



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 2, 2024 – 01:10 pm GMT

PDB ID : 5CQ1
Title : Disproportionating enzyme 1 from Arabidopsis - cycloamylose soak
Authors : O'Neill, E.C.; Stevenson, C.E.M.; Tantanarat, K.; Latousakis, D.; Donaldson, M.I.; Rejzek, M.; Limpaseni, T.; Smith, A.M.; Field, R.A.; Lawson, D.M.
Deposited on : 2015-07-21
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 : 1.8.4, CSD as541be (2020)
Xtrriage (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

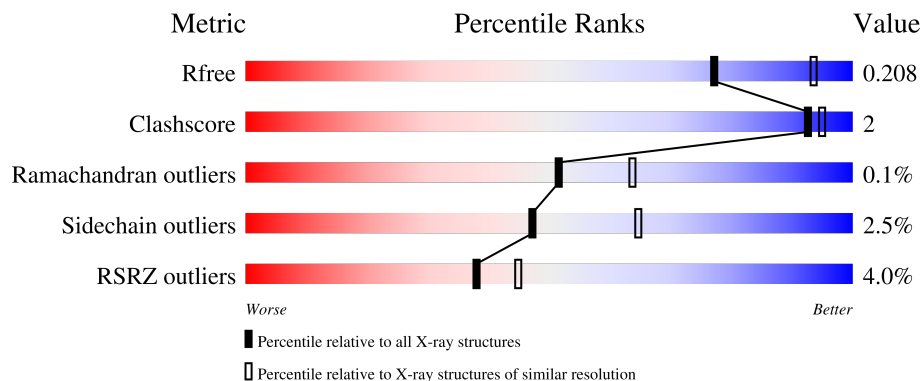
1 Overall quality at a glance

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	564	 2% 88% 9%
1	B	564	 5% 86% 5% 9%
2	C	2	 100%
2	D	2	 100%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8417 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 4-alpha-glucanotransferase DPE1, chloroplastic/amyloplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	514	4006	2570	678	747	11	0	1	0
1	B	511	3982	2557	678	735	12	0	1	0

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	MET	-	initiating methionine	UNP Q9LV91
A	14	HIS	-	expression tag	UNP Q9LV91
A	15	HIS	-	expression tag	UNP Q9LV91
A	16	HIS	-	expression tag	UNP Q9LV91
A	17	HIS	-	expression tag	UNP Q9LV91
A	18	HIS	-	expression tag	UNP Q9LV91
A	19	HIS	-	expression tag	UNP Q9LV91
A	20	GLY	-	expression tag	UNP Q9LV91
A	21	LYS	-	expression tag	UNP Q9LV91
A	22	PRO	-	expression tag	UNP Q9LV91
A	23	ILE	-	expression tag	UNP Q9LV91
A	24	PRO	-	expression tag	UNP Q9LV91
A	25	ASN	-	expression tag	UNP Q9LV91
A	26	PRO	-	expression tag	UNP Q9LV91
A	27	LEU	-	expression tag	UNP Q9LV91
A	28	LEU	-	expression tag	UNP Q9LV91
A	29	GLY	-	expression tag	UNP Q9LV91
A	30	LEU	-	expression tag	UNP Q9LV91
A	31	ASP	-	expression tag	UNP Q9LV91
A	32	SER	-	expression tag	UNP Q9LV91
A	33	THR	-	expression tag	UNP Q9LV91
A	34	GLU	-	expression tag	UNP Q9LV91
A	35	ASN	-	expression tag	UNP Q9LV91
A	36	LEU	-	expression tag	UNP Q9LV91
A	37	TYR	-	expression tag	UNP Q9LV91

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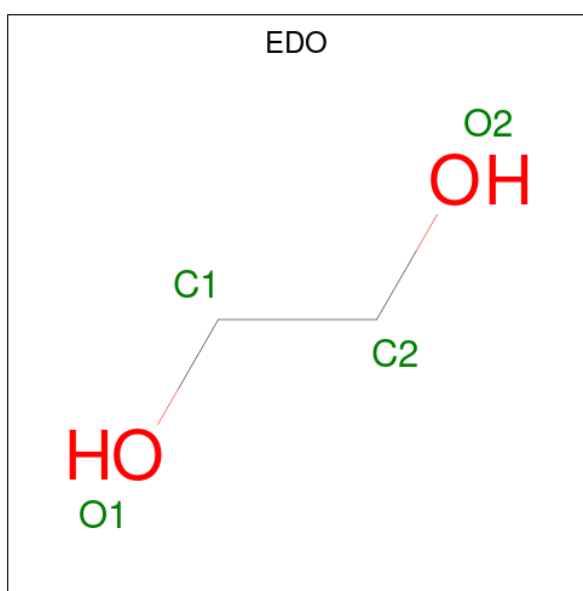
Chain	Residue	Modelled	Actual	Comment	Reference
A	38	PHE	-	expression tag	UNP Q9LV91
A	39	GLN	-	expression tag	UNP Q9LV91
A	40	GLY	-	expression tag	UNP Q9LV91
A	41	ILE	-	expression tag	UNP Q9LV91
A	42	ASP	-	expression tag	UNP Q9LV91
A	43	PRO	-	expression tag	UNP Q9LV91
A	44	PHE	-	expression tag	UNP Q9LV91
A	45	THR	-	expression tag	UNP Q9LV91
B	13	MET	-	initiating methionine	UNP Q9LV91
B	14	HIS	-	expression tag	UNP Q9LV91
B	15	HIS	-	expression tag	UNP Q9LV91
B	16	HIS	-	expression tag	UNP Q9LV91
B	17	HIS	-	expression tag	UNP Q9LV91
B	18	HIS	-	expression tag	UNP Q9LV91
B	19	HIS	-	expression tag	UNP Q9LV91
B	20	GLY	-	expression tag	UNP Q9LV91
B	21	LYS	-	expression tag	UNP Q9LV91
B	22	PRO	-	expression tag	UNP Q9LV91
B	23	ILE	-	expression tag	UNP Q9LV91
B	24	PRO	-	expression tag	UNP Q9LV91
B	25	ASN	-	expression tag	UNP Q9LV91
B	26	PRO	-	expression tag	UNP Q9LV91
B	27	LEU	-	expression tag	UNP Q9LV91
B	28	LEU	-	expression tag	UNP Q9LV91
B	29	GLY	-	expression tag	UNP Q9LV91
B	30	LEU	-	expression tag	UNP Q9LV91
B	31	ASP	-	expression tag	UNP Q9LV91
B	32	SER	-	expression tag	UNP Q9LV91
B	33	THR	-	expression tag	UNP Q9LV91
B	34	GLU	-	expression tag	UNP Q9LV91
B	35	ASN	-	expression tag	UNP Q9LV91
B	36	LEU	-	expression tag	UNP Q9LV91
B	37	TYR	-	expression tag	UNP Q9LV91
B	38	PHE	-	expression tag	UNP Q9LV91
B	39	GLN	-	expression tag	UNP Q9LV91
B	40	GLY	-	expression tag	UNP Q9LV91
B	41	ILE	-	expression tag	UNP Q9LV91
B	42	ASP	-	expression tag	UNP Q9LV91
B	43	PRO	-	expression tag	UNP Q9LV91
B	44	PHE	-	expression tag	UNP Q9LV91
B	45	THR	-	expression tag	UNP Q9LV91

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



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

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



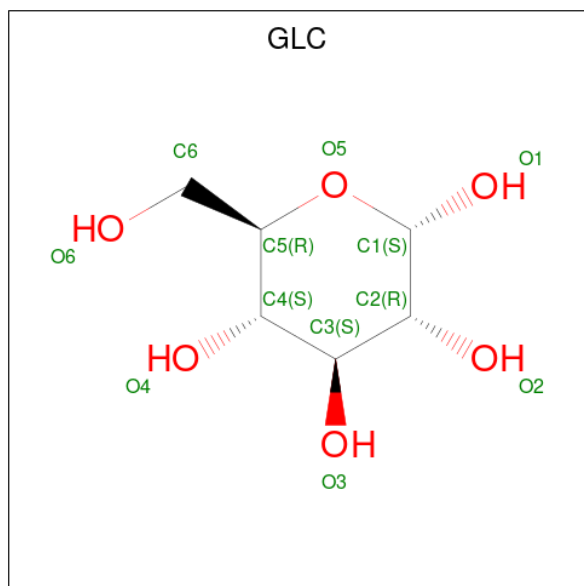
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 4 is alpha-D-glucopyranose (three-letter code: GLC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			12	6	6		

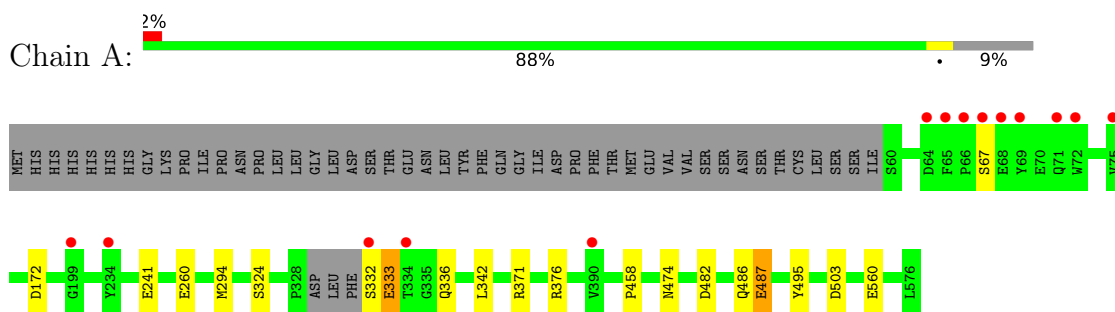
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	158	Total 158	O 158	0	0
5	B	151	Total 151	O 151	0	0

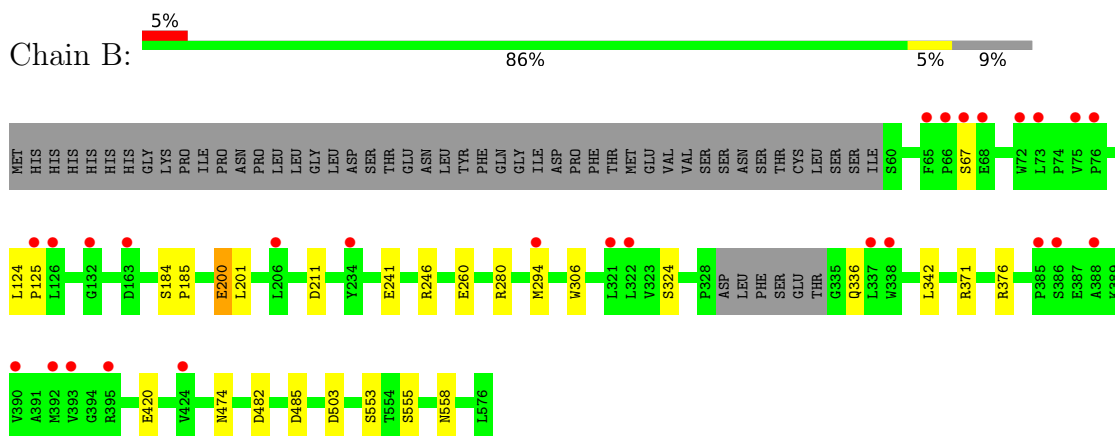
3 Residue-property plots [i](#)

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: 4-alpha-glucanotransferase DPE1, chloroplastic/amyloplastic



- Molecule 1: 4-alpha-glucanotransferase DPE1, chloroplastic/amyloplastic



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

Chain C: 

B6C1
GLC2

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

Chain D: 

B6C1
GLC2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	69.89Å 73.35Å 79.39Å 65.28° 69.49° 66.86°	Depositor
Resolution (Å)	32.03 – 2.30 32.03 – 2.30	Depositor EDS
% Data completeness (in resolution range)	92.4 (32.03-2.30) 92.5 (32.03-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.29Å)	Xtrriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.176 , 0.205 0.183 , 0.208	Depositor DCC
R_{free} test set	2688 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtrriage
Anisotropy	0.651	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8417	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 8.70% 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: BGC, GLC, EDO

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	0.59	0/4122	0.69	0/5615
1	B	0.59	0/4099	0.70	1/5584 (0.0%)
All	All	0.59	0/8221	0.70	1/11199 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	280	ARG	NE-CZ-NH2	-6.21	117.20	120.30

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	4006	0	3812	8	0
1	B	3982	0	3784	12	0
2	C	22	0	19	0	0
2	D	22	0	19	0	0
3	A	32	0	48	3	0
3	B	32	0	48	3	0
4	B	12	0	12	1	0
5	A	158	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	151	0	0	3	0
All	All	8417	0	7742	24	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 2.

All (24) 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)
3:A:610:EDO:H11	5:A:733:HOH:O	1.64	0.96
3:B:611:EDO:H22	5:B:754:HOH:O	1.69	0.90
3:A:610:EDO:C1	5:A:733:HOH:O	2.34	0.67
1:B:371:ARG:NH1	5:B:704:HOH:O	2.30	0.62
3:B:611:EDO:C2	5:B:754:HOH:O	2.36	0.60
1:B:474:ASN:O	3:B:609:EDO:H11	2.04	0.58
1:B:200:GLU:HG2	1:B:201:LEU:N	2.21	0.56
1:A:371:ARG:NH1	5:A:705:HOH:O	2.44	0.51
1:B:294:MET:O	1:B:294:MET:HG3	2.11	0.51
1:A:332:SER:O	1:A:333:GLU:HB2	2.12	0.50
1:A:486:GLN:O	1:A:487:GLU:HB2	2.13	0.49
1:A:458:PRO:HD2	1:A:495[A]:TYR:CE1	2.51	0.46
1:B:420:GLU:OE1	4:B:601:GLC:O3	2.14	0.45
1:B:324:SER:HA	1:B:342:LEU:HD21	2.00	0.42
1:A:294:MET:O	1:A:294:MET:HG3	2.18	0.42
1:B:324:SER:CA	1:B:342:LEU:HD21	2.49	0.42
1:A:324:SER:HA	1:A:342:LEU:HD21	2.01	0.42
1:A:324:SER:CA	1:A:342:LEU:HD21	2.51	0.41
1:B:555:SER:O	1:B:558:ASN:HB2	2.21	0.41
1:B:211:ASP:OD1	1:B:211:ASP:C	2.60	0.41
1:B:246:ARG:HG2	1:B:306:TRP:CZ2	2.56	0.41
1:A:474:ASN:O	3:A:609:EDO:H22	2.21	0.40
1:B:124:LEU:HB3	1:B:125:PRO:HD2	2.03	0.40
1:B:184:SER:N	1:B:185:PRO:CD	2.84	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	511/564 (91%)	501 (98%)	9 (2%)	1 (0%)	47	58
1	B	508/564 (90%)	500 (98%)	8 (2%)	0	100	100
All	All	1019/1128 (90%)	1001 (98%)	17 (2%)	1 (0%)	51	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	333	GLU

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	411/480 (86%)	401 (98%)	10 (2%)	49	66
1	B	405/480 (84%)	395 (98%)	10 (2%)	47	65
All	All	816/960 (85%)	796 (98%)	20 (2%)	47	65

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	SER
1	A	172	ASP
1	A	241	GLU
1	A	260	GLU
1	A	336	GLN

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Mol	Chain	Res	Type
1	A	376	ARG
1	A	482	ASP
1	A	487	GLU
1	A	503	ASP
1	A	560	GLU
1	B	67	SER
1	B	200	GLU
1	B	241	GLU
1	B	260	GLU
1	B	336	GLN
1	B	376	ARG
1	B	482	ASP
1	B	485	ASP
1	B	503	ASP
1	B	553	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	BGC	C	1	1,2	11,11,12	0.29	0	15,15,17	1.03	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GLC	C	2	2	11,11,12	0.45	0	15,15,17	1.23	2 (13%)
2	BGC	D	1	1,2	11,11,12	0.39	0	15,15,17	1.87	2 (13%)
2	GLC	D	2	2	11,11,12	0.35	0	15,15,17	0.85	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	C	1	1,2	-	0/2/19/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	BGC	D	1	1,2	-	2/2/19/22	0/1/1/1
2	GLC	D	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	BGC	C1-O5-C5	6.36	120.80	112.19
2	C	2	GLC	C1-O5-C5	3.06	116.34	112.19
2	C	1	BGC	C1-O5-C5	2.98	116.22	112.19
2	C	2	GLC	C1-C2-C3	2.57	112.82	109.67
2	D	2	GLC	C1-O5-C5	2.53	115.62	112.19
2	D	1	BGC	O4-C4-C5	-2.27	103.67	109.30

There are no chirality outliers.

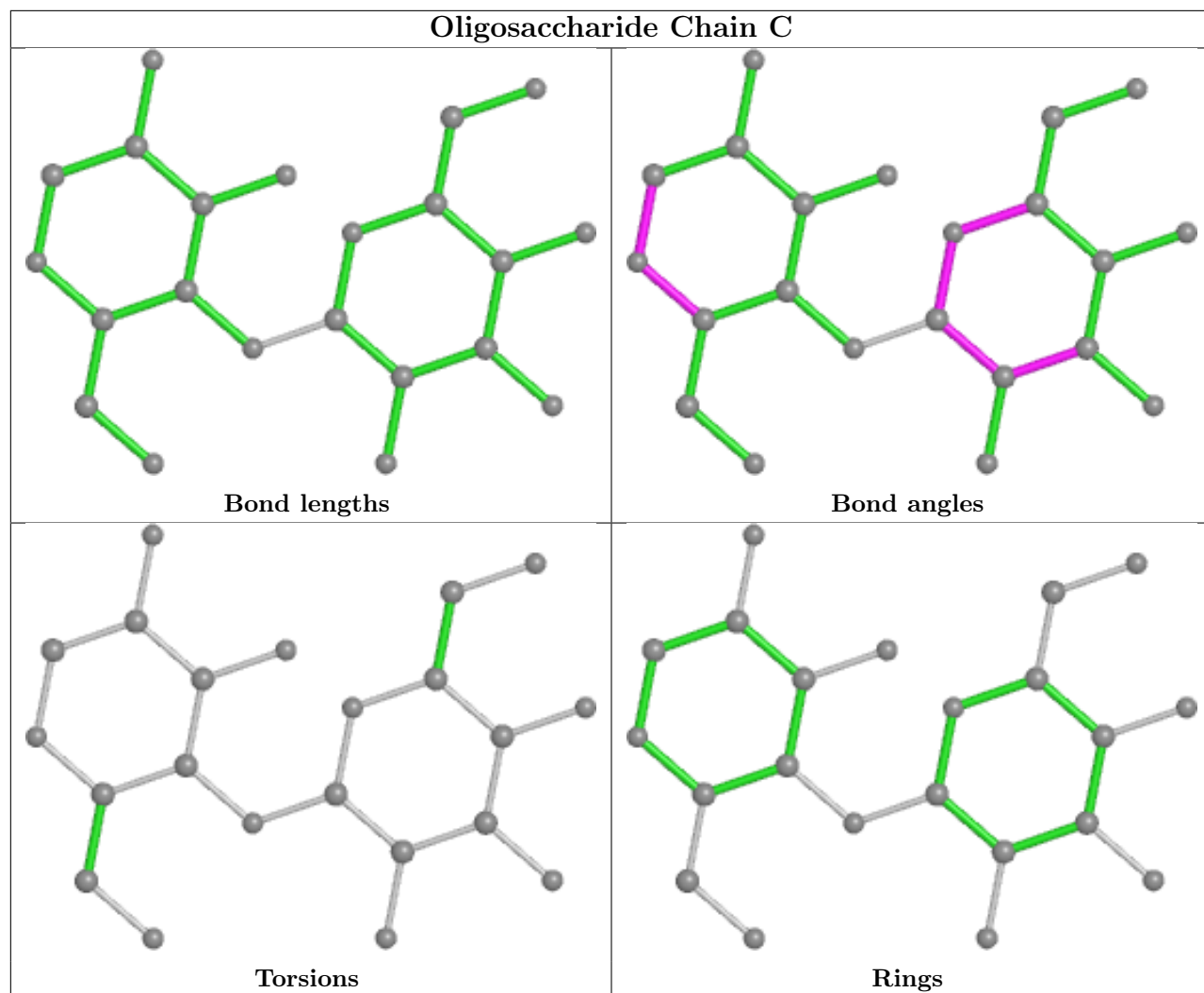
All (2) torsion outliers are listed below:

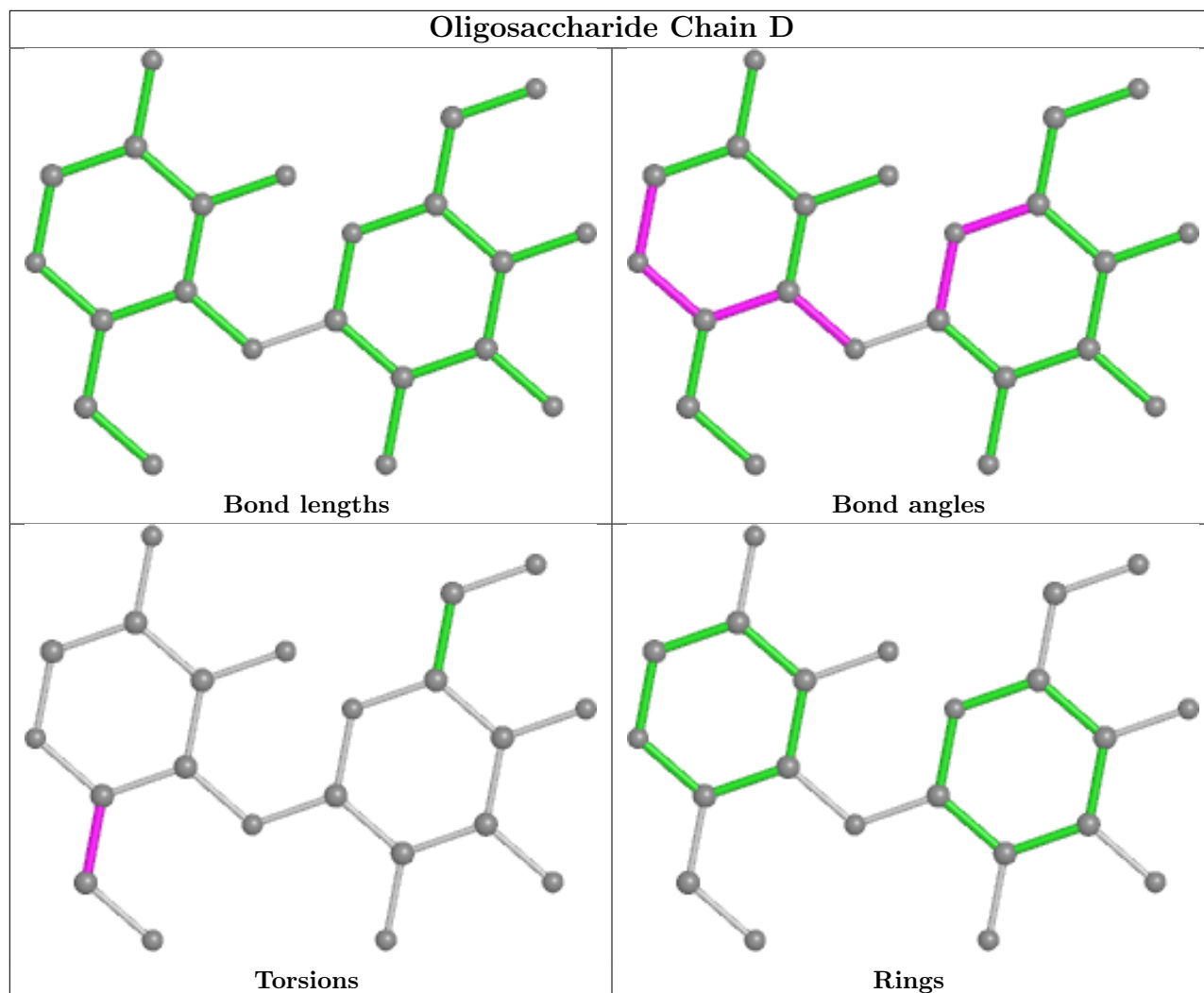
Mol	Chain	Res	Type	Atoms
2	D	1	BGC	O5-C5-C6-O6
2	D	1	BGC	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

17 ligands 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	$\# Z > 2$
3	EDO	A	604	-	3,3,3	0.88	0	2,2,2	0.47	0
3	EDO	B	609	-	3,3,3	0.96	0	2,2,2	0.28	0
3	EDO	B	607	-	3,3,3	0.77	0	2,2,2	0.28	0
3	EDO	A	606	-	3,3,3	0.51	0	2,2,2	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	A	609	-	3,3,3	0.76	0	2,2,2	0.34	0
3	EDO	A	605	-	3,3,3	0.45	0	2,2,2	0.17	0
3	EDO	A	603	-	3,3,3	0.58	0	2,2,2	0.47	0
3	EDO	B	604	-	3,3,3	0.51	0	2,2,2	0.41	0
3	EDO	B	611	-	3,3,3	0.33	0	2,2,2	0.41	0
4	GLC	B	601	-	12,12,12	0.58	0	17,17,17	0.89	0
3	EDO	B	608	-	3,3,3	0.74	0	2,2,2	0.22	0
3	EDO	A	608	-	3,3,3	0.73	0	2,2,2	0.13	0
3	EDO	A	607	-	3,3,3	0.67	0	2,2,2	0.24	0
3	EDO	B	610	-	3,3,3	0.56	0	2,2,2	0.45	0
3	EDO	A	610	-	3,3,3	0.75	0	2,2,2	0.22	0
3	EDO	B	606	-	3,3,3	1.04	0	2,2,2	0.04	0
3	EDO	B	605	-	3,3,3	0.66	0	2,2,2	0.33	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	EDO	A	604	-	-	1/1/1/1	-
3	EDO	B	609	-	-	0/1/1/1	-
3	EDO	B	607	-	-	1/1/1/1	-
3	EDO	A	606	-	-	1/1/1/1	-
3	EDO	A	609	-	-	0/1/1/1	-
3	EDO	A	605	-	-	1/1/1/1	-
3	EDO	A	603	-	-	0/1/1/1	-
3	EDO	B	604	-	-	1/1/1/1	-
3	EDO	B	611	-	-	0/1/1/1	-
4	GLC	B	601	-	-	2/2/22/22	0/1/1/1
3	EDO	B	608	-	-	1/1/1/1	-
3	EDO	A	608	-	-	1/1/1/1	-
3	EDO	A	607	-	-	1/1/1/1	-
3	EDO	B	610	-	-	1/1/1/1	-
3	EDO	A	610	-	-	0/1/1/1	-
3	EDO	B	606	-	-	1/1/1/1	-
3	EDO	B	605	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	601	GLC	O5-C5-C6-O6
4	B	601	GLC	C4-C5-C6-O6
3	B	610	EDO	O1-C1-C2-O2
3	A	608	EDO	O1-C1-C2-O2
3	B	608	EDO	O1-C1-C2-O2
3	A	605	EDO	O1-C1-C2-O2
3	B	607	EDO	O1-C1-C2-O2
3	B	604	EDO	O1-C1-C2-O2
3	B	606	EDO	O1-C1-C2-O2
3	A	607	EDO	O1-C1-C2-O2
3	A	604	EDO	O1-C1-C2-O2
3	A	606	EDO	O1-C1-C2-O2
3	B	605	EDO	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	609	EDO	1	0
3	A	609	EDO	1	0
3	B	611	EDO	2	0
4	B	601	GLC	1	0
3	A	610	EDO	2	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	514/564 (91%)	-0.27	14 (2%) 54 62	19, 38, 72, 111	0
1	B	511/564 (90%)	-0.03	27 (5%) 26 33	20, 40, 72, 105	0
All	All	1025/1128 (90%)	-0.15	41 (4%) 38 45	19, 39, 72, 111	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	393	VAL	6.2
1	A	67	SER	4.5
1	A	75	VAL	4.5
1	A	68	GLU	4.3
1	A	65	PHE	4.1
1	B	75	VAL	3.9
1	A	71	GLN	3.8
1	B	386	SER	3.8
1	B	337	LEU	3.4
1	B	67	SER	3.3
1	A	332	SER	3.2
1	A	334	THR	3.2
1	B	66	PRO	3.2
1	A	72	TRP	3.1
1	B	321	LEU	2.9
1	B	234	TYR	2.9
1	B	385	PRO	2.8
1	B	73	LEU	2.8
1	A	390	VAL	2.7
1	A	64	ASP	2.6
1	B	132	GLY	2.5
1	A	66	PRO	2.5
1	A	234	TYR	2.5
1	B	65	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	424	VAL	2.4
1	B	76	PRO	2.3
1	B	392	MET	2.3
1	B	390	VAL	2.2
1	B	72	TRP	2.2
1	B	388	ALA	2.2
1	B	294	MET	2.2
1	A	69	TYR	2.2
1	A	199	GLY	2.2
1	B	126	LEU	2.1
1	B	125	PRO	2.1
1	B	322	LEU	2.1
1	B	395	ARG	2.1
1	B	68	GLU	2.1
1	B	163	ASP	2.0
1	B	338[A]	TRP	2.0
1	B	206	LEU	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](#)

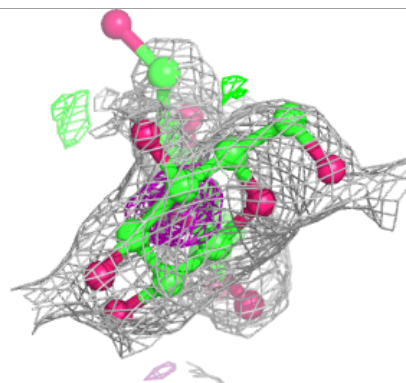
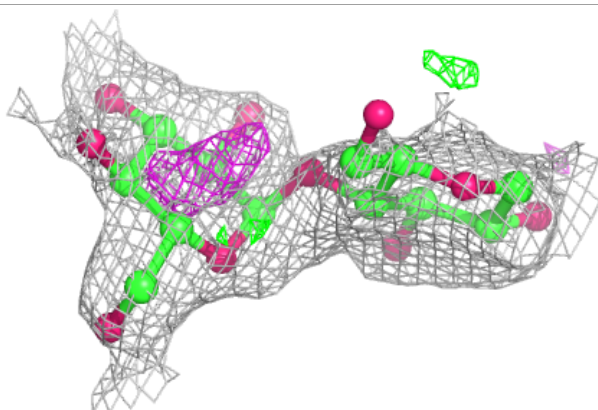
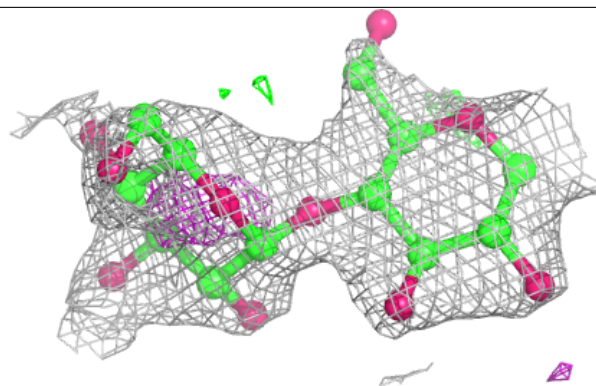
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	C	2	11/12	0.79	0.24	61,69,74,76	0
2	GLC	D	2	11/12	0.84	0.20	59,66,72,73	0
2	BGC	D	1	11/12	0.88	0.15	52,63,67,69	0
2	BGC	C	1	11/12	0.89	0.15	59,70,85,92	0

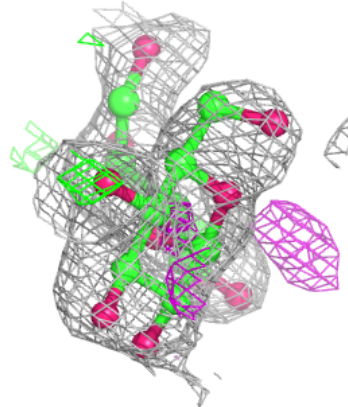
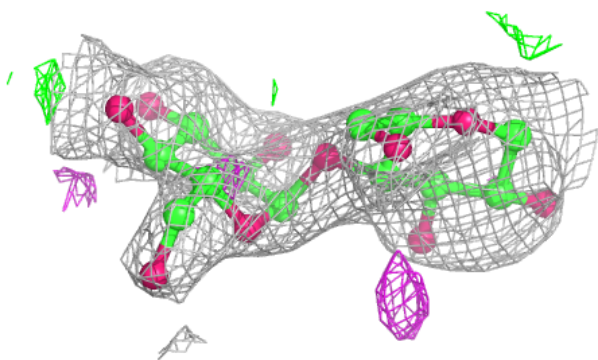
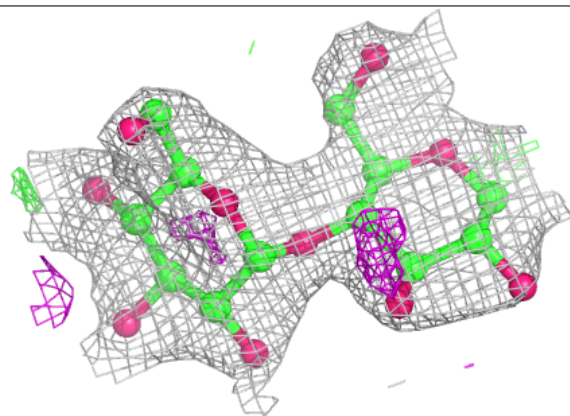
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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

**Electron density around Chain D:**

$2mF_o-DF_c$ (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(Å ²)	Q<0.9
3	EDO	B	608	4/4	0.74	0.24	66,66,68,68	0
4	GLC	B	601	12/12	0.82	0.18	64,71,80,83	0
3	EDO	B	606	4/4	0.86	0.27	47,49,50,50	0
3	EDO	A	604	4/4	0.87	0.16	48,49,53,53	0
3	EDO	B	611	4/4	0.89	0.20	48,48,51,53	0
3	EDO	A	608	4/4	0.89	0.15	47,52,54,55	0
3	EDO	B	605	4/4	0.90	0.22	54,62,63,68	0
3	EDO	A	605	4/4	0.93	0.19	52,53,55,57	0
3	EDO	B	609	4/4	0.93	0.16	34,40,40,43	0
3	EDO	B	607	4/4	0.94	0.13	44,46,46,47	0
3	EDO	B	604	4/4	0.95	0.10	42,43,44,45	0
3	EDO	A	610	4/4	0.95	0.15	43,44,45,45	0
3	EDO	B	610	4/4	0.96	0.16	52,57,60,65	0
3	EDO	A	609	4/4	0.96	0.10	38,43,43,45	0
3	EDO	A	606	4/4	0.96	0.26	62,62,64,65	0
3	EDO	A	607	4/4	0.97	0.13	44,44,45,48	0
3	EDO	A	603	4/4	0.98	0.09	39,40,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.