



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 6, 2023 – 06:00 pm BST

PDB ID : 4ZEB  
Title : PBP AccA from *A. tumefaciens* C58 in complex with agrocinopine A  
Authors : El Sahili, A.; Morera, S.  
Deposited on : 2015-04-20  
Resolution : 1.89 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](mailto: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>.

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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)  
Xtriage (Phenix) : 1.13  
EDS : 2.33  
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) : 2.33

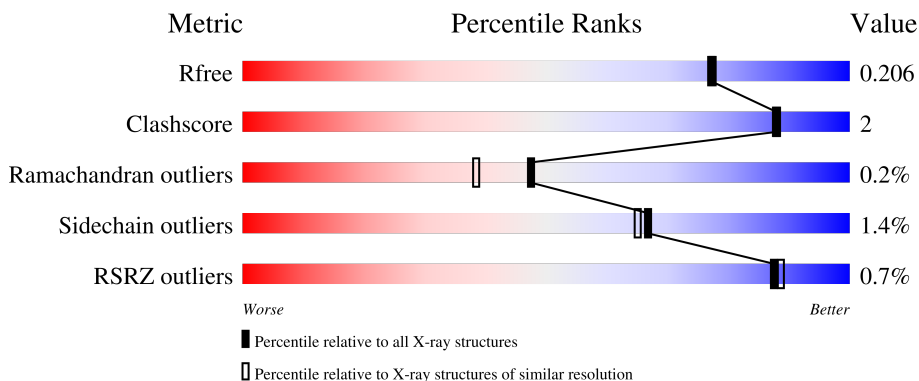
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.89 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	499	
1	B	499	
2	C	2	
2	D	2	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	C	1	X	-	-	-
2	GLC	D	1	X	-	-	-

## 2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 8520 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.

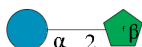
- Molecule 1 is a protein called ABC transporter substrate-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	490	3959	2555	662	723	19	0	4	0
1	B	489	3950	2549	661	720	20	0	2	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	MET	-	initiating methionine	UNP Q52012
A	522	HIS	-	expression tag	UNP Q52012
A	523	HIS	-	expression tag	UNP Q52012
A	524	HIS	-	expression tag	UNP Q52012
A	525	HIS	-	expression tag	UNP Q52012
A	526	HIS	-	expression tag	UNP Q52012
A	527	HIS	-	expression tag	UNP Q52012
B	29	MET	-	initiating methionine	UNP Q52012
B	522	HIS	-	expression tag	UNP Q52012
B	523	HIS	-	expression tag	UNP Q52012
B	524	HIS	-	expression tag	UNP Q52012
B	525	HIS	-	expression tag	UNP Q52012
B	526	HIS	-	expression tag	UNP Q52012
B	527	HIS	-	expression tag	UNP Q52012

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



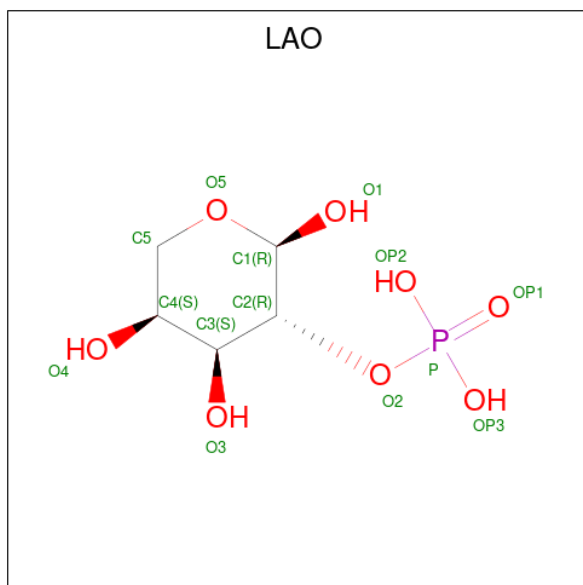
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	C	2	23	12	11	0	0	0

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

- Molecule 3 is 2-O-phosphono-alpha-L-arabinopyranose (three-letter code: LAO) (formula:  $C_5H_{11}O_8P$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			13	5	7	1		
3	B	1	Total	C	O	P	0	0
			13	5	7	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			7	4	3		

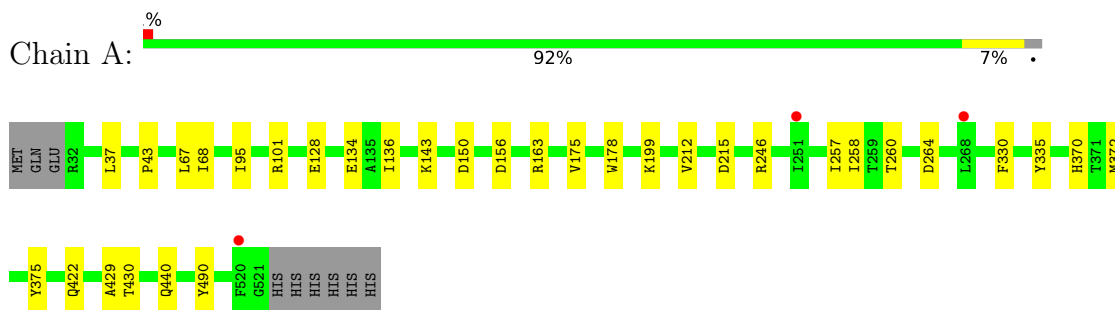
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	275	Total 275	O 275	0	0
6	B	253	Total 253	O 253	0	0

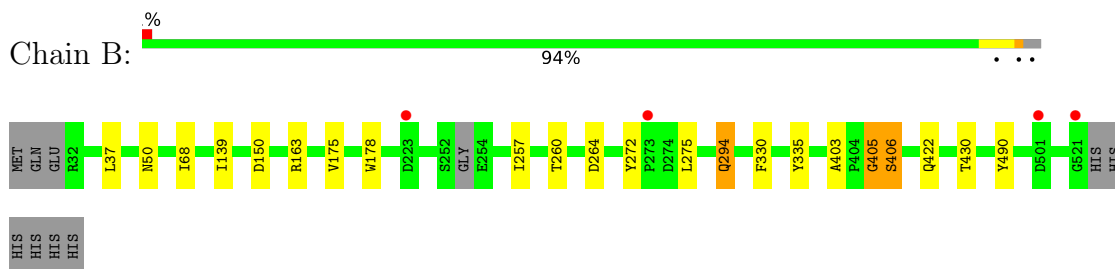
### 3 Residue-property plots [i](#)

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: ABC transporter substrate-binding protein



- Molecule 1: ABC transporter substrate-binding protein



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



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





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	177.21Å 51.24Å 119.95Å 90.00° 114.33° 90.00°	Depositor
Resolution (Å)	37.80 – 1.89 48.84 – 1.89	Depositor EDS
% Data completeness (in resolution range)	98.7 (37.80-1.89) 99.2 (48.84-1.89)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.22 (at 1.90Å)	Xtrriage
Refinement program	BUSTER 2.10.0	Depositor
R, $R_{free}$	0.172 , 0.200 0.177 , 0.206	Depositor DCC
$R_{free}$ test set	3903 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.7	Xtrriage
Anisotropy	0.363	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 42.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8520	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.49% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: EDO, FRU, PEG, LAO, 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	0.51	0/4080	0.65	0/5542
1	B	0.51	0/4064	0.66	0/5518
All	All	0.51	0/8144	0.65	0/11060

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	3959	0	3914	18	0
1	B	3950	0	3900	14	0
2	C	23	0	20	1	0
2	D	23	0	20	0	0
3	A	13	0	10	0	0
3	B	13	0	10	0	0
4	A	4	0	6	1	0
5	B	7	0	10	0	0
6	A	275	0	0	2	0
6	B	253	0	0	0	0
All	All	8520	0	7890	32	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 (32) 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:405:GLY:HA3	1:B:406:SER:HB2	1.45	0.98
1:B:403:ALA:O	1:B:406:SER:HB3	1.75	0.87
1:A:67:LEU:HD12	1:A:175[B]:VAL:HG21	1.68	0.74
1:B:405:GLY:CA	1:B:406:SER:HB2	2.18	0.72
1:A:67:LEU:HD12	1:A:175[B]:VAL:CG2	2.29	0.62
1:A:370:HIS:HD2	6:A:749:HOH:O	1.87	0.58
1:A:212[B]:VAL:HG13	1:A:215:ASP:HB3	1.86	0.57
1:A:372:MET:HE2	1:A:375:TYR:HB2	1.86	0.56
1:B:272:TYR:HB3	1:B:275:LEU:HD13	1.87	0.56
1:B:294:GLN:HE21	1:B:294:GLN:HA	1.73	0.54
1:B:403:ALA:O	1:B:406:SER:CB	2.53	0.53
1:B:50:ASN:HD21	1:B:139:ILE:HA	1.75	0.52
1:A:95:ILE:HD11	1:A:101:ARG:HB2	1.92	0.51
1:A:335:TYR:OH	1:A:422:GLN:NE2	2.45	0.50
1:B:403:ALA:O	1:B:405:GLY:N	2.44	0.50
1:B:68:ILE:HD12	1:B:175[B]:VAL:HG23	1.95	0.48
1:B:335:TYR:OH	1:B:422:GLN:NE2	2.46	0.48
1:B:422:GLN:HE22	1:B:430:THR:H	1.61	0.48
1:A:136:ILE:HD11	1:A:143:LYS:HD3	1.96	0.47
1:A:199:LYS:HE2	4:A:602:EDO:H12	1.96	0.47
1:A:150:ASP:HB3	1:A:163:ARG:HG3	1.98	0.46
1:A:422:GLN:HE22	1:A:430:THR:H	1.62	0.46
1:B:150:ASP:HB3	1:B:163:ARG:HG3	1.98	0.44
1:A:37:LEU:HD22	1:A:257:ILE:HB	2.00	0.43
1:B:405:GLY:CA	1:B:406:SER:CB	2.90	0.43
1:A:246:ARG:HG2	1:A:258:ILE:HB	2.01	0.43
1:A:440:GLN:NE2	6:A:701:HOH:O	2.51	0.43
1:A:128:GLU:HA	1:A:134:GLU:CG	2.49	0.42
1:B:37:LEU:HD22	1:B:257:ILE:HB	2.00	0.42
1:A:68:ILE:HD12	1:A:175[B]:VAL:HG13	2.01	0.42
1:A:429:ALA:HA	1:A:430:THR:HA	1.80	0.42
1:A:43:PRO:HA	2:C:1:GLC:H61	2.01	0.41

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	492/499 (99%)	475 (96%)	17 (4%)	0	100	100
1	B	487/499 (98%)	469 (96%)	16 (3%)	2 (0%)	34	24
All	All	979/998 (98%)	944 (96%)	33 (3%)	2 (0%)	47	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	406	SER
1	B	405	GLY

### 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	423/428 (99%)	417 (99%)	6 (1%)	67	65
1	B	421/428 (98%)	415 (99%)	6 (1%)	67	65
All	All	844/856 (99%)	832 (99%)	12 (1%)	67	65

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	156	ASP
1	A	178	TRP
1	A	260	THR
1	A	264	ASP

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Mol	Chain	Res	Type
1	A	330	PHE
1	A	490	TYR
1	B	178	TRP
1	B	260	THR
1	B	264	ASP
1	B	294	GLN
1	B	330	PHE
1	B	490	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	370	HIS
1	A	420	ASN
1	A	422	GLN
1	A	440	GLN
1	B	50	ASN
1	B	54	ASN
1	B	294	GLN
1	B	420	ASN
1	B	422	GLN
1	B	440	GLN
1	B	503	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

4 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	GLC	C	1	2	11,11,12	1.73	2 (18%)	15,15,17	2.60	8 (53%)
2	FRU	C	2	2,3	11,12,12	1.71	2 (18%)	10,18,18	1.37	2 (20%)
2	GLC	D	1	2	11,11,12	1.84	2 (18%)	15,15,17	2.46	5 (33%)
2	FRU	D	2	2,3	11,12,12	1.84	3 (27%)	10,18,18	1.34	1 (10%)

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	2/2/4/5	2/2/19/22	0/1/1/1
2	FRU	C	2	2,3	-	3/5/24/24	0/1/1/1
2	GLC	D	1	2	2/2/4/5	1/2/19/22	0/1/1/1
2	FRU	D	2	2,3	-	3/5/24/24	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	GLC	O5-C5	4.43	1.52	1.43
2	C	1	GLC	O5-C5	4.30	1.52	1.43
2	D	2	FRU	O2-C2	3.85	1.47	1.40
2	C	2	FRU	O2-C2	3.76	1.47	1.40
2	D	2	FRU	C4-C3	-3.12	1.39	1.52
2	C	2	FRU	C4-C3	-3.03	1.40	1.52
2	D	1	GLC	C4-C3	-2.72	1.45	1.52
2	C	1	GLC	C4-C3	-2.69	1.45	1.52
2	D	2	FRU	O5-C2	-2.19	1.39	1.43

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	GLC	C1-C2-C3	4.52	115.22	109.67
2	C	1	GLC	C3-C4-C5	4.50	118.28	110.24
2	C	1	GLC	C1-C2-C3	4.27	114.91	109.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	GLC	C2-C3-C4	4.21	118.19	110.89
2	D	1	GLC	C1-O5-C5	4.08	117.72	112.19
2	D	1	GLC	C2-C3-C4	3.91	117.66	110.89
2	D	1	GLC	C3-C4-C5	3.78	116.99	110.24
2	C	1	GLC	C1-O5-C5	3.41	116.81	112.19
2	C	1	GLC	O4-C4-C5	3.38	117.68	109.30
2	D	2	FRU	O1-C1-C2	3.08	118.41	111.86
2	D	1	GLC	O2-C2-C3	2.50	115.14	110.14
2	C	2	FRU	O1-C1-C2	2.45	117.06	111.86
2	C	2	FRU	O2-C2-O5	2.43	114.20	109.50
2	C	1	GLC	C6-C5-C4	2.27	118.32	113.00
2	C	1	GLC	O6-C6-C5	2.23	118.95	111.29
2	C	1	GLC	O2-C2-C3	2.04	114.22	110.14

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1	GLC	C3
2	C	1	GLC	C4
2	D	1	GLC	C3
2	D	1	GLC	C4

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	2	FRU	O1-C1-C2-C3
2	D	2	FRU	O1-C1-C2-O2
2	D	2	FRU	O1-C1-C2-O5
2	C	1	GLC	C4-C5-C6-O6
2	C	2	FRU	O1-C1-C2-O2
2	D	1	GLC	C4-C5-C6-O6
2	C	2	FRU	O1-C1-C2-C3
2	C	1	GLC	O5-C5-C6-O6
2	C	2	FRU	O1-C1-C2-O5

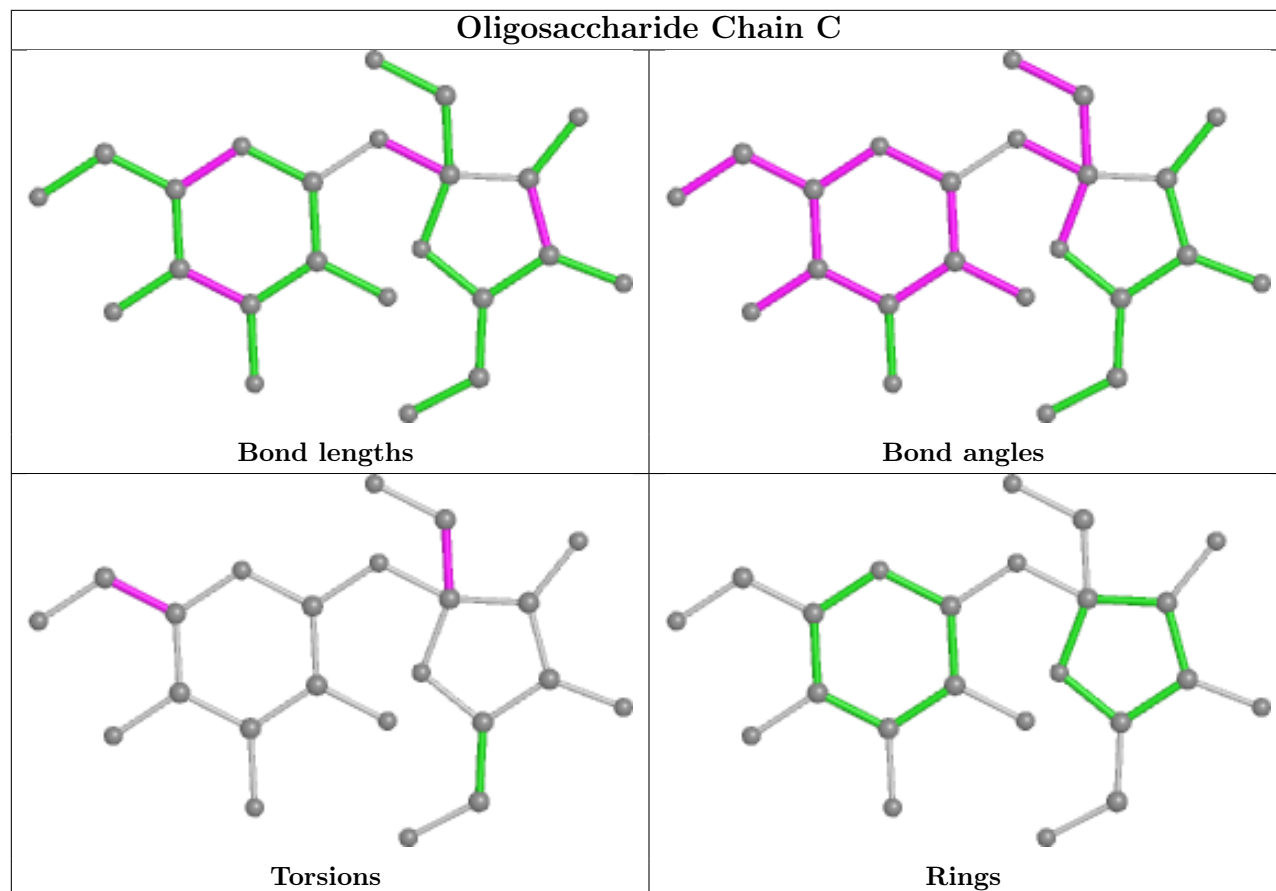
There are no ring outliers.

1 monomer is involved in 1 short contact:

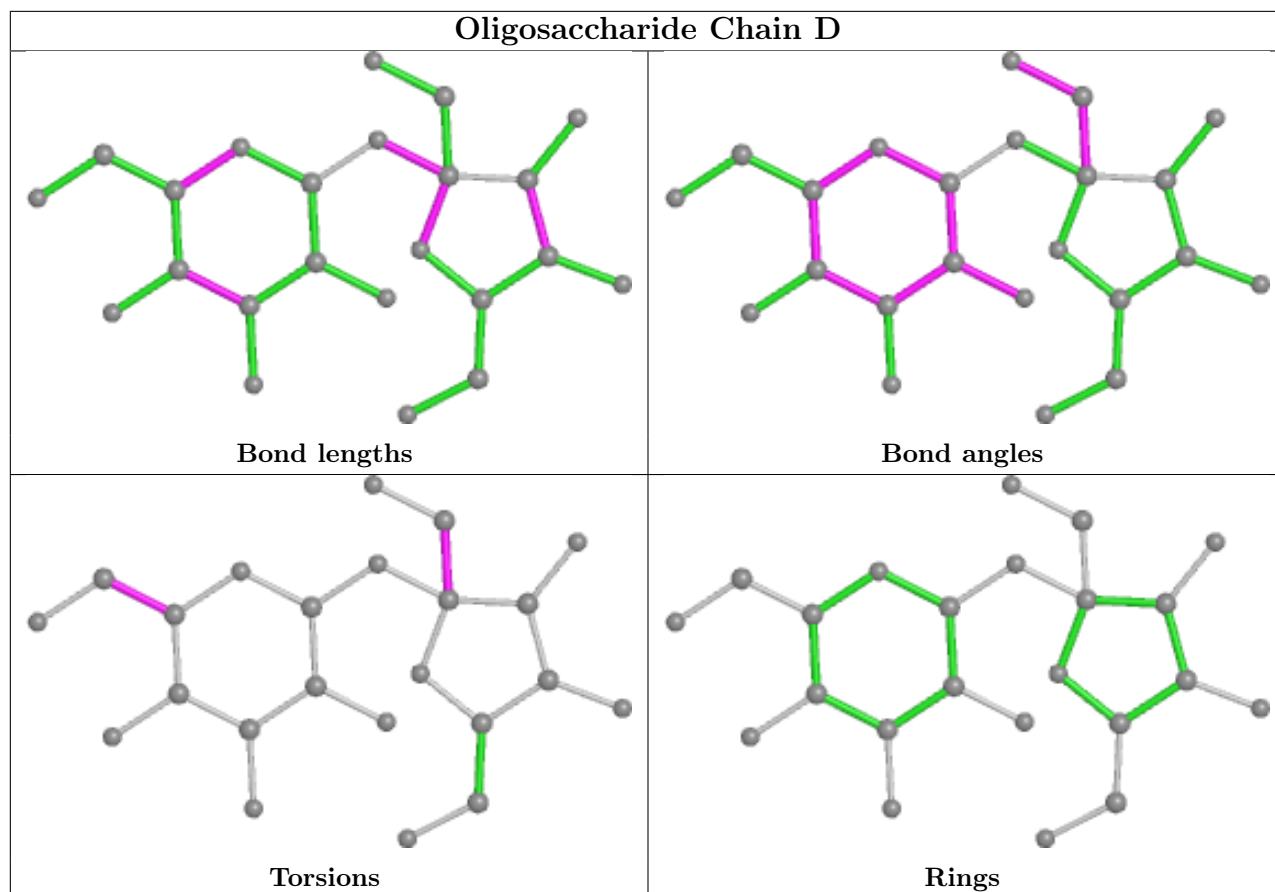
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	GLC	1	0

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](#)

4 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	$\# Z  > 2$
4	EDO	A	602	-	3,3,3	0.66	0	2,2,2	0.07	0
3	LAO	A	601	2	10,13,14	2.25	2 (20%)	12,18,21	1.36	2 (16%)
3	LAO	B	601	2	10,13,14	2.29	3 (30%)	12,18,21	1.70	4 (33%)
5	PEG	B	602	-	6,6,6	0.23	0	5,5,5	0.11	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
4	EDO	A	602	-	-	0/1/1/1	-
3	LAO	A	601	2	-	0/1/21/22	0/1/1/1
3	LAO	B	601	2	-	0/1/21/22	0/1/1/1
5	PEG	B	602	-	-	2/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	LAO	C4-C3	-6.21	1.43	1.52
3	B	601	LAO	C4-C3	-5.99	1.43	1.52
3	B	601	LAO	C3-C2	-2.93	1.44	1.52
3	A	601	LAO	C3-C2	-2.69	1.45	1.52
3	B	601	LAO	O3-C3	2.36	1.48	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	601	LAO	O5-C1-C2	3.26	114.23	109.51
3	A	601	LAO	C4-C3-C2	2.92	116.28	110.41
3	A	601	LAO	O5-C1-C2	2.44	113.04	109.51
3	B	601	LAO	O5-C5-C4	2.29	114.31	110.77
3	B	601	LAO	O4-C4-C5	-2.27	104.50	109.15
3	B	601	LAO	C4-C3-C2	2.22	114.87	110.41

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	602	PEG	C4-C3-O2-C2
5	B	602	PEG	C1-C2-O2-C3

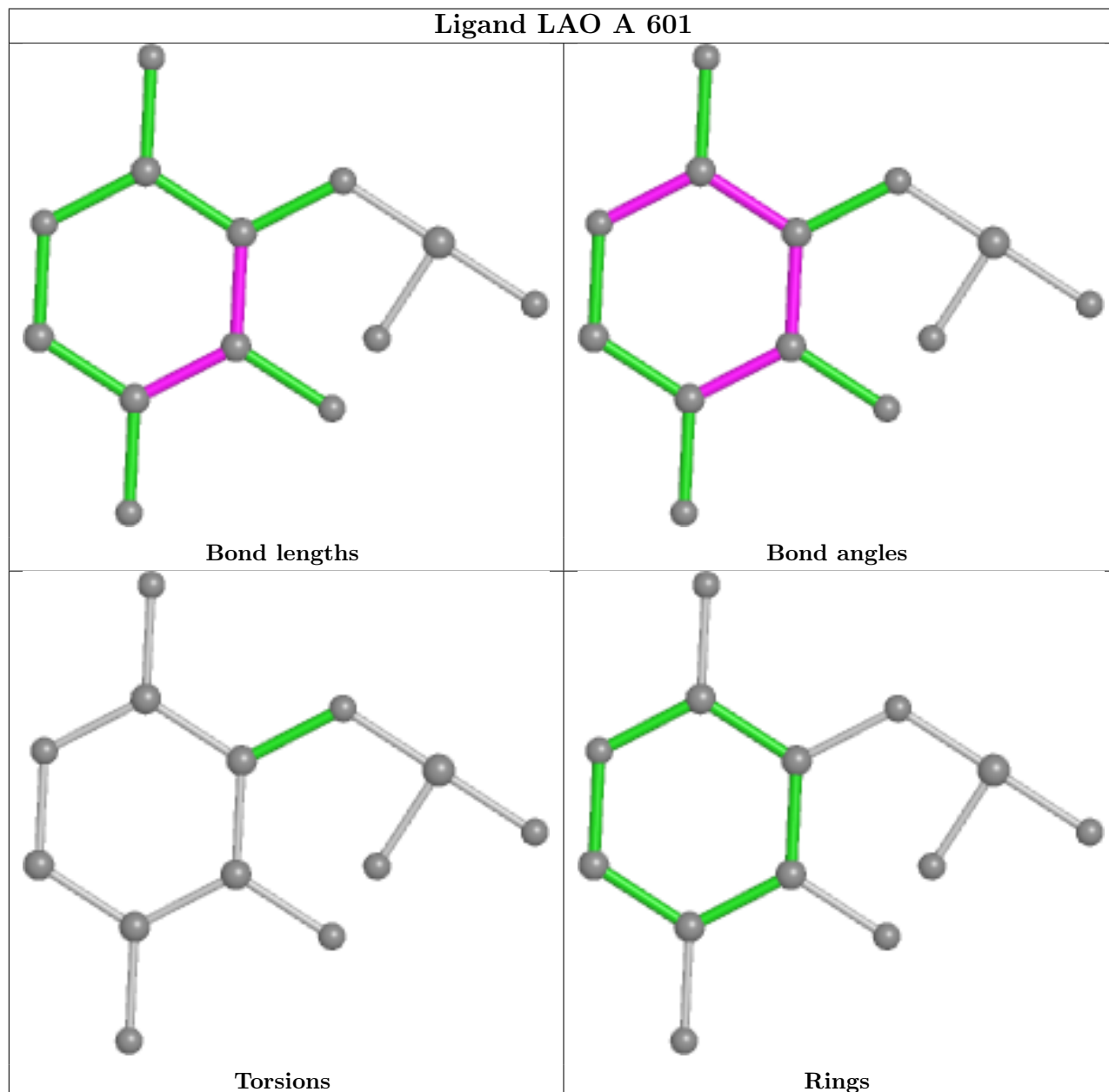
There are no ring outliers.

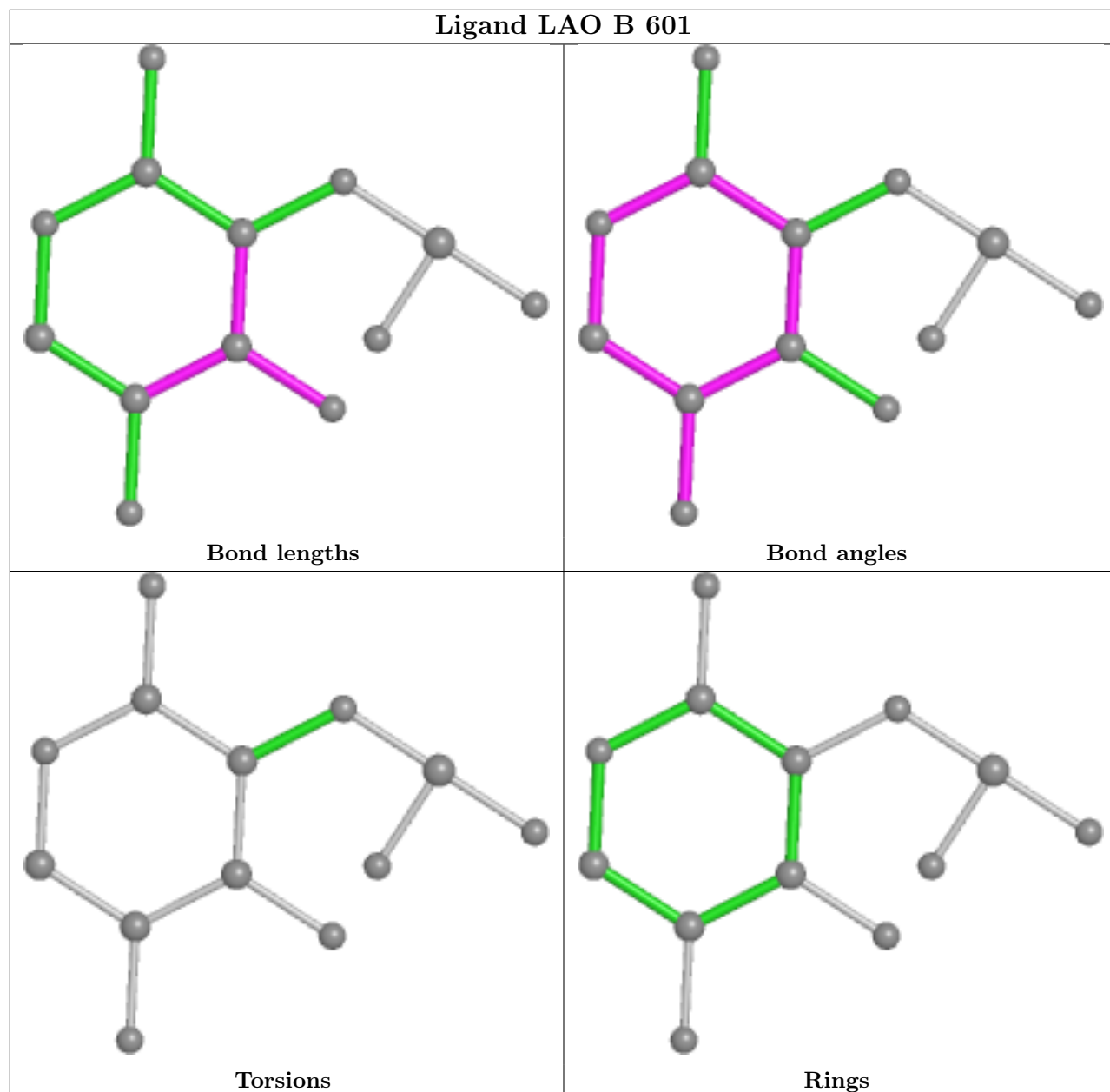
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	602	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	490/499 (98%)	-0.40	3 (0%) 89 90	21, 33, 56, 90	0
1	B	489/499 (97%)	-0.39	4 (0%) 86 87	21, 34, 58, 92	0
All	All	979/998 (98%)	-0.39	7 (0%) 87 88	21, 34, 57, 92	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	251	ILE	3.5
1	B	521	GLY	2.8
1	B	273	PRO	2.6
1	B	223	ASP	2.5
1	A	268	LEU	2.2
1	B	501	ASP	2.2
1	A	520	PHE	2.1

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

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

### 6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> 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.

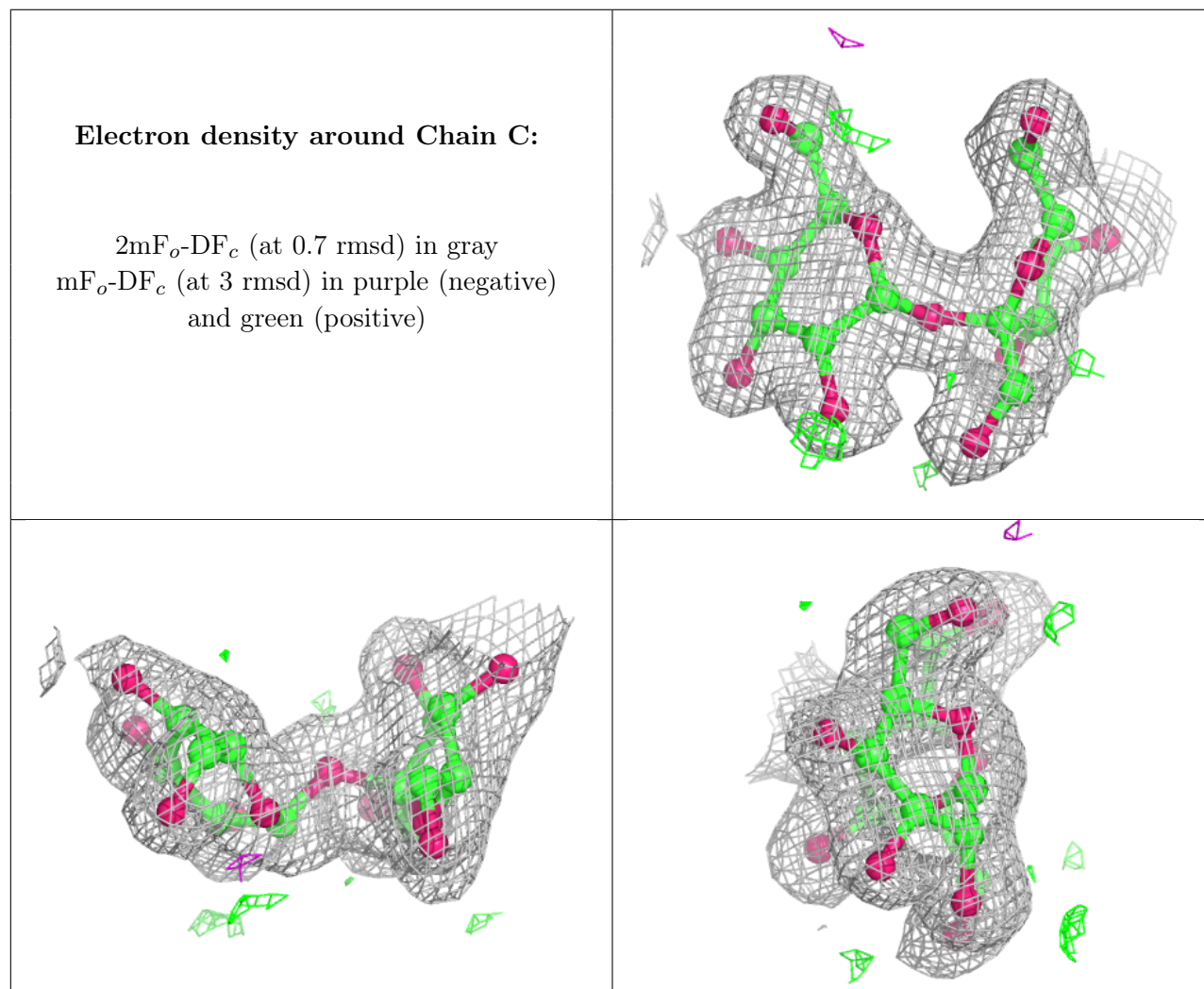
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	GLC	D	1	11/12	0.88	0.13	42,45,47,51	0
2	FRU	C	2	12/12	0.95	0.07	24,35,38,40	0

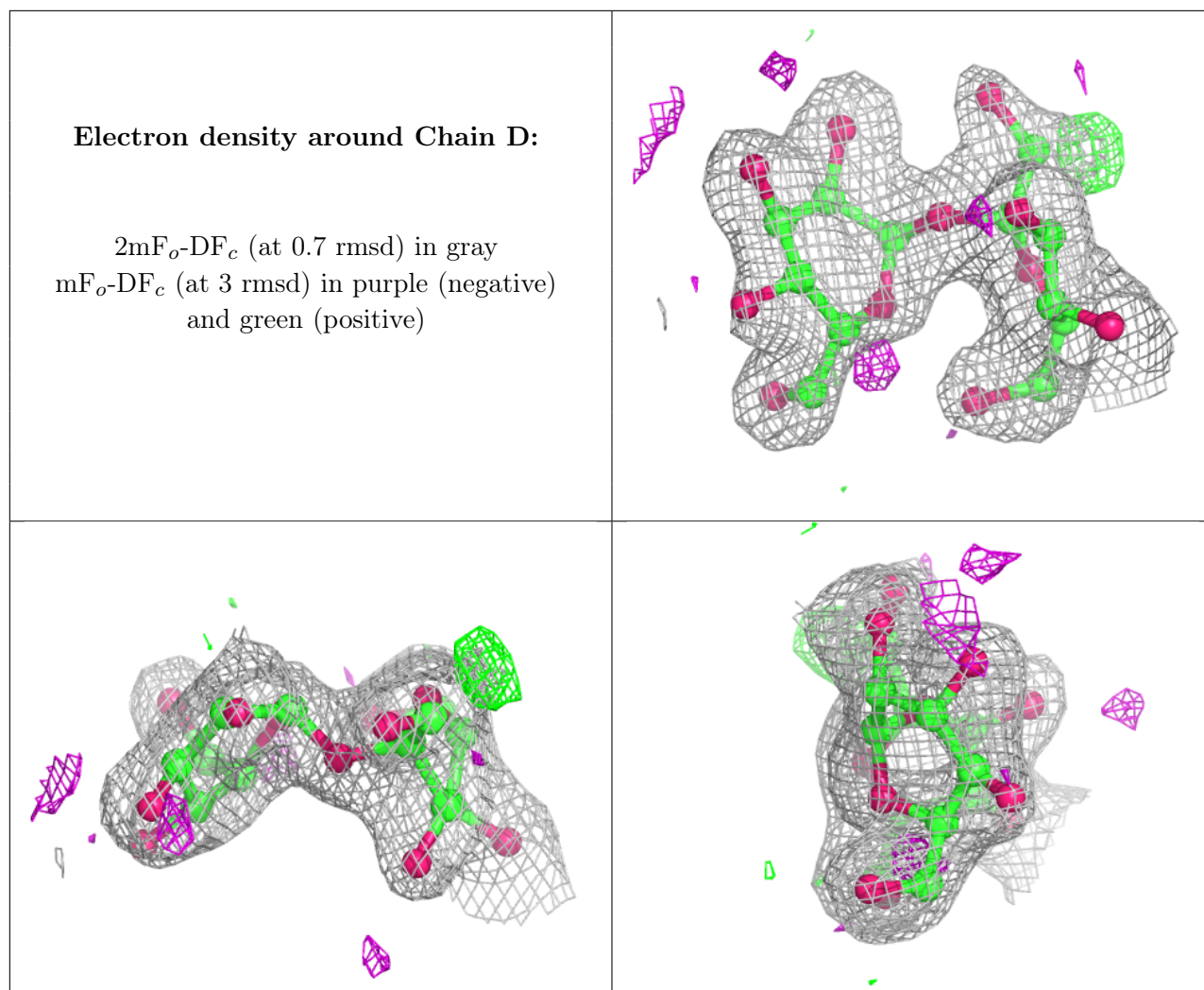
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GLC	C	1	11/12	0.95	0.08	38,41,45,46	0
2	FRU	D	2	12/12	0.97	0.09	26,39,45,48	0

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.



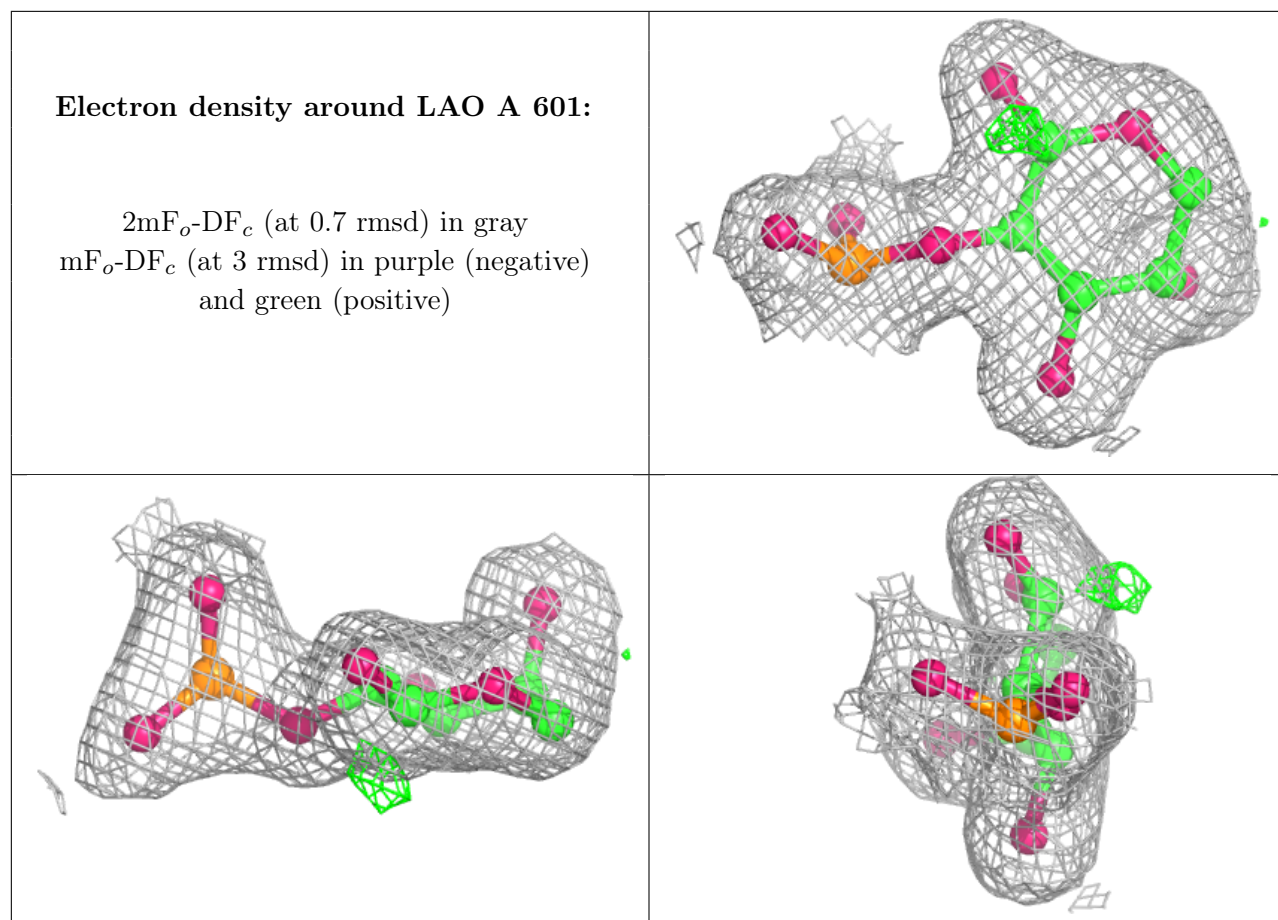


## 6.4 Ligands [i](#)

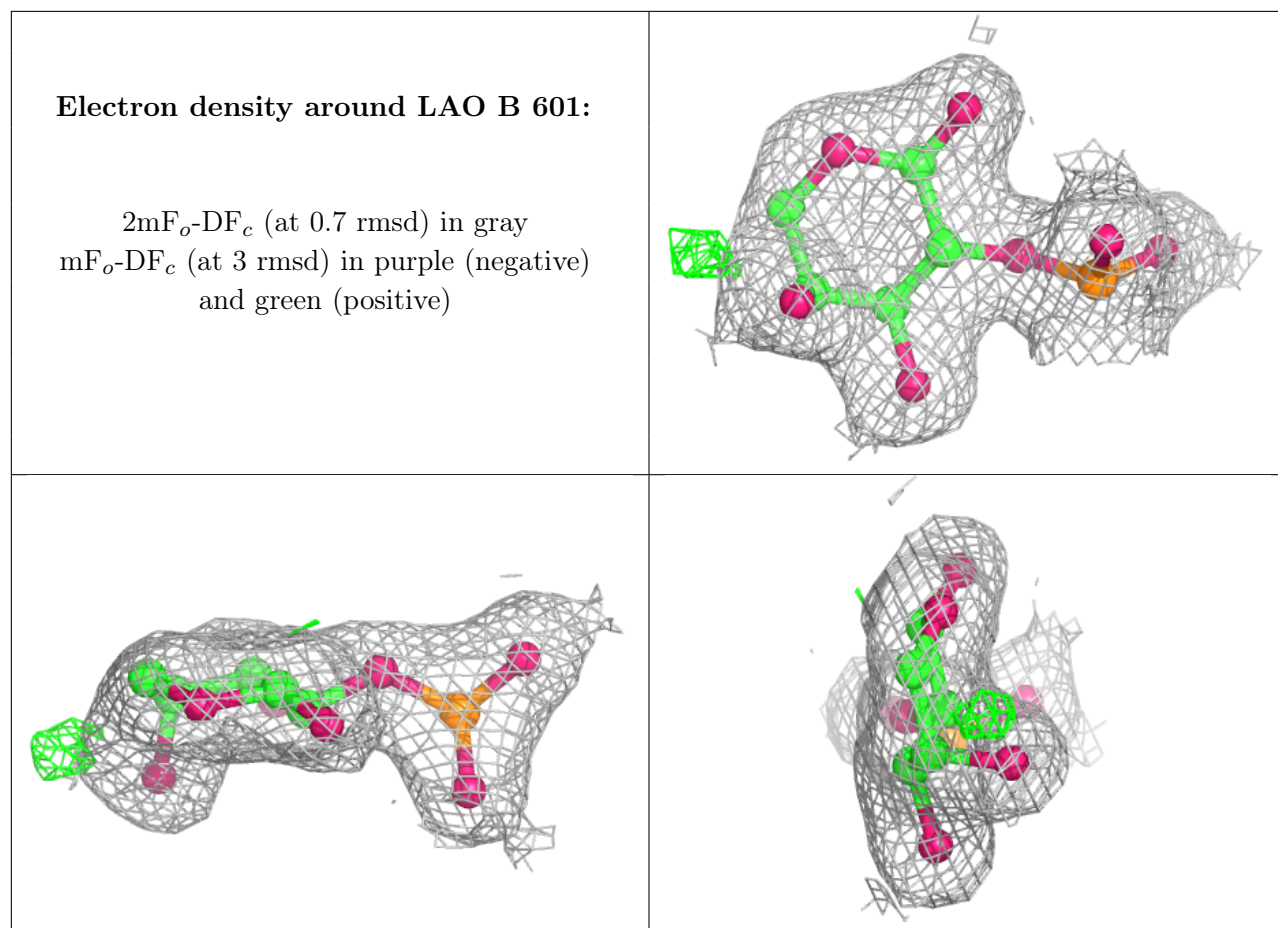
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	EDO	A	602	4/4	0.84	0.12	63,63,63,64	0
5	PEG	B	602	7/7	0.86	0.14	50,52,56,58	0
3	LAO	A	601	13/14	0.98	0.09	23,26,28,28	0
3	LAO	B	601	13/14	0.98	0.07	22,28,29,30	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.







## 6.5 Other polymers [i](#)

There are no such residues in this entry.