



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:25 PM UTC

PDB ID : 2D0F / pdb_00002d0f
Title : Crystal Structure of Thermoactinomyces vulgaris R-47 Alpha-Amylase 1 (TVAI) Mutant D356N complexed with P2, a pullulan model oligosaccharide
Authors : Abe, A.; Yoshida, H.; Tono-zuka, T.; Sakano, Y.; Kamitori, S.
Deposited on : 2005-08-02
Resolution : 2.08 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

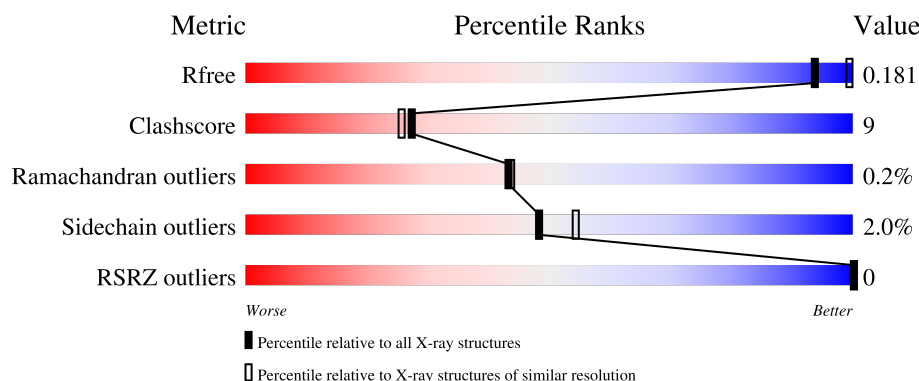
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	637	 84% 15% .
2	B	5	 40% 60%
3	C	3	 33% 67%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5615 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 alpha-amylase I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	637	Total	C	N	O	S	0	0	0
			5038	3192	843	993	10			

There is a discrepancy between the modelled and reference sequences:

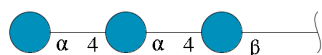
Chain	Residue	Modelled	Actual	Comment	Reference
A	356	ASN	ASP	engineered mutation	UNP Q60053

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	5	Total	C	O	0	0	0
			56	30	26			

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose.

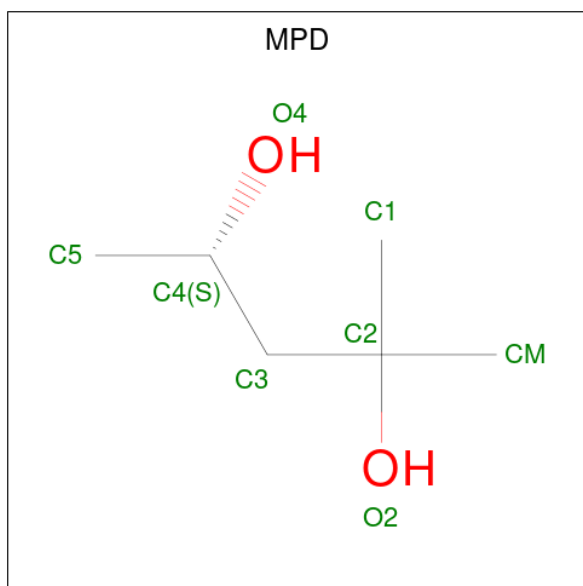


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	C	3	Total	C	O	0	0	0
			34	18	16			

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		

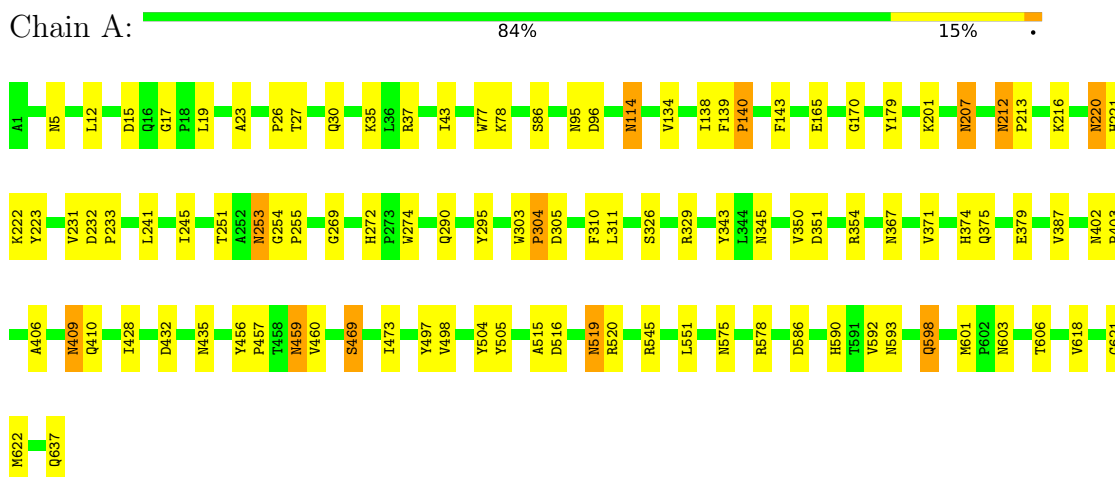
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	452	Total	O	0	0
			452	452		

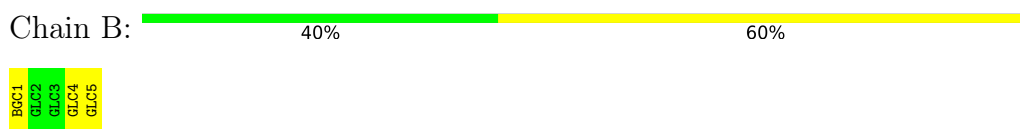
3 Residue-property plots

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 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: alpha-amylase I



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.53Å 50.43Å 107.83Å 90.00° 104.30° 90.00°	Depositor
Resolution (Å)	46.00 – 2.08 46.00 – 2.08	Depositor EDS
% Data completeness (in resolution range)	(Not available) (46.00-2.08) 99.8 (46.00-2.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	11.20 (at 2.08Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.154 , 0.187 0.149 , 0.181	Depositor DCC
R_{free} test set	3836 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	13.3	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 58.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5615	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% 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, BGC, GLC, MPD

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	0.35	0/5195	0.92	19/7108 (0.3%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	139	PHE	CA-C-N	7.45	129.15	119.84
1	A	139	PHE	C-N-CA	7.45	129.15	119.84
1	A	432	ASP	N-CA-C	-6.42	102.24	110.53
1	A	367	ASN	N-CA-C	6.18	119.85	112.93
1	A	428	ILE	N-CA-C	5.76	117.28	111.00
1	A	457	PRO	N-CA-C	-5.67	102.90	111.41
1	A	406	ALA	N-CA-C	5.53	119.51	112.87
1	A	212	ASN	N-CA-C	-5.43	102.15	110.07
1	A	231	VAL	N-CA-C	-5.41	102.60	109.58
1	A	86	SER	N-CA-C	5.39	118.45	110.48
1	A	17	GLY	CA-C-N	5.38	125.46	119.32
1	A	17	GLY	C-N-CA	5.38	125.46	119.32
1	A	15	ASP	N-CA-C	-5.36	101.14	108.38
1	A	456	TYR	N-CA-C	5.27	118.67	109.82
1	A	269	GLY	N-CA-C	-5.24	105.59	112.82
1	A	592	VAL	N-CA-C	5.23	116.12	108.53
1	A	222	LYS	N-CA-C	5.17	118.86	112.24
1	A	305	ASP	N-CA-C	5.16	119.29	112.89
1	A	351	ASP	N-CA-C	5.15	119.66	113.17

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	5038	0	4650	83	0
2	B	56	0	48	3	0
3	C	34	0	30	2	0
4	A	3	0	0	0	0
5	A	32	0	56	3	0
6	A	452	0	0	9	1
All	All	5615	0	4784	90	1

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 9.

All (90) 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:578:ARG:HE	1:A:637:GLN:HE21	1.17	0.93
1:A:272:HIS:HD2	1:A:274:TRP:H	1.17	0.90
1:A:603:ASN:HD21	1:A:621:GLY:H	1.16	0.88
3:C:2:GLC:H62	3:C:3:GLC:H5	1.62	0.80
1:A:37:ARG:HH22	1:A:598:GLN:HE22	1.30	0.76
1:A:603:ASN:ND2	1:A:621:GLY:H	1.85	0.74
1:A:207:ASN:H	1:A:207:ASN:HD22	1.33	0.74
1:A:603:ASN:HD21	1:A:621:GLY:N	1.85	0.73
1:A:221:HIS:CD2	1:A:223:TYR:H	2.08	0.72
1:A:221:HIS:HD2	1:A:223:TYR:H	1.36	0.71
1:A:220:ASN:HD22	1:A:221:HIS:N	1.92	0.67
1:A:519:ASN:ND2	1:A:520:ARG:HG3	2.10	0.66
1:A:606:THR:HG22	6:A:2350:HOH:O	1.97	0.63
1:A:201:LYS:HE3	1:A:255:PRO:O	1.97	0.63
2:B:4:GLC:H62	2:B:5:GLC:H5	1.81	0.61
3:C:2:GLC:H62	3:C:3:GLC:C5	2.30	0.60
1:A:586:ASP:OD1	1:A:590:HIS:HE1	1.83	0.60
1:A:578:ARG:HE	1:A:637:GLN:NE2	1.93	0.60
1:A:459:ASN:HD22	1:A:459:ASN:N	1.99	0.60
1:A:19:LEU:HD23	1:A:575:ASN:ND2	2.17	0.59
1:A:310:PHE:CE1	1:A:311:LEU:HD22	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:LYS:NZ	1:A:253:ASN:HD21	2.04	0.56
1:A:114:ASN:C	1:A:114:ASN:HD22	2.13	0.54
1:A:170:GLY:O	5:A:1002:MPD:HM1	2.08	0.54
1:A:345:ASN:ND2	1:A:387:VAL:HG13	2.22	0.54
1:A:409:ASN:HD22	1:A:409:ASN:H	1.56	0.54
1:A:272:HIS:CD2	1:A:274:TRP:H	2.10	0.53
1:A:143:PHE:HB3	6:A:2088:HOH:O	2.09	0.53
1:A:216:LYS:HE3	6:A:2370:HOH:O	2.08	0.53
1:A:578:ARG:NE	1:A:637:GLN:HE21	1.98	0.53
1:A:374:HIS:HE1	6:A:2028:HOH:O	1.92	0.52
1:A:37:ARG:NH2	1:A:598:GLN:HE22	2.04	0.51
1:A:37:ARG:HH22	1:A:598:GLN:NE2	2.05	0.51
1:A:78:LYS:NZ	1:A:78:LYS:HB3	2.27	0.49
1:A:253:ASN:ND2	1:A:253:ASN:H	2.11	0.49
1:A:354:ARG:C	1:A:354:ARG:HD3	2.38	0.49
1:A:519:ASN:ND2	1:A:519:ASN:H	2.12	0.48
1:A:23:ALA:O	1:A:26:PRO:HD3	2.14	0.47
1:A:343:TYR:O	1:A:350:VAL:HG22	2.14	0.47
1:A:303:TRP:CE2	1:A:304:PRO:HB3	2.49	0.47
1:A:545:ARG:HA	1:A:551:LEU:HD12	1.95	0.47
1:A:345:ASN:HD21	1:A:387:VAL:HG13	1.80	0.47
1:A:515:ALA:O	1:A:516:ASP:C	2.57	0.47
1:A:114:ASN:C	1:A:114:ASN:ND2	2.73	0.46
1:A:138:ILE:O	1:A:140:PRO:HD3	2.16	0.45
1:A:519:ASN:HD22	1:A:519:ASN:N	2.14	0.45
1:A:459:ASN:N	1:A:459:ASN:ND2	2.64	0.45
1:A:220:ASN:HD22	1:A:220:ASN:C	2.24	0.45
1:A:402:ASN:HB3	1:A:403:PRO:HD3	1.98	0.45
1:A:497:TYR:CG	1:A:498:VAL:N	2.85	0.45
5:A:1001:MPD:HM1	6:A:2408:HOH:O	2.16	0.45
1:A:134:VAL:H	1:A:207:ASN:ND2	2.15	0.44
1:A:216:LYS:HD3	1:A:233:PRO:CD	2.48	0.44
1:A:371:VAL:HG23	6:A:2294:HOH:O	2.16	0.44
1:A:326:SER:OG	1:A:329:ARG:HG3	2.16	0.44
6:A:2010:HOH:O	2:B:1:BGC:H2	2.17	0.44
1:A:253:ASN:H	1:A:253:ASN:HD22	1.66	0.43
1:A:603:ASN:HD22	1:A:618:VAL:HG12	1.83	0.43
1:A:290:GLN:HA	1:A:295:TYR:CD1	2.54	0.43
1:A:459:ASN:HD22	1:A:460:VAL:H	1.66	0.43
1:A:519:ASN:H	1:A:519:ASN:HD22	1.67	0.43
5:A:1004:MPD:H32	6:A:2442:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:GLN:O	1:A:379:GLU:HG3	2.18	0.43
1:A:603:ASN:HD22	1:A:603:ASN:HA	1.65	0.43
1:A:469:SER:HA	1:A:473:ILE:HD11	2.00	0.43
1:A:601:MET:HG2	1:A:637:GLN:NE2	2.34	0.43
1:A:43:ILE:HB	1:A:95:ASN:O	2.19	0.42
1:A:216:LYS:HD3	1:A:233:PRO:HD2	2.00	0.42
1:A:201:LYS:HZ1	1:A:253:ASN:HD21	1.67	0.42
1:A:207:ASN:HD22	1:A:207:ASN:N	2.02	0.42
1:A:251:THR:HA	1:A:254:GLY:O	2.19	0.42
1:A:212:ASN:HB3	1:A:213:PRO:HD2	2.02	0.42
1:A:253:ASN:HD22	1:A:253:ASN:C	2.27	0.42
1:A:27:THR:H	1:A:30:GLN:HE21	1.67	0.42
1:A:26:PRO:HA	1:A:30:GLN:NE2	2.35	0.42
1:A:593:ASN:HB3	1:A:622:MET:CE	2.50	0.42
1:A:374:HIS:HD2	1:A:410:GLN:OE1	2.03	0.41
1:A:12:LEU:HA	1:A:37:ARG:O	2.20	0.41
1:A:27:THR:H	1:A:30:GLN:NE2	2.19	0.41
1:A:216:LYS:HD2	1:A:232:ASP:OD2	2.21	0.41
1:A:35:LYS:HA	1:A:77:TRP:O	2.20	0.41
1:A:165:GLU:HB2	1:A:179:TYR:CD2	2.55	0.41
1:A:345:ASN:HD22	1:A:345:ASN:HA	1.74	0.41
2:B:4:GLC:C6	2:B:5:GLC:H5	2.48	0.41
1:A:402:ASN:N	1:A:403:PRO:CD	2.84	0.41
1:A:5:ASN:HA	1:A:96:ASP:O	2.21	0.40
1:A:409:ASN:H	1:A:409:ASN:ND2	2.17	0.40
1:A:504:TYR:O	1:A:505:TYR:C	2.64	0.40
1:A:621:GLY:HA2	6:A:2320:HOH:O	2.21	0.40
1:A:241:LEU:O	1:A:245:ILE:HG13	2.21	0.40

All (1) 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)
6:A:2178:HOH:O	6:A:2178:HOH:O[2_656]	1.10	1.10

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	635/637 (100%)	608 (96%)	26 (4%)	1 (0%)	43	44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	PRO

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	540/540 (100%)	529 (98%)	11 (2%)	48	54

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	ASN
1	A	207	ASN
1	A	220	ASN
1	A	253	ASN
1	A	304	PRO
1	A	409	ASN
1	A	435	ASN
1	A	459	ASN
1	A	469	SER
1	A	519	ASN

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Mol	Chain	Res	Type
1	A	598	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	10	ASN
1	A	14	HIS
1	A	30	GLN
1	A	56	ASN
1	A	95	ASN
1	A	114	ASN
1	A	153	GLN
1	A	160	ASN
1	A	195	GLN
1	A	207	ASN
1	A	212	ASN
1	A	220	ASN
1	A	221	HIS
1	A	242	GLN
1	A	253	ASN
1	A	266	ASN
1	A	272	HIS
1	A	314	ASN
1	A	345	ASN
1	A	364	ASN
1	A	367	ASN
1	A	374	HIS
1	A	402	ASN
1	A	409	ASN
1	A	410	GLN
1	A	435	ASN
1	A	455	ASN
1	A	459	ASN
1	A	494	GLN
1	A	512	GLN
1	A	519	ASN
1	A	539	GLN
1	A	546	ASN
1	A	575	ASN
1	A	585	ASN
1	A	590	HIS

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Mol	Chain	Res	Type
1	A	598	GLN
1	A	603	ASN
1	A	637	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 ⓘ

8 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	BGC	B	1	2	12,12,12	0.45	0	17,17,17	0.76	0
2	GLC	B	2	2	11,11,12	0.61	0	15,15,17	0.42	0
2	GLC	B	3	2	11,11,12	0.47	0	15,15,17	0.53	0
2	GLC	B	4	2	11,11,12	0.53	0	15,15,17	0.52	0
2	GLC	B	5	2	11,11,12	0.49	0	15,15,17	0.54	0
3	BGC	C	1	3	12,12,12	0.39	0	17,17,17	0.34	0
3	GLC	C	2	3	11,11,12	0.52	0	15,15,17	0.58	1 (6%)
3	GLC	C	3	3	11,11,12	0.50	0	15,15,17	0.59	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	BGC	B	1	2	-	0/2/22/22	0/1/1/1
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
2	GLC	B	5	2	-	2/2/19/22	0/1/1/1
3	BGC	C	1	3	-	0/2/22/22	0/1/1/1
3	GLC	C	2	3	-	0/2/19/22	0/1/1/1
3	GLC	C	3	3	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	3	GLC	C1-O5-C5	2.10	115.00	112.19
3	C	2	GLC	C1-O5-C5	2.04	114.92	112.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

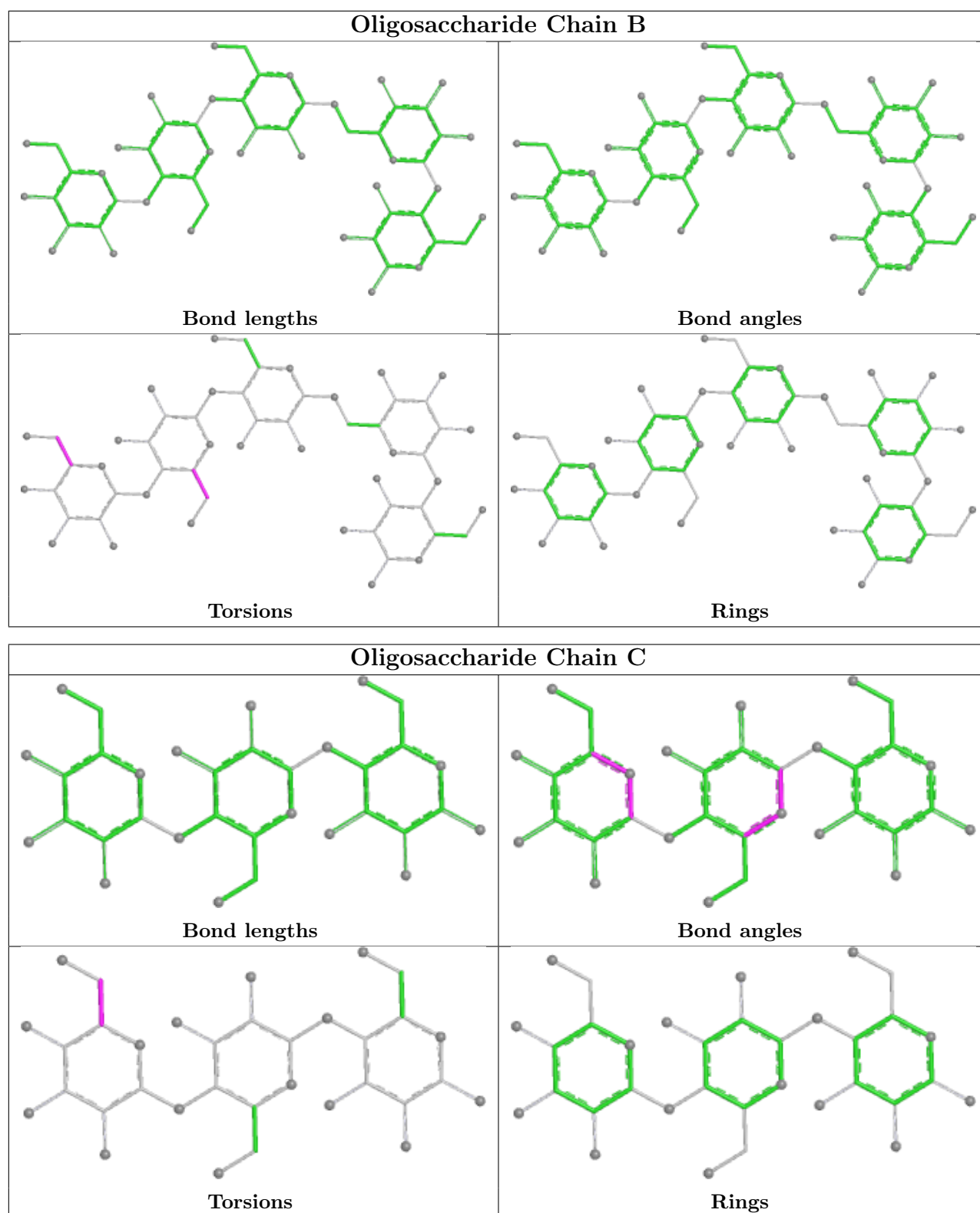
Mol	Chain	Res	Type	Atoms
2	B	5	GLC	O5-C5-C6-O6
2	B	5	GLC	C4-C5-C6-O6
3	C	3	GLC	O5-C5-C6-O6
2	B	4	GLC	O5-C5-C6-O6
3	C	3	GLC	C4-C5-C6-O6
2	B	4	GLC	C4-C5-C6-O6

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	4	GLC	2	0
3	C	2	GLC	2	0
2	B	1	BGC	1	0
3	C	3	GLC	2	0
2	B	5	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 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 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$
5	MPD	A	1002	-	7,7,7	0.53	0	9,10,10	0.50	0
5	MPD	A	1003	-	7,7,7	0.50	0	9,10,10	0.47	0
5	MPD	A	1001	-	7,7,7	0.58	0	9,10,10	0.63	0
5	MPD	A	1004	-	7,7,7	0.48	0	9,10,10	0.59	0

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
5	MPD	A	1002	-	-	3/5/5/5	-
5	MPD	A	1003	-	-	0/5/5/5	-
5	MPD	A	1001	-	-	3/5/5/5	-
5	MPD	A	1004	-	-	1/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1001	MPD	C1-C2-C3-C4
5	A	1001	MPD	O2-C2-C3-C4
5	A	1002	MPD	O2-C2-C3-C4
5	A	1004	MPD	O2-C2-C3-C4
5	A	1001	MPD	CM-C2-C3-C4
5	A	1002	MPD	C1-C2-C3-C4
5	A	1002	MPD	CM-C2-C3-C4

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1002	MPD	1	0
5	A	1001	MPD	1	0
5	A	1004	MPD	1	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	637/637 (100%)	-0.48	0 100 100	6, 11, 18, 24	0

There are no RSRZ outliers to report.

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](#)

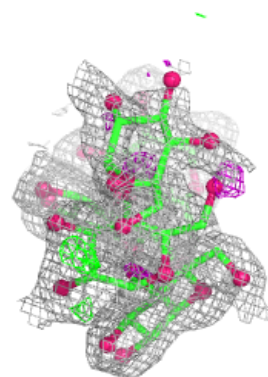
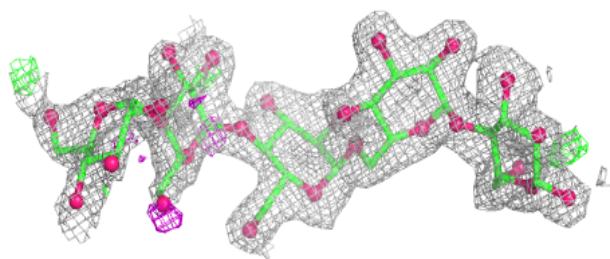
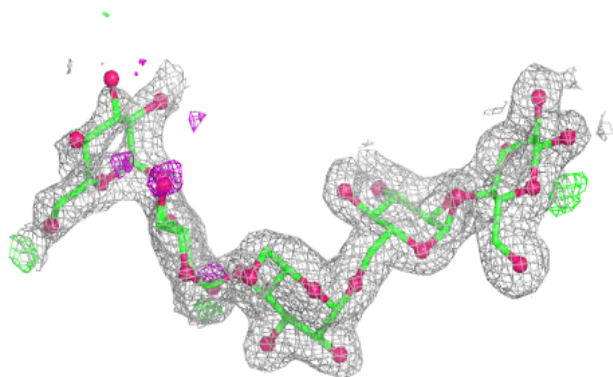
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GLC	C	3	11/12	0.46	0.21	60,62,63,64	0
3	BGC	C	1	12/12	0.64	0.16	52,54,55,56	0
2	GLC	B	5	11/12	0.77	0.15	44,45,46,48	0
2	GLC	B	4	11/12	0.79	0.13	33,37,41,45	0
3	GLC	C	2	11/12	0.80	0.17	52,54,56,57	0
2	GLC	B	3	11/12	0.92	0.08	18,20,22,29	0
2	BGC	B	1	12/12	0.94	0.07	12,16,17,17	0
2	GLC	B	2	11/12	0.95	0.06	11,14,15,16	0

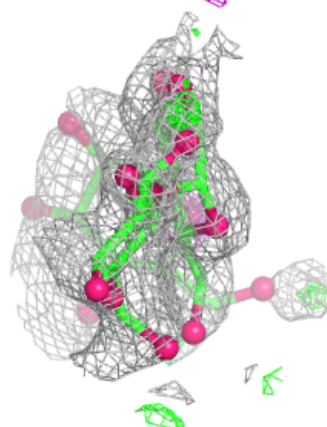
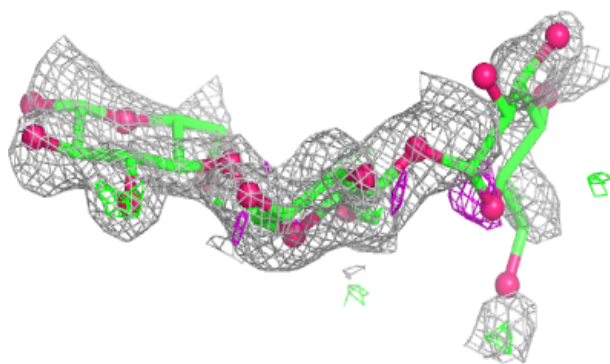
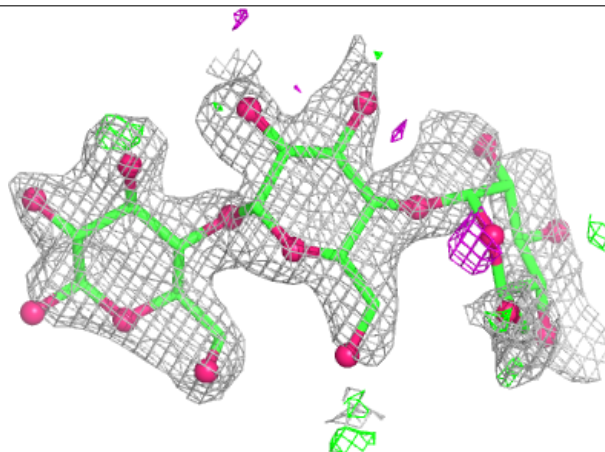
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain B:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around Chain C:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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
5	MPD	A	1004	8/8	0.59	0.34	71,71,71,72	0
5	MPD	A	1002	8/8	0.77	0.18	35,36,38,39	0
5	MPD	A	1001	8/8	0.80	0.16	36,37,38,39	0
5	MPD	A	1003	8/8	0.89	0.16	51,51,52,52	0
4	CA	A	2001	1/1	0.99	0.02	11,11,11,11	0
4	CA	A	2002	1/1	1.00	0.03	11,11,11,11	0
4	CA	A	2003	1/1	1.00	0.03	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.