:A:232:LYS:HG2	2.13	0.49
1:A:633:TYR:CG	1:A:634:PRO:HA	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:1:GLC:H62	2:D:2:GLC:H5	1.94	0.49
1:A:456:TYR:O	1:A:468:THR:HG23	2.13	0.48
2:D:1:GLC:H62	2:D:2:GLC:O5	2.13	0.48
1:A:558:GLY:HA2	1:A:581:VAL:O	2.14	0.47
1:A:524:SER:HB3	4:A:1152:HOH:O	2.13	0.47
1:A:479:ASN:ND2	1:A:479:ASN:N	2.59	0.47
1:A:194:LEU:HD23	1:A:194:LEU:HA	1.69	0.47
1:A:445:SER:HB3	1:A:479:ASN:ND2	2.29	0.47
1:A:652:PHE:C	1:A:653:LEU:HD23	2.39	0.47
1:A:12:PHE:HA	1:A:15:ASP:OD2	2.15	0.47
1:A:468:THR:HG22	1:A:469:LEU:N	2.30	0.47
1:A:141:THR:HG21	1:A:157:LEU:HB2	1.98	0.46
1:A:642:VAL:HB	1:A:643:PRO:HD2	1.97	0.46
1:A:121:LEU:HD12	1:A:121:LEU:C	2.39	0.46
1:A:505:GLY:HA2	1:A:506:PRO:C	2.40	0.46
1:A:34:PRO:HG2	1:A:49:TYR:CG	2.51	0.46
1:A:273:ALA:HB2	1:A:280:LEU:HD12	1.97	0.46
1:A:47:ARG:HH11	1:A:47:ARG:HD2	1.59	0.46
1:A:227:ARG:NH1	1:A:257:GLU:HB2	2.31	0.46
1:A:149:PRO:HA	1:A:154:ASN:ND2	2.30	0.46
1:A:617:ASP:OD2	1:A:620:LYS:HE2	2.15	0.46
1:A:325:ASP:HB2	4:A:1164:HOH:O	2.16	0.46
1:A:168:THR:O	1:A:169:ASN:HB2	2.16	0.45
1:A:413:TRP:CE2	2:B:1:GLC:H2	2.51	0.45
1:A:651:LYS:HE3	1:A:663:GLU:O	2.16	0.45
1:A:45:ASN:CG	1:A:48:LEU:HD22	2.41	0.45
1:A:87:ILE:HG21	1:A:89:TYR:CZ	2.52	0.45
1:A:445:SER:CB	1:A:479:ASN:ND2	2.80	0.45
1:A:582:LEU:HD23	1:A:582:LEU:HA	1.67	0.44
1:A:511:PRO:HB3	1:A:555:ALA:HA	1.99	0.44
1:A:627:ASN:ND2	2:D:2:GLC:H61	2.32	0.44
1:A:19:GLN:HG3	1:A:75:TRP:CD2	2.53	0.44
1:A:23:ASP:CG	1:A:50:CYS:H	2.25	0.44
1:A:60:LYS:HA	1:A:60:LYS:HD2	1.28	0.44
1:A:444:ILE:HD12	1:A:482:LEU:HB2	1.99	0.44
1:A:215:ILE:HD12	1:A:215:ILE:HA	1.83	0.44
1:A:266:SER:O	1:A:269:ASN:HB3	2.17	0.44
1:A:148:GLN:N	1:A:149:PRO:CD	2.79	0.44
1:A:607:THR:HG22	1:A:653:LEU:HD21	2.00	0.43
1:A:283:PHE:O	1:A:287:GLN:HG2	2.18	0.43
1:A:397:LEU:HA	1:A:400:CYS:SG	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:617:ASP:HB3	1:A:620:LYS:HE2	1.99	0.43
1:A:80:VAL:HA	1:A:107:LYS:O	2.18	0.43
1:A:412:ARG:O	2:B:2:GLC:H2	2.18	0.43
1:A:633:TYR:CD2	1:A:634:PRO:HA	2.53	0.43
1:A:89:TYR:O	1:A:90:SER:HB2	2.19	0.43
1:A:300:MET:HB2	1:A:415:ASN:O	2.18	0.43
1:A:147:ASP:C	1:A:149:PRO:HD3	2.44	0.43
1:A:290:ARG:O	1:A:294:ARG:HB3	2.18	0.43
1:A:468:THR:CG2	1:A:469:LEU:N	2.82	0.42
1:A:566:ALA:HA	1:A:571:THR:O	2.18	0.42
1:A:19:GLN:O	1:A:360:TYR:HB3	2.19	0.42
1:A:148:GLN:N	1:A:149:PRO:HD3	2.34	0.42
1:A:669:THR:CG2	1:A:670:PHE:N	2.82	0.42
1:A:320:GLN:HB3	4:A:1166:HOH:O	2.20	0.42
1:A:527:GLY:HA3	1:A:566:ALA:O	2.20	0.42
1:A:18:TYR:CB	1:A:71:VAL:HG11	2.49	0.42
1:A:605:TYR:CE1	1:A:655:LYS:HB2	2.55	0.42
1:A:143:PRO:HB3	1:A:196:ASP:OD2	2.19	0.42
1:A:249:TYR:CE2	1:A:250:LYS:HD2	2.54	0.42
1:A:643:PRO:HB2	1:A:646:LYS:HG3	2.02	0.42
1:A:87:ILE:HD13	1:A:143:PRO:HG2	2.01	0.42
1:A:187:GLU:CD	1:A:628:GLN:HB2	2.45	0.42
1:A:583:SER:HB2	1:A:612:GLU:OE1	2.20	0.42
1:A:133:ILE:CG2	1:A:134:ILE:N	2.83	0.41
1:A:290:ARG:NH2	1:A:330:GLU:O	2.53	0.41
1:A:361:GLY:HA3	1:A:366:MET:SD	2.60	0.41
1:A:413:TRP:CD1	2:B:1:GLC:H4	2.56	0.41
1:A:417:ASP:O	1:A:436:ARG:HG3	2.19	0.41
1:A:293:PHE:O	1:A:341:LYS:HD2	2.20	0.41
1:A:119:GLN:HE21	1:A:119:GLN:HB2	1.41	0.41
1:A:594:ASN:HB2	1:A:683:ASN:OD1	2.21	0.41
1:A:371:ASP:OD1	1:A:372:PRO:HA	2.21	0.41
1:A:116:ALA:HA	1:A:119:GLN:NE2	2.36	0.41
1:A:88:ASN:ND2	1:A:91:GLY:N	2.60	0.41
1:A:471:VAL:CG1	1:A:475:GLY:HA2	2.51	0.40
1:A:684:TRP:CD1	1:A:684:TRP:C	2.99	0.40
1:A:108:THR:O	1:A:110:PRO:HD3	2.21	0.40
1:A:258:TRP:CD1	1:A:258:TRP:C	2.99	0.40
1:A:308:GLU:HG2	1:A:308:GLU:H	1.77	0.40
1:A:664:GLY:O	1:A:685:GLN:HB2	2.21	0.40
1:A:340:ARG:HA	1:A:343:GLU:OE1	2.22	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	684/686 (100%)	647 (95%)	36 (5%)	1 (0%)	48 68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	629	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	556/556 (100%)	518 (93%)	38 (7%)	14 31

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	21	PHE
1	A	29	ASN
1	A	48	LEU
1	A	80	VAL
1	A	86	ILE
1	A	88	ASN

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Mol	Chain	Res	Type
1	A	92	VAL
1	A	103	ARG
1	A	107	LYS
1	A	115	ILE
1	A	128	LYS
1	A	146	SER
1	A	163	LEU
1	A	171	THR
1	A	172	GLN
1	A	173	ASN
1	A	227	ARG
1	A	308	GLU
1	A	318	ASP
1	A	330	GLU
1	A	341	LYS
1	A	375	ARG
1	A	380	SER
1	A	412	ARG
1	A	420	ILE
1	A	424	LYS
1	A	436	ARG
1	A	443	SER
1	A	455	SER
1	A	583	SER
1	A	595	ASN
1	A	613	LEU
1	A	653	LEU
1	A	660	VAL
1	A	670	PHE
1	A	680	ILE
1	A	682	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	88	ASN
1	A	119	GLN
1	A	120	ASN
1	A	169	ASN
1	A	269	ASN
1	A	392	GLN

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Mol	Chain	Res	Type
1	A	416	ASN
1	A	479	ASN
1	A	560	ASN
1	A	632	GLN
1	A	685	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 ⓘ

6 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	GLC	B	1	2	12,12,12	0.46	0	17,17,17	1.07	1 (5%)
2	GLC	B	2	2	11,11,12	0.43	0	15,15,17	1.16	3 (20%)
2	GLC	C	1	2	12,12,12	0.32	0	17,17,17	1.17	1 (5%)
2	GLC	C	2	2	11,11,12	0.50	0	15,15,17	1.04	1 (6%)
2	GLC	D	1	2	12,12,12	0.46	0	17,17,17	1.64	4 (23%)
2	GLC	D	2	2	11,11,12	0.24	0	15,15,17	1.72	3 (20%)

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	B	1	2	-	0/2/22/22	0/1/1/1
2	GLC	B	2	2	-	2/2/19/22	0/1/1/1
2	GLC	C	1	2	-	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	GLC	C1-O5-C5	4.93	118.79	112.19
2	D	1	GLC	O4-C4-C3	-3.45	102.24	110.38
2	D	1	GLC	O5-C1-C2	3.42	116.31	110.30
2	C	1	GLC	O5-C5-C4	-3.04	104.21	109.70
2	B	1	GLC	O5-C5-C4	-2.60	105.01	109.70
2	D	2	GLC	O5-C1-C2	2.58	116.95	110.79
2	D	2	GLC	O2-C2-C1	-2.55	103.39	109.22
2	D	1	GLC	C1-O5-C5	2.46	118.41	113.65
2	B	2	GLC	O4-C4-C5	-2.45	103.28	109.32
2	B	2	GLC	C1-O5-C5	2.24	115.19	112.19
2	C	2	GLC	O5-C5-C6	-2.18	103.41	107.66
2	B	2	GLC	O2-C2-C1	2.15	114.15	109.22
2	D	1	GLC	O4-C4-C5	-2.03	104.32	109.32

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	GLC	O5-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	D	2	GLC	O5-C5-C6-O6
2	D	2	GLC	C4-C5-C6-O6
2	B	2	GLC	C4-C5-C6-O6
2	B	2	GLC	O5-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 8 short contacts:

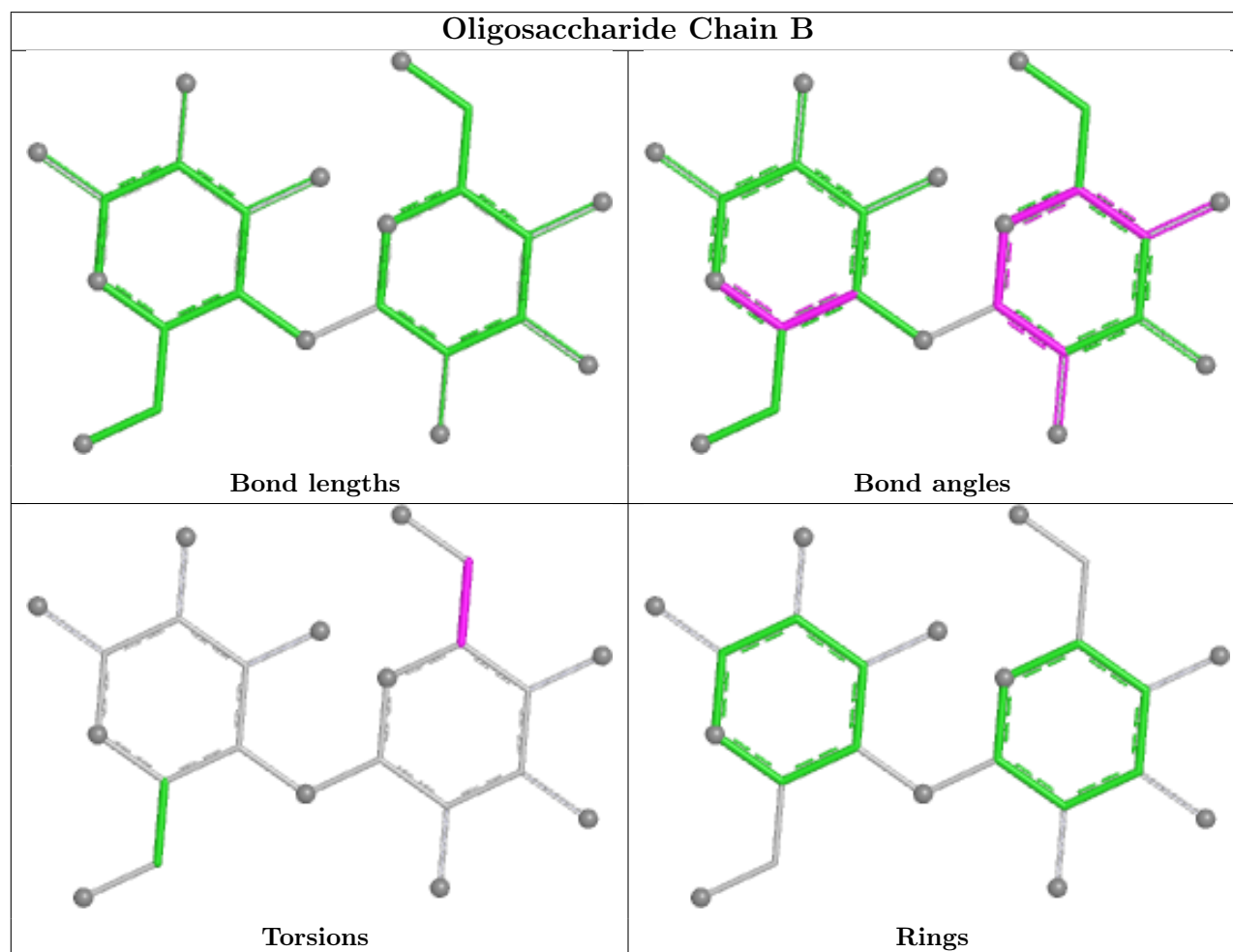
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	GLC	2	0

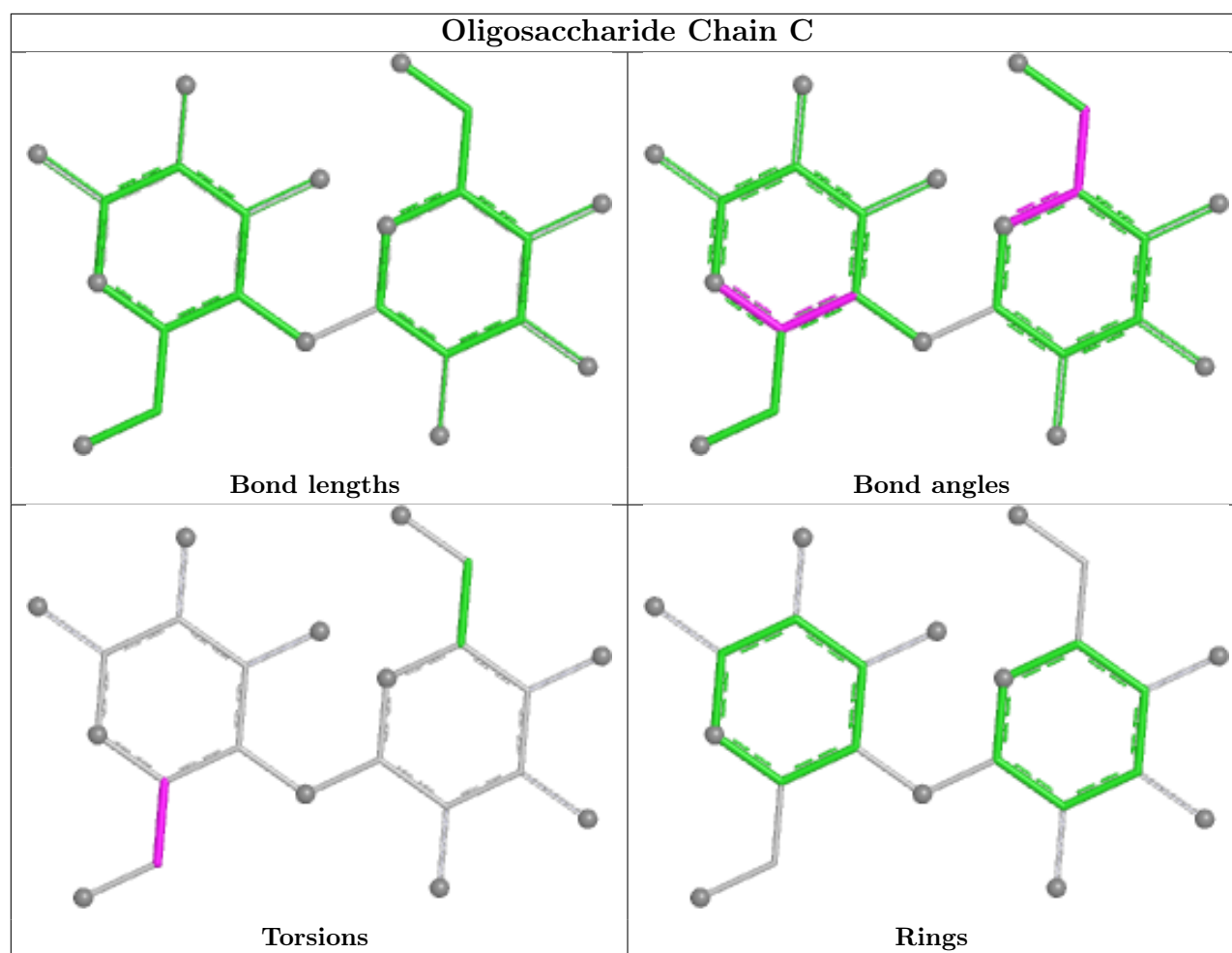
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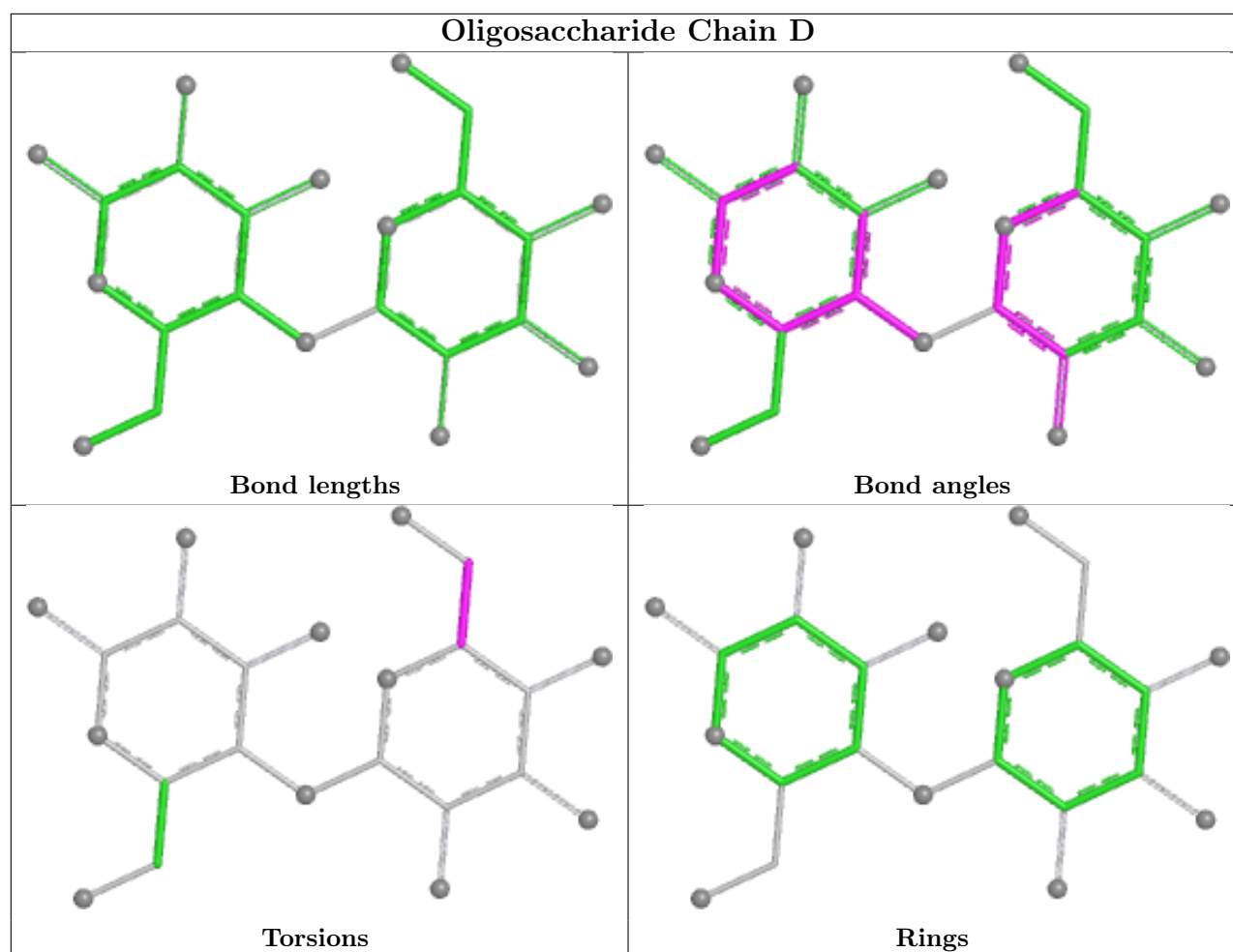
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	GLC	3	0
2	D	2	GLC	5	0
2	B	2	GLC	1	0

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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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.