



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:18 AM UTC

PDB ID : 2D2O / pdb_00002d2o
Title : Structure of a complex of Thermoactinomyces vulgaris R-47 alpha-amylase 2 with maltohexaose demonstrates the important role of aromatic residues at the reducing end of the substrate binding cleft
Authors : Ohtaki, A.; Mizuno, M.; Yoshida, H.; Tonozuka, T.; Sakano, Y.; Kamitori, S.
Deposited on : 2005-09-13
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

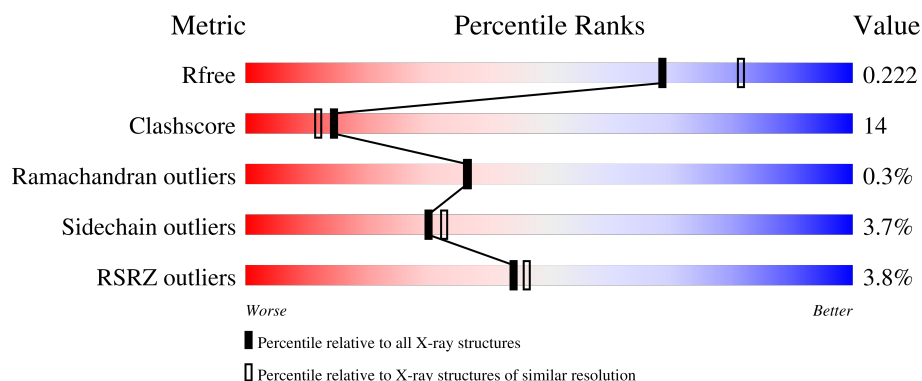
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	585	3% 74% 24% .
1	B	585	4% 73% 24% .
2	C	6	50% 50%
2	D	6	33% 17% 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	GLC	C	1	X	-	-	-
2	GLC	D	1	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10668 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 Neopullulanase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	585	Total	C	N	O	S	0	0	0
			4776	3056	832	873	15			
1	B	585	Total	C	N	O	S	0	0	0
			4776	3056	832	873	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	325	ASN	ASP	engineered mutation	UNP Q08751
B	325	ASN	ASP	engineered mutation	UNP Q08751

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	6	Total	C	O	0	0	0
			67	36	31			
2	D	6	Total	C	O	0	0	0
			67	36	31			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		

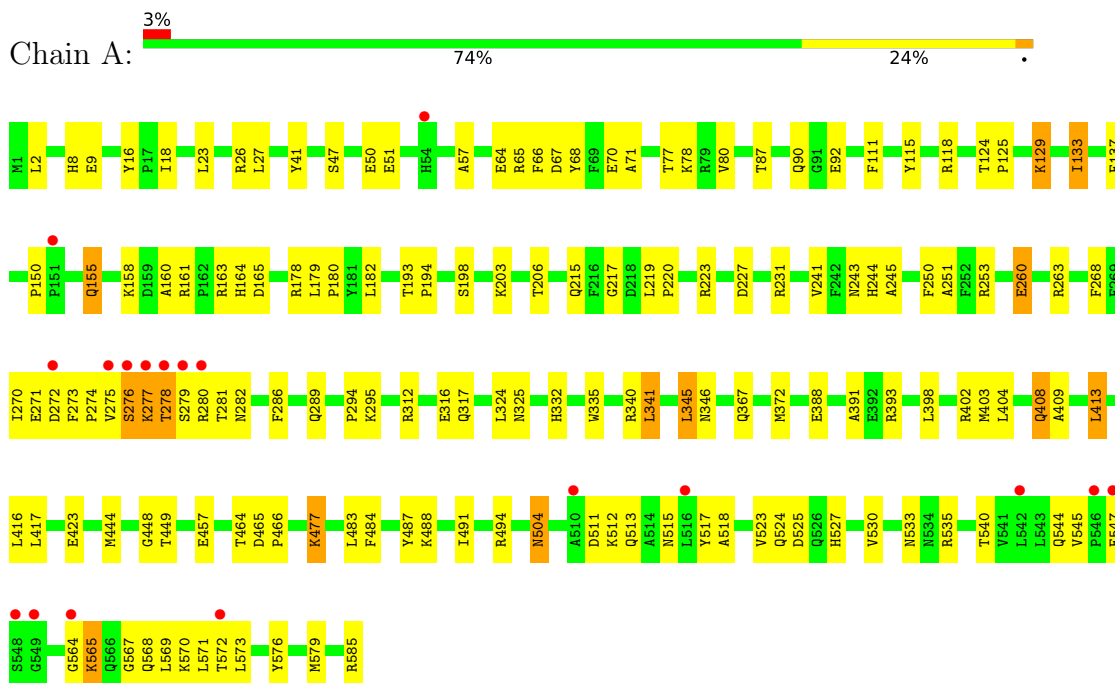
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	498	Total 498	O 498	0	0
4	B	482	Total 482	O 482	0	0

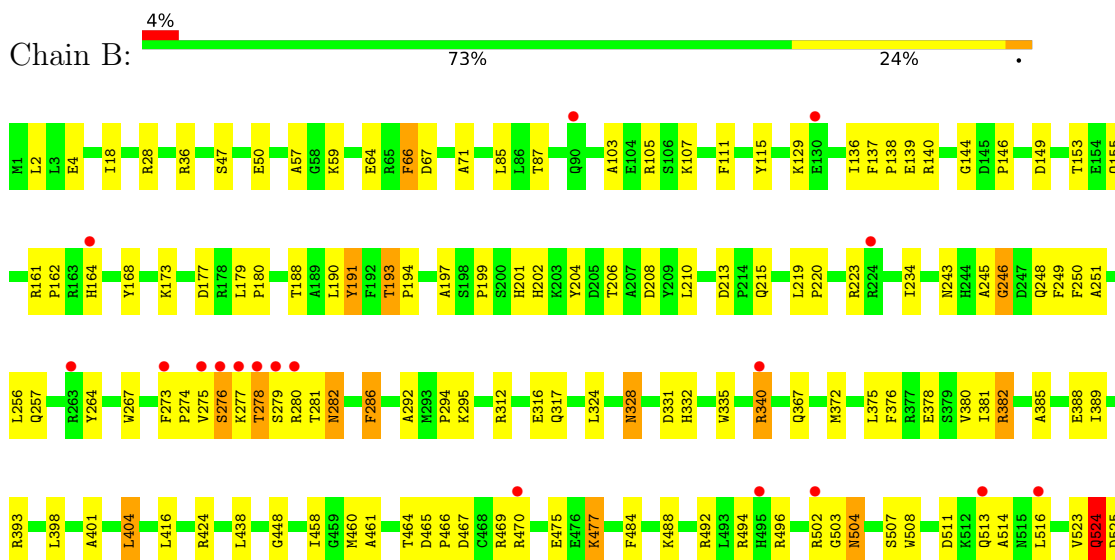
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Neopullulanase 2



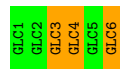
• Molecule 1: Neopullulanase 2





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

Chain C: 50% 50%



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

Chain D: 33% 17% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	113.69Å 118.71Å 112.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.00 – 2.10 46.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (46.00-2.10) 95.5 (46.00-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 1.98Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.184 , 0.217 0.191 , 0.222	Depositor DCC
R_{free} test set	6518 reflections (6.58%)	wwPDB-VP
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.329	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10668	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.37	0/4906	0.88	13/6641 (0.2%)
1	B	0.36	0/4906	0.89	18/6641 (0.3%)
All	All	0.37	0/9812	0.89	31/13282 (0.2%)

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	18	ILE	N-CA-C	-7.46	105.30	113.43
1	B	18	ILE	N-CA-C	-6.73	106.67	113.47
1	B	193	THR	N-CA-C	-6.65	100.36	110.07
1	B	66	PHE	N-CA-C	6.24	118.89	109.41
1	A	372	MET	N-CA-C	-6.23	100.17	109.15
1	B	67	ASP	N-CA-C	-6.14	100.19	109.95
1	B	161	ARG	CA-C-N	6.02	126.04	119.90
1	B	161	ARG	C-N-CA	6.02	126.04	119.90
1	A	66	PHE	N-CA-C	5.95	118.92	109.81
1	B	276	SER	N-CA-C	5.89	117.21	107.73
1	B	372	MET	N-CA-C	-5.86	100.71	109.15
1	A	206	THR	N-CA-C	5.85	118.18	109.59
1	A	137	PHE	N-CA-C	-5.83	96.92	109.81
1	A	524	GLN	CB-CA-C	-5.77	109.94	116.63
1	A	67	ASP	N-CA-C	-5.68	100.91	109.95
1	B	103	ALA	N-CA-C	-5.62	105.92	112.89
1	A	423	GLU	N-CA-C	-5.56	103.19	110.53
1	B	375	LEU	N-CA-C	-5.56	105.30	111.36
1	B	524	GLN	CB-CA-C	-5.53	108.60	116.34
1	B	206	THR	N-CA-C	5.50	117.67	109.59
1	B	137	PHE	CA-C-N	5.46	126.66	119.84
1	B	137	PHE	C-N-CA	5.46	126.66	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	111	PHE	N-CA-C	-5.45	102.81	110.50
1	A	80	VAL	N-CA-C	5.42	116.47	108.45
1	B	286	PHE	N-CA-C	-5.41	102.99	110.35
1	A	111	PHE	N-CA-C	-5.41	102.48	110.48
1	B	331	ASP	N-CA-C	5.30	117.56	110.35
1	B	246	GLY	N-CA-C	-5.13	104.87	113.02
1	A	276	SER	N-CA-C	5.10	121.66	110.80
1	A	41	TYR	N-CA-C	5.04	116.66	109.14
1	A	417	LEU	N-CA-C	-5.01	107.30	113.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4776	0	4609	136	0
1	B	4776	0	4609	133	0
2	C	67	0	57	5	0
2	D	67	0	57	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	498	0	0	15	0
4	B	482	0	0	15	0
All	All	10668	0	9332	266	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 14.

All (266) 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:2:LEU:HD21	1:B:2:LEU:HD21	1.25	1.14
1:A:129:LYS:H	1:A:129:LYS:HD2	1.24	1.02
1:A:525:ASP:HB3	1:A:585:ARG:HD3	1.49	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:THR:HG22	1:A:279:SER:H	1.35	0.91
1:A:544:GLN:HE22	1:A:568:GLN:HE21	1.20	0.90
1:A:408:GLN:H	1:A:408:GLN:HE21	1.16	0.89
1:B:511:ASP:OD2	1:B:513:GLN:HG2	1.71	0.89
1:A:129:LYS:HD2	1:A:129:LYS:N	1.91	0.83
1:B:28:ARG:HD3	1:B:66:PHE:CD2	2.12	0.83
1:B:552:THR:HB	4:B:1045:HOH:O	1.77	0.83
1:A:565:LYS:H	1:A:565:LYS:HD3	1.44	0.82
1:A:544:GLN:NE2	1:A:568:GLN:HE21	1.79	0.80
1:B:213:ASP:OD1	1:B:215:GLN:HG2	1.83	0.79
1:A:544:GLN:HE22	1:A:568:GLN:NE2	1.80	0.78
1:A:281:THR:H	1:A:289:GLN:HE22	1.32	0.76
1:B:129:LYS:HD2	1:B:502:ARG:NH2	2.00	0.76
1:B:533:ASN:HD21	1:B:539:GLN:HE21	1.31	0.75
1:A:2:LEU:CD2	1:B:2:LEU:HD21	2.13	0.75
1:A:57:ALA:HB2	1:A:71:ALA:HB2	1.68	0.75
1:A:271:GLU:HG2	1:A:282:ASN:O	1.86	0.75
1:B:202:HIS:HD2	1:B:204:TYR:H	1.35	0.74
1:A:223:ARG:HD3	1:A:317:GLN:HE21	1.53	0.74
1:A:408:GLN:H	1:A:408:GLN:NE2	1.87	0.73
1:A:477:LYS:HE3	1:A:477:LYS:HA	1.71	0.72
1:B:585:ARG:HB3	1:B:585:ARG:HH11	1.52	0.72
1:B:140:ARG:HH12	1:B:201:HIS:HB2	1.55	0.71
1:A:26:ARG:HD3	1:A:70:GLU:OE2	1.91	0.70
1:A:243:ASN:HD22	1:A:244:HIS:HD2	1.39	0.70
1:B:140:ARG:HG2	1:B:469:ARG:O	1.91	0.70
1:A:125:PRO:O	1:A:129:LYS:HE2	1.92	0.70
1:B:475:GLU:H	1:B:475:GLU:CD	2.00	0.70
1:B:484:PHE:CE1	1:B:488:LYS:HD2	2.28	0.68
1:A:2:LEU:HD21	1:B:2:LEU:CD2	2.14	0.68
1:B:243:ASN:HD21	1:B:295:LYS:NZ	1.91	0.68
1:B:542:LEU:HD21	1:B:570:LYS:HG2	1.76	0.67
1:A:312:ARG:O	1:A:316:GLU:HG3	1.94	0.67
1:B:278:THR:HG23	1:B:280:ARG:H	1.59	0.67
1:A:515:ASN:HB2	1:A:535:ARG:NH2	2.10	0.67
1:B:202:HIS:CD2	1:B:204:TYR:H	2.12	0.67
1:B:328:ASN:N	1:B:328:ASN:HD22	1.93	0.67
1:A:504:ASN:C	1:A:504:ASN:HD22	2.00	0.67
1:B:188:THR:HG21	4:B:1454:HOH:O	1.95	0.66
1:B:504:ASN:C	1:B:504:ASN:HD22	2.04	0.66
1:A:409:ALA:O	1:A:413:LEU:HD13	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:SER:HB3	1:B:50:GLU:HG3	1.77	0.66
1:A:341:LEU:HD22	1:A:345:LEU:HD22	1.77	0.65
1:A:540:THR:OG1	1:A:572:THR:HG22	1.97	0.64
1:B:256:LEU:HD23	1:B:275:VAL:HB	1.79	0.64
1:B:565:LYS:HE2	1:B:568:GLN:HB2	1.78	0.64
1:B:107:LYS:HG3	4:B:1399:HOH:O	1.98	0.64
1:A:133:ILE:HD12	1:A:133:ILE:N	2.12	0.64
2:C:3:GLC:H62	2:C:4:GLC:H5	1.80	0.63
1:A:388:GLU:OE1	1:B:107:LYS:HE3	1.98	0.63
1:A:525:ASP:HB2	4:A:1289:HOH:O	1.98	0.63
1:B:275:VAL:O	1:B:282:ASN:ND2	2.32	0.63
1:A:340:ARG:NH1	4:A:1365:HOH:O	2.31	0.62
1:A:465:ASP:OD2	2:C:6:GLC:O2	2.16	0.61
1:B:448:GLY:O	1:B:494:ARG:NH2	2.32	0.61
1:B:279:SER:HB3	4:B:1391:HOH:O	2.01	0.61
4:B:1204:HOH:O	2:D:6:GLC:H61	2.00	0.61
1:A:8:HIS:HD2	1:A:26:ARG:O	1.84	0.61
1:B:340:ARG:NH2	4:B:1415:HOH:O	2.30	0.61
1:B:328:ASN:HD22	1:B:328:ASN:H	1.49	0.60
1:B:340:ARG:HH21	1:B:340:ARG:HB3	1.66	0.60
1:A:391:ALA:H	1:A:515:ASN:ND2	2.00	0.60
1:B:585:ARG:HH11	1:B:585:ARG:CB	2.15	0.60
2:D:3:GLC:H62	2:D:4:GLC:H5	1.83	0.60
1:B:282:ASN:C	1:B:282:ASN:HD22	2.10	0.60
1:B:488:LYS:O	1:B:492:ARG:HG3	2.01	0.60
1:A:276:SER:O	1:A:277:LYS:HB2	2.02	0.59
1:B:173:LYS:HD3	1:B:177:ASP:OD2	2.02	0.59
1:A:408:GLN:HE21	1:A:408:GLN:N	1.96	0.58
1:B:467:ASP:OD1	1:B:470:ARG:HD3	2.03	0.58
1:A:260:GLU:HG3	1:A:273:PHE:CE2	2.38	0.58
1:A:564:GLY:HA2	1:A:568:GLN:O	2.03	0.58
1:B:514:ALA:HB1	1:B:516:LEU:HD13	1.85	0.57
1:A:219:LEU:HB3	1:A:220:PRO:HD3	1.87	0.57
1:A:504:ASN:C	1:A:504:ASN:ND2	2.61	0.57
1:B:382:ARG:HH21	1:B:382:ARG:HB2	1.69	0.57
4:A:1149:HOH:O	2:C:6:GLC:H61	2.03	0.57
1:B:525:ASP:HB3	1:B:585:ARG:HD3	1.86	0.57
1:B:57:ALA:HB2	1:B:71:ALA:HB2	1.86	0.56
1:B:164:HIS:CE1	1:B:466:PRO:HD3	2.41	0.56
1:B:547:GLU:HB2	4:B:1369:HOH:O	2.05	0.56
1:A:124:THR:OG1	1:A:129:LYS:HE3	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:332:HIS:HD2	1:B:367:GLN:OE1	1.89	0.56
1:A:332:HIS:HD2	1:A:367:GLN:OE1	1.90	0.55
1:A:565:LYS:HD3	1:A:565:LYS:N	2.18	0.55
1:A:241:VAL:HG13	1:A:325:ASN:HD22	1.72	0.55
1:A:545:VAL:HA	1:A:569:LEU:HD21	1.89	0.55
1:A:133:ILE:HD13	1:A:449:THR:CG2	2.37	0.55
1:A:160:ALA:O	1:A:161:ARG:HD2	2.06	0.54
1:B:328:ASN:H	1:B:328:ASN:ND2	2.03	0.54
1:A:164:HIS:CE1	2:C:6:GLC:H2	2.42	0.54
1:B:416:LEU:HD23	1:B:416:LEU:H	1.72	0.53
1:A:90:GLN:HA	1:A:90:GLN:HE21	1.74	0.53
1:A:115:TYR:CZ	1:B:295:LYS:HE2	2.43	0.53
1:A:26:ARG:HG2	1:A:70:GLU:HG3	1.91	0.53
1:B:382:ARG:HH21	1:B:382:ARG:CB	2.22	0.52
1:A:193:THR:HB	1:A:194:PRO:CD	2.40	0.52
1:A:271:GLU:HG3	1:A:272:ASP:OD2	2.10	0.52
1:B:136:ILE:O	1:B:138:PRO:HD3	2.08	0.52
1:B:470:ARG:HG3	4:B:1112:HOH:O	2.09	0.52
1:A:545:VAL:HA	1:A:569:LEU:CD2	2.39	0.52
1:B:153:THR:HG23	1:B:168:TYR:HA	1.90	0.52
1:A:448:GLY:O	1:A:494:ARG:NH2	2.41	0.52
1:B:28:ARG:HD3	1:B:66:PHE:CE2	2.45	0.52
1:B:105:ARG:HH11	1:B:105:ARG:HG2	1.74	0.52
1:B:492:ARG:O	1:B:496:ARG:HB2	2.09	0.52
1:B:542:LEU:HD12	1:B:568:GLN:HB3	1.90	0.52
1:B:282:ASN:ND2	1:B:282:ASN:C	2.67	0.52
1:A:515:ASN:HB2	1:A:535:ARG:HH21	1.74	0.52
1:A:268:PHE:HB2	1:A:270:ILE:CD1	2.39	0.52
1:B:273:PHE:HA	1:B:274:PRO:C	2.34	0.52
1:B:504:ASN:C	1:B:504:ASN:ND2	2.61	0.52
1:A:118:ARG:HG2	4:B:1007:HOH:O	2.10	0.51
1:A:295:LYS:HE2	1:B:115:TYR:CZ	2.46	0.51
1:B:275:VAL:O	1:B:276:SER:HB3	2.10	0.51
1:B:378:GLU:HG3	4:B:1362:HOH:O	2.09	0.51
1:A:158:LYS:HD2	1:A:158:LYS:O	2.11	0.51
1:B:64:GLU:OE1	1:B:393:ARG:HD3	2.11	0.51
1:A:47:SER:HB3	1:A:50:GLU:HG3	1.92	0.51
1:B:164:HIS:CE1	2:D:6:GLC:H2	2.46	0.51
1:A:150:PRO:HG3	1:A:215:GLN:CD	2.36	0.50
1:A:484:PHE:CE1	1:A:488:LYS:HD2	2.46	0.50
1:B:524:GLN:HB3	4:B:1430:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:546:PRO:HG2	1:B:553:TRP:CH2	2.47	0.50
1:A:276:SER:OG	1:A:277:LYS:N	2.44	0.50
1:A:540:THR:HA	1:A:571:LEU:O	2.11	0.50
1:B:219:LEU:HB3	1:B:220:PRO:HD3	1.93	0.50
1:A:281:THR:H	1:A:289:GLN:NE2	2.06	0.50
1:B:508:TRP:CZ2	1:B:546:PRO:HD3	2.47	0.50
1:B:542:LEU:CD2	1:B:570:LYS:HG2	2.42	0.49
1:A:227:ASP:O	1:A:231:ARG:HG2	2.12	0.49
1:B:139:GLU:OE2	1:B:140:ARG:NH1	2.45	0.49
1:B:507:SER:HB3	4:B:1120:HOH:O	2.13	0.49
1:A:133:ILE:HD13	1:A:449:THR:HG22	1.93	0.49
1:A:253:ARG:HG2	1:A:253:ARG:HH21	1.78	0.49
1:B:210:LEU:N	1:B:210:LEU:HD22	2.28	0.49
1:B:542:LEU:HD11	1:B:565:LYS:HZ3	1.76	0.49
1:A:65:ARG:HH21	1:A:65:ARG:HG2	1.78	0.49
1:A:178:ARG:NH2	4:A:1156:HOH:O	2.45	0.49
1:A:278:THR:HG22	1:A:279:SER:N	2.14	0.49
1:A:393:ARG:HH21	1:A:393:ARG:HG3	1.76	0.49
1:A:512:LYS:HG3	1:A:513:GLN:N	2.27	0.49
1:B:146:PRO:HA	1:B:149:ASP:OD1	2.12	0.49
1:A:215:GLN:HG3	4:A:1300:HOH:O	2.12	0.48
1:A:488:LYS:HB3	4:A:1385:HOH:O	2.13	0.48
1:B:278:THR:OG1	1:B:279:SER:N	2.46	0.48
1:B:551:LYS:HE3	4:B:1045:HOH:O	2.12	0.48
1:A:517:TYR:OH	4:A:1004:HOH:O	2.16	0.48
1:A:545:VAL:HG23	1:A:567:GLY:C	2.39	0.48
1:B:477:LYS:NZ	1:B:477:LYS:HB3	2.28	0.48
1:A:268:PHE:HB2	1:A:270:ILE:HD11	1.95	0.47
1:A:565:LYS:H	1:A:565:LYS:CD	2.20	0.47
1:A:27:LEU:C	1:A:27:LEU:HD23	2.39	0.47
1:A:155:GLN:HE21	1:A:155:GLN:HB3	1.52	0.47
1:A:263:ARG:HD3	4:A:1119:HOH:O	2.14	0.47
1:B:245:ALA:O	1:B:294:PRO:HD2	2.15	0.47
1:B:243:ASN:HD21	1:B:295:LYS:HZ2	1.62	0.47
1:B:199:PRO:HG3	1:B:248:GLN:NE2	2.30	0.47
1:B:424:ARG:HG3	1:B:461:ALA:C	2.39	0.47
1:A:87:THR:HA	1:A:92:GLU:O	2.16	0.46
1:A:324:LEU:HD13	1:A:335:TRP:CZ3	2.50	0.46
1:B:565:LYS:HD2	1:B:565:LYS:O	2.15	0.46
1:B:312:ARG:O	1:B:316:GLU:HG3	2.14	0.46
1:B:401:ALA:O	1:B:404:LEU:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:THR:OG1	1:A:289:GLN:NE2	2.49	0.46
1:B:144:GLY:C	1:B:173:LYS:HD2	2.41	0.46
1:A:64:GLU:HB2	4:A:1313:HOH:O	2.15	0.46
1:A:346:ASN:HD22	1:A:346:ASN:C	2.23	0.46
1:B:191:TYR:CD1	1:B:191:TYR:C	2.93	0.46
1:B:340:ARG:NH2	1:B:340:ARG:HB3	2.30	0.46
1:A:279:SER:O	1:A:289:GLN:NE2	2.49	0.46
1:A:51:GLU:OE1	1:A:51:GLU:HA	2.16	0.46
1:A:565:LYS:N	1:A:565:LYS:CD	2.79	0.45
1:B:376:PHE:O	1:B:380:VAL:HG23	2.17	0.45
1:A:57:ALA:CB	1:A:71:ALA:HB2	2.40	0.45
1:A:286:PHE:HZ	2:C:3:GLC:H61	1.80	0.45
1:A:16:TYR:C	1:A:23:LEU:HD12	2.40	0.45
1:B:140:ARG:NH1	1:B:201:HIS:HB2	2.27	0.45
1:A:163:ARG:NH1	1:A:165:ASP:OD1	2.49	0.45
1:A:272:ASP:O	1:A:282:ASN:ND2	2.50	0.45
1:B:551:LYS:HD2	1:B:551:LYS:O	2.17	0.45
1:A:244:HIS:CD2	1:A:286:PHE:HB2	2.51	0.45
1:A:273:PHE:HA	1:A:274:PRO:C	2.41	0.45
1:B:276:SER:O	1:B:277:LYS:HD2	2.17	0.45
1:A:564:GLY:HA3	1:A:569:LEU:HD13	1.99	0.45
1:B:146:PRO:HA	1:B:149:ASP:CG	2.41	0.45
1:B:164:HIS:ND1	4:B:1167:HOH:O	2.36	0.45
1:A:565:LYS:NZ	1:A:570:LYS:HE3	2.32	0.45
1:B:467:ASP:O	1:B:470:ARG:HG3	2.17	0.45
1:B:250:PHE:CG	1:B:251:ALA:N	2.85	0.44
1:B:324:LEU:HD13	1:B:335:TRP:CZ3	2.52	0.44
1:B:464:THR:O	1:B:465:ASP:C	2.61	0.44
1:A:402:ARG:HD2	4:A:1261:HOH:O	2.18	0.44
1:B:458:ILE:CD1	1:B:460:MET:HG3	2.47	0.44
1:A:64:GLU:HG3	4:A:1332:HOH:O	2.17	0.44
1:A:164:HIS:CE1	1:A:466:PRO:HD3	2.52	0.44
1:A:416:LEU:H	1:A:416:LEU:HD23	1.81	0.44
1:B:524:GLN:HB3	1:B:525:ASP:H	1.48	0.44
1:B:388:GLU:HB2	4:B:1187:HOH:O	2.17	0.44
1:B:542:LEU:HD11	1:B:565:LYS:NZ	2.32	0.44
1:A:124:THR:OG1	1:A:129:LYS:CE	2.66	0.44
1:B:193:THR:HB	1:B:194:PRO:CD	2.47	0.43
1:A:564:GLY:CA	1:A:568:GLN:O	2.67	0.43
1:A:281:THR:N	1:A:289:GLN:HE22	2.10	0.43
1:A:527:HIS:HD2	4:A:1315:HOH:O	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:533:ASN:O	1:A:576:TYR:HA	2.17	0.43
1:B:249:PHE:CE2	1:B:251:ALA:HB3	2.53	0.43
1:B:475:GLU:CD	1:B:475:GLU:N	2.74	0.43
1:A:65:ARG:HB3	1:B:4:GLU:HG3	1.99	0.43
1:B:105:ARG:HG2	1:B:105:ARG:NH1	2.34	0.43
1:A:245:ALA:O	1:A:294:PRO:HD2	2.19	0.43
1:B:197:ALA:HB3	1:B:208:ASP:HB3	2.01	0.43
1:A:133:ILE:CD1	1:A:449:THR:CG2	2.97	0.43
1:A:518:ALA:HA	1:A:530:VAL:O	2.19	0.43
1:B:243:ASN:ND2	1:B:295:LYS:NZ	2.64	0.43
1:B:275:VAL:O	1:B:275:VAL:HG12	2.18	0.42
1:A:275:VAL:HG12	1:A:276:SER:N	2.34	0.42
1:B:416:LEU:H	1:B:416:LEU:CD2	2.32	0.42
1:B:565:LYS:HD2	1:B:565:LYS:C	2.44	0.42
1:A:270:ILE:N	1:A:270:ILE:HD12	2.34	0.42
1:A:9:GLU:HA	4:A:1205:HOH:O	2.18	0.42
1:A:569:LEU:N	1:A:569:LEU:HD22	2.35	0.42
1:B:190:LEU:HG	1:B:234:ILE:CG2	2.50	0.42
1:B:193:THR:HB	1:B:194:PRO:HD2	2.02	0.42
1:B:516:LEU:HD11	1:B:539:GLN:NE2	2.33	0.42
1:B:162:PRO:HG2	1:B:470:ARG:HA	2.00	0.42
1:B:546:PRO:CG	1:B:553:TRP:CH2	3.02	0.42
1:A:444:MET:CE	1:A:491:ILE:HG12	2.49	0.42
1:A:523:VAL:O	1:A:523:VAL:HG13	2.20	0.42
1:A:444:MET:HE2	1:A:491:ILE:CG1	2.49	0.42
1:B:286:PHE:HZ	2:D:3:GLC:H61	1.83	0.42
1:B:381:ILE:O	1:B:385:ALA:HB3	2.20	0.42
1:A:77:THR:O	1:A:78:LYS:HB2	2.20	0.42
1:A:179:LEU:N	1:A:180:PRO:CD	2.83	0.41
1:A:133:ILE:N	1:A:133:ILE:CD1	2.80	0.41
1:A:457:GLU:HA	1:A:487:TYR:CE1	2.55	0.41
1:B:577:GLN:HG2	1:B:578:GLY:N	2.36	0.41
1:A:281:THR:HG22	4:A:1419:HOH:O	2.20	0.41
1:B:36:ARG:HB3	1:B:87:THR:HB	2.03	0.41
1:B:264:TYR:O	1:B:267:TRP:HB2	2.20	0.41
1:A:217:GLY:HA2	4:A:1015:HOH:O	2.19	0.41
1:B:465:ASP:HA	1:B:466:PRO:HA	1.88	0.41
1:B:546:PRO:HG2	1:B:553:TRP:CZ2	2.55	0.41
1:A:68:TYR:CD2	1:A:403:MET:HG3	2.55	0.41
1:A:250:PHE:CG	1:A:251:ALA:N	2.88	0.41
1:B:279:SER:O	1:B:281:THR:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:523:VAL:O	1:B:523:VAL:HG13	2.20	0.41
1:A:444:MET:HE2	1:A:491:ILE:HG12	2.02	0.41
1:B:223:ARG:CZ	1:B:317:GLN:OE1	2.68	0.41
1:B:382:ARG:HG3	1:B:388:GLU:HB3	2.02	0.41
1:B:541:VAL:HG22	1:B:542:LEU:N	2.36	0.41
1:A:198:SER:OG	1:A:203:LYS:HA	2.21	0.41
1:A:345:LEU:HD12	1:A:345:LEU:HA	1.96	0.41
1:B:179:LEU:N	1:B:180:PRO:CD	2.84	0.41
1:B:574:ARG:HG3	1:B:577:GLN:OE1	2.21	0.41
1:B:503:GLY:HA2	1:B:523:VAL:HB	2.03	0.41
1:A:573:LEU:HD11	1:A:579:MET:HG3	2.03	0.40
1:A:464:THR:O	1:A:465:ASP:C	2.64	0.40
1:A:511:ASP:OD2	1:A:511:ASP:C	2.64	0.40
1:B:246:GLY:HA2	1:B:292:ALA:O	2.22	0.40
1:B:382:ARG:HB3	1:B:389:ILE:HG12	2.03	0.40
1:A:223:ARG:HH11	1:A:223:ARG:HA	1.85	0.40
1:A:477:LYS:HE3	1:A:477:LYS:CA	2.47	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	583/585 (100%)	555 (95%)	26 (4%)	2 (0%)	36	36
1	B	583/585 (100%)	551 (94%)	31 (5%)	1 (0%)	43	44
All	All	1166/1170 (100%)	1106 (95%)	57 (5%)	3 (0%)	36	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	LYS

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Mol	Chain	Res	Type
1	A	278	THR
1	B	278	THR

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	493/493 (100%)	476 (97%)	17 (3%)	32	35
1	B	493/493 (100%)	474 (96%)	19 (4%)	28	31
All	All	986/986 (100%)	950 (96%)	36 (4%)	30	33

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	LYS
1	A	133	ILE
1	A	155	GLN
1	A	182	LEU
1	A	260	GLU
1	A	280	ARG
1	A	341	LEU
1	A	345	LEU
1	A	398	LEU
1	A	404	LEU
1	A	408	GLN
1	A	413	LEU
1	A	477	LYS
1	A	483	LEU
1	A	504	ASN
1	A	547	GLU
1	A	565	LYS
1	B	59	LYS
1	B	85	LEU
1	B	155	GLN
1	B	191	TYR

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Mol	Chain	Res	Type
1	B	257	GLN
1	B	282	ASN
1	B	328	ASN
1	B	340	ARG
1	B	382	ARG
1	B	398	LEU
1	B	404	LEU
1	B	438	LEU
1	B	477	LYS
1	B	504	ASN
1	B	524	GLN
1	B	542	LEU
1	B	565	LYS
1	B	573	LEU
1	B	585	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	HIS
1	A	22	GLN
1	A	90	GLN
1	A	135	GLN
1	A	155	GLN
1	A	244	HIS
1	A	248	GLN
1	A	289	GLN
1	A	317	GLN
1	A	325	ASN
1	A	332	HIS
1	A	346	ASN
1	A	367	GLN
1	A	373	ASN
1	A	390	HIS
1	A	408	GLN
1	A	415	ASN
1	A	443	GLN
1	A	504	ASN
1	A	509	HIS
1	A	515	ASN
1	A	527	HIS
1	A	563	HIS

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Mol	Chain	Res	Type
1	A	568	GLN
1	B	135	GLN
1	B	164	HIS
1	B	202	HIS
1	B	243	ASN
1	B	244	HIS
1	B	248	GLN
1	B	282	ASN
1	B	328	ASN
1	B	332	HIS
1	B	357	HIS
1	B	415	ASN
1	B	504	ASN
1	B	509	HIS
1	B	533	ASN
1	B	539	GLN
1	B	544	GLN
1	B	563	HIS
1	B	566	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

12 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	C	1	2	12,12,12	0.43	0	17,17,17	0.73	0
2	GLC	C	2	2	11,11,12	0.67	0	15,15,17	0.56	0
2	GLC	C	3	2	11,11,12	0.64	0	15,15,17	0.81	1 (6%)
2	GLC	C	4	2	11,11,12	0.62	0	15,15,17	1.12	2 (13%)
2	GLC	C	5	2	11,11,12	0.63	0	15,15,17	0.57	0
2	GLC	C	6	2	11,11,12	0.47	0	15,15,17	0.74	1 (6%)
2	GLC	D	1	2	12,12,12	0.42	0	17,17,17	0.46	0
2	GLC	D	2	2	11,11,12	0.56	0	15,15,17	0.65	1 (6%)
2	GLC	D	3	2	11,11,12	0.65	0	15,15,17	0.78	1 (6%)
2	GLC	D	4	2	11,11,12	0.58	0	15,15,17	1.00	1 (6%)
2	GLC	D	5	2	11,11,12	0.57	0	15,15,17	0.55	0
2	GLC	D	6	2	11,11,12	0.47	0	15,15,17	0.68	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	GLC	C	1	2	1/1/5/5	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	GLC	C	3	2	-	2/2/19/22	0/1/1/1
2	GLC	C	4	2	-	0/2/19/22	0/1/1/1
2	GLC	C	5	2	-	0/2/19/22	0/1/1/1
2	GLC	C	6	2	-	2/2/19/22	0/1/1/1
2	GLC	D	1	2	1/1/5/5	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	0/2/19/22	0/1/1/1
2	GLC	D	3	2	-	2/2/19/22	0/1/1/1
2	GLC	D	4	2	-	0/2/19/22	0/1/1/1
2	GLC	D	5	2	-	0/2/19/22	0/1/1/1
2	GLC	D	6	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4	GLC	C1-O5-C5	2.85	116.01	112.19
2	C	3	GLC	C1-O5-C5	2.75	115.87	112.19
2	D	3	GLC	C1-O5-C5	2.63	115.71	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4	GLC	C1-O5-C5	2.51	115.54	112.19
2	C	6	GLC	C1-O5-C5	2.46	115.49	112.19
2	D	6	GLC	C1-O5-C5	2.19	115.11	112.19
2	D	2	GLC	C1-O5-C5	2.04	114.93	112.19
2	C	4	GLC	C1-C2-C3	-2.04	106.67	109.64

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1	GLC	C1
2	D	1	GLC	C1

All (10) torsion outliers are listed below:

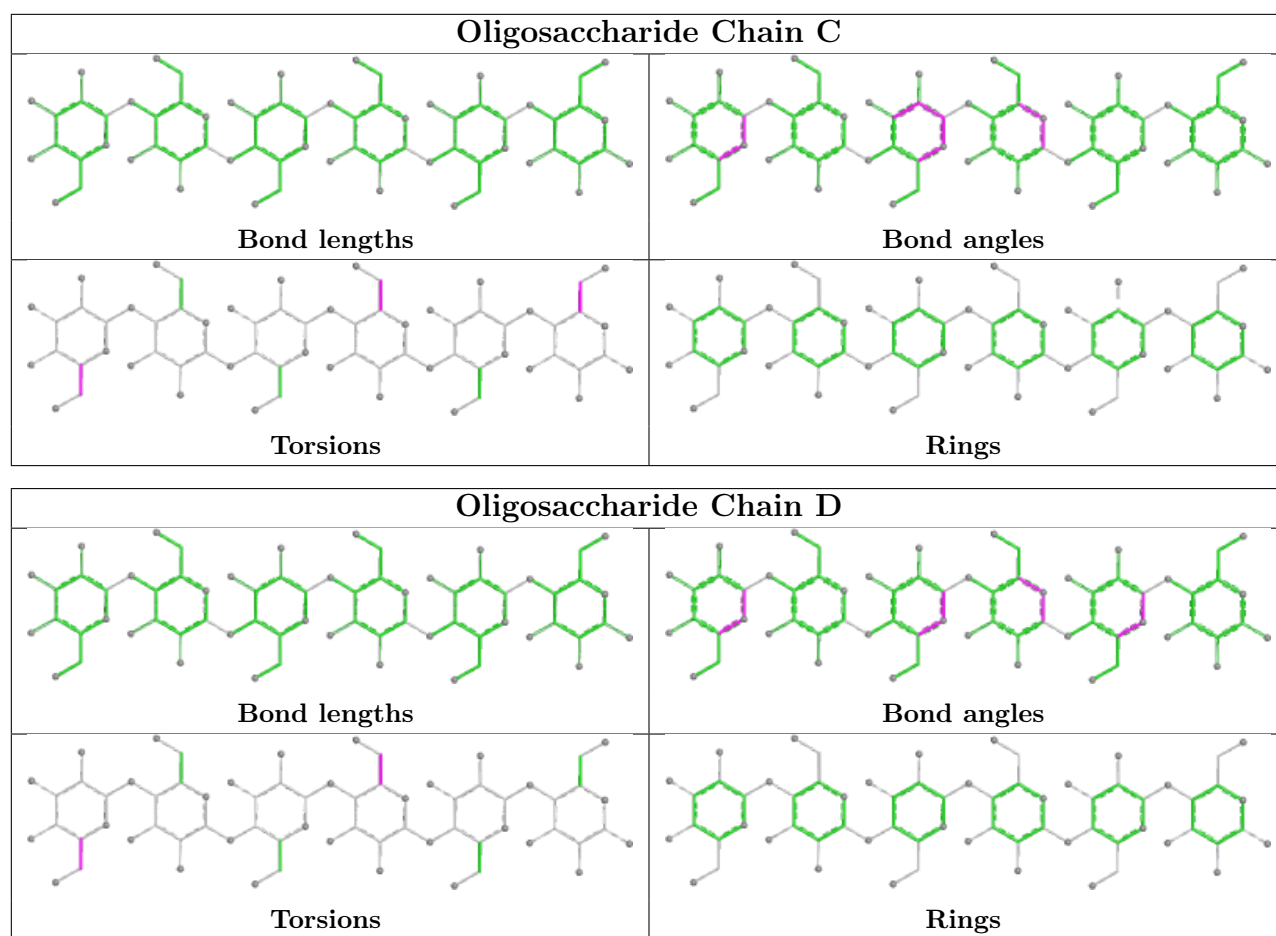
Mol	Chain	Res	Type	Atoms
2	C	6	GLC	O5-C5-C6-O6
2	D	6	GLC	C4-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	D	6	GLC	O5-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6
2	C	3	GLC	C4-C5-C6-O6
2	D	3	GLC	C4-C5-C6-O6
2	C	6	GLC	C4-C5-C6-O6
2	C	3	GLC	O5-C5-C6-O6
2	D	3	GLC	O5-C5-C6-O6

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	4	GLC	1	0
2	D	3	GLC	2	0
2	D	6	GLC	2	0
2	C	4	GLC	1	0
2	C	6	GLC	3	0
2	C	3	GLC	2	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

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	585/585 (100%)	0.05	18 (3%)	51	54	15, 25, 47, 72	0
1	B	585/585 (100%)	0.11	26 (4%)	39	41	17, 27, 50, 83	0
All	All	1170/1170 (100%)	0.08	44 (3%)	44	46	15, 26, 48, 83	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	549	GLY	4.7
1	B	275	VAL	4.2
1	A	276	SER	4.2
1	A	275	VAL	4.1
1	B	278	THR	3.7
1	B	276	SER	3.6
1	B	90	GLN	3.3
1	B	224	ARG	3.2
1	B	552	THR	3.2
1	B	516	LEU	3.2
1	A	564	GLY	3.2
1	A	279	SER	3.1
1	B	565	LYS	2.9
1	A	277	LYS	2.8
1	A	548	SER	2.8
1	B	279	SER	2.8
1	B	536	GLY	2.8
1	A	280	ARG	2.7
1	B	548	SER	2.7
1	A	151	PRO	2.7
1	A	547	GLU	2.6
1	B	280	ARG	2.6
1	A	572	THR	2.5
1	A	278	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	546	PRO	2.5
1	B	164	HIS	2.4
1	B	566	GLN	2.3
1	B	502	ARG	2.3
1	B	550	GLY	2.3
1	B	495	HIS	2.2
1	B	570	LYS	2.2
1	B	513	GLN	2.1
1	B	340	ARG	2.1
1	B	546	PRO	2.1
1	A	516	LEU	2.1
1	B	130	GLU	2.1
1	A	272	ASP	2.1
1	B	273	PHE	2.1
1	A	510	ALA	2.1
1	B	277	LYS	2.0
1	A	542	LEU	2.0
1	B	263	ARG	2.0
1	B	470	ARG	2.0
1	A	54	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

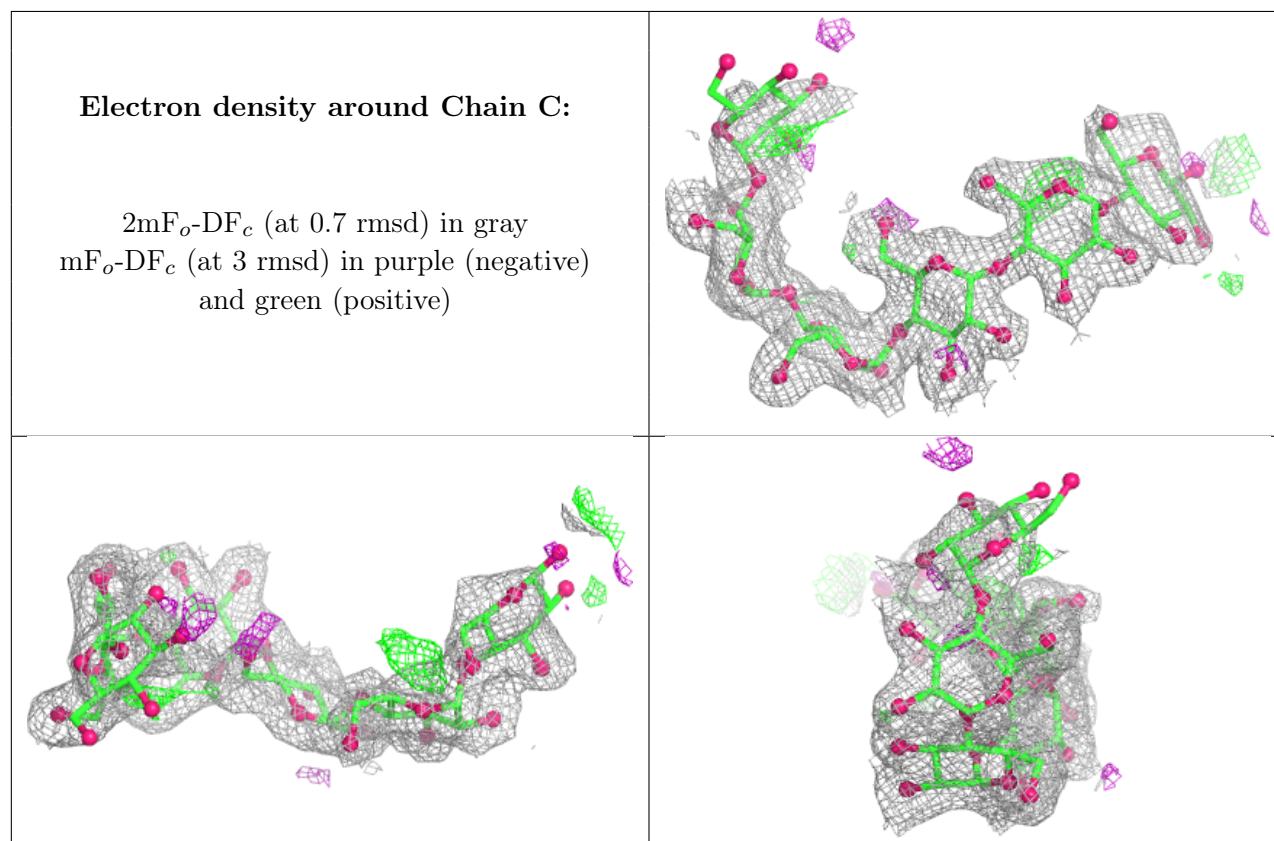
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	D	1	12/12	0.63	0.18	51,64,66,67	0
2	GLC	D	6	11/12	0.65	0.17	43,47,49,51	11
2	GLC	C	6	11/12	0.66	0.23	42,47,50,51	11
2	GLC	C	1	12/12	0.72	0.19	40,48,52,55	11
2	GLC	D	2	11/12	0.76	0.18	33,38,40,43	11
2	GLC	C	2	11/12	0.86	0.14	23,29,34,37	11
2	GLC	D	3	11/12	0.92	0.09	24,28,31,33	0
2	GLC	C	3	11/12	0.93	0.08	20,24,27,33	0

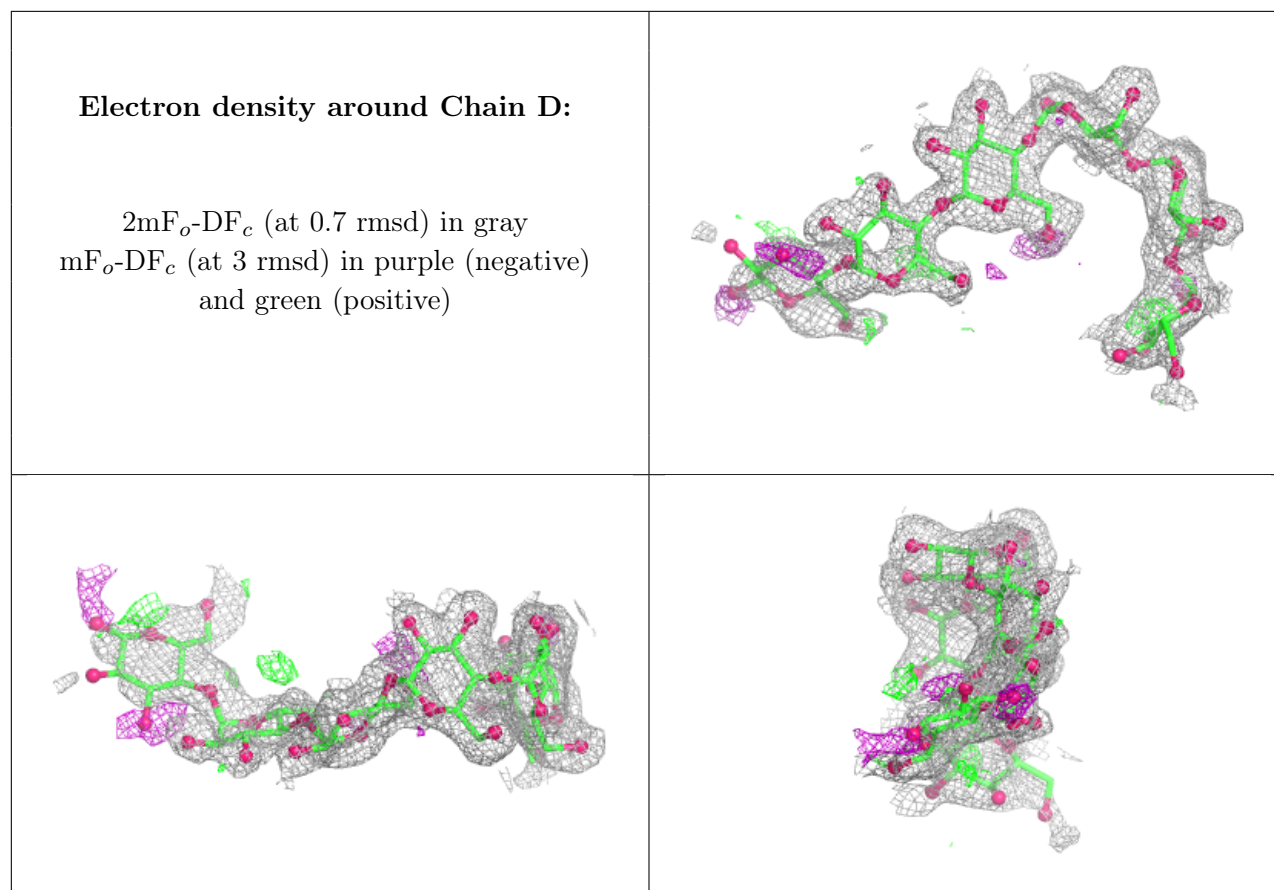
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GLC	D	4	11/12	0.94	0.07	23,26,27,28	0
2	GLC	C	4	11/12	0.96	0.05	17,21,22,23	0
2	GLC	D	5	11/12	0.97	0.06	27,29,31,37	0
2	GLC	C	5	11/12	0.97	0.05	21,23,26,33	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.





6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CA	B	1002	1/1	0.99	0.03	32,32,32,32	0
3	CA	A	1001	1/1	1.00	0.01	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.