



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:11 PM UTC

PDB ID : 4ZWC / pdb_00004zwc
Title : Crystal structure of maltose-bound human GLUT3 in the outward-open conformation at 2.6 angstrom
Authors : Deng, D.; Sun, P.C.; Yan, C.Y.; Yan, N.
Deposited on : 2015-05-19
Resolution : 2.60 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

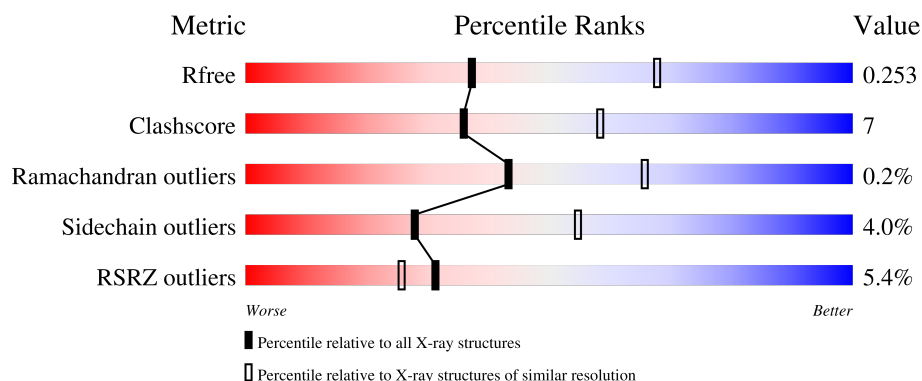
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	518	5% 74% 15% 10%
1	B	518	5% 73% 16% 10%
2	C	2	50% 50%
2	D	2	100%

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
3	OLC	A	501	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7359 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 Solute carrier family 2, facilitated glucose transporter member 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	468	Total	C	N	O	S	0	0	0
			3593	2361	581	631	20			
1	B	468	Total	C	N	O	S	0	0	0
			3593	2361	581	631	20			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP P11169
A	-20	HIS	-	expression tag	UNP P11169
A	-19	HIS	-	expression tag	UNP P11169
A	-18	HIS	-	expression tag	UNP P11169
A	-17	HIS	-	expression tag	UNP P11169
A	-16	HIS	-	expression tag	UNP P11169
A	-15	HIS	-	expression tag	UNP P11169
A	-14	HIS	-	expression tag	UNP P11169
A	-13	HIS	-	expression tag	UNP P11169
A	-12	HIS	-	expression tag	UNP P11169
A	-11	HIS	-	expression tag	UNP P11169
A	-10	SER	-	expression tag	UNP P11169
A	-9	GLY	-	expression tag	UNP P11169
A	-8	ASP	-	expression tag	UNP P11169
A	-7	GLU	-	expression tag	UNP P11169
A	-6	VAL	-	expression tag	UNP P11169
A	-5	ASP	-	expression tag	UNP P11169
A	-4	ALA	-	expression tag	UNP P11169
A	-3	GLY	-	expression tag	UNP P11169
A	-2	SER	-	expression tag	UNP P11169
A	-1	GLY	-	expression tag	UNP P11169
A	0	HIS	-	expression tag	UNP P11169
A	43	THR	ASN	engineered mutation	UNP P11169
B	-21	MET	-	expression tag	UNP P11169

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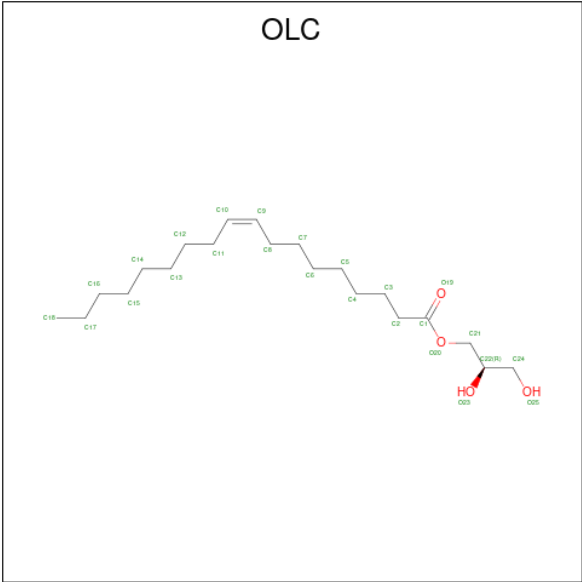
Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	HIS	-	expression tag	UNP P11169
B	-19	HIS	-	expression tag	UNP P11169
B	-18	HIS	-	expression tag	UNP P11169
B	-17	HIS	-	expression tag	UNP P11169
B	-16	HIS	-	expression tag	UNP P11169
B	-15	HIS	-	expression tag	UNP P11169
B	-14	HIS	-	expression tag	UNP P11169
B	-13	HIS	-	expression tag	UNP P11169
B	-12	HIS	-	expression tag	UNP P11169
B	-11	HIS	-	expression tag	UNP P11169
B	-10	SER	-	expression tag	UNP P11169
B	-9	GLY	-	expression tag	UNP P11169
B	-8	ASP	-	expression tag	UNP P11169
B	-7	GLU	-	expression tag	UNP P11169
B	-6	VAL	-	expression tag	UNP P11169
B	-5	ASP	-	expression tag	UNP P11169
B	-4	ALA	-	expression tag	UNP P11169
B	-3	GLY	-	expression tag	UNP P11169
B	-2	SER	-	expression tag	UNP P11169
B	-1	GLY	-	expression tag	UNP P11169
B	0	HIS	-	expression tag	UNP P11169
B	43	THR	ASN	engineered mutation	UNP P11169

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is (2R)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (CCD ID: OLC) (formula: C₂₁H₄₀O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			25	21	4		
3	A	1	Total	C	O	0	0
			25	21	4		
3	B	1	Total	C	O	0	0
			25	21	4		

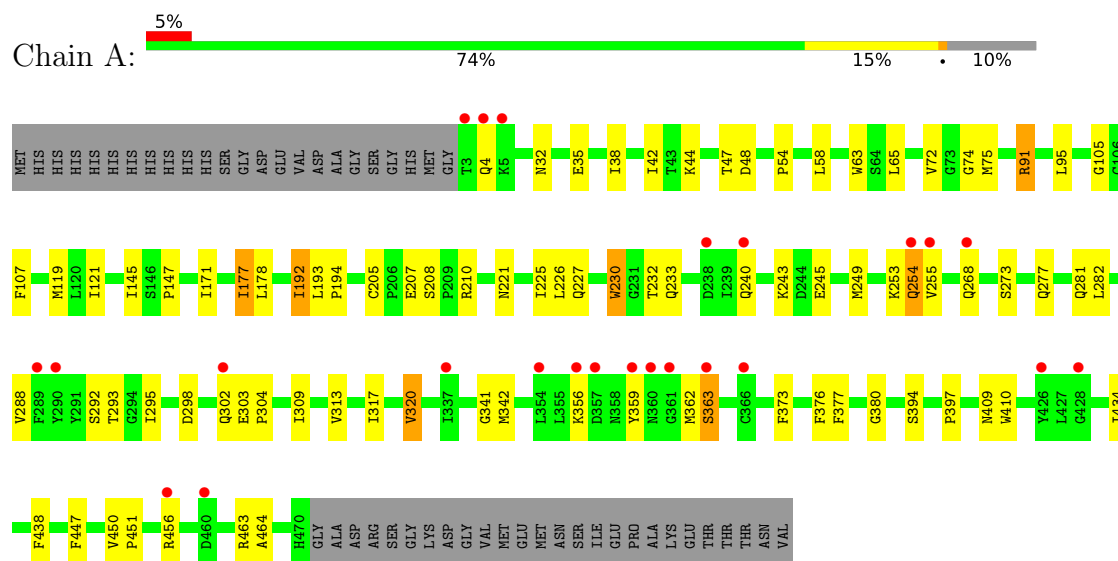
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	26	Total	O	0	0
			26	26		
4	B	26	Total	O	0	0
			26	26		

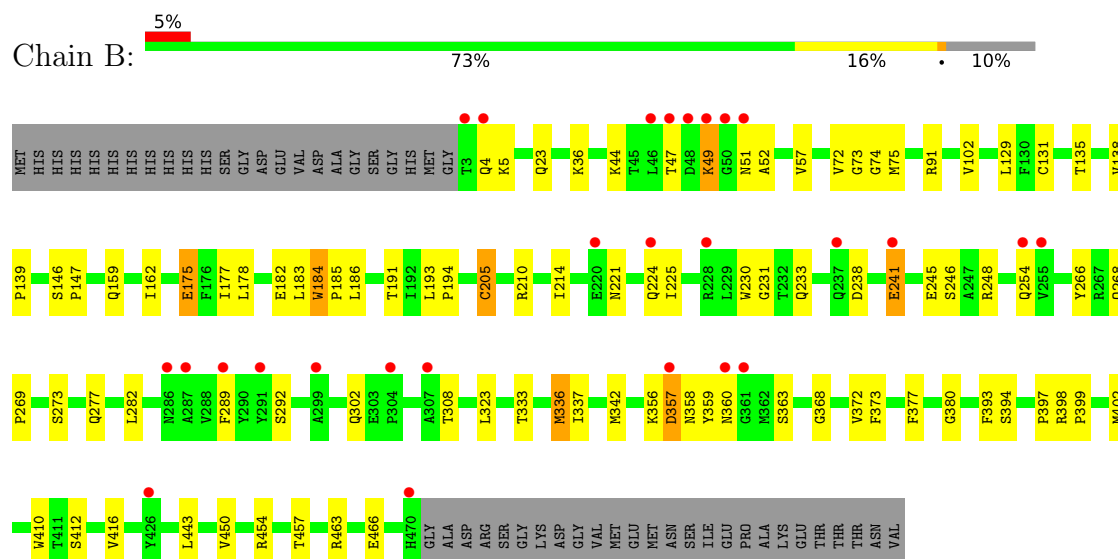
3 Residue-property plots [i](#)

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: Solute carrier family 2, facilitated glucose transporter member 3



- Molecule 1: Solute carrier family 2, facilitated glucose transporter member 3



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

Chain C:  50% 50%

GLC1
GLC2

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

Chain D:  100%

GLC1
GLC2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.11Å 121.88Å 96.08Å 90.00° 108.14° 90.00°	Depositor
Resolution (Å)	31.70 – 2.60 31.70 – 2.60	Depositor EDS
% Data completeness (in resolution range)	94.5 (31.70-2.60) 89.7 (31.70-2.60)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.62 (at 2.61Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE: 1.9_1692)	Depositor
R, R_{free}	0.215 , 0.251 0.224 , 0.253	Depositor DCC
R_{free} test set	2479 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	42.5	Xtrriage
Anisotropy	0.380	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.3	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	0.92	EDS
Total number of atoms	7359	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.66 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.1475e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: OLC, GLC

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.63	0/3674	0.91	2/4986 (0.0%)
1	B	0.67	1/3674 (0.0%)	0.95	11/4986 (0.2%)
All	All	0.65	1/7348 (0.0%)	0.93	13/9972 (0.1%)

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
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	138	VAL	CA-C	5.05	1.57	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	360	ASN	N-CA-C	6.60	118.56	111.36
1	A	450	VAL	N-CA-C	6.38	113.37	107.56
1	B	146	SER	CA-C-N	6.37	126.75	119.93
1	B	146	SER	C-N-CA	6.37	126.75	119.93
1	B	358	ASN	N-CA-C	-6.14	105.05	112.54
1	B	450	VAL	N-CA-C	5.79	113.28	107.55
1	B	184	TRP	CA-C-N	-5.62	113.10	119.28
1	B	184	TRP	C-N-CA	-5.62	113.10	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	32	ASN	N-CA-C	5.62	117.21	111.14
1	B	205	CYS	CA-C-N	-5.53	114.27	119.85
1	B	205	CYS	C-N-CA	-5.53	114.27	119.85
1	B	138	VAL	CA-C-N	-5.48	114.03	119.56
1	B	138	VAL	C-N-CA	-5.48	114.03	119.56

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	254	GLN	Peptide
1	B	357	ASP	Peptide

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	3593	0	3691	53	0
1	B	3593	0	3691	50	0
2	C	23	0	21	1	0
2	D	23	0	21	0	0
3	A	50	0	80	13	0
3	B	25	0	40	1	0
4	A	26	0	0	2	0
4	B	26	0	0	1	0
All	All	7359	0	7544	108	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 7.

All (108) 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:B:241:GLU:OE2	1:B:457:THR:HG21	1.43	1.17
1:A:464:ALA:HB3	3:A:502:OLC:H21	1.13	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:ILE:HG23	1:A:208:SER:HB3	1.46	0.97
1:A:464:ALA:CB	3:A:502:OLC:H21	1.96	0.96
1:A:145:ILE:CG2	1:A:208:SER:HB3	1.95	0.96
1:B:72:VAL:HA	1:B:75:MET:HE3	1.58	0.84
1:A:447:PHE:HZ	3:A:502:OLC:H17	1.42	0.82
1:A:320:VAL:CG2	3:A:501:OLC:H8A	2.11	0.81
1:A:447:PHE:CZ	3:A:502:OLC:H17	2.16	0.78
1:A:207:GLU:OE2	1:A:208:SER:N	2.27	0.67
1:B:5:LYS:NZ	1:B:231:GLY:O	2.31	0.64
3:A:502:OLC:H4A	3:A:502:OLC:O19	1.97	0.64
1:A:72:VAL:HA	1:A:75:MET:HE3	1.79	0.63
1:B:182:GLU:N	1:B:182:GLU:OE1	2.31	0.63
1:B:342:MET:HE1	1:B:377:PHE:HD1	1.64	0.63
1:A:145:ILE:HG22	1:A:208:SER:HB3	1.80	0.62
1:A:249:MET:HG2	1:A:254:GLN:HE21	1.65	0.62
1:B:210:ARG:NH1	1:B:245:GLU:OE2	2.34	0.61
1:A:221:ASN:O	1:A:225:ILE:HG12	2.01	0.60
3:A:502:OLC:H10	3:A:502:OLC:H14	1.83	0.60
1:A:210:ARG:NH1	1:A:245:GLU:OE2	2.34	0.60
1:A:356:LYS:NZ	1:A:363:SER:OG	2.36	0.59
1:A:447:PHE:CZ	3:A:502:OLC:H15A	2.37	0.59
1:A:451:PRO:HB3	3:A:502:OLC:H5A	1.85	0.58
1:A:245:GLU:OE1	4:A:601:HOH:O	2.17	0.58
1:B:241:GLU:OE2	1:B:457:THR:CG2	2.35	0.57
1:A:447:PHE:CE2	3:A:502:OLC:H15A	2.40	0.57
1:A:44:LYS:NZ	1:A:48:ASP:OD2	2.38	0.57
1:B:336:MET:HG2	1:B:443:LEU:HD12	1.86	0.57
1:B:51:ASN:OD1	1:B:52:ALA:N	2.30	0.57
1:A:145:ILE:CG2	1:A:208:SER:CB	2.78	0.56
1:B:248:ARG:NH1	1:B:466:GLU:OE2	2.37	0.56
1:B:51:ASN:CG	1:B:52:ALA:H	2.11	0.56
1:A:320:VAL:HG22	3:A:501:OLC:H8A	1.85	0.56
1:A:145:ILE:HG22	1:A:208:SER:CB	2.36	0.56
1:A:91:ARG:HD2	1:A:205:CYS:O	2.07	0.55
1:A:65:LEU:HD23	1:A:121:ILE:HD13	1.89	0.55
1:B:44:LYS:HA	1:B:47:THR:HG22	1.91	0.53
1:B:214:ILE:HG23	1:B:246:SER:HA	1.89	0.53
1:A:281:GLN:NE2	1:A:409:ASN:OD1	2.37	0.53
1:A:171:ILE:HD13	1:A:309:ILE:HD13	1.91	0.52
1:B:356:LYS:HE2	1:B:363:SER:HB3	1.91	0.52
1:B:175:GLU:H	1:B:175:GLU:CD	2.18	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:MET:HG2	1:B:443:LEU:CD1	2.40	0.51
1:B:193:LEU:HB2	1:B:194:PRO:HD3	1.93	0.51
1:B:342:MET:HE3	1:B:377:PHE:HB2	1.93	0.51
1:B:4:GLN:OE1	1:B:147:PRO:HB3	2.11	0.51
1:A:288:VAL:O	1:A:292:SER:HB3	2.11	0.51
1:B:159:GLN:HA	1:B:162:ILE:HD12	1.93	0.50
1:B:221:ASN:O	1:B:225:ILE:HG12	2.11	0.50
1:A:74:GLY:HA3	1:A:410:TRP:CE3	2.46	0.50
1:A:253:LYS:O	4:A:602:HOH:O	2.19	0.50
1:A:342:MET:HE1	1:A:377:PHE:HD1	1.77	0.50
1:A:313:VAL:O	1:A:317:ILE:HG13	2.12	0.49
1:B:183:LEU:HD22	1:B:186:LEU:HD12	1.95	0.49
1:A:342:MET:HE3	1:A:377:PHE:HB2	1.95	0.49
1:B:23:GLN:OE1	1:B:191:THR:OG1	2.20	0.49
1:B:182:GLU:H	1:B:182:GLU:CD	2.21	0.49
1:B:266:TYR:C	1:B:269:PRO:HD2	2.38	0.48
3:A:501:OLC:H11	3:A:501:OLC:H8	1.69	0.47
1:B:398:ARG:HB3	1:B:399:PRO:HD3	1.97	0.47
1:A:107:PHE:HA	1:A:119:MET:HE3	1.97	0.47
1:B:273:SER:O	1:B:277:GLN:HG2	2.15	0.46
1:B:289:PHE:O	1:B:292:SER:OG	2.32	0.46
1:A:298:ASP:OD2	1:A:356:LYS:HE3	2.16	0.46
1:A:359:TYR:O	1:A:362:MET:HG2	2.16	0.46
1:B:131:CYS:O	1:B:135:THR:HG23	2.17	0.45
1:A:227:GLN:HG2	1:A:233:GLN:HA	1.99	0.45
1:A:249:MET:HG2	1:A:254:GLN:NE2	2.32	0.45
1:B:394:SER:O	1:B:397:PRO:HD2	2.18	0.44
3:B:501:OLC:H2A	3:B:501:OLC:H5	1.70	0.44
1:A:240:GLN:HA	1:A:243:LYS:HB2	1.99	0.44
1:A:394:SER:O	1:A:397:PRO:HD2	2.17	0.44
1:B:269:PRO:HB2	1:B:393:PHE:CZ	2.53	0.44
1:A:342:MET:HE2	1:A:373:PHE:CZ	2.53	0.44
1:B:91:ARG:HD3	1:B:205:CYS:O	2.16	0.44
1:A:273:SER:O	1:A:277:GLN:HG2	2.18	0.44
1:B:210:ARG:HD3	4:B:605:HOH:O	2.16	0.44
1:A:38:ILE:O	1:A:42:ILE:HG13	2.18	0.44
1:B:230:TRP:HZ3	1:B:238:ASP:CG	2.26	0.44
1:B:356:LYS:HD2	1:B:356:LYS:O	2.17	0.44
1:A:341:GLY:HA3	1:A:376:PHE:CD1	2.52	0.43
1:B:342:MET:HE2	1:B:373:PHE:CZ	2.53	0.43
1:B:412:SER:O	1:B:416:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:2:GLC:HO4	2:C:2:GLC:HO6	1.59	0.43
1:A:4:GLN:OE1	1:A:147:PRO:HB3	2.19	0.43
1:A:451:PRO:HG3	3:A:502:OLC:H3	2.00	0.43
1:B:73:GLY:O	1:B:129:LEU:HD23	2.18	0.43
1:B:333:THR:O	1:B:337:ILE:HG13	2.19	0.43
1:B:268:GLN:HB3	1:B:269:PRO:HD3	2.02	0.42
1:A:105:GLY:HA3	1:A:192:ILE:HG22	2.01	0.42
1:B:356:LYS:HD2	1:B:356:LYS:C	2.44	0.42
1:B:139:PRO:HG2	1:B:402:MET:HE1	2.02	0.42
1:A:303:GLU:HA	1:A:304:PRO:HD3	1.83	0.42
1:B:74:GLY:HA3	1:B:410:TRP:CE3	2.55	0.41
1:B:184:TRP:CG	1:B:185:PRO:HD3	2.55	0.41
1:B:357:ASP:HA	1:B:359:TYR:O	2.20	0.41
1:B:282:LEU:HD23	1:B:282:LEU:HA	1.85	0.41
1:B:368:GLY:O	1:B:372:VAL:HG23	2.21	0.41
1:B:224:GLN:HG3	1:B:225:ILE:N	2.35	0.41
1:A:230:TRP:HB3	1:A:232:THR:HG22	2.02	0.41
1:A:434:ILE:HG22	1:A:438:PHE:CE2	2.55	0.41
1:A:177:ILE:HB	1:A:178:LEU:H	1.74	0.41
1:B:177:ILE:O	1:B:178:LEU:HD23	2.21	0.41
1:A:193:LEU:HB2	1:A:194:PRO:HD3	2.03	0.40
1:B:49:LYS:HA	1:B:49:LYS:HD2	1.71	0.40
1:A:35:GLU:HG3	1:A:63:TRP:CE2	2.56	0.40
1:A:54:PRO:HB3	1:A:58:LEU:HD23	2.03	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	466/518 (90%)	455 (98%)	10 (2%)	1 (0%)	43 66

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	466/518 (90%)	452 (97%)	13 (3%)	1 (0%)	43 66
All	All	932/1036 (90%)	907 (97%)	23 (2%)	2 (0%)	43 66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	380	GLY
1	A	380	GLY

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	387/428 (90%)	370 (96%)	17 (4%)	25 50
1	B	387/428 (90%)	373 (96%)	14 (4%)	31 58
All	All	774/856 (90%)	743 (96%)	31 (4%)	28 55

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	THR
1	A	91	ARG
1	A	95	LEU
1	A	177	ILE
1	A	192	ILE
1	A	226	LEU
1	A	230	TRP
1	A	255	VAL
1	A	268	GLN
1	A	282	LEU
1	A	293	THR
1	A	295	ILE
1	A	302	GLN
1	A	320	VAL

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Mol	Chain	Res	Type
1	A	363	SER
1	A	456	ARG
1	A	463	ARG
1	B	36	LYS
1	B	49	LYS
1	B	57	VAL
1	B	102	VAL
1	B	175	GLU
1	B	233	GLN
1	B	241	GLU
1	B	254	GLN
1	B	302	GLN
1	B	308	THR
1	B	323	LEU
1	B	336	MET
1	B	454	ARG
1	B	463	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	221	ASN
1	A	224	GLN
1	A	286	ASN
1	A	413	ASN
1	B	254	GLN
1	B	286	ASN

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 ⓘ

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	C	1	2	12,12,12	0.73	0	17,17,17	1.36	2 (11%)
2	GLC	C	2	2	11,11,12	0.71	0	15,15,17	1.64	3 (20%)
2	GLC	D	1	2	12,12,12	0.71	0	17,17,17	1.31	3 (17%)
2	GLC	D	2	2	11,11,12	0.59	0	15,15,17	1.93	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	C	1	2	-	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	1/2/19/22	0/1/1/1
2	GLC	D	1	2	-	2/2/22/22	0/1/1/1
2	GLC	D	2	2	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	GLC	C1-O5-C5	4.51	118.23	112.19
2	C	1	GLC	C1-O5-C5	3.60	120.61	113.65
2	D	2	GLC	C2-C3-C4	-3.12	105.37	110.86
2	D	2	GLC	C1-C2-C3	-2.88	105.45	109.64
2	C	2	GLC	C1-O5-C5	2.83	115.98	112.19
2	D	1	GLC	C1-O5-C5	2.45	118.40	113.65
2	C	2	GLC	O3-C3-C2	2.33	114.80	110.05
2	D	1	GLC	O5-C1-C2	2.27	114.29	110.30
2	D	1	GLC	C3-C4-C5	-2.19	106.27	110.23
2	C	2	GLC	C2-C3-C4	-2.17	107.05	110.86
2	C	1	GLC	O3-C3-C2	-2.06	105.53	110.38

There are no chirality outliers.

All (6) torsion outliers are listed below:

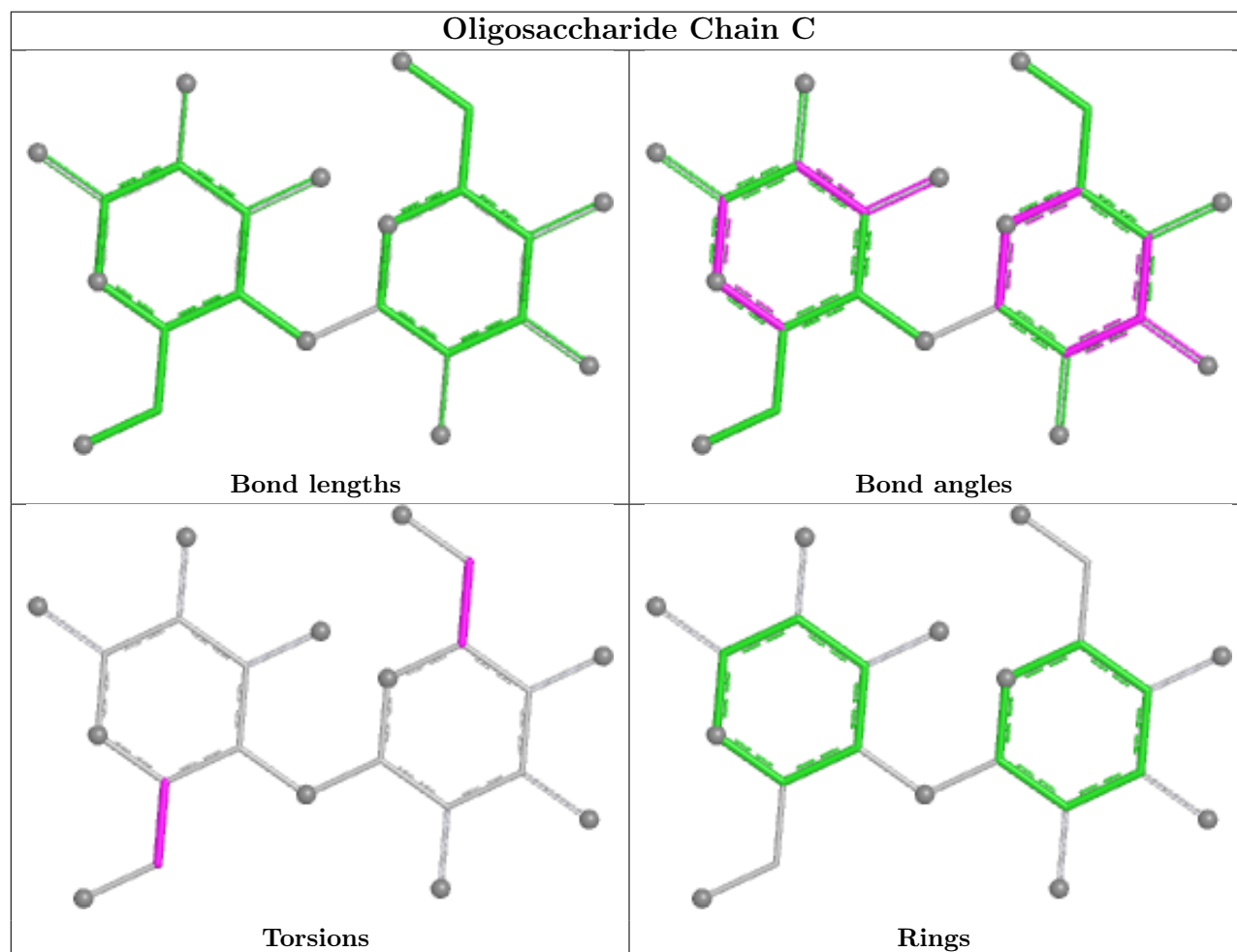
Mol	Chain	Res	Type	Atoms
2	D	1	GLC	O5-C5-C6-O6
2	D	2	GLC	O5-C5-C6-O6
2	C	2	GLC	O5-C5-C6-O6
2	D	1	GLC	C4-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6

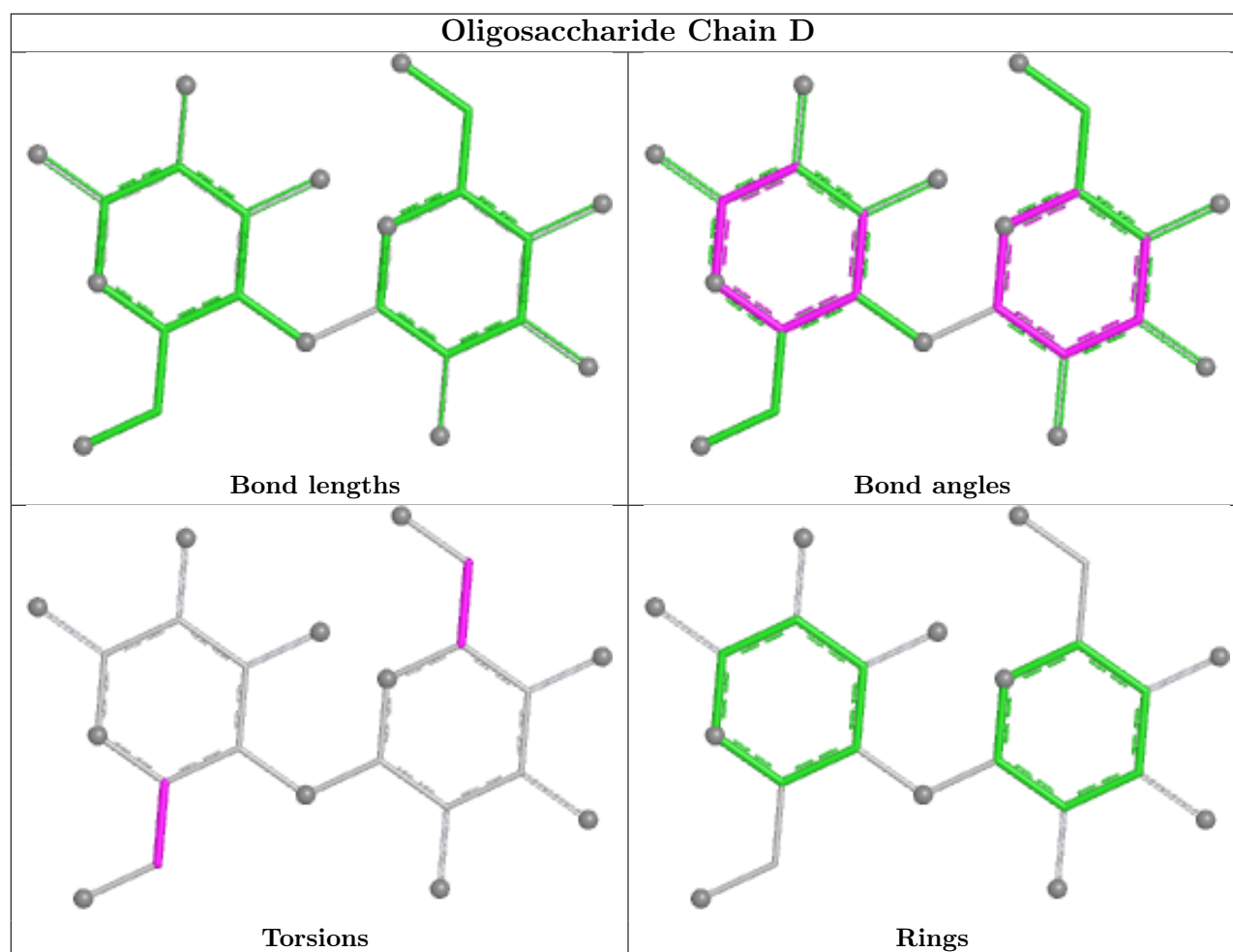
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	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.





5.6 Ligand geometry [i](#)

3 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	OLC	A	501	-	24,24,24	0.88	1 (4%)	25,25,25	0.94	1 (4%)
3	OLC	B	501	-	24,24,24	0.88	1 (4%)	25,25,25	1.03	2 (8%)
3	OLC	A	502	-	24,24,24	0.86	1 (4%)	25,25,25	1.07	2 (8%)

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	OLC	A	501	-	-	14/24/24/24	-
3	OLC	B	501	-	-	17/24/24/24	-
3	OLC	A	502	-	-	8/24/24/24	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	OLC	O20-C1	4.13	1.45	1.33
3	B	501	OLC	O20-C1	4.08	1.45	1.33
3	A	502	OLC	O20-C1	3.72	1.44	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	OLC	O20-C1-C2	3.41	122.23	111.83
3	B	501	OLC	O20-C1-C2	3.15	121.43	111.83
3	A	501	OLC	O20-C1-C2	2.81	120.39	111.83
3	B	501	OLC	O20-C1-O19	-2.36	117.73	123.63
3	A	502	OLC	O20-C1-O19	-2.04	118.52	123.63

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	501	OLC	C21-C22-C24-O25
3	B	501	OLC	O20-C21-C22-C24
3	B	501	OLC	O20-C21-C22-O23
3	A	501	OLC	C2-C1-O20-C21
3	A	502	OLC	C2-C1-O20-C21
3	A	501	OLC	O19-C1-O20-C21
3	A	501	OLC	O20-C21-C22-O23
3	A	501	OLC	C1-C2-C3-C4
3	A	502	OLC	O19-C1-O20-C21
3	A	501	OLC	C5-C6-C7-C8
3	B	501	OLC	O23-C22-C24-O25
3	A	501	OLC	C12-C13-C14-C15
3	B	501	OLC	C2-C1-O20-C21
3	A	502	OLC	C11-C12-C13-C14

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Mol	Chain	Res	Type	Atoms
3	B	501	OLC	C12-C13-C14-C15
3	B	501	OLC	C4-C5-C6-C7
3	B	501	OLC	O19-C1-O20-C21
3	A	501	OLC	O20-C21-C22-C24
3	B	501	OLC	C14-C15-C16-C17
3	A	501	OLC	C6-C7-C8-C9
3	B	501	OLC	C6-C7-C8-C9
3	B	501	OLC	C11-C12-C13-C14
3	A	502	OLC	C5-C6-C7-C8
3	A	501	OLC	C11-C12-C13-C14
3	B	501	OLC	C10-C11-C12-C13
3	A	501	OLC	C3-C4-C5-C6
3	A	501	OLC	C2-C3-C4-C5
3	A	502	OLC	C2-C3-C4-C5
3	A	502	OLC	C15-C16-C17-C18
3	B	501	OLC	C3-C4-C5-C6
3	A	501	OLC	C10-C11-C12-C13
3	A	501	OLC	C15-C16-C17-C18
3	A	501	OLC	C13-C14-C15-C16
3	B	501	OLC	C5-C6-C7-C8
3	B	501	OLC	C2-C3-C4-C5
3	B	501	OLC	C15-C16-C17-C18
3	B	501	OLC	C7-C8-C9-C10
3	A	502	OLC	C3-C4-C5-C6
3	A	502	OLC	C13-C14-C15-C16

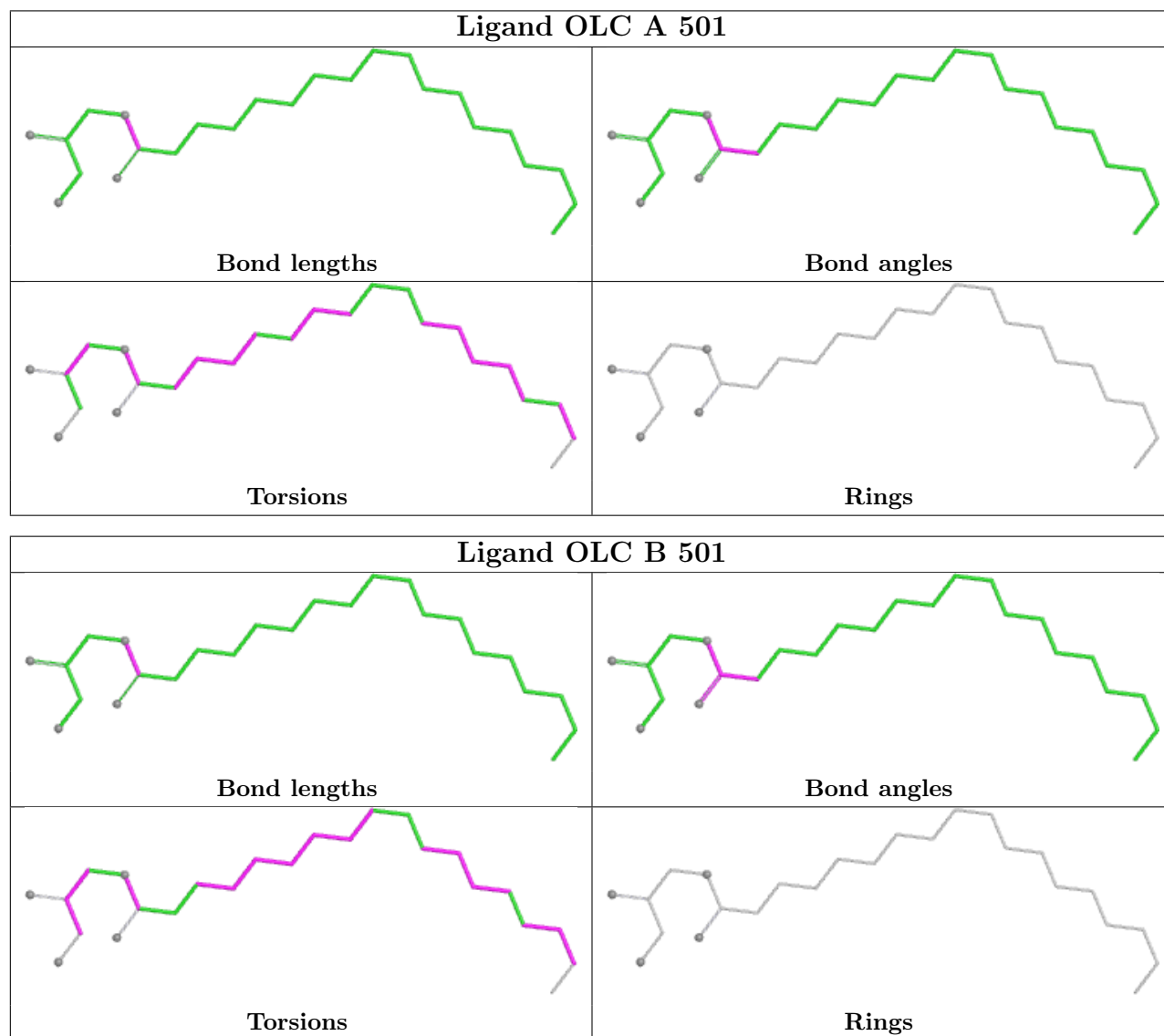
There are no ring outliers.

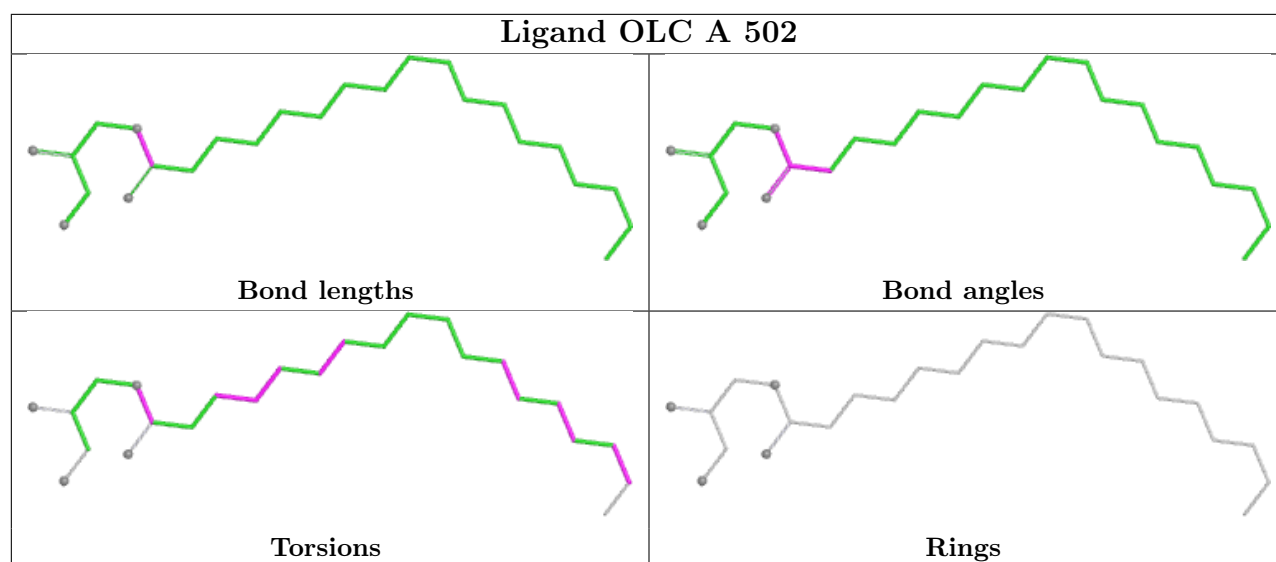
3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	OLC	3	0
3	B	501	OLC	1	0
3	A	502	OLC	10	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	468/518 (90%)	0.39	24 (5%) 33 28	34, 45, 71, 94	0
1	B	468/518 (90%)	0.46	27 (5%) 29 23	35, 47, 77, 94	0
All	All	936/1036 (90%)	0.42	51 (5%) 31 26	34, 46, 74, 94	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	THR	6.7
1	B	3	THR	6.5
1	B	50	GLY	5.6
1	A	302	GLN	4.9
1	A	254	GLN	4.3
1	B	51	ASN	4.2
1	A	4	GLN	4.1
1	B	47	THR	4.1
1	B	254	GLN	4.0
1	A	240	GLN	3.9
1	B	357	ASP	3.9
1	B	286	ASN	3.7
1	B	287	ALA	3.5
1	B	48	ASP	3.5
1	B	304	PRO	3.5
1	B	360	ASN	3.2
1	B	4	GLN	3.1
1	B	49	LYS	3.1
1	B	299	ALA	3.1
1	A	255	VAL	3.1
1	B	241	GLU	3.1
1	B	470	HIS	3.1
1	B	237	GLN	3.0
1	A	456	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	426	TYR	2.9
1	B	361	GLY	2.9
1	B	426	TYR	2.8
1	A	460	ASP	2.8
1	A	5	LYS	2.7
1	A	359	TYR	2.7
1	A	289	PHE	2.7
1	B	224	GLN	2.6
1	A	357	ASP	2.6
1	B	255	VAL	2.5
1	B	307	ALA	2.4
1	A	354	LEU	2.4
1	A	356	LYS	2.4
1	A	361	GLY	2.4
1	B	228	ARG	2.3
1	B	291	TYR	2.3
1	A	363	SER	2.3
1	A	360	ASN	2.2
1	B	289	PHE	2.2
1	A	290	TYR	2.2
1	A	268	GLN	2.2
1	A	428	GLY	2.2
1	A	366	CYS	2.1
1	B	220	GLU	2.1
1	B	46	LEU	2.1
1	A	337	ILE	2.0
1	A	238	ASP	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.

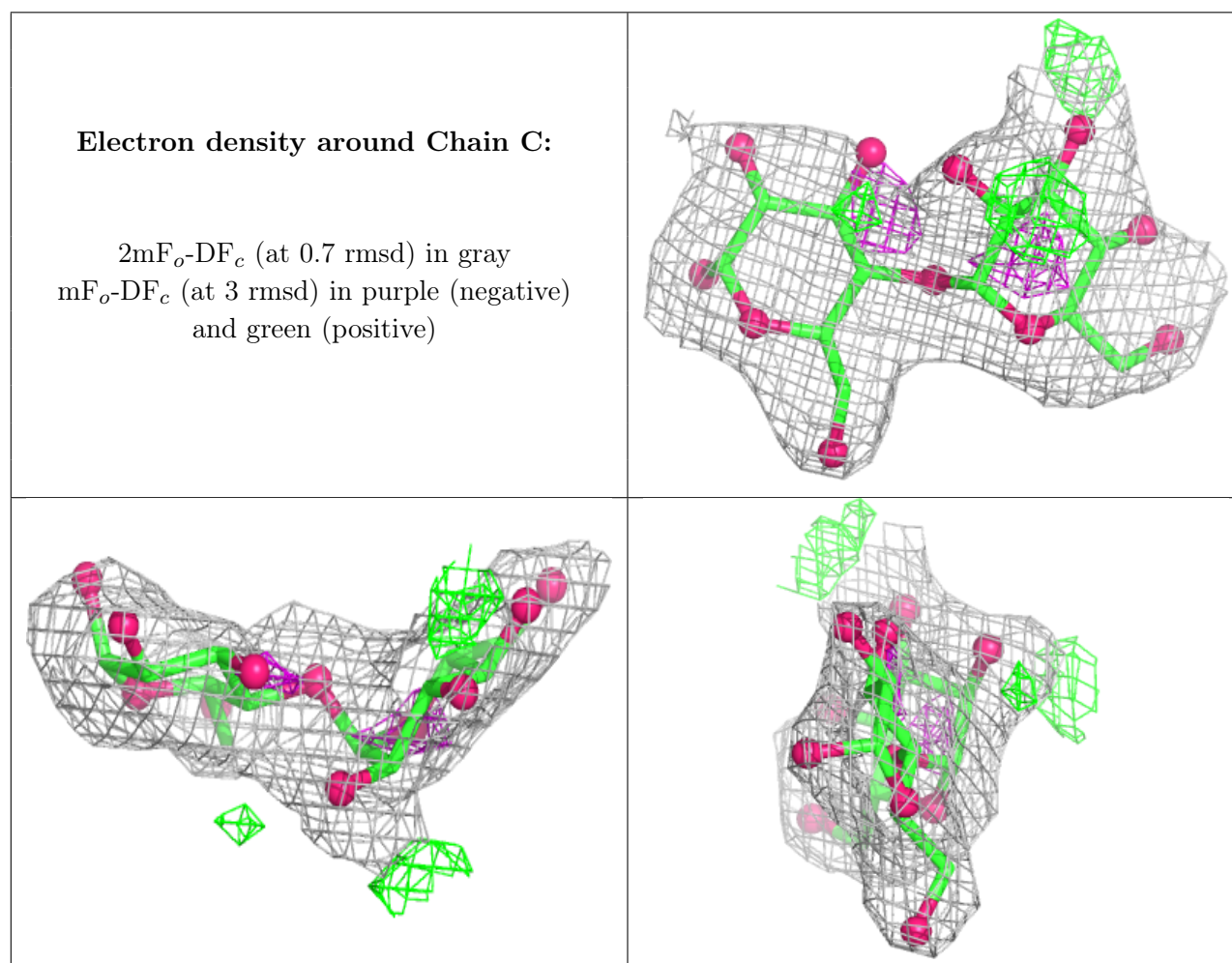
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GLC	C	2	11/12	0.81	0.14	58,60,60,61	0

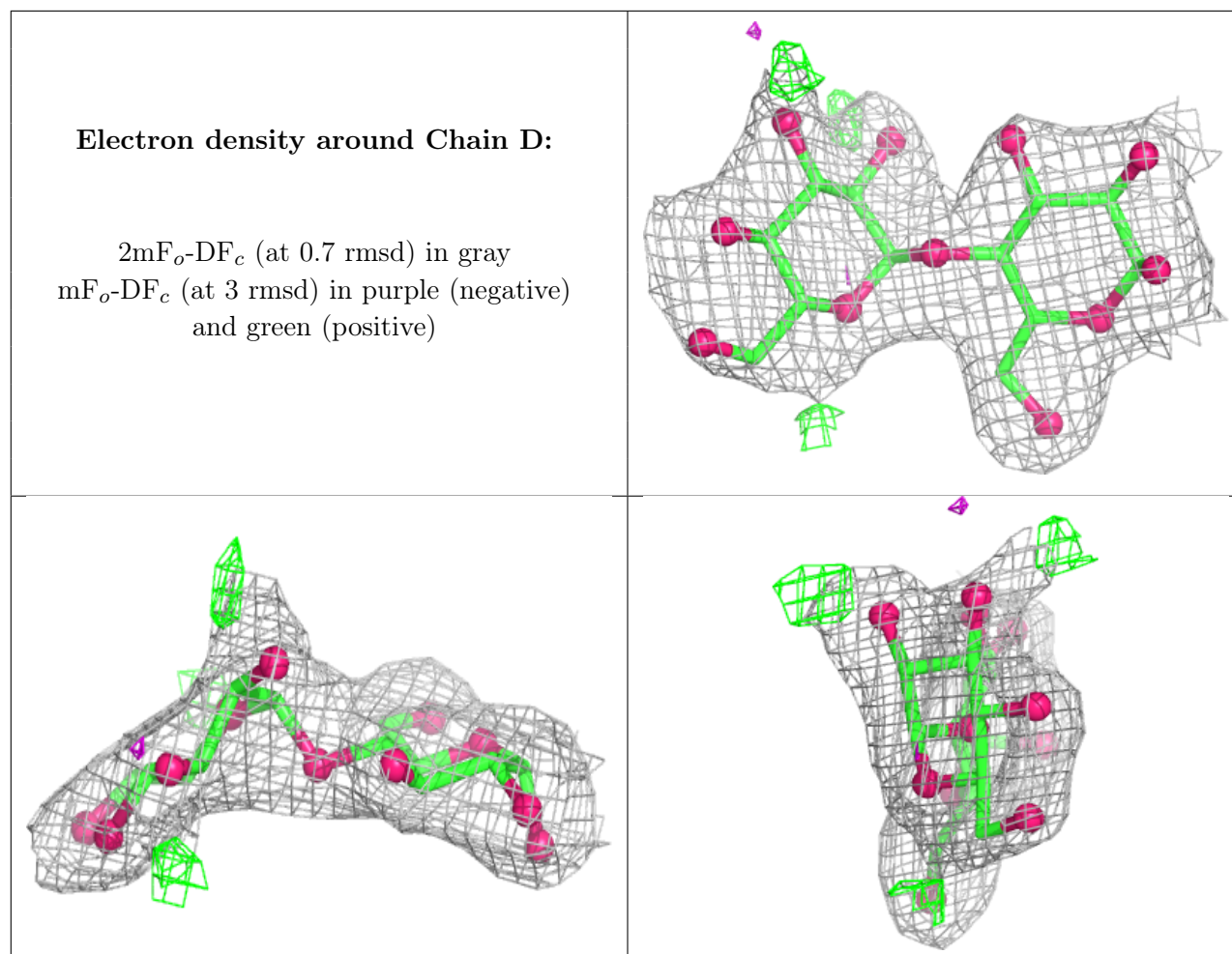
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GLC	D	2	11/12	0.85	0.13	57,59,60,60	0
2	GLC	D	1	12/12	0.91	0.12	56,56,58,58	0
2	GLC	C	1	12/12	0.91	0.12	58,59,60,60	0

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](#)

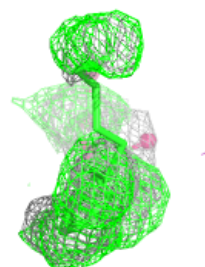
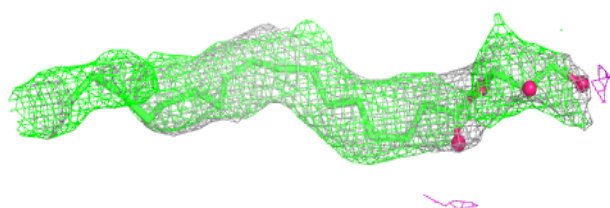
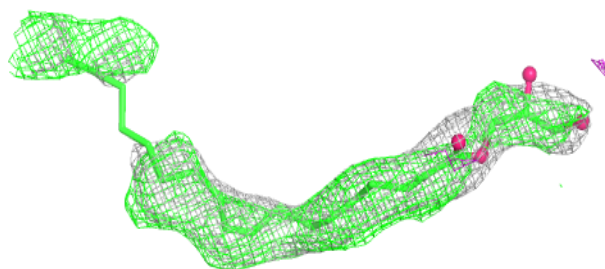
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	OLC	A	501	25/25	0.77	0.57	63,65,66,66	25
3	OLC	A	502	25/25	0.82	0.22	54,54,55,55	25
3	OLC	B	501	25/25	0.89	0.40	59,60,60,60	25

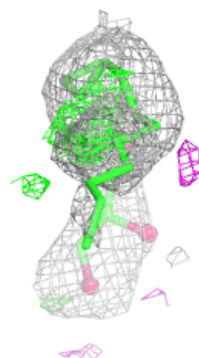
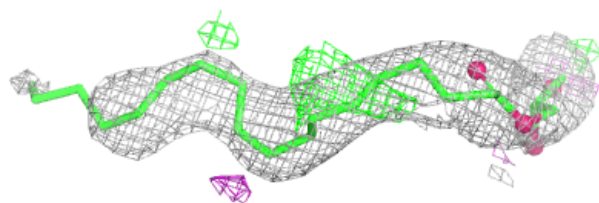
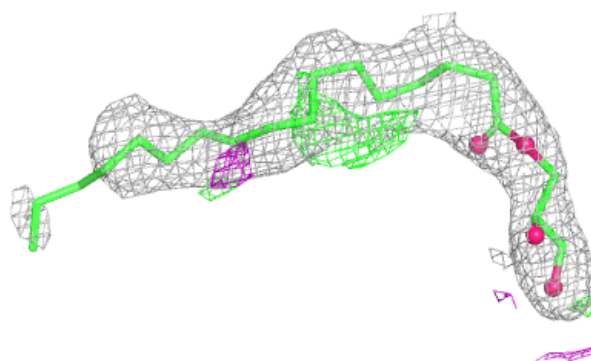
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

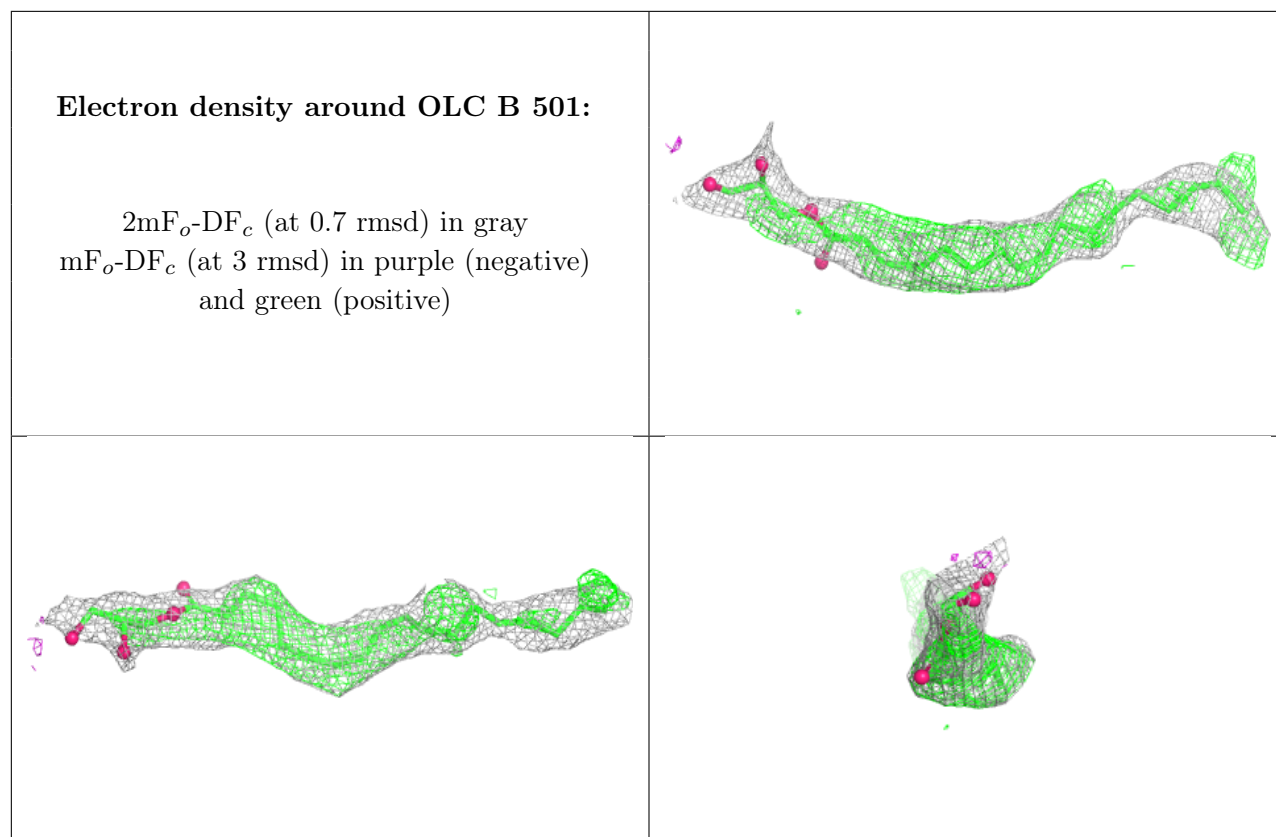
Electron density around OLC A 501:

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

**Electron density around OLC A 502:**

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





6.5 Other polymers [i](#)

There are no such residues in this entry.