



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 29, 2024 – 08:39 PM EST

PDB ID : 1FA2
Title : CRYSTAL STRUCTURE OF BETA-AMYLASE FROM SWEET POTATO
Authors : Lee, B.I.; Cheong, C.G.; Suh, S.W.
Deposited on : 2000-07-12
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

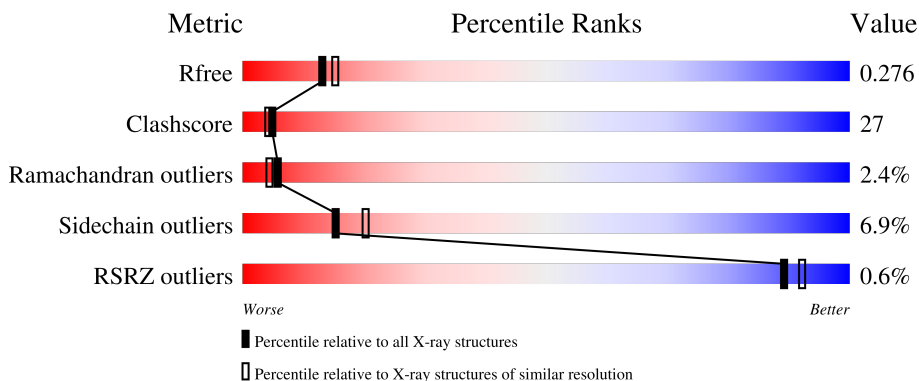
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	498	
2	B	2	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4072 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.

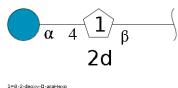
- Molecule 1 is a protein called BETA-AMYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	498	3904	2488	664	729	23	52	0	0

There are 2 discrepancies between the modelled and reference sequences:

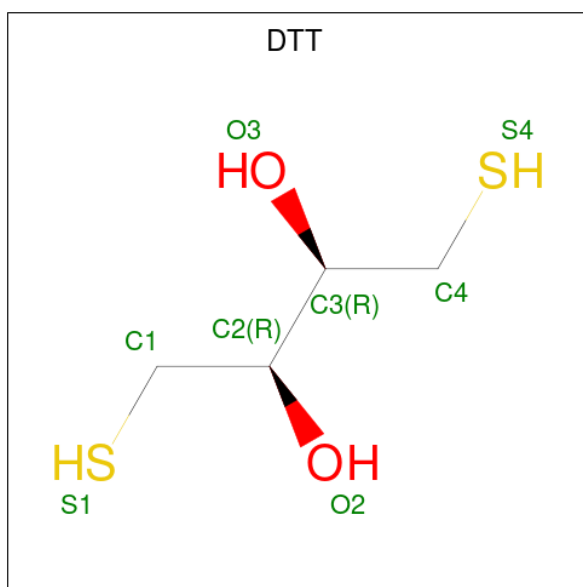
Chain	Residue	Modelled	Actual	Comment	Reference
A	399	LYS	ASN	SEE REMARK 999	UNP P10537
A	400	LEU	VAL	SEE REMARK 999	UNP P10537

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	B	2	22	12	10	0	0	0

- Molecule 3 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (three-letter code: DTT) (formula: C₄H₁₀O₂S₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			8	4	2	2		

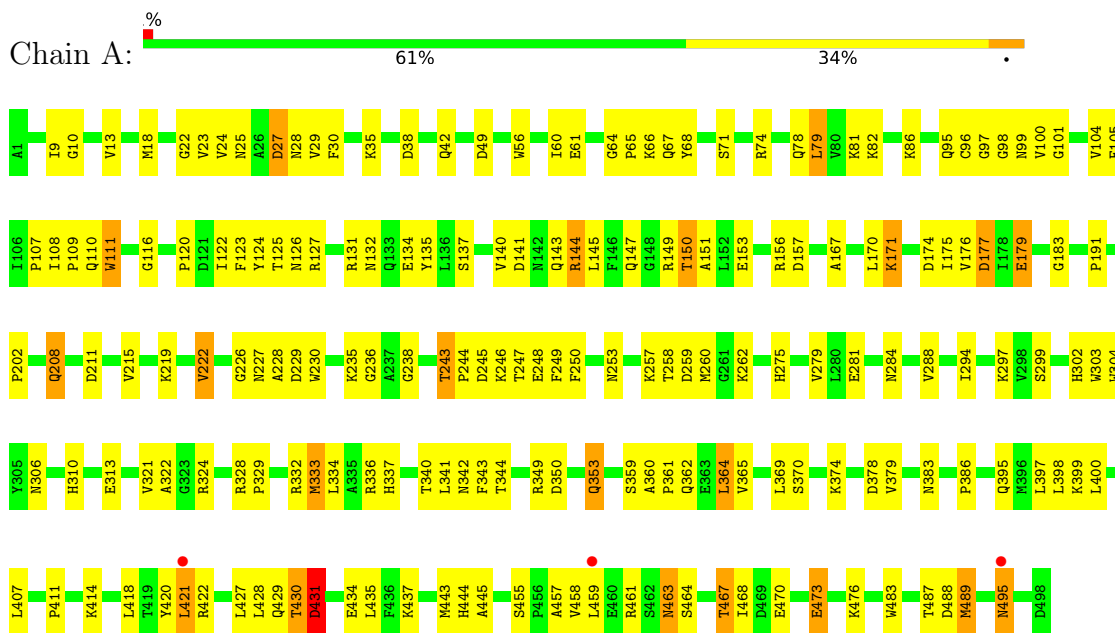
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	138	Total	O	0	0
			138	138		

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: BETA-AMYLASE



- Molecule 2: alpha-D-glucopyranose-(1-4)-2-deoxy-beta-D-arabino-hexopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	129.63Å 129.63Å 68.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.66 – 2.30 91.66 – 1.81	Depositor EDS
% Data completeness (in resolution range)	87.9 (91.66-2.30) 65.5 (91.66-1.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.12 (at 1.81Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.213 , 0.281 0.214 , 0.276	Depositor DCC
R_{free} test set	3475 reflections (9.93%)	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtrriage
Anisotropy	0.738	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 58.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4072	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: RR7, GLC, DTT

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	0.38	0/4006	0.63	2/5426 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	27	ASP	N-CA-C	-5.52	96.11	111.00
1	A	177	ASP	CB-CG-OD2	5.38	123.14	118.30

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	3904	0	3751	210	0
2	B	22	0	10	2	0
3	A	8	0	8	0	0
4	A	138	0	0	8	0
All	All	4072	0	3769	210	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 27.

All (210) 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:386:PRO:HA	1:A:421:LEU:HD22	1.25	1.19
1:A:383:ASN:HD21	1:A:421:LEU:HB3	1.16	1.10
1:A:171:LYS:NZ	1:A:171:LYS:HA	1.80	0.95
1:A:65:PRO:HG3	1:A:147:GLN:HE22	1.28	0.95
1:A:336:ARG:HD2	1:A:467:THR:HG23	1.48	0.94
1:A:332:ARG:HH12	1:A:473:GLU:HG2	1.34	0.92
1:A:421:LEU:HD12	1:A:422:ARG:H	1.35	0.92
1:A:421:LEU:HD12	1:A:422:ARG:N	1.85	0.91
1:A:65:PRO:HG3	1:A:147:GLN:NE2	1.87	0.89
1:A:350:ASP:H	1:A:362:GLN:NE2	1.70	0.88
1:A:350:ASP:N	1:A:362:GLN:HE21	1.72	0.87
1:A:461:ARG:HH21	1:A:461:ARG:HG3	1.41	0.85
1:A:383:ASN:ND2	1:A:421:LEU:HB3	1.90	0.85
1:A:350:ASP:H	1:A:362:GLN:HE21	0.88	0.84
1:A:332:ARG:NH1	1:A:473:GLU:HG2	1.92	0.84
1:A:150:THR:HG21	4:A:1100:HOH:O	1.76	0.83
1:A:258:THR:HG22	1:A:260:MET:H	1.44	0.83
1:A:284:ASN:ND2	1:A:461:ARG:HH22	1.76	0.82
1:A:171:LYS:HA	1:A:171:LYS:HZ3	1.44	0.80
1:A:349:ARG:HE	1:A:395:GLN:HG2	1.47	0.80
1:A:321:VAL:HG12	1:A:322:ALA:H	1.47	0.79
1:A:247:THR:HG22	1:A:249:PHE:H	1.47	0.79
1:A:321:VAL:HG12	1:A:322:ALA:N	1.98	0.79
1:A:95:GLN:HG3	1:A:107:PRO:HA	1.66	0.77
1:A:176:VAL:HG12	1:A:177:ASP:OD2	1.84	0.76
1:A:332:ARG:HH12	1:A:473:GLU:CG	2.02	0.73
1:A:27:ASP:O	1:A:29:VAL:N	2.21	0.71
1:A:369:LEU:HD22	1:A:379:VAL:HG11	1.71	0.71
1:A:13:VAL:HG11	1:A:400:LEU:HD22	1.73	0.71
1:A:78:GLN:O	1:A:82:LYS:HG3	1.91	0.70
1:A:421:LEU:CD1	1:A:422:ARG:H	2.05	0.70
1:A:284:ASN:ND2	1:A:461:ARG:HH12	1.89	0.70
1:A:333:MET:HE1	1:A:334:LEU:HA	1.74	0.70
1:A:61:GLU:HG2	1:A:111:TRP:CZ3	2.27	0.70
1:A:284:ASN:HB2	1:A:461:ARG:HH12	1.53	0.70
1:A:284:ASN:CB	1:A:461:ARG:HH12	2.04	0.69
1:A:284:ASN:ND2	1:A:461:ARG:NH2	2.41	0.68
1:A:470:GLU:OE1	1:A:470:GLU:HA	1.94	0.68
1:A:284:ASN:HD22	1:A:461:ARG:HH12	1.41	0.68
1:A:167:ALA:O	1:A:171:LYS:HG2	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:THR:HG22	1:A:249:PHE:N	2.10	0.67
1:A:132:ASN:OD1	1:A:134:GLU:HG2	1.95	0.67
1:A:98:GLY:HA3	1:A:105:PHE:CD1	2.30	0.67
1:A:141:ASP:OD2	1:A:275:HIS:HD2	1.78	0.66
1:A:56:TRP:CH2	1:A:96:CYS:HB2	2.30	0.66
1:A:243:THR:HG22	1:A:246:LYS:HG3	1.77	0.66
1:A:156:ARG:HG3	1:A:157:ASP:N	2.11	0.65
1:A:294:ILE:CD1	1:A:461:ARG:HH11	2.10	0.65
1:A:445:ALA:HA	1:A:489:MET:HE2	1.77	0.65
1:A:463:ASN:HD22	1:A:464:SER:N	1.95	0.65
1:A:171:LYS:HA	1:A:171:LYS:HZ2	1.62	0.65
1:A:222:VAL:HG21	1:A:230:TRP:HB2	1.79	0.65
1:A:179:GLU:OE2	1:A:297:LYS:HE2	1.97	0.64
1:A:461:ARG:HG3	1:A:461:ARG:NH2	2.12	0.64
1:A:150:THR:HG22	1:A:153:GLU:H	1.63	0.64
1:A:302:HIS:HE1	2:B:2:GLC:O2	1.80	0.63
1:A:430:THR:O	1:A:434:GLU:HG2	1.98	0.63
1:A:302:HIS:HD2	1:A:359:SER:OG	1.81	0.63
1:A:343:PHE:HZ	1:A:364:LEU:HD13	1.63	0.63
1:A:397:LEU:HA	1:A:400:LEU:HD12	1.80	0.63
1:A:150:THR:HG22	1:A:153:GLU:HG3	1.81	0.62
1:A:141:ASP:HA	1:A:151:ALA:HB3	1.82	0.62
1:A:321:VAL:CG1	1:A:322:ALA:H	2.13	0.62
1:A:383:ASN:HD21	1:A:421:LEU:CB	2.03	0.62
1:A:463:ASN:HD22	1:A:464:SER:H	1.46	0.62
1:A:400:LEU:HD13	1:A:443:MET:SD	2.40	0.61
1:A:61:GLU:HG2	1:A:111:TRP:HZ3	1.65	0.61
1:A:495:ASN:N	1:A:495:ASN:HD22	1.96	0.61
1:A:222:VAL:CG2	1:A:230:TRP:HB2	2.31	0.60
1:A:400:LEU:CD1	1:A:443:MET:SD	2.90	0.60
1:A:284:ASN:CG	1:A:461:ARG:HH22	2.03	0.60
1:A:140:VAL:CG2	1:A:143:GLN:HB2	2.31	0.59
1:A:258:THR:HG22	1:A:260:MET:N	2.15	0.59
1:A:341:LEU:HD23	1:A:342:ASN:N	2.17	0.59
1:A:116:GLY:HA3	1:A:122:ILE:HD11	1.83	0.59
1:A:191:PRO:HG2	1:A:208:GLN:HG2	1.83	0.59
1:A:294:ILE:HD11	1:A:461:ARG:HH11	1.67	0.59
1:A:468:ILE:HG22	1:A:470:GLU:N	2.16	0.59
1:A:258:THR:O	1:A:262:LYS:HG3	2.03	0.59
1:A:341:LEU:HD23	1:A:341:LEU:C	2.24	0.57
1:A:284:ASN:ND2	1:A:461:ARG:NH1	2.51	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:ASN:HD21	1:A:421:LEU:HD23	1.70	0.57
1:A:383:ASN:ND2	1:A:421:LEU:HD23	2.19	0.57
1:A:13:VAL:HG22	1:A:414:LYS:HD3	1.86	0.57
1:A:468:ILE:C	1:A:470:GLU:H	2.07	0.57
1:A:22:GLY:O	1:A:24:VAL:N	2.38	0.56
1:A:211:ASP:O	1:A:215:VAL:HG23	2.05	0.56
1:A:333:MET:CE	1:A:337:HIS:HD2	2.18	0.56
1:A:365:VAL:O	1:A:369:LEU:HG	2.06	0.56
1:A:284:ASN:HD22	1:A:461:ARG:NH1	2.01	0.55
1:A:143:GLN:HB3	1:A:145:LEU:HD13	1.88	0.55
1:A:110:GLN:HG3	4:A:1040:HOH:O	2.06	0.55
1:A:257:LYS:O	1:A:262:LYS:HE3	2.07	0.55
1:A:140:VAL:HG23	1:A:143:GLN:HB2	1.88	0.55
1:A:127:ARG:NH1	1:A:235:LYS:HA	2.22	0.55
1:A:421:LEU:CD1	1:A:422:ARG:N	2.66	0.54
1:A:81:LYS:NZ	1:A:174:ASP:OD2	2.39	0.54
1:A:227:ASN:O	1:A:229:ASP:N	2.40	0.54
1:A:243:THR:HG22	1:A:246:LYS:H	1.72	0.54
1:A:420:TYR:HE2	1:A:427:LEU:HD22	1.71	0.54
1:A:64:GLY:O	1:A:67:GLN:HG2	2.08	0.54
1:A:353:GLN:HG2	1:A:359:SER:OG	2.09	0.53
1:A:487:THR:HG22	1:A:488:ASP:N	2.23	0.53
1:A:243:THR:CG2	1:A:246:LYS:HG3	2.39	0.53
1:A:253:ASN:ND2	1:A:257:LYS:HE2	2.24	0.52
1:A:49:ASP:OD1	1:A:444:HIS:HE1	1.92	0.52
1:A:247:THR:HG21	1:A:250:PHE:HD1	1.74	0.52
1:A:18:MET:HG2	4:A:1075:HOH:O	2.10	0.52
1:A:306:ASN:ND2	1:A:360:ALA:HB2	2.25	0.52
1:A:495:ASN:N	1:A:495:ASN:ND2	2.58	0.52
1:A:247:THR:HG22	1:A:248:GLU:N	2.25	0.51
1:A:383:ASN:HD21	1:A:421:LEU:CD2	2.23	0.51
1:A:395:GLN:OE1	1:A:399:LYS:HE3	2.11	0.51
1:A:310:HIS:HD2	1:A:313:GLU:OE1	1.93	0.51
1:A:66:LYS:HE2	1:A:149:ARG:HD3	1.93	0.51
1:A:333:MET:HE2	1:A:337:HIS:HD2	1.75	0.51
1:A:179:GLU:HG2	1:A:297:LYS:CE	2.41	0.50
1:A:238:GLY:HA2	4:A:1087:HOH:O	2.10	0.50
1:A:86:LYS:HG2	1:A:176:VAL:HG21	1.93	0.50
1:A:236:GLY:O	1:A:247:THR:HG23	2.11	0.50
1:A:247:THR:CG2	1:A:249:PHE:H	2.22	0.50
1:A:468:ILE:CG2	1:A:470:GLU:HB2	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:ASN:HB2	1:A:461:ARG:NH1	2.26	0.50
1:A:400:LEU:HD21	1:A:418:LEU:HB2	1.92	0.50
1:A:22:GLY:C	1:A:24:VAL:N	2.65	0.50
1:A:35:LYS:O	1:A:38:ASP:HB2	2.12	0.49
1:A:431:ASP:O	1:A:434:GLU:HB2	2.12	0.49
1:A:302:HIS:CE1	2:B:2:GLC:O2	2.62	0.49
1:A:61:GLU:HG2	1:A:111:TRP:CH2	2.48	0.49
1:A:324:ARG:HG2	1:A:324:ARG:HH11	1.78	0.49
1:A:18:MET:SD	1:A:421:LEU:O	2.71	0.49
1:A:284:ASN:ND2	1:A:464:SER:OG	2.45	0.49
1:A:321:VAL:CG1	1:A:322:ALA:N	2.66	0.49
1:A:247:THR:HG21	1:A:250:PHE:CD1	2.47	0.49
1:A:284:ASN:O	1:A:288:VAL:HG13	2.13	0.49
1:A:60:ILE:HG22	1:A:68:TYR:HB3	1.94	0.48
1:A:42:GLN:OE1	1:A:428:LEU:HG	2.13	0.48
1:A:61:GLU:O	1:A:111:TRP:HZ3	1.96	0.48
1:A:340:THR:HG23	1:A:378:ASP:O	2.13	0.48
1:A:99:ASN:ND2	1:A:100:VAL:H	2.11	0.48
1:A:411:PRO:HG2	1:A:414:LYS:HB2	1.96	0.48
1:A:120:PRO:HB2	1:A:131:ARG:CZ	2.43	0.48
1:A:445:ALA:HA	1:A:489:MET:CE	2.43	0.48
1:A:473:GLU:O	1:A:476:LYS:HG2	2.13	0.48
1:A:258:THR:CG2	1:A:259:ASP:N	2.77	0.48
1:A:258:THR:HG22	1:A:259:ASP:N	2.29	0.48
1:A:253:ASN:HD21	1:A:257:LYS:HE2	1.77	0.47
1:A:468:ILE:HG22	1:A:470:GLU:HB2	1.95	0.47
1:A:226:GLY:C	1:A:227:ASN:HD22	2.18	0.47
1:A:144:ARG:C	1:A:145:LEU:HD12	2.35	0.47
1:A:365:VAL:HG11	1:A:399:LYS:HD3	1.95	0.47
1:A:341:LEU:C	1:A:341:LEU:CD2	2.82	0.47
1:A:333:MET:HE2	1:A:337:HIS:CD2	2.50	0.47
1:A:421:LEU:CG	1:A:422:ARG:H	2.28	0.47
1:A:123:PHE:HB3	1:A:132:ASN:O	2.15	0.47
1:A:343:PHE:CZ	1:A:364:LEU:HD13	2.48	0.46
1:A:22:GLY:C	1:A:24:VAL:H	2.19	0.46
1:A:65:PRO:HG2	4:A:1081:HOH:O	2.15	0.46
1:A:170:LEU:HD23	1:A:175:ILE:HB	1.97	0.46
1:A:243:THR:HG23	1:A:245:ASP:H	1.80	0.46
1:A:9:ILE:HG13	1:A:10:GLY:N	2.29	0.46
1:A:328:ARG:N	1:A:329:PRO:HD2	2.31	0.46
1:A:71:SER:O	1:A:74:ARG:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:SER:OG	1:A:361:PRO:HD3	2.17	0.45
1:A:140:VAL:HG22	1:A:143:GLN:HB2	1.97	0.45
1:A:400:LEU:HD11	1:A:443:MET:SD	2.56	0.45
1:A:124:TYR:HA	4:A:1022:HOH:O	2.16	0.45
1:A:284:ASN:CG	1:A:461:ARG:HH12	2.19	0.45
1:A:349:ARG:NE	1:A:395:GLN:HG2	2.25	0.44
1:A:420:TYR:CZ	1:A:421:LEU:HD21	2.52	0.44
1:A:487:THR:HG22	1:A:489:MET:H	1.82	0.44
1:A:284:ASN:ND2	1:A:461:ARG:CZ	2.81	0.44
1:A:370:SER:O	1:A:374:LYS:HG3	2.18	0.44
1:A:470:GLU:OE1	1:A:470:GLU:CA	2.63	0.44
1:A:333:MET:HE1	1:A:334:LEU:HD23	2.00	0.43
1:A:467:THR:O	1:A:468:ILE:HG13	2.19	0.43
1:A:247:THR:CG2	1:A:248:GLU:N	2.81	0.43
1:A:303:TRP:O	1:A:304:TRP:HB2	2.18	0.43
1:A:116:GLY:HA3	1:A:122:ILE:CD1	2.48	0.43
1:A:179:GLU:HG2	1:A:297:LYS:HE2	2.00	0.43
1:A:281:GLU:HG3	1:A:337:HIS:HE1	1.84	0.43
1:A:30:PHE:CZ	1:A:79:LEU:HD12	2.54	0.42
1:A:299:SER:HA	1:A:344:THR:OG1	2.19	0.42
1:A:353:GLN:HG2	1:A:359:SER:CB	2.49	0.42
1:A:111:TRP:CD1	1:A:111:TRP:C	2.92	0.42
1:A:10:GLY:O	1:A:414:LYS:HE2	2.19	0.42
1:A:49:ASP:OD1	1:A:444:HIS:CE1	2.73	0.42
1:A:125:THR:HG22	1:A:126:ASN:O	2.19	0.42
1:A:294:ILE:O	1:A:294:ILE:HG13	2.19	0.42
1:A:461:ARG:NH2	1:A:461:ARG:CG	2.77	0.42
1:A:468:ILE:C	1:A:470:GLU:N	2.73	0.42
1:A:144:ARG:O	1:A:145:LEU:HD12	2.18	0.42
1:A:226:GLY:C	1:A:227:ASN:ND2	2.73	0.42
1:A:275:HIS:O	1:A:279:VAL:HG23	2.20	0.42
1:A:108:ILE:O	1:A:109:PRO:C	2.58	0.41
1:A:179:GLU:HG2	1:A:297:LYS:HE3	2.02	0.41
1:A:332:ARG:HH12	1:A:473:GLU:CD	2.23	0.41
1:A:468:ILE:HG22	1:A:470:GLU:H	1.85	0.41
1:A:333:MET:CE	1:A:337:HIS:CD2	3.01	0.41
1:A:253:ASN:HA	1:A:257:LYS:HD3	2.03	0.41
1:A:18:MET:HB2	1:A:421:LEU:O	2.21	0.41
1:A:60:ILE:O	1:A:68:TYR:HA	2.20	0.41
1:A:135:TYR:CD2	1:A:183:GLY:HA2	2.56	0.41
1:A:244:PRO:HD3	1:A:304:TRP:CZ2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ILE:HG23	4:A:1093:HOH:O	2.21	0.41
1:A:137:SER:O	1:A:140:VAL:HG12	2.20	0.40
1:A:284:ASN:CB	1:A:461:ARG:NH1	2.78	0.40
1:A:294:ILE:HD13	1:A:461:ARG:HH11	1.85	0.40
1:A:153:GLU:HG2	1:A:156:ARG:HH22	1.86	0.40
1:A:122:ILE:CG2	4:A:1093:HOH:O	2.69	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	496/498 (100%)	449 (90%)	35 (7%)	12 (2%)	6 4

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	97	GLY
1	A	104	VAL
1	A	228	ALA
1	A	421	LEU
1	A	455	SER
1	A	457	ALA
1	A	101	GLY
1	A	430	THR
1	A	458	VAL
1	A	23	VAL
1	A	431	ASP

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	406/416 (98%)	378 (93%)	28 (7%)	15 20

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	79	LEU
1	A	111	TRP
1	A	144	ARG
1	A	150	THR
1	A	171	LYS
1	A	179	GLU
1	A	202	PRO
1	A	208	GLN
1	A	219	LYS
1	A	222	VAL
1	A	243	THR
1	A	333	MET
1	A	353	GLN
1	A	364	LEU
1	A	398	LEU
1	A	407	LEU
1	A	429	GLN
1	A	431	ASP
1	A	435	LEU
1	A	437	LYS
1	A	459	LEU
1	A	463	ASN
1	A	467	THR
1	A	473	GLU
1	A	483	TRP
1	A	489	MET
1	A	495	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	67	GLN
1	A	99	ASN
1	A	110	GLN
1	A	143	GLN
1	A	147	GLN
1	A	208	GLN
1	A	227	ASN
1	A	253	ASN
1	A	275	HIS
1	A	278	GLN
1	A	284	ASN
1	A	302	HIS
1	A	306	ASN
1	A	310	HIS
1	A	337	HIS
1	A	362	GLN
1	A	366	GLN
1	A	429	GLN
1	A	432	ASN
1	A	444	HIS
1	A	463	ASN
1	A	495	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](#)

2 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	RR7	B	1	2	11,11,11	1.63	2 (18%)	13,15,15	3.47	2 (15%)
2	GLC	B	2	2	11,11,12	2.04	2 (18%)	15,15,17	1.49	4 (26%)

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	RR7	B	1	2	-	2/2/18/18	0/1/1/1
2	GLC	B	2	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	2	GLC	O5-C1	5.09	1.51	1.43
2	B	1	RR7	C3-C4	3.93	1.58	1.52
2	B	2	GLC	C1-C2	2.49	1.57	1.52
2	B	1	RR7	O4-C4	2.18	1.48	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	RR7	C3-C4-C5	-11.58	98.43	109.97
2	B	1	RR7	C2-C3-C4	3.17	115.29	110.69
2	B	2	GLC	O5-C5-C6	-2.77	102.87	107.20
2	B	2	GLC	C1-C2-C3	2.63	112.90	109.67
2	B	2	GLC	O5-C1-C2	-2.15	107.46	110.77
2	B	2	GLC	C1-O5-C5	2.09	115.02	112.19

There are no chirality outliers.

All (2) torsion outliers are listed below:

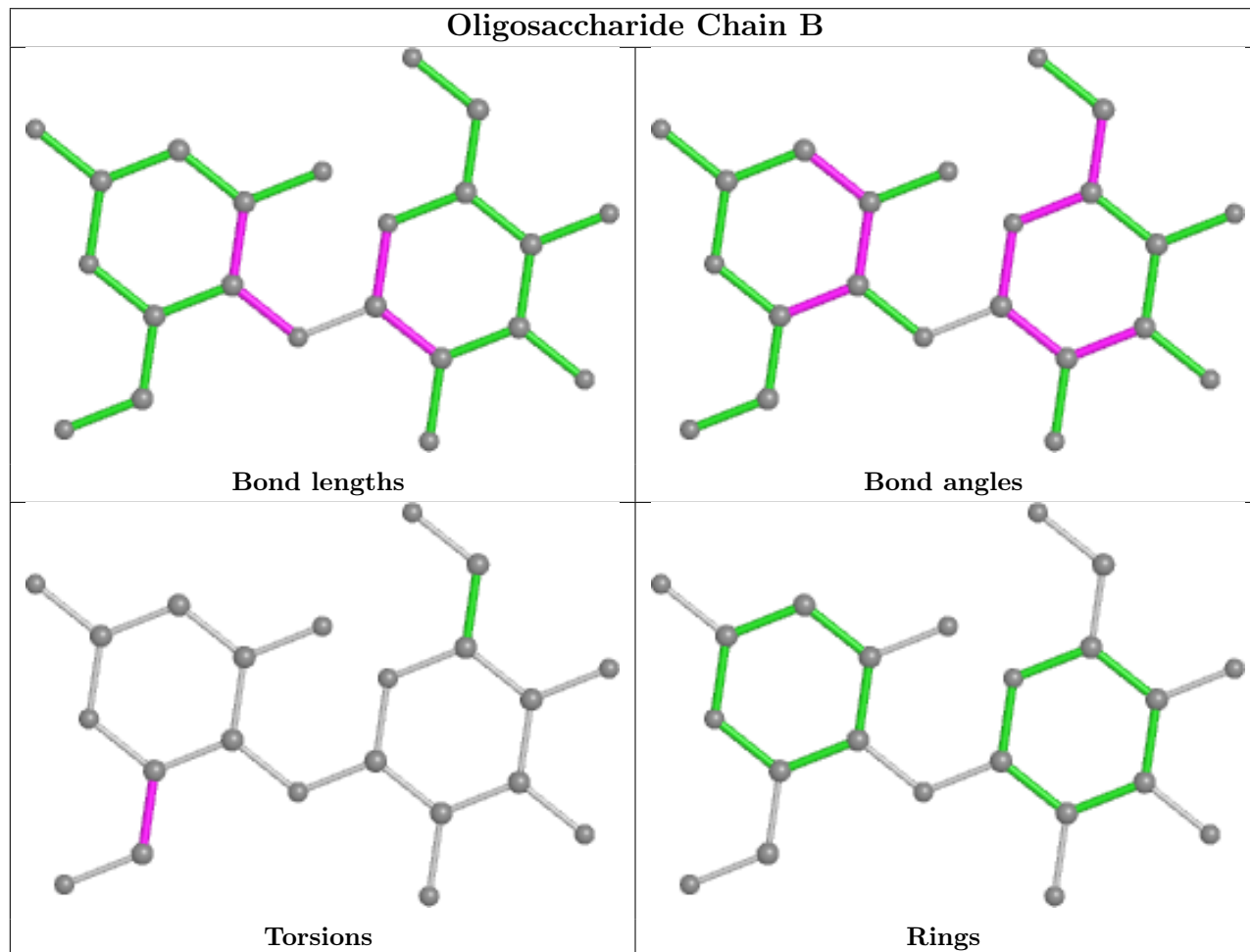
Mol	Chain	Res	Type	Atoms
2	B	1	RR7	O5-C5-C6-O6
2	B	1	RR7	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	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](#)

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	DTT	A	1409	-	7,7,7	2.16	2 (28%)	4,8,8	0.92	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
3	DTT	A	1409	-	-	0/8/8/8	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1409	DTT	C3-C2	4.13	1.63	1.52
3	A	1409	DTT	C4-S4	3.45	1.88	1.81

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](#)

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	485/498 (97%)	-0.42	3 (0%) 89 92	25, 41, 63, 76	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	459	LEU	2.7
1	A	495	ASN	2.1
1	A	421	LEU	2.1

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

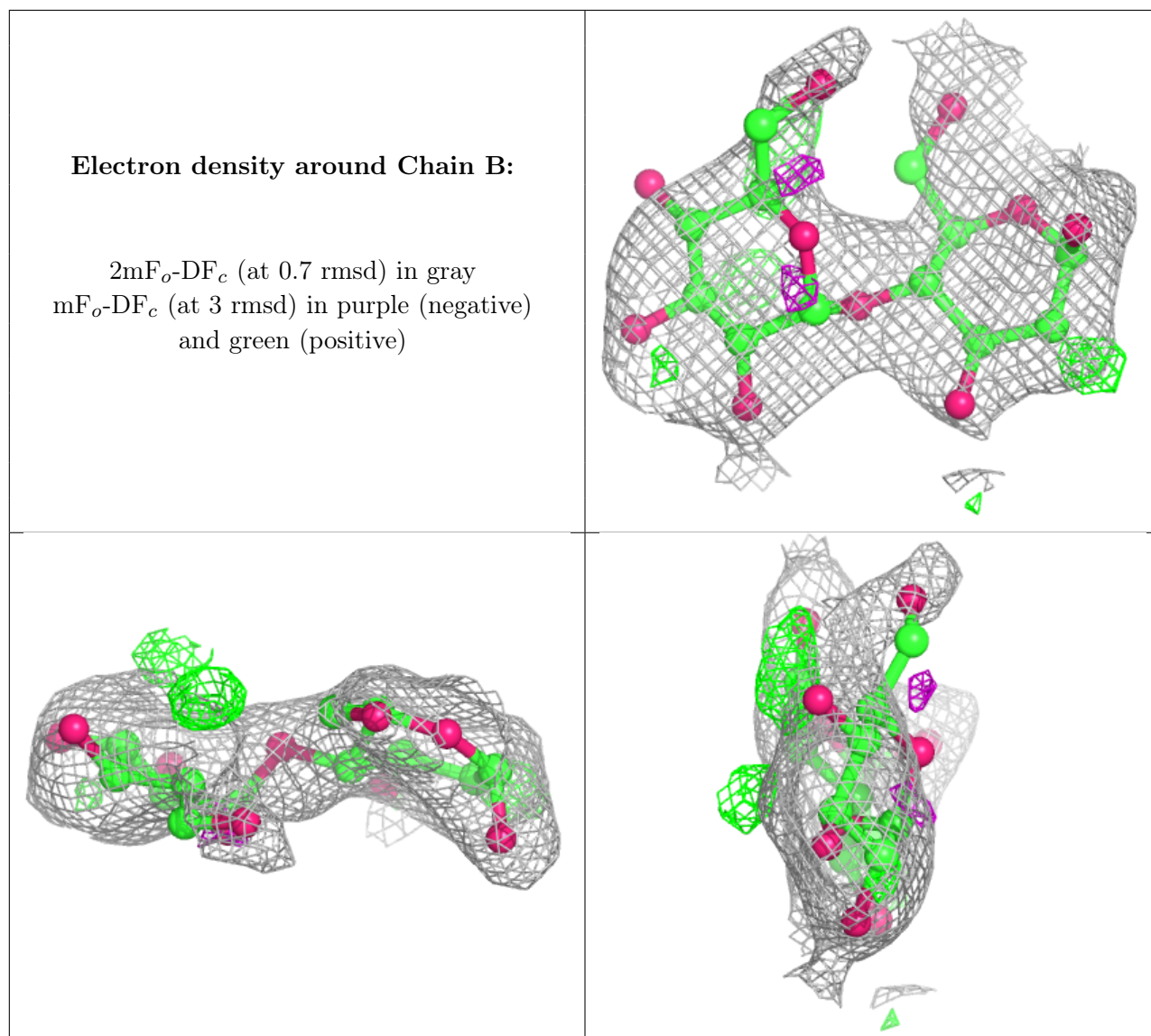
There are no non-standard protein/DNA/RNA residues in this entry.

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	B	2	11/12	0.67	0.27	75,77,81,82	0
2	RR7	B	1	11/11	0.73	0.17	57,67,73,73	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(Å ²)	Q<0.9
3	DTT	A	1409	8/8	0.78	0.14	80,82,83,87	0

6.5 Other polymers [i](#)

There are no such residues in this entry.