



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 12:02 AM BST

PDB ID : 5KOR
Title : Arabidopsis thaliana fucosyltransferase 1 (FUT1) in complex with GDP and a xylo-oligosaccharide
Authors : Rocha, J.; de Sanctis, D.; Breton, C.
Deposited on : 2016-07-01
Resolution : 2.20 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 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 : 2.13.1
buster-report : 1.1.7 (2018)
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) : 2.13.1

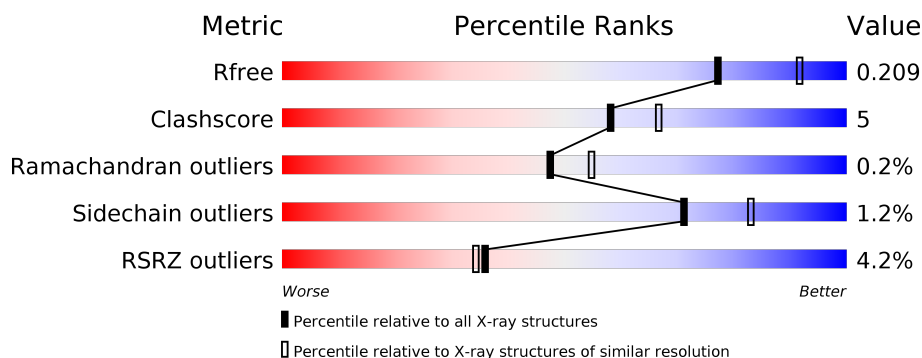
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 on 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	521	<div> <div>5%</div> <div> <div></div> <div>76%</div> <div>10%</div> <div>14%</div> </div> </div>
1	B	521	<div> <div>3%</div> <div> <div></div> <div>78%</div> <div>8%</div> <div>14%</div> </div> </div>
1	C	521	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>11%</div> <div>14%</div> </div> </div>
1	D	521	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>11%</div> <div>13%</div> </div> </div>
2	E	9	<div> <div></div> <div> <div>11%</div> <div>89%</div> </div> </div>
2	F	9	<div> <div></div> <div>100%</div> </div>

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	XYS	E	6	-	-	-	X
2	GAL	E	7	-	-	-	X
2	GAL	F	7	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 15538 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 Galactoside 2-alpha-L-fucosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	2	0
			3638	2345	601	669	23			
1	B	450	Total	C	N	O	S	0	3	0
			3647	2345	608	671	23			
1	C	449	Total	C	N	O	S	0	4	0
			3629	2334	600	671	24			
1	D	452	Total	C	N	O	S	0	2	0
			3659	2352	612	672	23			

There are 124 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	38	HIS	-	expression tag	UNP Q9SWH5
A	39	HIS	-	expression tag	UNP Q9SWH5
A	40	HIS	-	expression tag	UNP Q9SWH5
A	41	HIS	-	expression tag	UNP Q9SWH5
A	42	HIS	-	expression tag	UNP Q9SWH5
A	43	HIS	-	expression tag	UNP Q9SWH5
A	44	GLY	-	expression tag	UNP Q9SWH5
A	45	MET	-	expression tag	UNP Q9SWH5
A	46	ALA	-	expression tag	UNP Q9SWH5
A	47	SER	-	expression tag	UNP Q9SWH5
A	48	MET	-	expression tag	UNP Q9SWH5
A	49	THR	-	expression tag	UNP Q9SWH5
A	50	GLY	-	expression tag	UNP Q9SWH5
A	51	GLY	-	expression tag	UNP Q9SWH5
A	52	GLN	-	expression tag	UNP Q9SWH5
A	53	GLN	-	expression tag	UNP Q9SWH5
A	54	MET	-	expression tag	UNP Q9SWH5
A	55	GLY	-	expression tag	UNP Q9SWH5
A	56	ARG	-	expression tag	UNP Q9SWH5
A	57	ASP	-	expression tag	UNP Q9SWH5
A	58	LEU	-	expression tag	UNP Q9SWH5

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	59	TYR	-	expression tag	UNP Q9SWH5
A	60	ASP	-	expression tag	UNP Q9SWH5
A	61	ASP	-	expression tag	UNP Q9SWH5
A	62	ASP	-	expression tag	UNP Q9SWH5
A	63	ASP	-	expression tag	UNP Q9SWH5
A	64	LYS	-	expression tag	UNP Q9SWH5
A	65	SER	-	expression tag	UNP Q9SWH5
A	66	ARG	-	expression tag	UNP Q9SWH5
A	67	LEU	-	expression tag	UNP Q9SWH5
A	68	GLN	-	expression tag	UNP Q9SWH5
B	38	HIS	-	expression tag	UNP Q9SWH5
B	39	HIS	-	expression tag	UNP Q9SWH5
B	40	HIS	-	expression tag	UNP Q9SWH5
B	41	HIS	-	expression tag	UNP Q9SWH5
B	42	HIS	-	expression tag	UNP Q9SWH5
B	43	HIS	-	expression tag	UNP Q9SWH5
B	44	GLY	-	expression tag	UNP Q9SWH5
B	45	MET	-	expression tag	UNP Q9SWH5
B	46	ALA	-	expression tag	UNP Q9SWH5
B	47	SER	-	expression tag	UNP Q9SWH5
B	48	MET	-	expression tag	UNP Q9SWH5
B	49	THR	-	expression tag	UNP Q9SWH5
B	50	GLY	-	expression tag	UNP Q9SWH5
B	51	GLY	-	expression tag	UNP Q9SWH5
B	52	GLN	-	expression tag	UNP Q9SWH5
B	53	GLN	-	expression tag	UNP Q9SWH5
B	54	MET	-	expression tag	UNP Q9SWH5
B	55	GLY	-	expression tag	UNP Q9SWH5
B	56	ARG	-	expression tag	UNP Q9SWH5
B	57	ASP	-	expression tag	UNP Q9SWH5
B	58	LEU	-	expression tag	UNP Q9SWH5
B	59	TYR	-	expression tag	UNP Q9SWH5
B	60	ASP	-	expression tag	UNP Q9SWH5
B	61	ASP	-	expression tag	UNP Q9SWH5
B	62	ASP	-	expression tag	UNP Q9SWH5
B	63	ASP	-	expression tag	UNP Q9SWH5
B	64	LYS	-	expression tag	UNP Q9SWH5
B	65	SER	-	expression tag	UNP Q9SWH5
B	66	ARG	-	expression tag	UNP Q9SWH5
B	67	LEU	-	expression tag	UNP Q9SWH5
B	68	GLN	-	expression tag	UNP Q9SWH5
C	38	HIS	-	expression tag	UNP Q9SWH5

Continued on next page...

Continued from previous page...

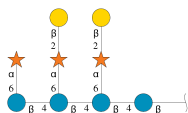
Chain	Residue	Modelled	Actual	Comment	Reference
C	39	HIS	-	expression tag	UNP Q9SWH5
C	40	HIS	-	expression tag	UNP Q9SWH5
C	41	HIS	-	expression tag	UNP Q9SWH5
C	42	HIS	-	expression tag	UNP Q9SWH5
C	43	HIS	-	expression tag	UNP Q9SWH5
C	44	GLY	-	expression tag	UNP Q9SWH5
C	45	MET	-	expression tag	UNP Q9SWH5
C	46	ALA	-	expression tag	UNP Q9SWH5
C	47	SER	-	expression tag	UNP Q9SWH5
C	48	MET	-	expression tag	UNP Q9SWH5
C	49	THR	-	expression tag	UNP Q9SWH5
C	50	GLY	-	expression tag	UNP Q9SWH5
C	51	GLY	-	expression tag	UNP Q9SWH5
C	52	GLN	-	expression tag	UNP Q9SWH5
C	53	GLN	-	expression tag	UNP Q9SWH5
C	54	MET	-	expression tag	UNP Q9SWH5
C	55	GLY	-	expression tag	UNP Q9SWH5
C	56	ARG	-	expression tag	UNP Q9SWH5
C	57	ASP	-	expression tag	UNP Q9SWH5
C	58	LEU	-	expression tag	UNP Q9SWH5
C	59	TYR	-	expression tag	UNP Q9SWH5
C	60	ASP	-	expression tag	UNP Q9SWH5
C	61	ASP	-	expression tag	UNP Q9SWH5
C	62	ASP	-	expression tag	UNP Q9SWH5
C	63	ASP	-	expression tag	UNP Q9SWH5
C	64	LYS	-	expression tag	UNP Q9SWH5
C	65	SER	-	expression tag	UNP Q9SWH5
C	66	ARG	-	expression tag	UNP Q9SWH5
C	67	LEU	-	expression tag	UNP Q9SWH5
C	68	GLN	-	expression tag	UNP Q9SWH5
D	38	HIS	-	expression tag	UNP Q9SWH5
D	39	HIS	-	expression tag	UNP Q9SWH5
D	40	HIS	-	expression tag	UNP Q9SWH5
D	41	HIS	-	expression tag	UNP Q9SWH5
D	42	HIS	-	expression tag	UNP Q9SWH5
D	43	HIS	-	expression tag	UNP Q9SWH5
D	44	GLY	-	expression tag	UNP Q9SWH5
D	45	MET	-	expression tag	UNP Q9SWH5
D	46	ALA	-	expression tag	UNP Q9SWH5
D	47	SER	-	expression tag	UNP Q9SWH5
D	48	MET	-	expression tag	UNP Q9SWH5
D	49	THR	-	expression tag	UNP Q9SWH5

Continued on next page...

Continued from previous page...

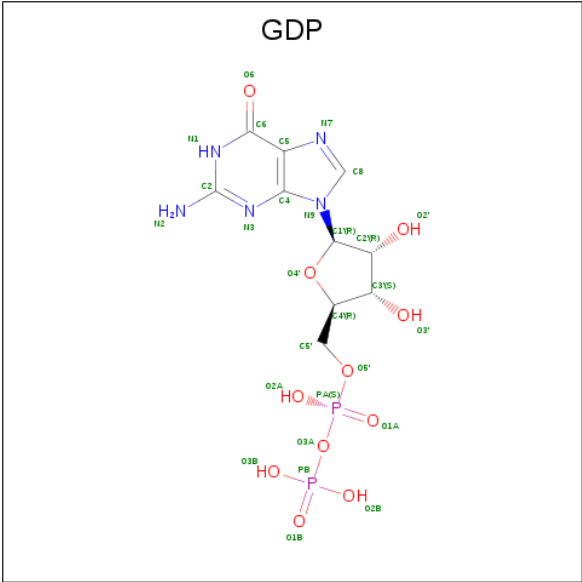
Chain	Residue	Modelled	Actual	Comment	Reference
D	50	GLY	-	expression tag	UNP Q9SWH5
D	51	GLY	-	expression tag	UNP Q9SWH5
D	52	GLN	-	expression tag	UNP Q9SWH5
D	53	GLN	-	expression tag	UNP Q9SWH5
D	54	MET	-	expression tag	UNP Q9SWH5
D	55	GLY	-	expression tag	UNP Q9SWH5
D	56	ARG	-	expression tag	UNP Q9SWH5
D	57	ASP	-	expression tag	UNP Q9SWH5
D	58	LEU	-	expression tag	UNP Q9SWH5
D	59	TYR	-	expression tag	UNP Q9SWH5
D	60	ASP	-	expression tag	UNP Q9SWH5
D	61	ASP	-	expression tag	UNP Q9SWH5
D	62	ASP	-	expression tag	UNP Q9SWH5
D	63	ASP	-	expression tag	UNP Q9SWH5
D	64	LYS	-	expression tag	UNP Q9SWH5
D	65	SER	-	expression tag	UNP Q9SWH5
D	66	ARG	-	expression tag	UNP Q9SWH5
D	67	LEU	-	expression tag	UNP Q9SWH5
D	68	GLN	-	expression tag	UNP Q9SWH5

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



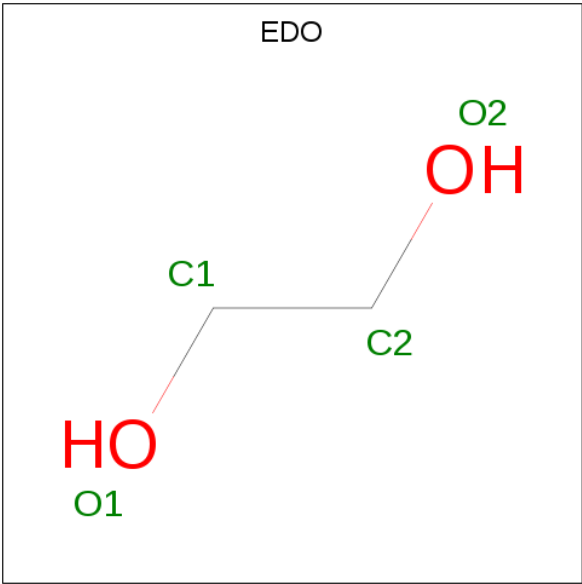
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	9	Total	C	O	0	0	0
			94	51	43			
2	F	9	Total	C	O	0	0	0
			94	51	43			

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	C	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	D	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Cl 1 1	0	0
5	D	1	Total Cl 1 1	0	0
5	C	1	Total Cl 1 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	129	Total O 129 129	0	0

Continued on next page...

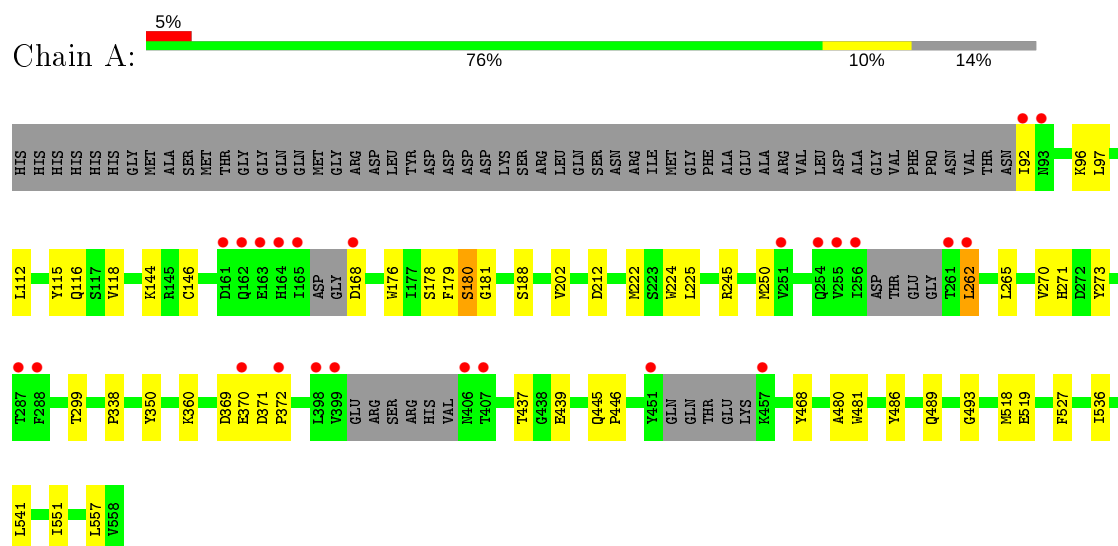
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	152	Total 152	O 152	0	1
6	C	187	Total 187	O 187	0	1
6	D	138	Total 138	O 138	0	0

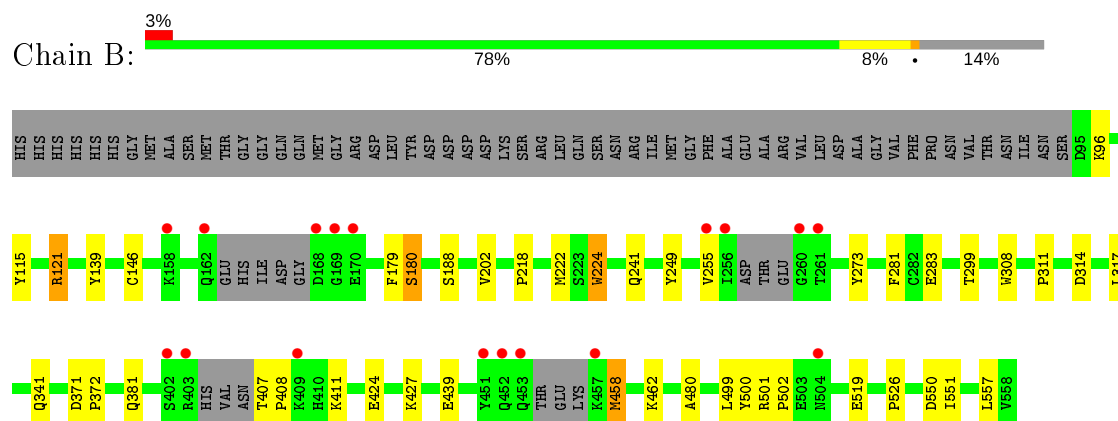
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

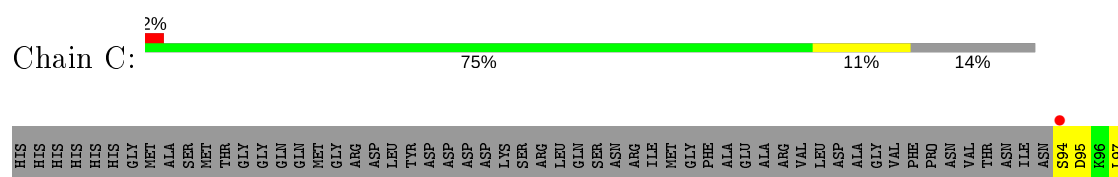
- Molecule 1: Galactoside 2-alpha-L-fucosyltransferase

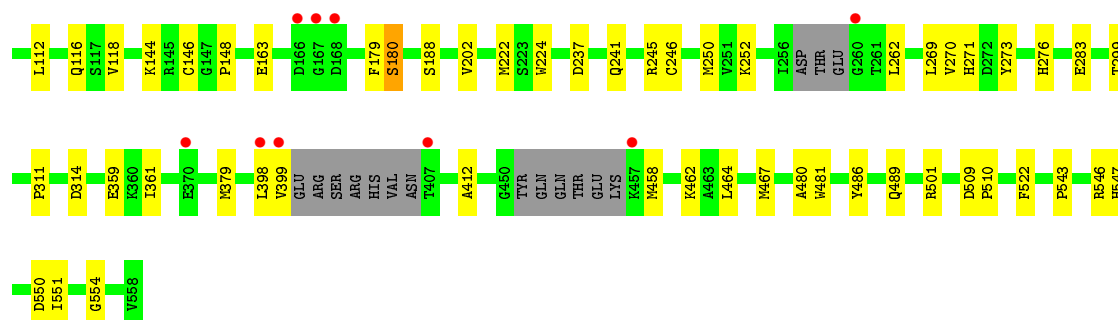


- Molecule 1: Galactoside 2-alpha-L-fucosyltransferase

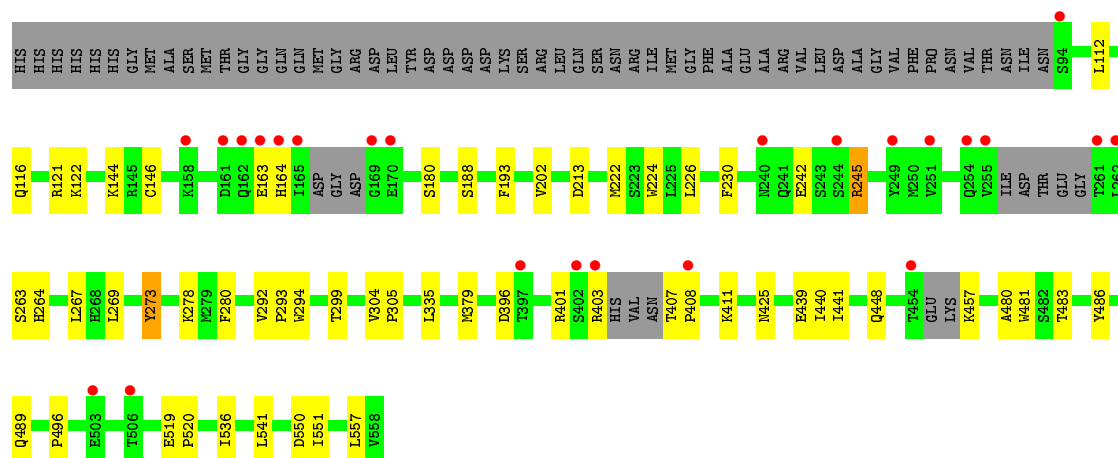
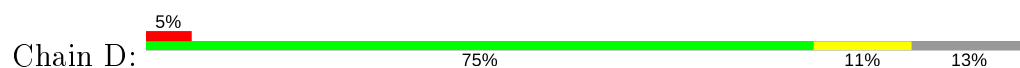


- Molecule 1: Galactoside 2-alpha-L-fucosyltransferase





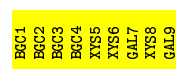
- Molecule 1: Galactoside 2-alpha-L-fucosyltransferase



- Molecule 2: alpha-D-xylopyranose-(1-6)-beta-D-glucopyranose-(1-4)-[beta-D-galactopyranose-(1-2)-alpha-D-xylopyranose-(1-6)]beta-D-glucopyranose-(1-4)-[beta-D-galactopyranose-(1-2)-alpha-D-xylopyranose-(1-6)]beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



- Molecule 2: alpha-D-xylopyranose-(1-6)-beta-D-glucopyranose-(1-4)-[beta-D-galactopyranose-(1-2)-alpha-D-xylopyranose-(1-6)]beta-D-glucopyranose-(1-4)-[beta-D-galactopyranose-(1-2)-alpha-D-xylopyranose-(1-6)]beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.97Å 85.81Å 150.72Å 90.00° 96.19° 90.00°	Depositor
Resolution (Å)	39.68 – 2.20 49.28 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (39.68-2.20) 99.8 (49.28-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.20Å)	Xtriage
Refinement program	PHENIX (1.10 _2155: ???)	Depositor
R, R_{free}	0.179 , 0.215 0.172 , 0.209	Depositor DCC
R_{free} test set	5666 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	29.2	Xtriage
Anisotropy	0.714	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15538	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.28% 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: GDP, BGC, CL, EDO, GAL, XYS

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	0.48	0/3741	0.57	0/5071
1	B	0.49	0/3749	0.58	0/5078
1	C	0.51	0/3731	0.61	0/5056
1	D	0.49	0/3761	0.58	0/5094
All	All	0.49	0/14982	0.59	0/20299

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	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	379[B]	MET	Mainchain

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	3638	0	3532	32	0
1	B	3647	0	3539	28	0
1	C	3629	0	3520	40	0
1	D	3659	0	3558	42	0
2	E	94	0	78	0	0
2	F	94	0	78	0	0
3	A	28	0	12	0	0
3	B	28	0	12	0	0
3	C	28	0	12	0	0
3	D	28	0	12	0	0
4	A	12	0	18	0	0
4	B	8	0	12	1	0
4	C	24	0	36	4	0
4	D	12	0	18	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	A	129	0	0	0	0
6	B	152	0	0	0	0
6	C	187	0	0	0	0
6	D	138	0	0	1	0
All	All	15538	0	14437	142	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 5.

All (142) 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:C:398:LEU:O	1:C:399:VAL:CG1	1.73	1.34
1:C:398:LEU:O	1:C:399:VAL:HG12	1.01	1.18
1:C:398:LEU:C	1:C:399:VAL:HG12	1.87	0.96
1:B:255:VAL:O	1:B:255:VAL:HG12	1.65	0.94
1:D:396:ASP:OD1	1:D:401[A]:ARG:NH2	2.11	0.83
1:B:121:ARG:HB3	1:B:121:ARG:HH21	1.44	0.80
1:B:121:ARG:CB	1:B:121:ARG:HH21	1.97	0.78
1:D:163:GLU:HB3	1:D:164:HIS:CE1	2.21	0.76
1:C:398:LEU:O	1:C:399:VAL:HG13	1.86	0.74
1:C:163:GLU:HB3	1:C:245:ARG:HH12	1.52	0.73
1:B:255:VAL:O	1:B:255:VAL:CG1	2.34	0.71
1:B:188:SER:HB2	1:B:299:THR:HG23	1.73	0.70
1:A:188:SER:HB2	1:A:299:THR:HG23	1.73	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:LEU:O	1:C:116:GLN:HB3	1.92	0.70
1:D:188:SER:HB2	1:D:299:THR:HG23	1.75	0.69
1:A:551:ILE:HG23	1:A:551:ILE:O	1.94	0.68
1:D:407:THR:HB	1:D:408:PRO:CD	2.25	0.66
1:D:121:ARG:HH21	1:D:121:ARG:HG3	1.60	0.66
1:A:536:ILE:HD11	1:A:541:LEU:HD21	1.76	0.65
1:D:163:GLU:HB3	1:D:164:HIS:ND1	2.10	0.65
1:C:188:SER:HB2	1:C:299:THR:HG23	1.79	0.65
1:C:250:MET:HE1	1:C:262:LEU:HD11	1.78	0.64
1:B:458:MET:O	1:B:462:LYS:HG3	1.96	0.64
1:B:371:ASP:OD1	1:B:372:PRO:HD2	1.96	0.64
1:A:480:ALA:O	1:A:481:TRP:HB2	1.99	0.63
1:C:144:LYS:HA	1:C:222[A]:MET:HG2	1.80	0.62
1:B:411:LYS:HE3	1:B:439:GLU:OE1	2.00	0.62
1:A:369:ASP:OD1	1:A:370:GLU:N	2.34	0.60
1:C:241:GLN:OE1	4:C:613:EDO:H22	2.01	0.60
1:D:411:LYS:HE3	1:D:439:GLU:OE1	2.01	0.60
1:A:338:PRO:HD3	1:A:468[A]:TYR:CZ	2.36	0.59
1:D:242:GLU:HA	1:D:242:GLU:OE1	2.03	0.59
1:C:543:PRO:O	1:C:546:ARG:HD3	2.04	0.58
1:B:121:ARG:NH2	1:B:121:ARG:HB3	2.17	0.57
1:B:96:LYS:HD3	1:B:115:TYR:HB3	1.86	0.56
1:D:401[B]:ARG:NH1	6:D:701:HOH:O	2.38	0.56
1:D:403:ARG:N	1:D:403:ARG:HD3	2.21	0.56
1:D:267:LEU:HD13	1:D:280:PHE:CE1	2.41	0.55
1:D:481:TRP:CH2	1:D:551:ILE:HD12	2.42	0.55
1:C:144:LYS:HA	1:C:222[B]:MET:HG3	1.89	0.54
1:B:551:ILE:O	1:B:551:ILE:HG23	2.08	0.54
1:C:412:ALA:HB2	4:C:614:EDO:H12	1.90	0.54
1:A:144:LYS:HA	1:A:222:MET:HG3	1.91	0.53
1:A:146:CYS:HB2	1:A:202:VAL:HG21	1.91	0.53
1:B:311:PRO:HA	1:B:314:ASP:OD2	2.09	0.52
1:D:112:LEU:O	1:D:116:GLN:HB3	2.09	0.52
1:A:527:PHE:HB3	1:A:536:ILE:O	2.10	0.51
1:C:547:HIS:HA	1:C:554:GLY:O	2.10	0.51
1:C:311:PRO:HA	1:C:314:ASP:OD2	2.10	0.51
1:A:486:TYR:O	1:A:489:GLN:HG2	2.10	0.51
1:B:480:ALA:HB1	1:B:550:ASP:HB2	1.93	0.51
1:D:213:ASP:OD2	1:D:457:LYS:HE2	2.11	0.51
1:D:263:SER:O	1:D:293:PRO:HD2	2.11	0.50
1:D:146:CYS:HB2	1:D:202:VAL:HG21	1.94	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:407:THR:HB	1:D:408:PRO:HD3	1.93	0.49
1:A:179:PHE:O	1:A:180:SER:HB3	2.12	0.49
1:C:269:LEU:HB2	1:C:299:THR:HB	1.94	0.49
1:B:139:TYR:C	1:B:139:TYR:CD2	2.86	0.49
1:B:146:CYS:HB2	1:B:202:VAL:HG21	1.94	0.49
1:A:92:ILE:HA	1:A:97:LEU:HD12	1.95	0.48
1:D:163:GLU:CB	1:D:164:HIS:ND1	2.77	0.48
1:B:283:GLU:HB2	1:B:311:PRO:HD2	1.94	0.48
1:A:245:ARG:HA	1:A:250:MET:HE3	1.95	0.48
1:C:509:ASP:HA	1:C:510:PRO:C	2.35	0.47
1:C:94:SER:HB3	1:C:97:LEU:HB2	1.96	0.47
1:D:411:LYS:HB2	1:D:441:ILE:HD13	1.96	0.47
1:C:359:GLU:OE2	1:C:361:ILE:HD11	2.14	0.47
1:C:95:ASP:HB2	1:C:118:VAL:HG21	1.97	0.46
1:A:116:GLN:C	1:A:118:VAL:H	2.19	0.46
1:B:407:THR:HG23	1:B:408:PRO:HD2	1.98	0.46
1:D:121:ARG:HG2	1:D:122:LYS:O	2.14	0.46
1:D:480:ALA:HB1	1:D:550:ASP:HB2	1.96	0.46
1:C:250:MET:CE	1:C:262:LEU:HD11	2.45	0.46
1:C:252:LYS:HB3	1:C:276:HIS:CG	2.50	0.46
1:D:144:LYS:HA	1:D:222:MET:HG3	1.97	0.46
1:D:226:LEU:HD11	1:D:230:PHE:CG	2.51	0.46
1:D:273:TYR:CE1	1:D:278:LYS:HE2	2.51	0.46
1:A:551:ILE:CG2	1:A:551:ILE:O	2.62	0.46
1:B:281:PHE:CD2	1:B:526:PRO:HG3	2.50	0.46
1:A:536:ILE:CD1	1:A:541:LEU:HD21	2.46	0.45
1:C:458:MET:O	1:C:462:LYS:HG3	2.16	0.45
1:B:427:LYS:HE2	4:B:612:EDO:O1	2.17	0.45
1:C:252:LYS:HB3	1:C:276:HIS:CD2	2.51	0.45
1:C:480:ALA:O	1:C:481:TRP:HB2	2.17	0.45
1:A:116:GLN:C	1:A:118:VAL:N	2.69	0.45
1:C:464:LEU:HA	1:C:467:MET:CE	2.47	0.45
1:C:551:ILE:O	1:C:551:ILE:HG23	2.16	0.45
1:C:146:CYS:HB2	1:C:202:VAL:HG21	1.98	0.45
1:C:148:PRO:O	1:C:222[B]:MET:HE1	2.17	0.45
1:D:121:ARG:NH2	1:D:121:ARG:HG3	2.27	0.44
1:D:264:HIS:HA	1:D:294:TRP:O	2.18	0.44
1:A:116:GLN:O	1:A:118:VAL:N	2.51	0.44
1:B:218:PRO:HG3	1:B:224:TRP:CE3	2.52	0.44
1:B:381:GLN:HG2	1:B:499:LEU:HD22	1.98	0.44
1:D:489:GLN:OE1	1:D:496:PRO:HG3	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:481:TRP:CZ3	1:D:551:ILE:HD12	2.52	0.44
1:B:241:GLN:O	1:B:249:TYR:HB2	2.18	0.44
1:A:245:ARG:HA	1:A:250:MET:CE	2.48	0.44
1:B:500:TYR:O	1:B:502:PRO:HD3	2.17	0.43
1:C:283:GLU:OE1	1:C:311:PRO:HG2	2.19	0.43
1:A:180:SER:OG	1:A:181:GLY:N	2.51	0.43
1:D:483:THR:HA	1:D:486:TYR:CD2	2.53	0.43
1:C:252:LYS:NZ	4:C:613:EDO:H21	2.34	0.43
1:C:311:PRO:HA	1:C:314:ASP:CG	2.39	0.43
1:A:96:LYS:HG2	1:A:115:TYR:HB3	1.99	0.43
1:D:519:GLU:HG3	1:D:557:LEU:HB2	2.01	0.43
1:C:486:TYR:O	1:C:489:GLN:HG2	2.18	0.43
1:C:480:ALA:HB1	1:C:550:ASP:HB2	2.01	0.43
1:A:437:THR:OG1	1:A:439:GLU:HG3	2.20	0.42
1:D:486:TYR:HB3	1:D:520:PRO:HB3	2.01	0.42
1:C:95:ASP:N	1:C:95:ASP:OD1	2.52	0.42
1:B:308:TRP:CG	1:B:317:LEU:HD13	2.54	0.42
1:C:179:PHE:O	1:C:180:SER:HB3	2.20	0.42
1:A:112:LEU:O	1:A:116:GLN:HB2	2.19	0.42
1:D:193:PHE:CD2	1:D:335:LEU:HD11	2.55	0.42
1:D:408:PRO:HB2	1:D:440:ILE:HD12	2.02	0.41
1:A:176:TRP:CH2	1:A:178:SER:HA	2.56	0.41
1:A:519:GLU:HG3	1:A:557:LEU:HB2	2.03	0.41
1:B:519:GLU:HG3	1:B:557:LEU:HB2	2.02	0.41
1:D:536:ILE:HD11	1:D:541:LEU:HD21	2.03	0.41
1:B:179:PHE:O	1:B:180:SER:CB	2.68	0.41
1:A:493:GLY:HA2	1:A:518:MET:SD	2.59	0.41
1:D:245:ARG:HA	1:D:245:ARG:HD2	1.62	0.41
1:B:458:MET:HG3	1:B:462:LYS:HE3	2.02	0.41
1:C:241:GLN:HB2	4:C:613:EDO:O1	2.21	0.41
1:C:270:VAL:HG12	1:C:271:HIS:N	2.36	0.41
1:A:445:GLN:HA	1:A:446:PRO:HD3	1.94	0.41
1:B:188:SER:HB2	1:B:299:THR:CG2	2.47	0.41
1:D:480:ALA:O	1:D:481:TRP:HB2	2.21	0.41
1:A:212:ASP:HB3	1:A:225:LEU:HD11	2.02	0.40
1:A:270:VAL:HG12	1:A:271:HIS:N	2.36	0.40
1:D:269:LEU:HB2	1:D:299:THR:HB	2.03	0.40
1:D:448:GLN:O	1:D:448:GLN:HG3	2.21	0.40
1:D:292:VAL:HA	1:D:293:PRO:HD3	1.96	0.40
1:D:541:LEU:HD23	1:D:541:LEU:HA	1.86	0.40
1:A:350:TYR:CE2	1:A:360:LYS:HE2	2.56	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:304:VAL:N	1:D:305:PRO:CD	2.83	0.40
1:D:379[B]:MET:HG2	1:D:425:ASN:CG	2.42	0.40
1:C:241:GLN:HA	1:C:246:CYS:SG	2.62	0.40
1:A:262:LEU:HD23	1:A:265:LEU:HD13	2.03	0.40
1:A:371:ASP:CG	1:A:372:PRO:HD2	2.41	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	442/521 (85%)	427 (97%)	14 (3%)	1 (0%)	47	55
1	B	443/521 (85%)	427 (96%)	15 (3%)	1 (0%)	47	55
1	C	445/521 (85%)	433 (97%)	11 (2%)	1 (0%)	47	55
1	D	444/521 (85%)	428 (96%)	15 (3%)	1 (0%)	47	55
All	All	1774/2084 (85%)	1715 (97%)	55 (3%)	4 (0%)	47	55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	180	SER
1	A	180	SER
1	C	180	SER
1	D	180	SER

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	403/460 (88%)	399 (99%)	4 (1%)	76	86
1	B	403/460 (88%)	394 (98%)	9 (2%)	52	65
1	C	402/460 (87%)	397 (99%)	5 (1%)	71	83
1	D	405/460 (88%)	402 (99%)	3 (1%)	84	91
All	All	1613/1840 (88%)	1592 (99%)	21 (1%)	71	81

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	168	ASP
1	A	224	TRP
1	A	262	LEU
1	A	273	TYR
1	B	121	ARG
1	B	222	MET
1	B	224	TRP
1	B	273	TYR
1	B	341	GLN
1	B	424[A]	GLU
1	B	424[B]	GLU
1	B	458	MET
1	B	501	ARG
1	C	224	TRP
1	C	237	ASP
1	C	273	TYR
1	C	501	ARG
1	C	522	PHE
1	D	224	TRP
1	D	245	ARG
1	D	273	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	410	HIS

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 ⓘ

18 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	E	1	2	12,12,12	0.67	0	17,17,17	1.22	3 (17%)
2	BGC	E	2	2	11,11,12	0.38	0	15,15,17	1.29	2 (13%)
2	BGC	E	3	2	11,11,12	0.48	0	15,15,17	1.01	0
2	BGC	E	4	2	11,11,12	0.66	0	15,15,17	1.10	1 (6%)
2	XYS	E	5	2	9,9,10	0.88	1 (11%)	10,12,14	2.07	2 (20%)
2	XYS	E	6	2	9,9,10	0.75	0	10,12,14	2.35	3 (30%)
2	GAL	E	7	2	11,11,12	1.47	2 (18%)	15,15,17	0.77	0
2	XYS	E	8	2	9,9,10	1.02	1 (11%)	10,12,14	1.42	2 (20%)
2	GAL	E	9	2	11,11,12	1.34	2 (18%)	15,15,17	1.07	1 (6%)
2	BGC	F	1	2	12,12,12	0.50	0	17,17,17	0.91	1 (5%)
2	BGC	F	2	2	11,11,12	0.58	0	15,15,17	1.29	2 (13%)
2	BGC	F	3	2	11,11,12	0.53	0	15,15,17	1.31	3 (20%)
2	BGC	F	4	2	11,11,12	0.56	0	15,15,17	1.35	1 (6%)
2	XYS	F	5	2	9,9,10	1.04	1 (11%)	10,12,14	1.91	3 (30%)
2	XYS	F	6	2	9,9,10	0.91	1 (11%)	10,12,14	1.86	3 (30%)
2	GAL	F	7	2	11,11,12	1.57	2 (18%)	15,15,17	1.12	1 (6%)
2	XYS	F	8	2	9,9,10	1.08	1 (11%)	10,12,14	2.00	2 (20%)
2	GAL	F	9	2	11,11,12	1.46	3 (27%)	15,15,17	1.08	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
2	BGC	E	1	2	-	1/2/22/22	0/1/1/1
2	BGC	E	2	2	-	0/2/19/22	0/1/1/1
2	BGC	E	3	2	-	1/2/19/22	0/1/1/1
2	BGC	E	4	2	-	0/2/19/22	0/1/1/1
2	XYS	E	5	2	-	-	0/1/1/1
2	XYS	E	6	2	-	-	0/1/1/1
2	GAL	E	7	2	-	2/2/19/22	0/1/1/1
2	XYS	E	8	2	-	-	0/1/1/1
2	GAL	E	9	2	-	1/2/19/22	0/1/1/1
2	BGC	F	1	2	-	2/2/22/22	0/1/1/1
2	BGC	F	2	2	-	0/2/19/22	0/1/1/1
2	BGC	F	3	2	-	2/2/19/22	0/1/1/1
2	BGC	F	4	2	-	0/2/19/22	0/1/1/1
2	XYS	F	5	2	-	-	0/1/1/1
2	XYS	F	6	2	-	-	0/1/1/1
2	GAL	F	7	2	-	2/2/19/22	0/1/1/1
2	XYS	F	8	2	-	-	0/1/1/1
2	GAL	F	9	2	-	1/2/19/22	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	7	GAL	C4-C3	2.75	1.59	1.52
2	E	9	GAL	C1-C2	2.72	1.58	1.52
2	F	7	GAL	O5-C5	2.58	1.48	1.43
2	F	9	GAL	O5-C5	2.40	1.48	1.43
2	E	9	GAL	C2-C3	2.39	1.56	1.52
2	F	8	XYS	O5-C1	-2.37	1.38	1.42
2	F	9	GAL	C2-C3	2.30	1.55	1.52
2	F	7	GAL	C2-C3	2.23	1.55	1.52
2	E	8	XYS	O5-C1	-2.17	1.38	1.42
2	F	5	XYS	O5-C1	-2.16	1.38	1.42
2	F	9	GAL	O5-C1	2.15	1.47	1.43
2	F	6	XYS	O5-C1	-2.14	1.38	1.42
2	E	5	XYS	O5-C1	-2.10	1.38	1.42
2	E	7	GAL	O5-C1	2.09	1.47	1.43

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	6	XYS	C5-O5-C1	6.49	121.50	111.52
2	E	5	XYS	C5-O5-C1	5.31	119.69	111.52
2	F	8	XYS	C5-O5-C1	5.26	119.61	111.52
2	F	5	XYS	C5-O5-C1	4.67	118.70	111.52
2	F	6	XYS	C5-O5-C1	4.07	117.78	111.52
2	F	4	BGC	C1-C2-C3	3.55	114.03	109.67
2	E	8	XYS	C5-O5-C1	2.87	115.94	111.52
2	F	7	GAL	C1-O5-C5	2.65	115.78	112.19
2	F	2	BGC	O4-C4-C5	-2.61	102.82	109.30
2	F	3	BGC	O4-C4-C5	-2.59	102.86	109.30
2	E	1	BGC	O4-C4-C5	-2.54	103.00	109.30
2	E	4	BGC	O3-C3-C2	-2.53	105.16	109.99
2	F	3	BGC	C6-C5-C4	-2.52	107.11	113.00
2	E	8	XYS	C1-C2-C3	2.37	112.58	109.67
2	E	2	BGC	C1-C2-C3	2.33	112.53	109.67
2	E	1	BGC	O5-C5-C6	2.29	112.12	106.44
2	E	2	BGC	O4-C4-C5	-2.29	103.62	109.30
2	F	1	BGC	O5-C5-C6	2.28	112.10	106.44
2	F	6	XYS	C1-C2-C3	2.27	112.45	109.67
2	E	6	XYS	C1-C2-C3	2.26	112.44	109.67
2	E	6	XYS	C5-C4-C3	2.25	112.43	109.67
2	F	6	XYS	C4-C3-C2	-2.22	108.28	110.92
2	E	9	GAL	C1-C2-C3	2.22	112.39	109.67
2	F	5	XYS	O3-C3-C4	2.20	114.21	109.99
2	E	1	BGC	O5-C5-C4	2.18	113.64	109.69
2	F	5	XYS	C4-C3-C2	-2.02	108.52	110.92
2	F	3	BGC	C3-C4-C5	2.02	113.84	110.24
2	F	8	XYS	O4-C4-C3	-2.01	106.10	110.14
2	E	5	XYS	C5-C4-C3	2.01	112.14	109.67
2	F	2	BGC	C1-C2-C3	2.00	112.12	109.67

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	7	GAL	O5-C5-C6-O6
2	E	7	GAL	C4-C5-C6-O6
2	F	3	BGC	O5-C5-C6-O6
2	F	9	GAL	O5-C5-C6-O6
2	E	9	GAL	O5-C5-C6-O6
2	F	1	BGC	C4-C5-C6-O6
2	E	3	BGC	O5-C5-C6-O6
2	E	1	BGC	C4-C5-C6-O6

Continued on next page...

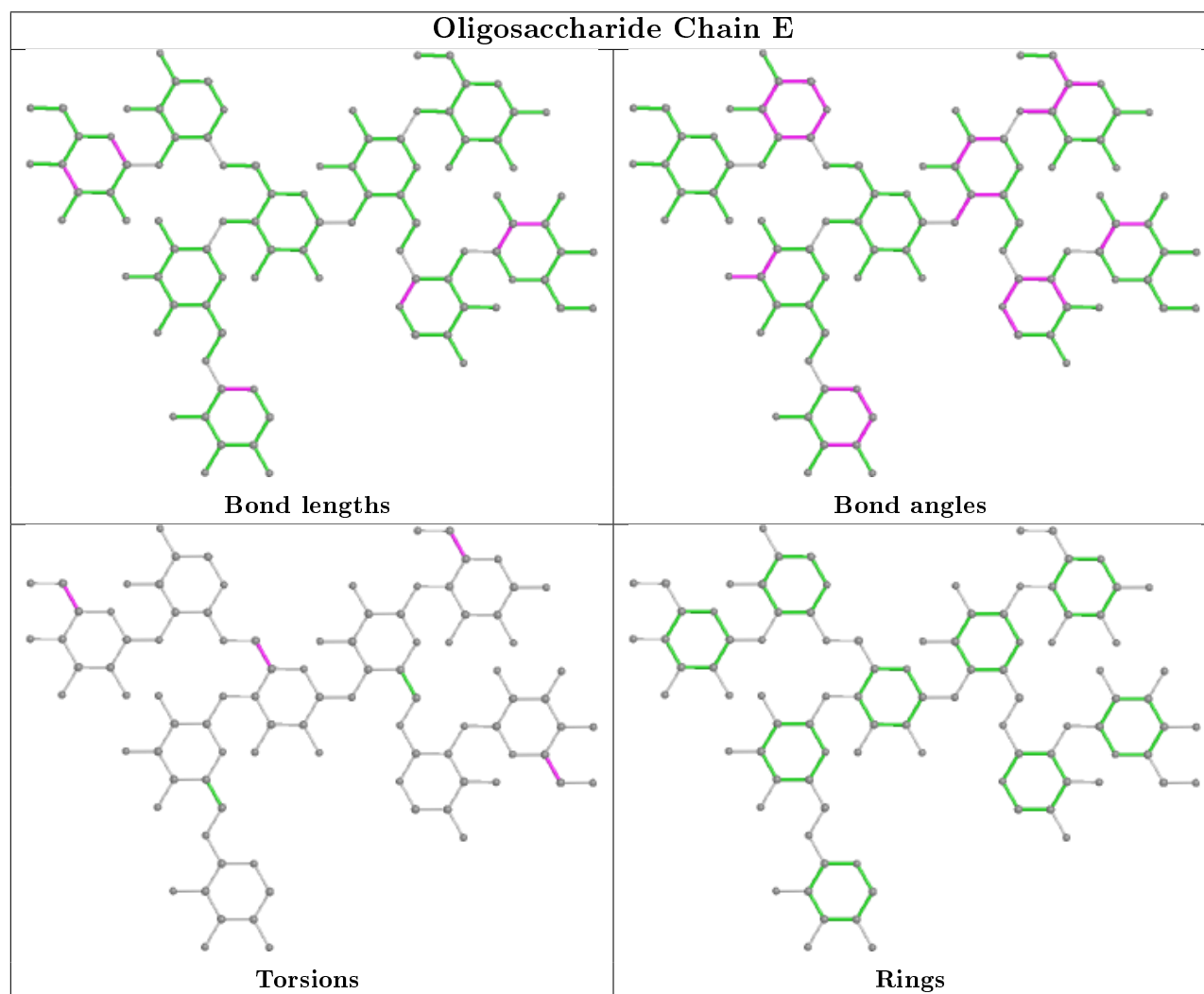
Continued from previous page...

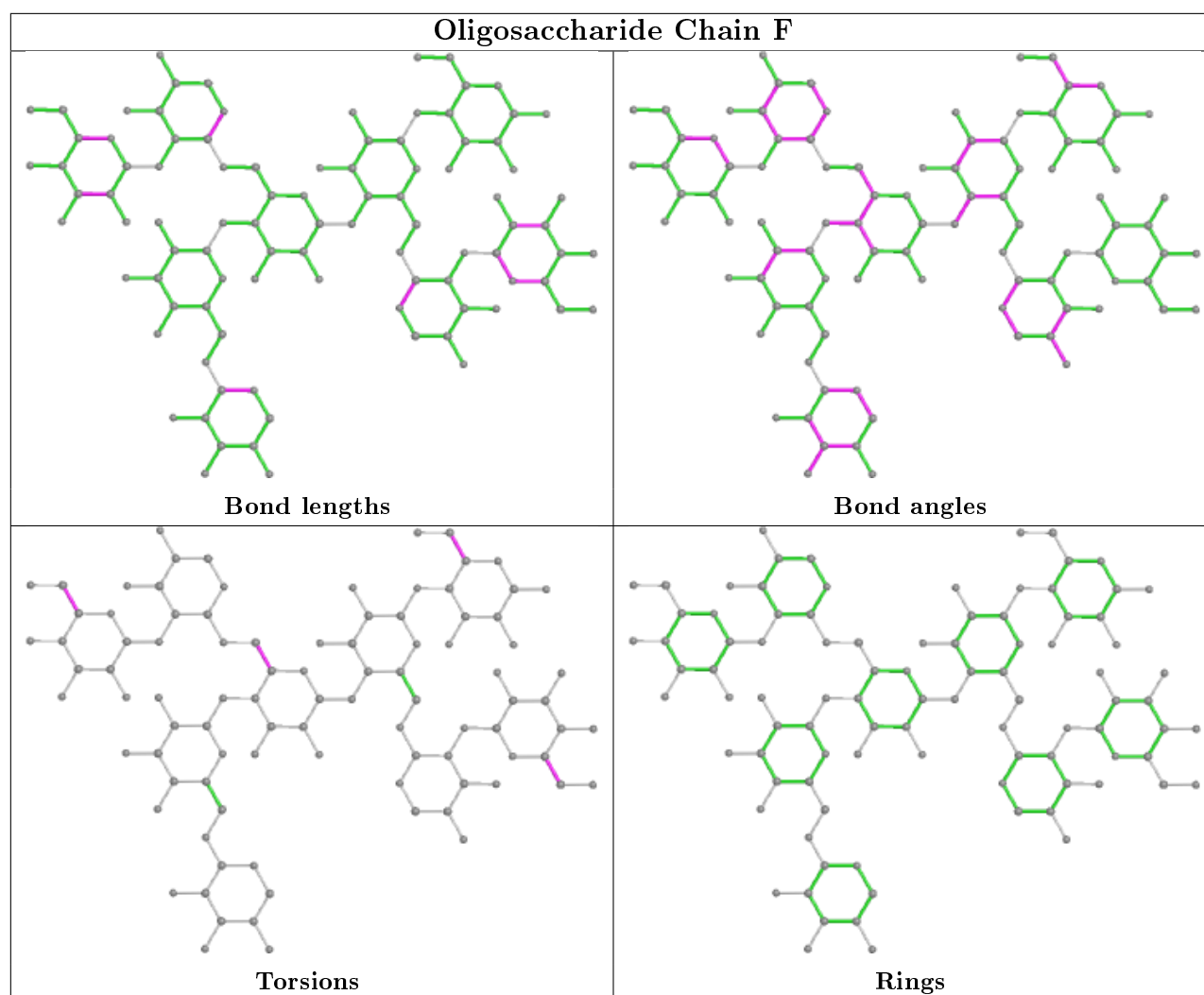
Mol	Chain	Res	Type	Atoms
2	F	3	BGC	C4-C5-C6-O6
2	F	1	BGC	O5-C5-C6-O6
2	F	7	GAL	O5-C5-C6-O6
2	F	7	GAL	C4-C5-C6-O6

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](#)

Of 21 ligands modelled in this entry, 3 are monoatomic - leaving 18 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GDP	D	601	-	24,30,30	1.22	3 (12%)	31,47,47	1.42	4 (12%)
3	GDP	B	601	-	24,30,30	1.20	3 (12%)	31,47,47	1.34	5 (16%)
4	EDO	D	603	-	3,3,3	0.54	0	2,2,2	0.21	0
4	EDO	C	616	-	3,3,3	0.59	0	2,2,2	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	A	604	-	3,3,3	0.39	0	2,2,2	0.83	0
4	EDO	D	604	-	3,3,3	0.47	0	2,2,2	0.31	0
4	EDO	C	612	-	3,3,3	0.34	0	2,2,2	0.40	0
4	EDO	B	612	-	3,3,3	0.45	0	2,2,2	0.41	0
4	EDO	A	603	-	3,3,3	0.52	0	2,2,2	0.31	0
3	GDP	A	601	-	24,30,30	1.14	2 (8%)	31,47,47	1.46	4 (12%)
3	GDP	C	601	-	24,30,30	1.17	2 (8%)	31,47,47	1.51	5 (16%)
4	EDO	B	611	-	3,3,3	0.46	0	2,2,2	0.29	0
4	EDO	C	613	-	3,3,3	0.74	0	2,2,2	0.94	0
4	EDO	C	614	-	3,3,3	0.32	0	2,2,2	0.25	0
4	EDO	A	602	-	3,3,3	0.44	0	2,2,2	0.49	0
4	EDO	D	602	-	3,3,3	0.41	0	2,2,2	0.52	0
4	EDO	C	615	-	3,3,3	0.46	0	2,2,2	0.31	0
4	EDO	C	611	-	3,3,3	0.56	0	2,2,2	0.49	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	GDP	D	601	-	-	5/12/32/32	0/3/3/3
3	GDP	B	601	-	-	5/12/32/32	0/3/3/3
4	EDO	D	603	-	-	1/1/1/1	-
4	EDO	C	616	-	-	0/1/1/1	-
4	EDO	A	604	-	-	1/1/1/1	-
4	EDO	D	604	-	-	0/1/1/1	-
4	EDO	C	612	-	-	1/1/1/1	-
4	EDO	B	612	-	-	1/1/1/1	-
4	EDO	A	603	-	-	0/1/1/1	-
3	GDP	A	601	-	-	7/12/32/32	0/3/3/3
3	GDP	C	601	-	-	7/12/32/32	0/3/3/3
4	EDO	B	611	-	-	0/1/1/1	-
4	EDO	C	613	-	-	1/1/1/1	-
4	EDO	C	614	-	-	1/1/1/1	-
4	EDO	A	602	-	-	0/1/1/1	-
4	EDO	D	602	-	-	0/1/1/1	-
4	EDO	C	615	-	-	0/1/1/1	-
4	EDO	C	611	-	-	0/1/1/1	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	601	GDP	C6-C5	3.38	1.47	1.41
3	A	601	GDP	C6-C5	3.21	1.46	1.41
3	D	601	GDP	C6-C5	3.11	1.46	1.41
3	B	601	GDP	C6-C5	2.85	1.46	1.41
3	C	601	GDP	C5-C4	2.48	1.47	1.40
3	B	601	GDP	PB-O2B	-2.43	1.45	1.54
3	D	601	GDP	C5-C4	2.37	1.47	1.40
3	D	601	GDP	PB-O2B	-2.33	1.45	1.54
3	B	601	GDP	C5-C4	2.23	1.46	1.40
3	A	601	GDP	C5-C4	2.20	1.46	1.40

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	601	GDP	C5-C6-N1	-4.65	117.07	123.43
3	A	601	GDP	C6-C5-C4	-4.20	116.79	120.80
3	D	601	GDP	C5-C6-N1	-4.05	117.89	123.43
3	B	601	GDP	C6-C5-C4	-3.70	117.26	120.80
3	C	601	GDP	C6-N1-C2	3.43	121.38	115.93
3	D	601	GDP	C6-C5-C4	-3.42	117.54	120.80
3	B	601	GDP	C2-N3-C4	3.27	119.09	115.36
3	A	601	GDP	C5-C6-N1	-3.24	119.00	123.43
3	A	601	GDP	C2-N3-C4	3.22	119.03	115.36
3	C	601	GDP	C6-C5-C4	-3.18	117.76	120.80
3	D	601	GDP	C6-N1-C2	3.01	120.71	115.93
3	B	601	GDP	C5-C6-N1	-2.79	119.61	123.43
3	A	601	GDP	C6-N1-C2	2.78	120.35	115.93
3	D	601	GDP	C2-N3-C4	2.72	118.46	115.36
3	C	601	GDP	C2-N3-C4	2.63	118.37	115.36
3	B	601	GDP	PA-O3A-PB	-2.57	124.02	132.83
3	B	601	GDP	C6-N1-C2	2.23	119.47	115.93
3	C	601	GDP	C4-C5-N7	-2.18	107.13	109.40

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	601	GDP	PA-O3A-PB-O2B
3	B	601	GDP	PA-O3A-PB-O2B
3	B	601	GDP	O4'-C4'-C5'-O5'
3	C	601	GDP	PA-O3A-PB-O3B
3	A	601	GDP	PA-O3A-PB-O3B

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	A	601	GDP	O4'-C4'-C5'-O5'
3	D	601	GDP	O4'-C4'-C5'-O5'
3	C	601	GDP	O4'-C4'-C5'-O5'
3	C	601	GDP	C3'-C4'-C5'-O5'
3	D	601	GDP	C3'-C4'-C5'-O5'
3	B	601	GDP	C3'-C4'-C5'-O5'
3	A	601	GDP	C3'-C4'-C5'-O5'
4	A	604	EDO	O1-C1-C2-O2
3	C	601	GDP	PB-O3A-PA-O1A
3	D	601	GDP	C4'-C5'-O5'-PA
3	A	601	GDP	C4'-C5'-O5'-PA
4	C	613	EDO	O1-C1-C2-O2
3	D	601	GDP	PB-O3A-PA-O5'
3	B	601	GDP	PB-O3A-PA-O5'
3	C	601	GDP	PB-O3A-PA-O5'
3	A	601	GDP	PB-O3A-PA-O5'
3	B	601	GDP	C4'-C5'-O5'-PA
3	C	601	GDP	C4'-C5'-O5'-PA
4	C	612	EDO	O1-C1-C2-O2
4	D	603	EDO	O1-C1-C2-O2
4	C	614	EDO	O1-C1-C2-O2
3	A	601	GDP	PA-O3A-PB-O1B
4	B	612	EDO	O1-C1-C2-O2
3	C	601	GDP	PA-O3A-PB-O2B
3	A	601	GDP	PB-O3A-PA-O1A

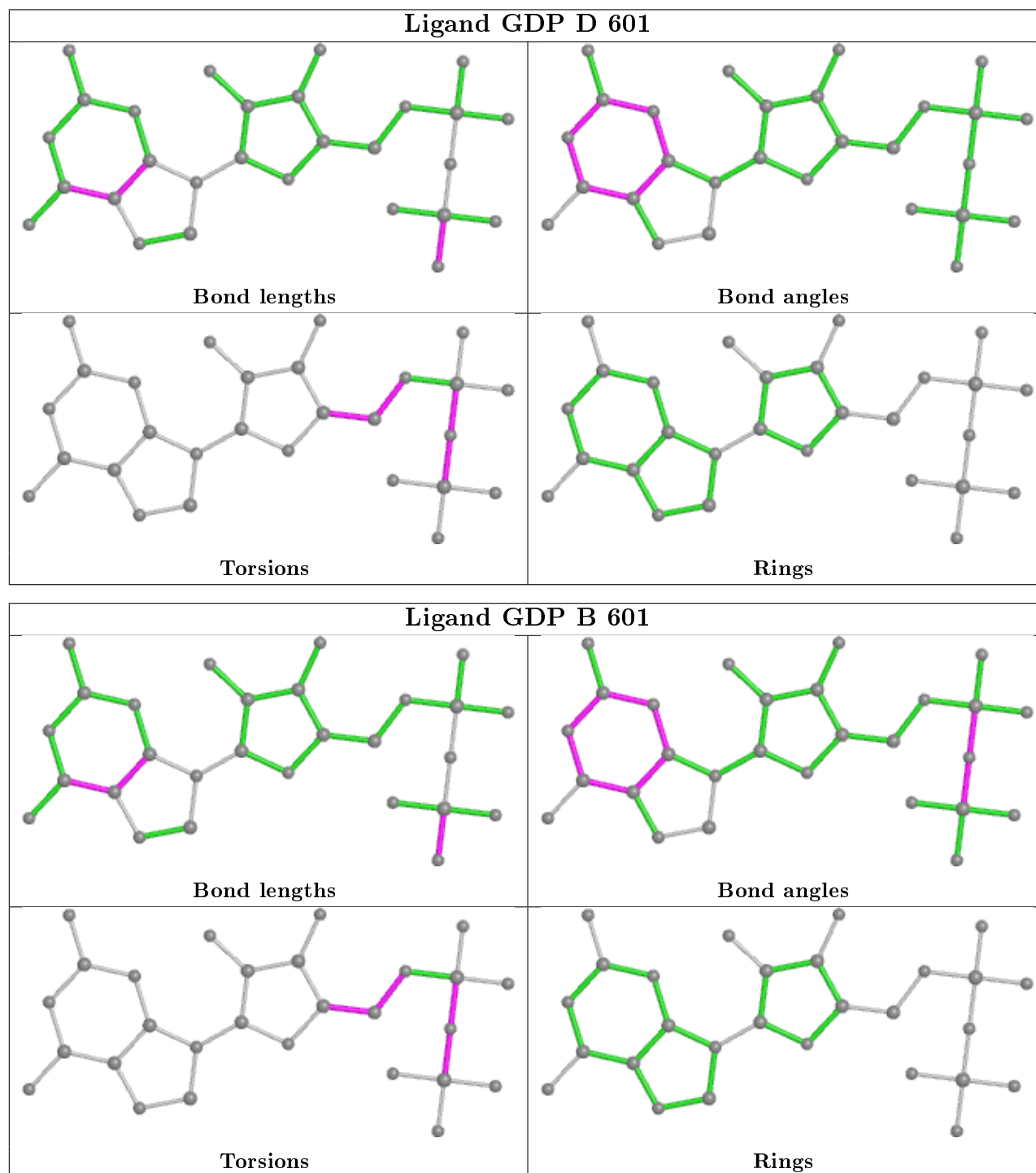
There are no ring outliers.

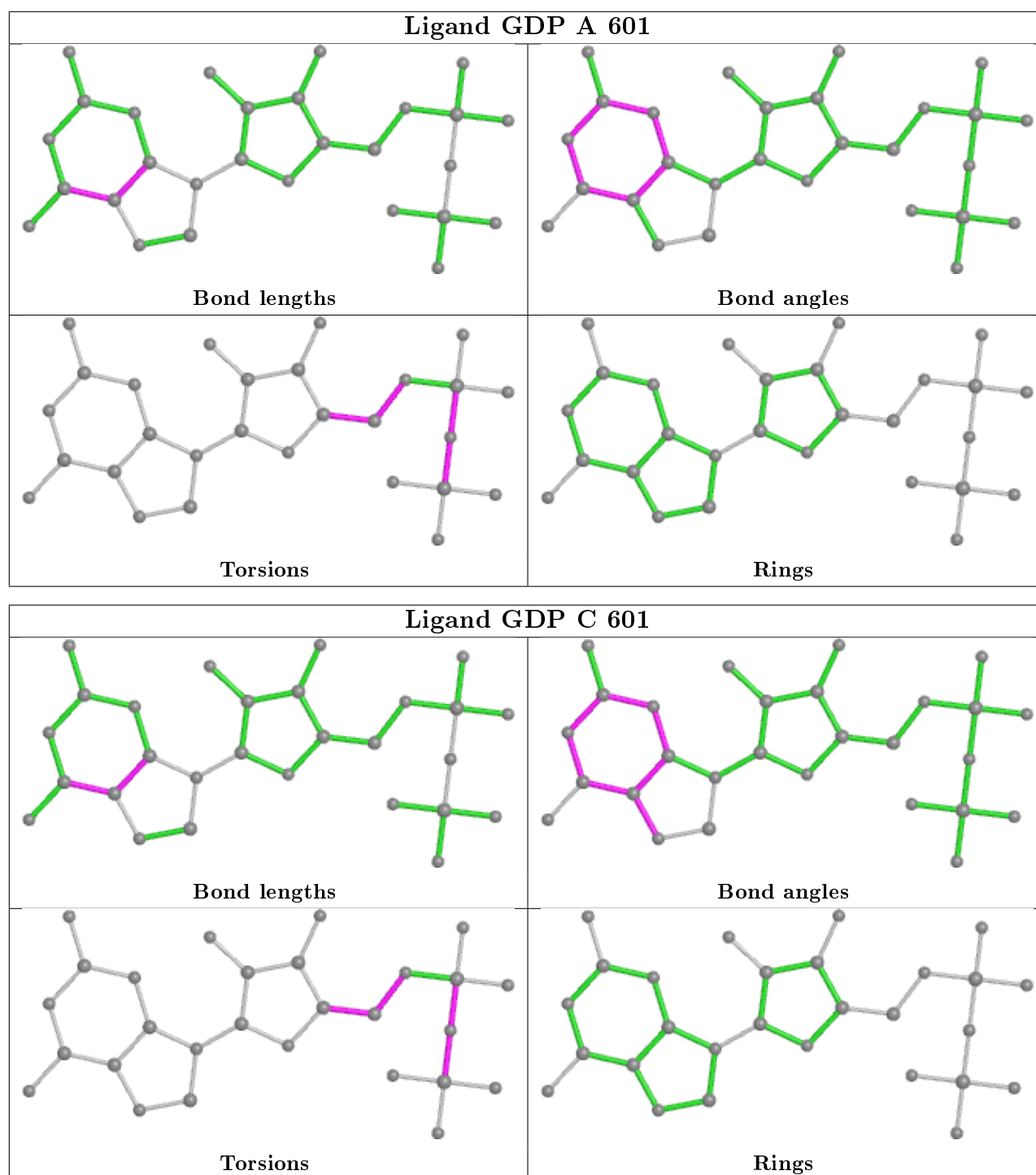
3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	612	EDO	1	0
4	C	613	EDO	3	0
4	C	614	EDO	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 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	450/521 (86%)	-0.04	24 (5%) 26 25	20, 35, 73, 120	0
1	B	450/521 (86%)	-0.11	17 (3%) 40 38	19, 35, 71, 111	0
1	C	449/521 (86%)	-0.18	10 (2%) 62 59	18, 30, 58, 99	0
1	D	452/521 (86%)	-0.02	24 (5%) 26 25	18, 36, 74, 119	0
All	All	1801/2084 (86%)	-0.09	75 (4%) 36 34	18, 34, 71, 120	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	164	HIS	6.4
1	A	406	ASN	5.7
1	D	255	VAL	5.7
1	B	256	ILE	5.1
1	D	261	THR	5.0
1	C	166	ASP	4.7
1	D	403	ARG	4.7
1	D	164	HIS	4.6
1	C	457	LYS	4.5
1	A	92	ILE	4.3
1	A	372	PRO	4.2
1	A	255	VAL	4.2
1	D	169	GLY	4.2
1	A	451	TYR	4.1
1	B	261	THR	4.1
1	A	398	LEU	4.1
1	A	163	GLU	4.1
1	A	261	THR	4.0
1	D	503	GLU	3.9
1	A	457	LYS	3.9
1	D	163	GLU	3.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	403	ARG	3.7
1	A	256	ILE	3.5
1	B	168	ASP	3.3
1	A	262	LEU	3.3
1	D	454	THR	3.3
1	D	165	ILE	3.2
1	A	399	VAL	3.2
1	D	262	LEU	3.1
1	D	162	GLN	3.1
1	B	162	GLN	3.1
1	B	451	TYR	3.0
1	C	167	GLY	3.0
1	B	402	SER	3.0
1	D	161	ASP	3.0
1	B	260	GLY	2.9
1	B	457	LYS	2.9
1	A	93	ASN	2.9
1	D	402	SER	2.9
1	B	170	GLU	2.7
1	D	158	LYS	2.7
1	C	407	THR	2.6
1	A	165	ILE	2.6
1	A	161	ASP	2.6
1	D	244	SER	2.6
1	C	370	GLU	2.6
1	A	288	PHE	2.6
1	B	453	GLN	2.5
1	D	249	TYR	2.5
1	D	397	THR	2.5
1	C	398	LEU	2.5
1	C	399	VAL	2.5
1	A	162	GLN	2.5
1	A	407	THR	2.4
1	A	370	GLU	2.4
1	B	169	GLY	2.4
1	C	168	ASP	2.4
1	D	170	GLU	2.4
1	B	504	ASN	2.3
1	A	251	VAL	2.3
1	D	94	SER	2.3
1	D	254	GLN	2.3
1	B	409	LYS	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	452	GLN	2.2
1	D	506	THR	2.2
1	A	168	ASP	2.2
1	D	408	PRO	2.1
1	A	287	THR	2.1
1	C	94	SER	2.1
1	B	255	VAL	2.1
1	B	158	LYS	2.1
1	A	254	GLN	2.1
1	D	240	ASN	2.0
1	D	251	VAL	2.0
1	C	260	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

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.

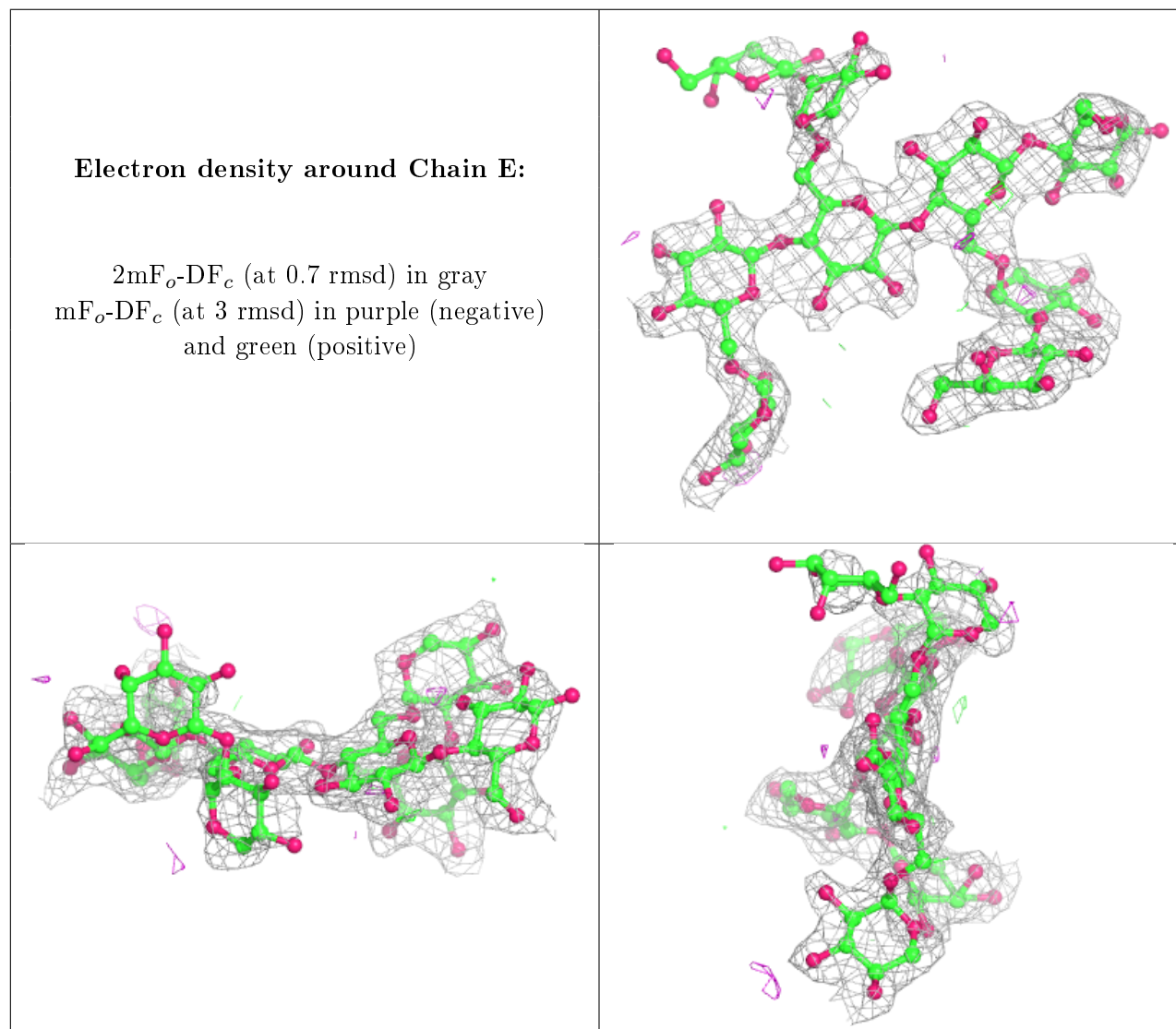
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GAL	F	7	11/12	0.52	0.65	101,107,108,109	0
2	GAL	E	7	11/12	0.62	0.62	103,105,106,106	0
2	XYS	E	6	9/10	0.65	0.42	91,94,96,100	0
2	XYS	F	6	9/10	0.80	0.25	81,83,86,93	0
2	BGC	E	4	11/12	0.81	0.22	64,72,75,77	0
2	XYS	E	5	9/10	0.82	0.17	59,61,62,62	0
2	BGC	E	3	11/12	0.84	0.17	62,65,80,86	0
2	BGC	E	1	12/12	0.87	0.23	64,69,70,72	0
2	XYS	E	8	9/10	0.88	0.16	51,54,56,56	0
2	BGC	F	1	12/12	0.88	0.21	53,59,63,66	0
2	XYS	F	5	9/10	0.91	0.11	42,43,44,44	0
2	BGC	E	2	11/12	0.92	0.14	56,58,61,61	0
2	BGC	F	3	11/12	0.92	0.19	41,50,65,73	0
2	BGC	F	4	11/12	0.93	0.23	45,49,55,57	0
2	XYS	F	8	9/10	0.93	0.11	39,46,49,53	0
2	GAL	E	9	11/12	0.94	0.12	40,42,47,48	0

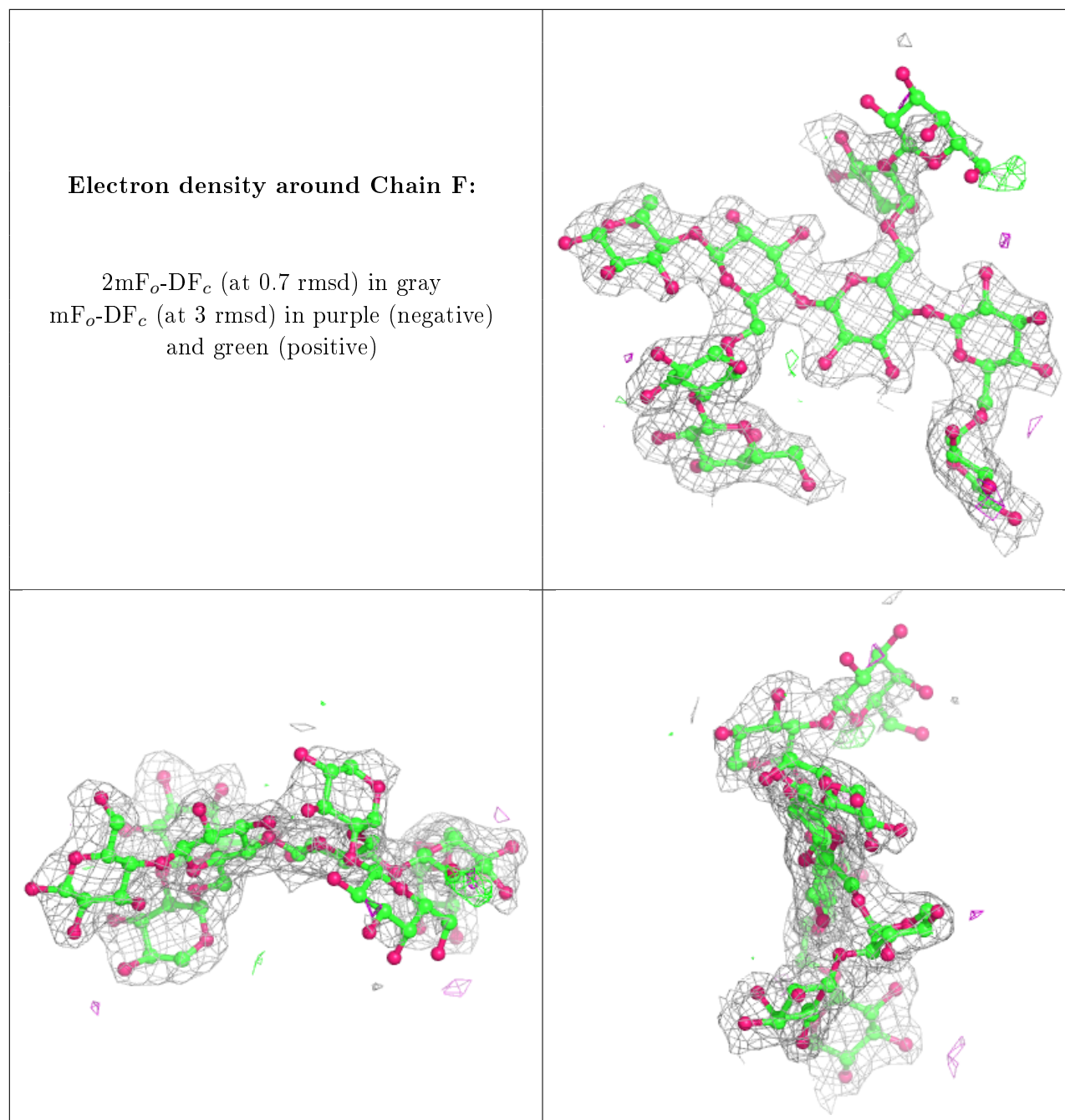
Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BGC	F	2	11/12	0.95	0.17	42,45,48,48	0
2	GAL	F	9	11/12	0.96	0.10	29,34,36,36	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(\AA^2)	Q<0.9
4	EDO	A	603	4/4	0.79	0.18	59,61,63,64	0
4	EDO	C	614	4/4	0.80	0.18	58,58,59,59	0

Continued on next page...

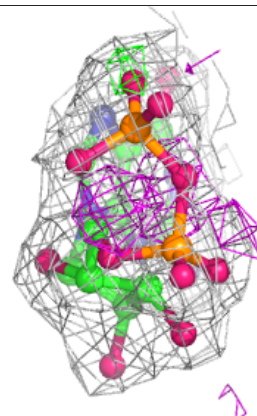
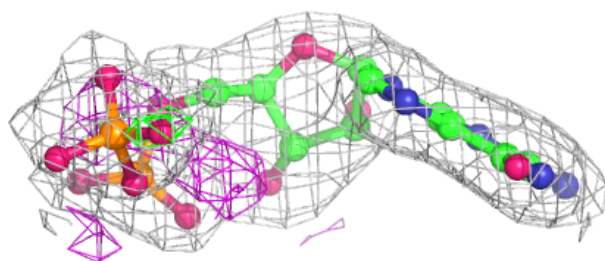
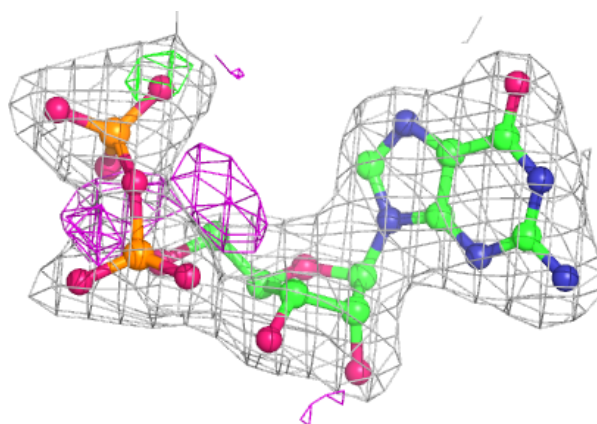
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	EDO	C	613	4/4	0.85	0.28	38,43,44,49	0
4	EDO	C	612	4/4	0.87	0.13	65,66,67,68	0
4	EDO	C	616	4/4	0.88	0.14	63,64,64,65	0
4	EDO	D	604	4/4	0.89	0.24	60,61,63,64	0
4	EDO	C	615	4/4	0.89	0.19	63,64,65,66	0
4	EDO	A	604	4/4	0.90	0.21	49,54,58,62	0
3	GDP	A	601	28/28	0.90	0.15	38,51,67,67	0
5	CL	B	613	1/1	0.92	0.09	57,57,57,57	0
3	GDP	B	601	28/28	0.94	0.12	21,34,41,45	0
4	EDO	A	602	4/4	0.94	0.15	36,37,37,38	0
3	GDP	C	601	28/28	0.94	0.12	24,36,49,51	0
4	EDO	D	603	4/4	0.94	0.18	57,57,58,61	0
3	GDP	D	601	28/28	0.95	0.12	27,34,40,43	0
4	EDO	D	602	4/4	0.96	0.13	42,43,43,44	0
4	EDO	B	612	4/4	0.96	0.16	56,56,58,60	0
5	CL	D	605	1/1	0.96	0.11	58,58,58,58	0
5	CL	C	617	1/1	0.98	0.09	51,51,51,51	0
4	EDO	B	611	4/4	0.98	0.14	37,37,38,40	0
4	EDO	C	611	4/4	0.98	0.09	31,32,33,34	0

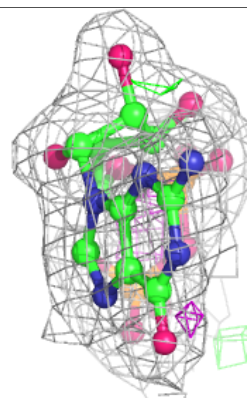
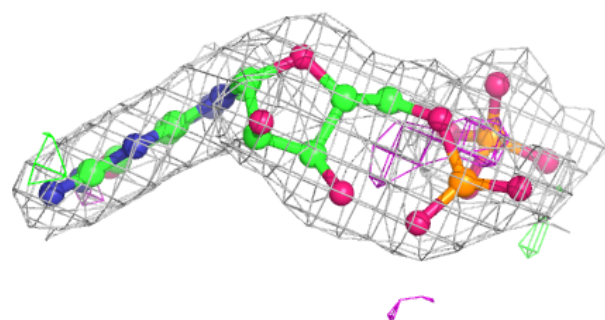
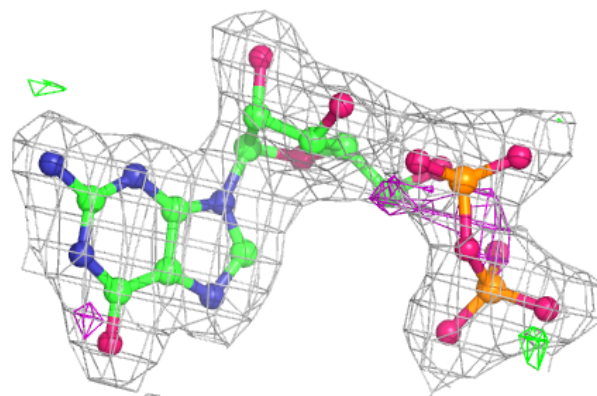
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GDP A 601:

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

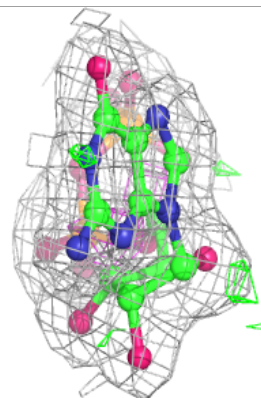
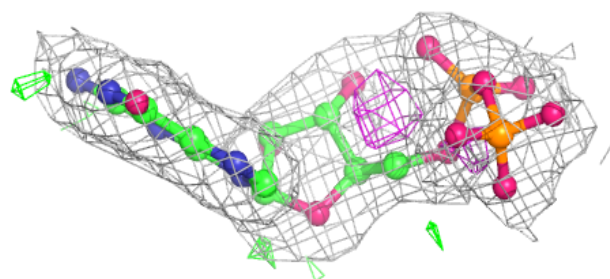
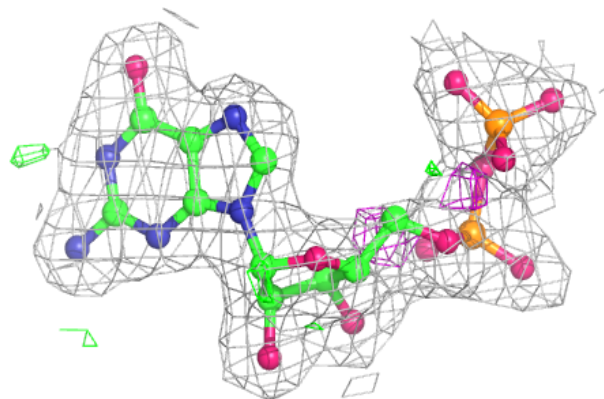
**Electron density around GDP B 601:**

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

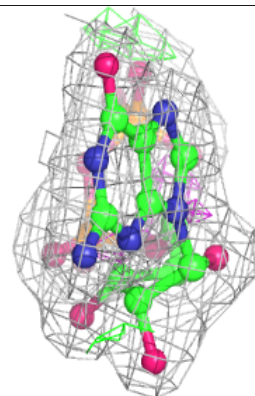
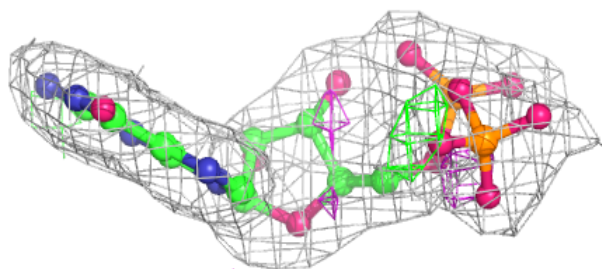
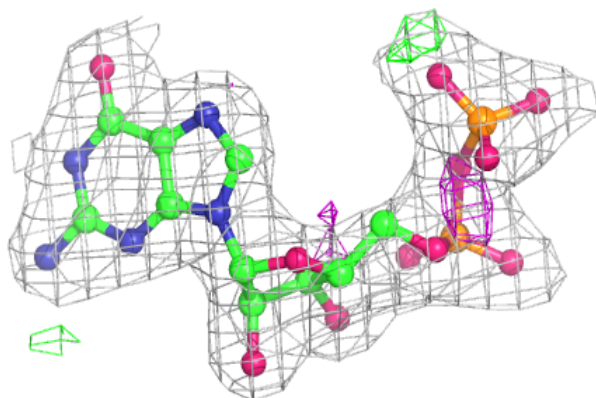


Electron density around GDP C 601:

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

**Electron density around GDP D 601:**

$2mF_o-DF_c$ (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.