Full wwPDB X-ray Structure Validation Report (i)

Jan 27, 2024 – 01:35 PM EST

PDB ID	:	1BYB
Title	:	CRYSTAL STRUCTURES OF SOYBEAN BETA-AMYLASE REACTED WITH BETA-MALTOSE AND MALTAL: ACTIVE SITE COMPONENTS AND THEIR APPARENT ROLE IN CATALYSIS
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Deposited on	:	1994-01-25
Resolution	:	1.90 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 (i) 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 (i)) were used in the production of this report:

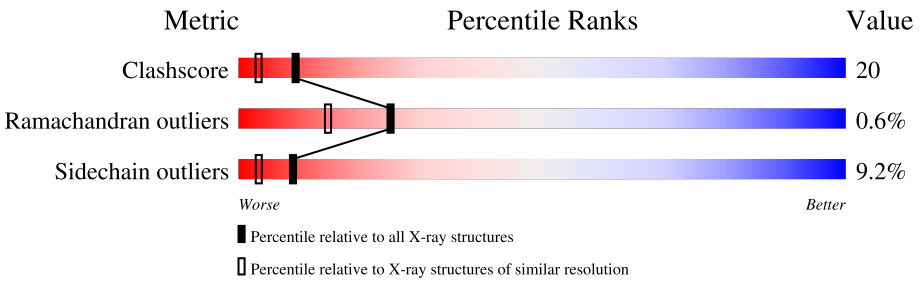
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	495	64% 28% 6% ••
2	B	4	25% 75%

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	B	1	X	-	-	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4307 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.

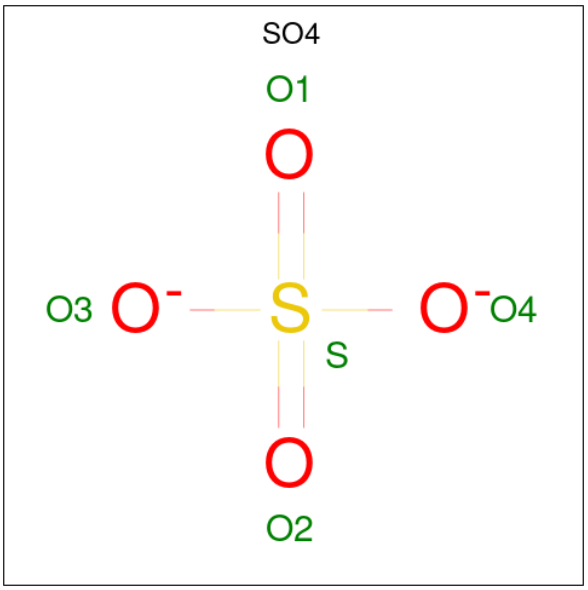
- Molecule 1 is a protein called BETA-AMYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	491	3925	2518	662	728	17	0	0	1

- Molecule 2 is an oligosaccharide called 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
			Total	C	O			
2	B	4	45	24	21	0	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
3	A	1	5	4	1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	332	332	332	0	0

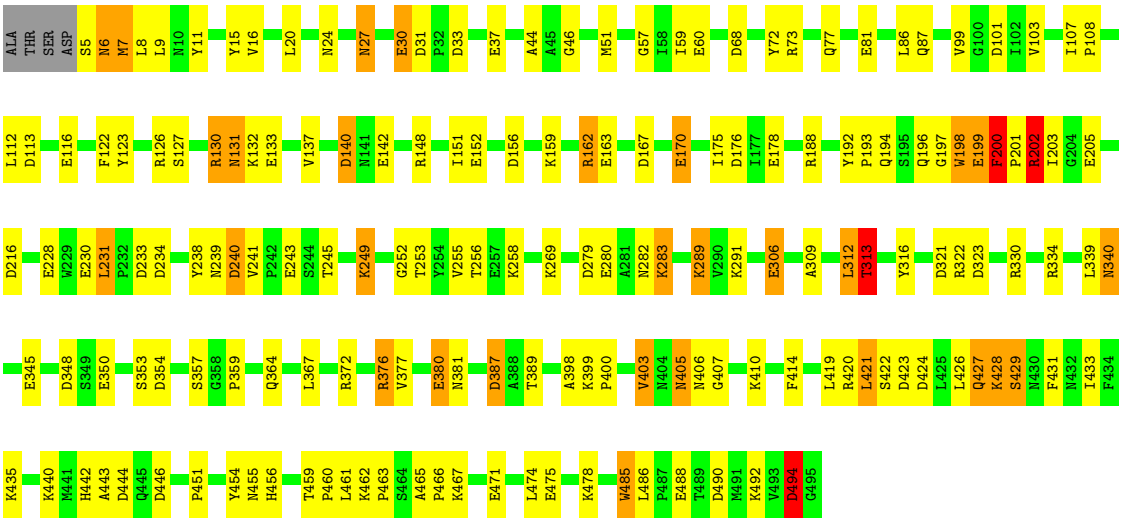
3 Residue-property plots (i)

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. 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: BETA-AMYLASE

Chain A: 64% 28% 6% ••



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

Chain B: 25% 75%

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	86.20Å 86.20Å 144.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	9.00 – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) (9.00-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.149 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4307	wwPDB-VP
Average B, all atoms ($Å^2$)	22.0	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, SO4

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	1.17	22/4032 (0.5%)	1.31	57/5479 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	37	GLU	CD-OE2	9.85	1.36	1.25
1	A	170	GLU	CD-OE1	9.23	1.35	1.25
1	A	306	GLU	CD-OE2	9.14	1.35	1.25
1	A	116	GLU	CD-OE2	8.87	1.35	1.25
1	A	488	GLU	CD-OE2	8.19	1.34	1.25
1	A	163	GLU	CD-OE1	8.14	1.34	1.25
1	A	475	GLU	CD-OE1	8.07	1.34	1.25
1	A	228	GLU	CD-OE2	7.99	1.34	1.25
1	A	243	GLU	CD-OE1	7.72	1.34	1.25
1	A	60	GLU	CD-OE2	7.19	1.33	1.25
1	A	30	GLU	CD-OE2	6.68	1.33	1.25
1	A	230	GLU	CD-OE1	6.50	1.32	1.25
1	A	471	GLU	CD-OE2	6.29	1.32	1.25
1	A	380	GLU	CD-OE2	-5.81	1.19	1.25
1	A	81	GLU	CD-OE1	-5.74	1.19	1.25
1	A	142	GLU	CD-OE1	5.64	1.31	1.25
1	A	178	GLU	CD-OE1	5.51	1.31	1.25
1	A	205	GLU	CD-OE1	5.32	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	345	GLU	CD-OE2	5.25	1.31	1.25
1	A	350	GLU	CD-OE1	5.10	1.31	1.25
1	A	280	GLU	CD-OE1	5.09	1.31	1.25
1	A	152	GLU	CD-OE2	5.07	1.31	1.25

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	313	THR	N-CA-CB	11.49	132.13	110.30
1	A	240	ASP	CB-CG-OD2	10.02	127.32	118.30
1	A	321	ASP	CB-CG-OD1	9.89	127.20	118.30
1	A	162	ARG	NE-CZ-NH1	9.64	125.12	120.30
1	A	240	ASP	CB-CG-OD1	-9.17	110.05	118.30
1	A	233	ASP	CB-CG-OD2	-8.67	110.49	118.30
1	A	188	ARG	NE-CZ-NH1	8.31	124.46	120.30
1	A	31	ASP	CB-CG-OD1	8.18	125.66	118.30
1	A	323	ASP	CB-CG-OD1	7.89	125.40	118.30
1	A	130	ARG	NE-CZ-NH1	7.86	124.23	120.30
1	A	156	ASP	CB-CG-OD1	-7.84	111.24	118.30
1	A	200	PHE	CB-CA-C	7.55	125.51	110.40
1	A	33	ASP	CB-CG-OD1	-7.41	111.63	118.30
1	A	446	ASP	CB-CG-OD2	-7.40	111.64	118.30
1	A	202	ARG	NE-CZ-NH2	7.38	123.99	120.30
1	A	321	ASP	CB-CG-OD2	-7.30	111.73	118.30
1	A	420	ARG	NE-CZ-NH2	7.30	123.95	120.30
1	A	372	ARG	NE-CZ-NH2	-7.24	116.68	120.30
1	A	233	ASP	CB-CG-OD1	7.05	124.64	118.30
1	A	279	ASP	CB-CG-OD1	-6.93	112.06	118.30
1	A	31	ASP	CB-CG-OD2	-6.75	112.22	118.30
1	A	216	ASP	CB-CG-OD2	-6.74	112.23	118.30
1	A	322	ARG	NE-CZ-NH2	-6.70	116.95	120.30
1	A	494	ASP	CB-CG-OD1	-6.69	112.28	118.30
1	A	323	ASP	CB-CG-OD2	-6.67	112.30	118.30
1	A	140	ASP	CB-CG-OD2	6.62	124.26	118.30
1	A	376	ARG	NE-CZ-NH1	6.59	123.60	120.30
1	A	156	ASP	CB-CG-OD2	6.56	124.20	118.30
1	A	216	ASP	CB-CG-OD1	6.52	124.17	118.30
1	A	279	ASP	CB-CG-OD2	6.32	123.99	118.30
1	A	316	TYR	CB-CG-CD2	-6.28	117.23	121.00
1	A	140	ASP	CB-CG-OD1	-6.22	112.70	118.30
1	A	354	ASP	CB-CG-OD2	-6.06	112.85	118.30
1	A	387	ASP	CB-CG-OD2	5.99	123.69	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	322	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	A	334	ARG	NE-CZ-NH1	5.96	123.28	120.30
1	A	68	ASP	CB-CG-OD1	5.87	123.58	118.30
1	A	446	ASP	CB-CG-OD1	5.86	123.57	118.30
1	A	423	ASP	CB-CG-OD2	-5.84	113.04	118.30
1	A	167	ASP	CB-CG-OD1	5.69	123.42	118.30
1	A	312	LEU	O-C-N	-5.69	113.60	122.70
1	A	490	ASP	CB-CG-OD2	-5.67	113.19	118.30
1	A	424	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	316	TYR	CB-CG-CD1	5.49	124.30	121.00
1	A	348	ASP	CB-CG-OD1	-5.47	113.38	118.30
1	A	113	ASP	CB-CG-OD2	-5.35	113.48	118.30
1	A	176	ASP	CB-CG-OD1	5.31	123.08	118.30
1	A	176	ASP	CB-CG-OD2	-5.28	113.55	118.30
1	A	148	ARG	NE-CZ-NH1	5.25	122.92	120.30
1	A	387	ASP	CB-CG-OD1	-5.22	113.60	118.30
1	A	348	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	73	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	A	410	LYS	N-CA-CB	5.05	119.69	110.60
1	A	20	LEU	C-N-CA	-5.05	111.70	122.30
1	A	33	ASP	CB-CG-OD2	5.04	122.83	118.30
1	A	101	ASP	CB-CG-OD2	5.04	122.83	118.30
1	A	330	ARG	NE-CZ-NH2	-5.01	117.80	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	11	TYR	Sidechain

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	3925	0	3824	150	0
2	B	45	0	40	3	0
3	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	332	0	0	19	1
All	All	4307	0	3864	153	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 20.

All (153) 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:202:ARG:HG3	1:A:239:ASN:CA	1.40	1.47
1:A:202:ARG:CG	1:A:239:ASN:HA	1.55	1.35
1:A:201:PRO:O	1:A:202:ARG:CD	1.83	1.27
1:A:200:PHE:CG	1:A:201:PRO:HD3	1.71	1.26
1:A:201:PRO:O	1:A:202:ARG:HD3	1.31	1.20
1:A:202:ARG:CG	1:A:239:ASN:CA	2.18	1.09
1:A:421:LEU:C	1:A:421:LEU:HD22	1.72	1.09
1:A:202:ARG:HG3	1:A:239:ASN:C	1.75	1.06
1:A:202:ARG:HG2	1:A:239:ASN:HD22	1.20	1.06
1:A:99:VAL:HG12	4:A:868:HOH:O	1.56	1.05
1:A:202:ARG:HG2	1:A:239:ASN:ND2	1.72	1.04
1:A:200:PHE:CD2	1:A:201:PRO:HD3	1.92	1.04
1:A:198:TRP:CH2	1:A:200:PHE:HA	1.93	1.02
1:A:202:ARG:NE	1:A:239:ASN:O	1.96	0.97
1:A:203:ILE:HG23	1:A:240:ASP:O	1.64	0.97
1:A:202:ARG:CG	1:A:239:ASN:HD22	1.77	0.96
1:A:202:ARG:HG3	1:A:239:ASN:HA	0.99	0.96
1:A:202:ARG:HG2	1:A:239:ASN:CB	1.97	0.94
1:A:202:ARG:CG	1:A:239:ASN:CB	2.44	0.94
1:A:367:LEU:HD22	1:A:377:VAL:HG11	1.53	0.90
1:A:405:ASN:HD22	1:A:405:ASN:H	1.18	0.89
1:A:199:GLU:H	1:A:239:ASN:HD21	1.15	0.89
1:A:201:PRO:O	1:A:202:ARG:HD2	1.73	0.89
1:A:421:LEU:C	1:A:421:LEU:CD2	2.42	0.87
1:A:200:PHE:CG	1:A:201:PRO:CD	2.57	0.87
1:A:202:ARG:CB	1:A:239:ASN:HA	2.04	0.87
1:A:289:LYS:HZ1	1:A:460:PRO:CD	1.90	0.84
4:A:868:HOH:O	2:B:2:GLC:H62	1.77	0.84
1:A:5:SER:HB3	1:A:9:LEU:CD1	2.07	0.83
1:A:202:ARG:HG3	1:A:239:ASN:CB	2.08	0.82
1:A:200:PHE:CD1	1:A:201:PRO:HD3	2.16	0.81
1:A:202:ARG:CB	1:A:239:ASN:HD22	1.93	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:459:THR:HG22	1:A:460:PRO:HD2	1.61	0.80
1:A:289:LYS:HZ1	1:A:460:PRO:N	1.78	0.80
1:A:381:ASN:HD21	1:A:419:LEU:HB3	1.46	0.79
1:A:6:ASN:HD21	1:A:8:LEU:HB2	1.47	0.79
1:A:421:LEU:HD22	1:A:421:LEU:O	1.86	0.75
1:A:6:ASN:ND2	1:A:8:LEU:HB2	2.01	0.75
1:A:357:SER:HB3	1:A:359:PRO:HD3	1.69	0.75
1:A:202:ARG:HG2	1:A:239:ASN:CG	2.06	0.74
1:A:245:THR:O	1:A:249:LYS:HG3	1.86	0.73
1:A:202:ARG:HB3	1:A:238:TYR:O	1.89	0.72
1:A:313:THR:N	4:A:566:HOH:O	2.20	0.72
1:A:405:ASN:H	1:A:405:ASN:ND2	1.87	0.72
1:A:202:ARG:CG	1:A:239:ASN:HB3	2.18	0.71
1:A:459:THR:CG2	1:A:460:PRO:HD2	2.23	0.68
1:A:282:ASN:ND2	4:A:697:HOH:O	2.28	0.66
1:A:367:LEU:HD22	1:A:377:VAL:CG1	2.24	0.66
1:A:202:ARG:HB3	1:A:239:ASN:HA	1.77	0.66
1:A:159:LYS:HE3	4:A:556:HOH:O	1.94	0.66
1:A:462:LYS:HB3	1:A:463:PRO:HD2	1.77	0.65
1:A:202:ARG:HG2	1:A:239:ASN:HB3	1.75	0.65
1:A:289:LYS:HZ1	1:A:460:PRO:HD3	1.62	0.63
1:A:398:ALA:HB2	4:A:658:HOH:O	1.99	0.62
2:B:3:GLC:H61	2:B:4:GLC:H5	1.81	0.62
1:A:201:PRO:C	1:A:202:ARG:HD2	2.19	0.62
1:A:387:ASP:OD2	1:A:389:THR:HB	1.99	0.62
1:A:376:ARG:NH2	4:A:514:HOH:O	2.32	0.62
1:A:131:ASN:HD22	1:A:132:LYS:N	1.98	0.61
1:A:5:SER:HB3	1:A:9:LEU:HD12	1.82	0.60
1:A:269:LYS:HE2	4:A:808:HOH:O	2.00	0.60
1:A:131:ASN:HD22	1:A:131:ASN:C	2.04	0.59
2:B:2:GLC:HO4	2:B:3:GLC:C1	2.08	0.59
1:A:253:THR:O	1:A:256:THR:HB	2.02	0.59
1:A:202:ARG:HG3	1:A:239:ASN:O	2.03	0.59
1:A:199:GLU:H	1:A:239:ASN:ND2	1.93	0.59
1:A:427:GLN:HG2	1:A:428:LYS:N	2.17	0.59
1:A:256:THR:HG22	1:A:258:LYS:H	1.69	0.57
1:A:199:GLU:O	1:A:200:PHE:C	2.43	0.57
1:A:381:ASN:ND2	1:A:419:LEU:HB3	2.17	0.56
1:A:459:THR:HG22	1:A:460:PRO:CD	2.32	0.55
1:A:131:ASN:ND2	1:A:133:GLU:H	2.05	0.55
1:A:7:MET:CE	1:A:403:VAL:HG13	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:TRP:CZ2	1:A:200:PHE:O	2.60	0.54
1:A:122:PHE:HB3	1:A:131:ASN:O	2.06	0.54
1:A:197:GLY:O	1:A:198:TRP:C	2.45	0.54
1:A:309:ALA:O	1:A:312:LEU:O	2.25	0.54
1:A:192:TYR:HB2	1:A:198:TRP:CD2	2.43	0.54
1:A:421:LEU:HD22	1:A:422:SER:N	2.22	0.54
1:A:57:GLY:HA2	1:A:108:PRO:HD3	1.91	0.53
1:A:289:LYS:HZ2	1:A:289:LYS:HB3	1.73	0.53
1:A:193:PRO:HG2	1:A:196:GLN:HB2	1.88	0.53
1:A:249:LYS:C	1:A:249:LYS:HD3	2.28	0.53
1:A:201:PRO:C	1:A:202:ARG:CD	2.72	0.52
1:A:202:ARG:HB2	1:A:239:ASN:HD22	1.72	0.52
1:A:7:MET:HE3	1:A:403:VAL:CG1	2.39	0.51
1:A:123:TYR:OH	1:A:137:VAL:HG23	2.09	0.51
1:A:289:LYS:CB	1:A:289:LYS:NZ	2.73	0.51
1:A:399:LYS:NZ	4:A:782:HOH:O	2.20	0.51
1:A:405:ASN:HD22	1:A:405:ASN:N	1.97	0.50
1:A:306:GLU:HG2	4:A:731:HOH:O	2.11	0.50
1:A:7:MET:HE3	1:A:403:VAL:HG13	1.93	0.50
1:A:201:PRO:HG3	4:A:506:HOH:O	2.11	0.50
1:A:381:ASN:HD21	1:A:419:LEU:CB	2.19	0.50
1:A:364:GLN:HG2	1:A:485:TRP:CE2	2.46	0.49
1:A:421:LEU:O	1:A:421:LEU:HD13	2.11	0.49
1:A:289:LYS:NZ	1:A:460:PRO:N	2.57	0.49
1:A:381:ASN:ND2	1:A:419:LEU:CB	2.75	0.49
1:A:289:LYS:HZ2	1:A:289:LYS:CB	2.26	0.48
1:A:339:LEU:HD23	1:A:339:LEU:C	2.33	0.48
1:A:202:ARG:CG	1:A:239:ASN:ND2	2.49	0.48
1:A:126:ARG:HH11	1:A:126:ARG:HG3	1.78	0.48
1:A:414:PHE:O	1:A:456:HIS:HE1	1.97	0.47
1:A:465:ALA:HB1	1:A:466:PRO:HD2	1.96	0.47
1:A:5:SER:HB3	1:A:9:LEU:HD11	1.93	0.47
1:A:15:TYR:CE1	1:A:175:ILE:HD11	2.49	0.47
1:A:431:PHE:O	1:A:435:LYS:HG3	2.15	0.47
1:A:140:ASP:HB3	1:A:151:ILE:CD1	2.45	0.47
1:A:399:LYS:CE	4:A:782:HOH:O	2.63	0.47
1:A:44:ALA:HB3	4:A:594:HOH:O	2.15	0.47
1:A:443:ALA:O	1:A:444:ASP:HB2	2.14	0.46
1:A:200:PHE:CD2	1:A:201:PRO:CD	2.81	0.46
1:A:291:LYS:HD2	1:A:376:ARG:NH1	2.31	0.46
1:A:27:ASN:HB2	4:A:702:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:340:ASN:HD21	1:A:380:GLU:HG3	1.81	0.46
1:A:340:ASN:HD22	1:A:340:ASN:C	2.20	0.45
1:A:252:GLY:O	1:A:255:VAL:HG22	2.14	0.45
1:A:241:VAL:HG22	4:A:609:HOH:O	2.16	0.45
1:A:107:ILE:HB	1:A:108:PRO:HD2	1.97	0.45
1:A:192:TYR:HB2	1:A:198:TRP:CE2	2.52	0.45
1:A:494:ASP:OD1	1:A:494:ASP:N	2.50	0.44
1:A:5:SER:HB2	1:A:407:GLY:HA2	1.99	0.44
1:A:427:GLN:OE1	1:A:429:SER:N	2.51	0.44
1:A:429:SER:O	1:A:433:ILE:HG12	2.17	0.44
1:A:122:PHE:CD1	1:A:132:LYS:HA	2.52	0.44
1:A:256:THR:CG2	4:A:723:HOH:O	2.66	0.44
1:A:200:PHE:CD1	1:A:201:PRO:CD	2.93	0.44
1:A:231:LEU:HD22	4:A:692:HOH:O	2.18	0.44
1:A:51:MET:HB2	1:A:87:GLN:HE21	1.83	0.44
1:A:399:LYS:HE3	1:A:403:VAL:HG23	2.00	0.43
1:A:283:LYS:HB3	1:A:283:LYS:HE3	1.33	0.43
1:A:24:ASN:HD22	1:A:30:GLU:CD	2.21	0.43
1:A:59:ILE:HD11	1:A:72:TYR:CD2	2.54	0.43
1:A:198:TRP:CZ3	1:A:200:PHE:HA	2.45	0.42
1:A:16:VAL:CG1	1:A:421:LEU:HB2	2.49	0.42
1:A:159:LYS:HD3	4:A:653:HOH:O	2.19	0.42
1:A:202:ARG:NH2	1:A:241:VAL:HG13	2.35	0.42
1:A:381:ASN:OD1	1:A:419:LEU:N	2.52	0.42
1:A:400:PRO:HB3	1:A:485:TRP:CD2	2.55	0.42
1:A:7:MET:CE	1:A:403:VAL:CG1	2.97	0.42
1:A:234:ASP:OD1	1:A:234:ASP:N	2.53	0.42
1:A:462:LYS:HB3	1:A:463:PRO:CD	2.46	0.41
1:A:198:TRP:CH2	1:A:200:PHE:CA	2.84	0.41
1:A:7:MET:HE1	1:A:403:VAL:HG13	2.01	0.41
1:A:194:GLN:HE21	1:A:194:GLN:HB2	1.56	0.41
1:A:440:LYS:HA	1:A:440:LYS:HD3	1.82	0.41
1:A:289:LYS:HZ1	1:A:460:PRO:CA	2.34	0.41
1:A:256:THR:HG23	4:A:723:HOH:O	2.20	0.41
1:A:451:PRO:HA	1:A:454:TYR:CE2	2.56	0.41
1:A:465:ALA:HB1	1:A:466:PRO:CD	2.50	0.40
1:A:16:VAL:HG11	1:A:421:LEU:HB2	2.02	0.40
1:A:46:GLY:O	1:A:442:HIS:HE1	2.04	0.40
1:A:312:LEU:HD23	1:A:312:LEU:N	2.36	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)
4:A:647:HOH:O	4:A:829:HOH:O[2_665]	2.19	0.01

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	489/495 (99%)	465 (95%)	21 (4%)	3 (1%)	25 15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	200	PHE
1	A	313	THR
1	A	198	TRP

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	422/425 (99%)	383 (91%)	39 (9%)	9 3

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	7	MET
1	A	27	ASN

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Mol	Chain	Res	Type
1	A	77	GLN
1	A	86	LEU
1	A	103	VAL
1	A	112	LEU
1	A	127	SER
1	A	130	ARG
1	A	131	ASN
1	A	162	ARG
1	A	170	GLU
1	A	199	GLU
1	A	200	PHE
1	A	202	ARG
1	A	231	LEU
1	A	249	LYS
1	A	283	LYS
1	A	289	LYS
1	A	313	THR
1	A	340	ASN
1	A	353	SER
1	A	403	VAL
1	A	405	ASN
1	A	406	ASN
1	A	421	LEU
1	A	426	LEU
1	A	427	GLN
1	A	428	LYS
1	A	429	SER
1	A	455	ASN
1	A	461	LEU
1	A	467	LYS
1	A	474	LEU
1	A	478	LYS
1	A	485	TRP
1	A	486	LEU
1	A	492	LYS
1	A	494	ASP

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	6	ASN
1	A	27	ASN

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Mol	Chain	Res	Type
1	A	87	GLN
1	A	131	ASN
1	A	194	GLN
1	A	239	ASN
1	A	268	ASN
1	A	307	ASN
1	A	340	ASN
1	A	405	ASN
1	A	432	ASN
1	A	455	ASN

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)

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	GLC	B	1	2	12,12,12	1.04	1 (8%)	17,17,17	1.64	4 (23%)
2	GLC	B	2	2	11,11,12	0.98	1 (9%)	15,15,17	1.51	4 (26%)
2	GLC	B	3	2	11,11,12	0.87	1 (9%)	15,15,17	1.87	4 (26%)
2	GLC	B	4	2	11,11,12	1.07	1 (9%)	15,15,17	1.45	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	1/1/5/5	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	-	0/2/19/22	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	4	GLC	O5-C1	2.52	1.47	1.43
2	B	1	GLC	C1-C2	-2.40	1.46	1.52
2	B	2	GLC	C4-C3	2.09	1.57	1.52
2	B	3	GLC	O2-C2	-2.06	1.39	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	3	GLC	O5-C1-C2	-4.59	103.69	110.77
2	B	4	GLC	C1-O5-C5	3.27	116.62	112.19
2	B	1	GLC	C1-O5-C5	-3.20	107.62	113.66
2	B	1	GLC	C1-C2-C3	-3.09	103.91	110.31
2	B	3	GLC	C1-C2-C3	-2.93	106.07	109.67
2	B	1	GLC	O5-C1-C2	-2.84	105.22	110.28
2	B	4	GLC	O2-C2-C1	2.71	114.70	109.15
2	B	2	GLC	O3-C3-C2	-2.63	104.96	109.99
2	B	1	GLC	O1-C1-O5	2.61	118.22	110.38
2	B	2	GLC	O4-C4-C3	2.51	116.15	110.35
2	B	3	GLC	C3-C4-C5	-2.36	106.03	110.24
2	B	2	GLC	C1-C2-C3	2.30	112.50	109.67
2	B	3	GLC	C1-O5-C5	-2.13	109.31	112.19
2	B	4	GLC	O6-C6-C5	2.11	118.53	111.29
2	B	2	GLC	C6-C5-C4	-2.04	108.23	113.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	1	GLC	C1

There are no torsion outliers.

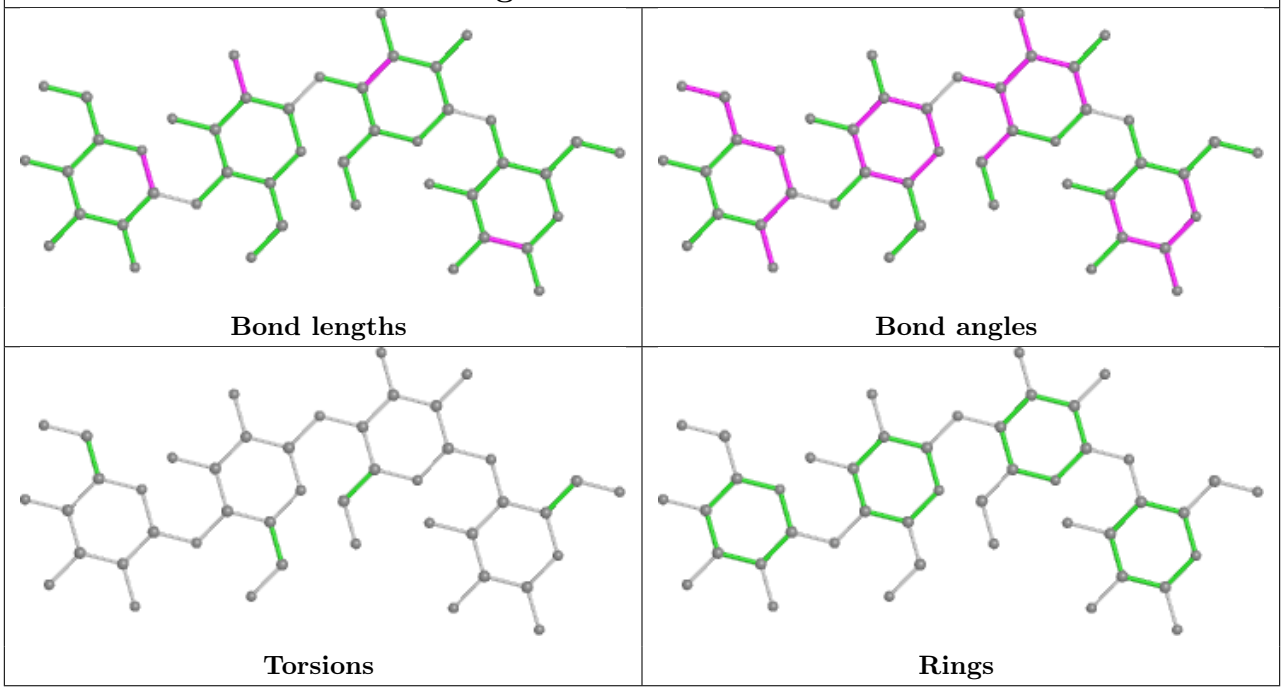
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3	GLC	2	0
2	B	2	GLC	2	0
2	B	4	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.

Oligosaccharide Chain B



5.6 Ligand geometry (i)

1 ligand is 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	$\#\|Z\| > 2$
3	SO4	A	860	-	4,4,4	0.38	0	6,6,6	0.42	0

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 (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.