



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 18, 2021 – 12:13 PM JST

PDB ID : 7DDH  
Title : Crystal structures of Na<sup>+</sup>,K<sup>+</sup>-ATPase in complex with digoxin  
Authors : Ogawa, H.; Cornelius, F.; Kanai, R.; Motoyama, K.; Vilsen, B.; Toyoshima, C.  
Deposited on : 2020-10-29  
Resolution : 3.46 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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.

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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	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.16
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)	:	2.16

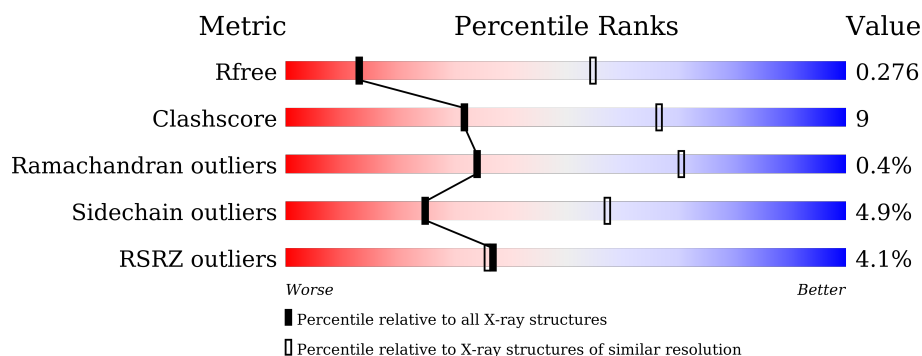
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.46 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	1291 (3.52-3.40)
Clashscore	141614	1372 (3.52-3.40)
Ramachandran outliers	138981	1337 (3.52-3.40)
Sidechain outliers	138945	1338 (3.52-3.40)
RSRZ outliers	127900	1205 (3.52-3.40)

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	1016	<div> <div>4%</div> <div> <div></div> <div>75%</div> <div>21%</div> <div>..</div> </div> </div>
1	C	1016	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>21%</div> <div>..</div> </div> </div>
2	B	303	<div> <div>7%</div> <div> <div></div> <div>69%</div> <div>25%</div> <div>..</div> </div> </div>
2	D	303	<div> <div>4%</div> <div> <div></div> <div>64%</div> <div>28%</div> <div>6%</div> </div> </div>
3	E	65	<div> <div></div> <div> <div>40%</div> <div>9%</div> <div>51%</div> </div> </div>
3	G	65	<div> <div></div> <div> <div>48%</div> <div></div> <div>51%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
4	F	2	 100%
4	H	2	 50% 50%
4	I	2	 50% 50%
4	J	2	 50% 50%

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
4	NAG	H	2	-	-	-	X
8	PCW	A	1106	-	-	-	X
8	PCW	A	1107	-	-	-	X
8	PCW	A	1108	-	-	-	X
8	PCW	C	1105	-	-	-	X

## 2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 21344 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 Sodium/potassium-transporting ATPase subunit alpha-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	996	Total	C	N	O	P	S	0	0	0
			7730	4922	1301	1459	1	47			
1	C	996	Total	C	N	O	P	S	0	0	0
			7730	4922	1301	1459	1	47			

- Molecule 2 is a protein called Sodium/potassium-transporting ATPase subunit beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	291	Total	C	N	O	S	0	0	0
			2386	1546	390	437	13			
2	D	285	Total	C	N	O	S	0	0	0
			2334	1514	383	424	13			

- Molecule 3 is a protein called FXYD domain-containing ion transport regulator.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	32	Total	C	N	O	0	0	0
			255	174	37	44			
3	E	32	Total	C	N	O	0	0	0
			255	174	37	44			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	J	2	Total	C	N	O	0	0	0
			28	16	2	10			

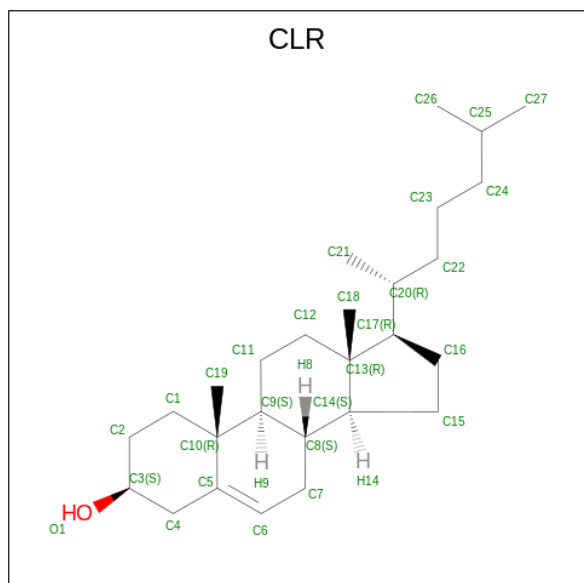
- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Mg	0	0
			2	2		
5	C	2	Total	Mg	0	0
			2	2		

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

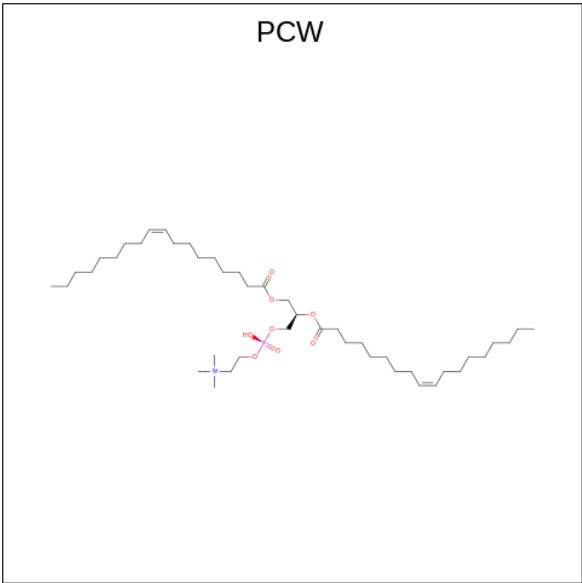
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		
6	C	1	Total	Na	0	0
			1	1		

- Molecule 7 is CHOLESTEROL (three-letter code: CLR) (formula: C<sub>27</sub>H<sub>46</sub>O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			28	27	1		
7	B	1	Total	C	O	0	0
			28	27	1		
7	G	1	Total	C	O	0	0
			28	27	1		
7	C	1	Total	C	O	0	0
			28	27	1		
7	D	1	Total	C	O	0	0
			28	27	1		
7	E	1	Total	C	O	0	0
			28	27	1		

- Molecule 8 is 1,2-DIOLEOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: PCW) (formula: C<sub>44</sub>H<sub>85</sub>NO<sub>8</sub>P).



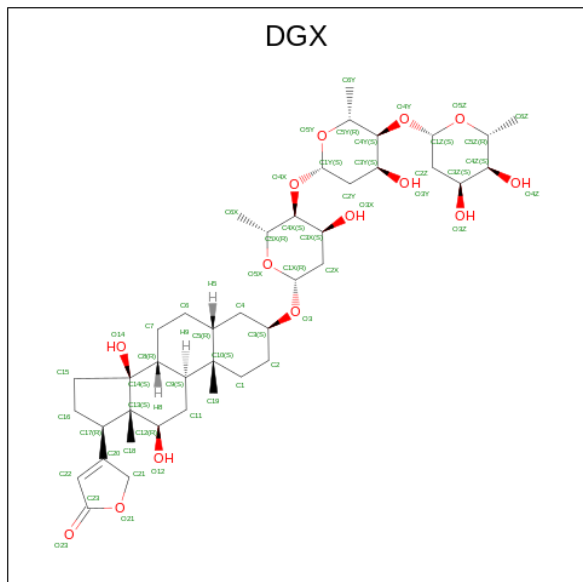
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	N	O	P	0	0
			22	12	1	8	1		
8	A	1	Total	C	N	O	P	0	0
			22	12	1	8	1		
8	A	1	Total	C	N	O	P	0	0
			22	12	1	8	1		
8	A	1	Total	C	N	O	P	0	0
			22	12	1	8	1		
8	A	1	Total	C	N	O	P	0	0
			22	12	1	8	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	B	1	Total	C	N	O	P	0	0
			22	12	1	8	1		
8	C	1	Total	C	N	O	P	0	0
			22	12	1	8	1		
8	C	1	Total	C	N	O	P	0	0
			22	12	1	8	1		
8	C	1	Total	C	N	O	P	0	0
			22	12	1	8	1		
8	C	1	Total	C	N	O	P	0	0
			22	12	1	8	1		

- Molecule 9 is DIGOXIN (three-letter code: DGX) (formula:  $C_{41}H_{64}O_{14}$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			55	41	14		
9	C	1	Total	C	O	0	0
			55	41	14		

- Molecule 10 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	B	1	Total	C	N	O	0	0
			14	8	1	5		
10	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 11 is water.

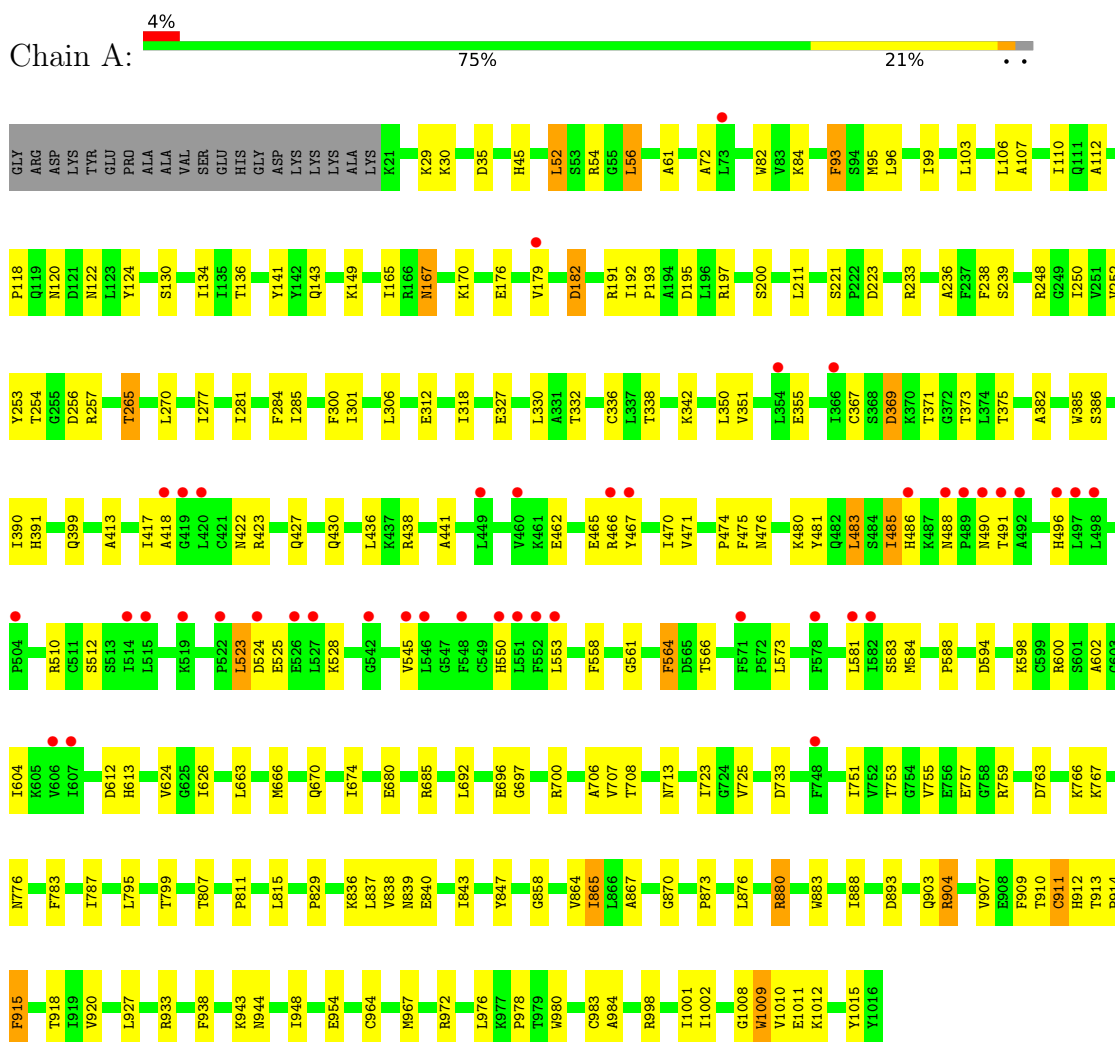
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	5	Total	O	0	0
			5	5		
11	C	5	Total	O	0	0
			5	5		



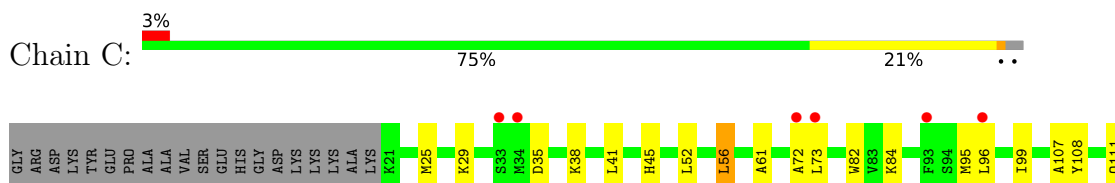
### 3 Residue-property plots

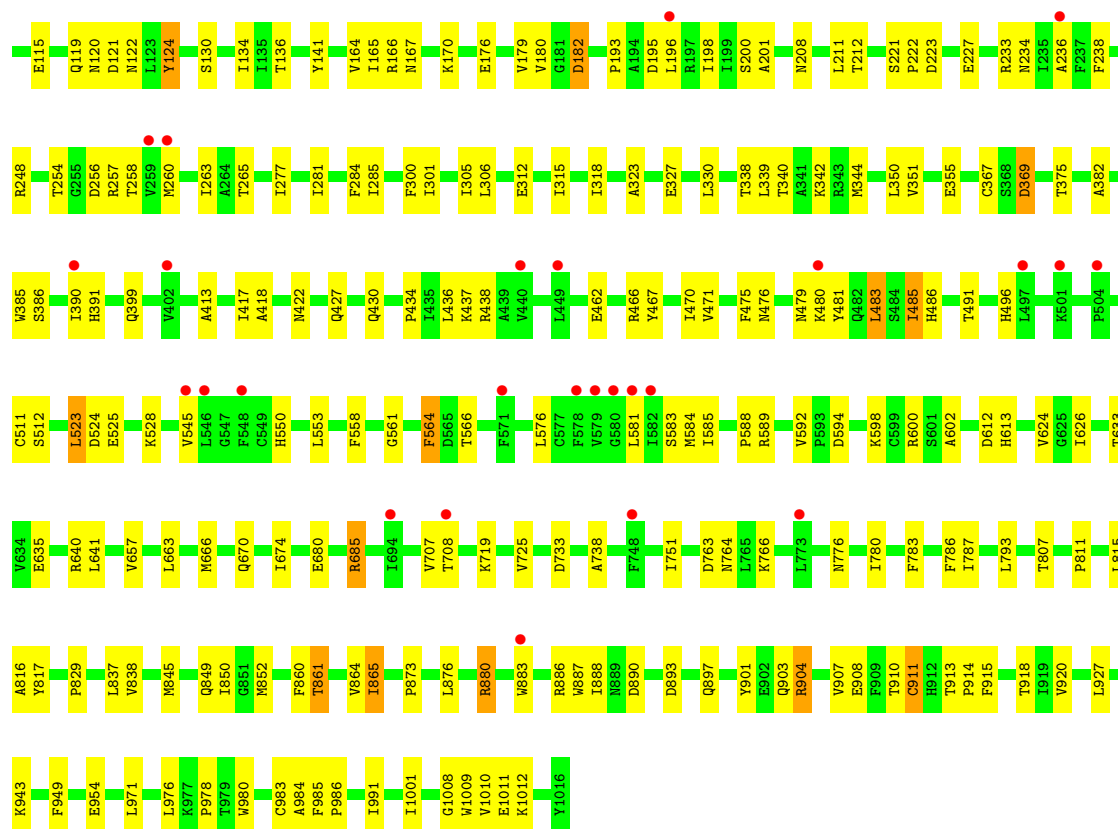
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Sodium/potassium-transporting ATPase subunit alpha-1

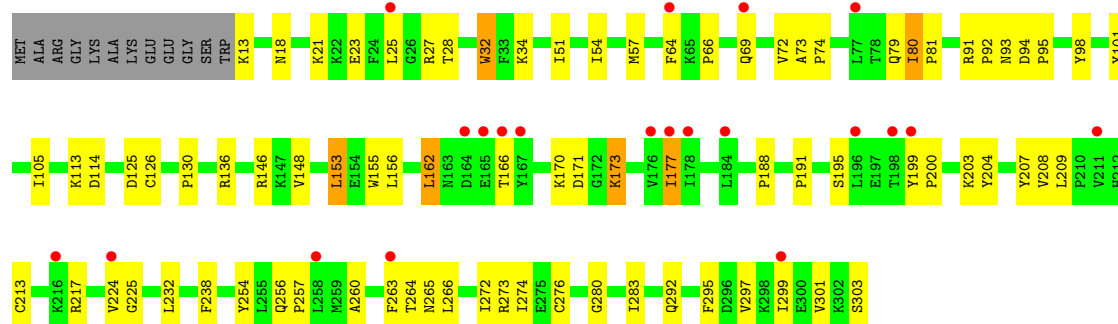


- Molecule 1: Sodium/potassium-transporting ATPase subunit alpha-1

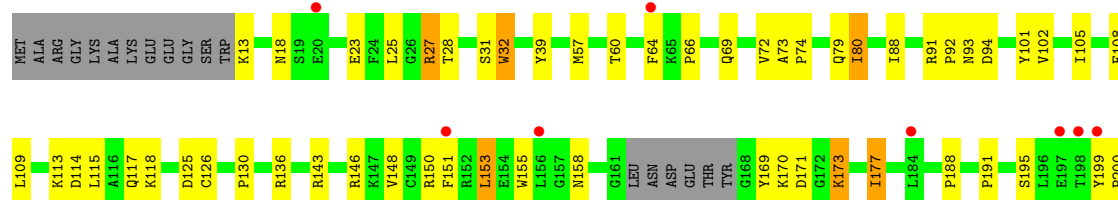


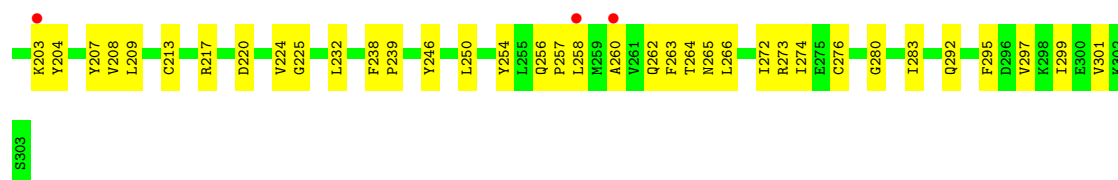


• Molecule 2: Sodium/potassium-transporting ATPase subunit beta-1



• Molecule 2: Sodium/potassium-transporting ATPase subunit beta-1

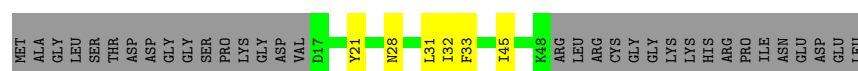




- Molecule 3: FXYD domain-containing ion transport regulator



- Molecule 3: FXYD domain-containing ion transport regulator



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	114.97Å 118.49Å 495.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.99 – 3.46 49.72 – 3.46	Depositor EDS
% Data completeness (in resolution range)	55.1 (15.99-3.46) 55.5 (49.72-3.46)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.70 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, $R_{free}$	0.241 , 0.272 0.252 , 0.276	Depositor DCC
$R_{free}$ test set	2477 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	108.1	Xtriage
Anisotropy	0.321	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 48.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.38$ , $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.068 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	21344	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	119.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, NAG, NA, PCW, PHD, DGX, CLR

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	0.29	0/7867	0.50	0/10674
1	C	0.29	0/7867	0.51	0/10674
2	B	0.29	0/2449	0.54	0/3301
2	D	0.30	0/2395	0.54	0/3225
3	E	0.32	0/261	0.51	0/354
3	G	0.33	0/261	0.47	0/354
All	All	0.29	0/21100	0.51	0/28582

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 [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	7730	0	7777	137	0
1	C	7730	0	7777	135	0
2	B	2386	0	2361	49	0
2	D	2334	0	2317	55	0
3	E	255	0	259	6	0
3	G	255	0	259	0	0
4	F	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	28	0	25	0	0
4	I	28	0	25	1	0
4	J	28	0	25	0	0
5	A	2	0	0	0	0
5	C	2	0	0	0	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
7	A	28	0	46	2	0
7	B	28	0	46	0	0
7	C	28	0	46	3	0
7	D	28	0	46	1	0
7	E	28	0	46	6	0
7	G	28	0	46	1	0
8	A	110	0	90	9	0
8	B	22	0	18	2	0
8	C	88	0	72	4	0
9	A	55	0	64	1	0
9	C	55	0	64	2	0
10	B	14	0	13	0	0
10	D	14	0	13	0	0
11	A	5	0	0	0	0
11	C	5	0	0	0	0
All	All	21344	0	21460	380	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 9.

The worst 5 of 380 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:LYS:HG3	1:A:141:TYR:HE1	1.41	0.85
1:C:978:PRO:HB3	7:E:101:CLR:H192	1.62	0.82
1:C:864:VAL:HG22	2:D:57:MET:HG3	1.63	0.81
1:C:430:GLN:HG3	1:C:438:ARG:HB2	1.65	0.78
2:B:80:ILE:HG12	2:B:177:ILE:HG12	1.66	0.78

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	993/1016 (98%)	907 (91%)	83 (8%)	3 (0%)	41	75
1	C	993/1016 (98%)	906 (91%)	84 (8%)	3 (0%)	41	75
2	B	289/303 (95%)	260 (90%)	27 (9%)	2 (1%)	22	60
2	D	281/303 (93%)	252 (90%)	27 (10%)	2 (1%)	22	60
3	E	30/65 (46%)	28 (93%)	2 (7%)	0	100	100
3	G	30/65 (46%)	26 (87%)	4 (13%)	0	100	100
All	All	2616/2768 (94%)	2379 (91%)	227 (9%)	10 (0%)	34	70

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	200	PRO
1	A	193	PRO
1	C	193	PRO
1	C	306	LEU
2	D	200	PRO

### 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	846/861 (98%)	813 (96%)	33 (4%)	32	64
1	C	846/861 (98%)	812 (96%)	34 (4%)	31	63
2	B	261/269 (97%)	240 (92%)	21 (8%)	12	40

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	255/269 (95%)	234 (92%)	21 (8%)	11	39
3	E	26/52 (50%)	25 (96%)	1 (4%)	33	64
3	G	26/52 (50%)	25 (96%)	1 (4%)	33	64
All	All	2260/2364 (96%)	2149 (95%)	111 (5%)	25	57

5 of 111 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	203	LYS
1	C	180	VAL
2	D	170	LYS
2	B	256	GLN
1	C	56	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	935	ASN
1	C	111	GLN
1	C	482	GLN
1	A	488	ASN
1	A	897	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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
1	PHD	A	369	1,5	9,11,12	0.93	0	10,15,17	1.47	2 (20%)
1	PHD	C	369	1,5	9,11,12	0.92	0	10,15,17	1.39	2 (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
1	PHD	A	369	1,5	-	0/8/11/13	-
1	PHD	C	369	1,5	-	0/8/11/13	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	369	PHD	CA-CB-CG	2.91	118.95	112.86
1	A	369	PHD	OD1-CG-CB	2.78	118.76	111.11
1	A	369	PHD	CA-CB-CG	2.71	118.54	112.86
1	C	369	PHD	OD1-CG-CB	2.38	117.66	111.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	369	PHD	2	0
1	C	369	PHD	1	0

## 5.5 Carbohydrates [i](#)

8 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
4	NAG	F	1	2,4	14,14,15	0.35	0	17,19,21	0.43	0
4	NAG	F	2	4	14,14,15	0.28	0	17,19,21	0.49	0
4	NAG	H	1	2,4	14,14,15	0.62	1 (7%)	17,19,21	0.70	0
4	NAG	H	2	4	14,14,15	0.32	0	17,19,21	0.36	0
4	NAG	I	1	2,4	14,14,15	0.39	0	17,19,21	0.46	0
4	NAG	I	2	4	14,14,15	0.31	0	17,19,21	0.48	0
4	NAG	J	1	2,4	14,14,15	0.61	1 (7%)	17,19,21	0.63	0
4	NAG	J	2	4	14,14,15	0.31	0	17,19,21	0.34	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	NAG	F	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1
4	NAG	H	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	H	2	4	-	0/6/23/26	0/1/1/1
4	NAG	I	1	2,4	-	2/6/23/26	0/1/1/1
4	NAG	I	2	4	-	0/6/23/26	0/1/1/1
4	NAG	J	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	1	NAG	O5-C1	-2.18	1.40	1.43
4	J	1	NAG	O5-C1	-2.09	1.40	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

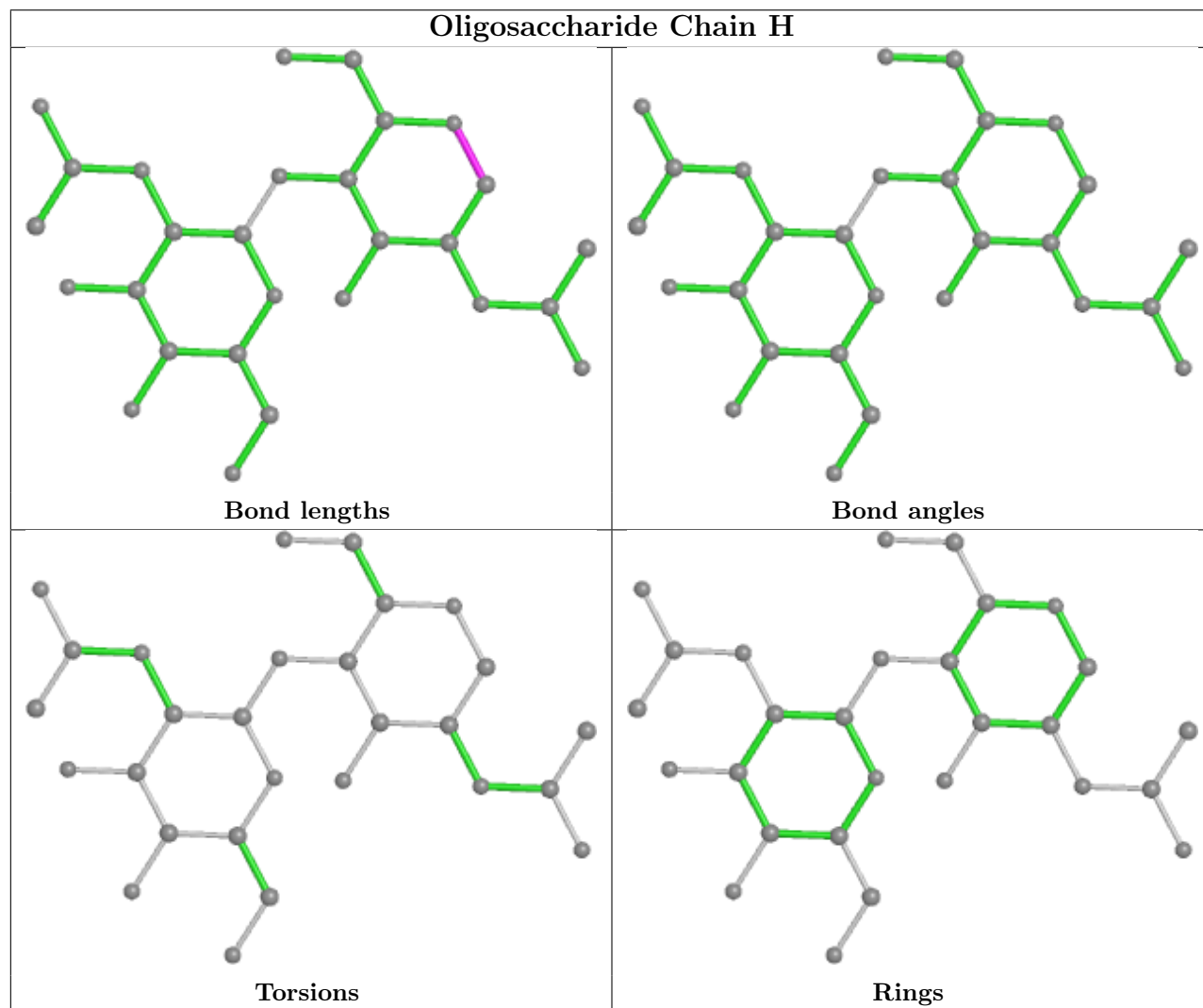
Mol	Chain	Res	Type	Atoms
4	I	1	NAG	C4-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6

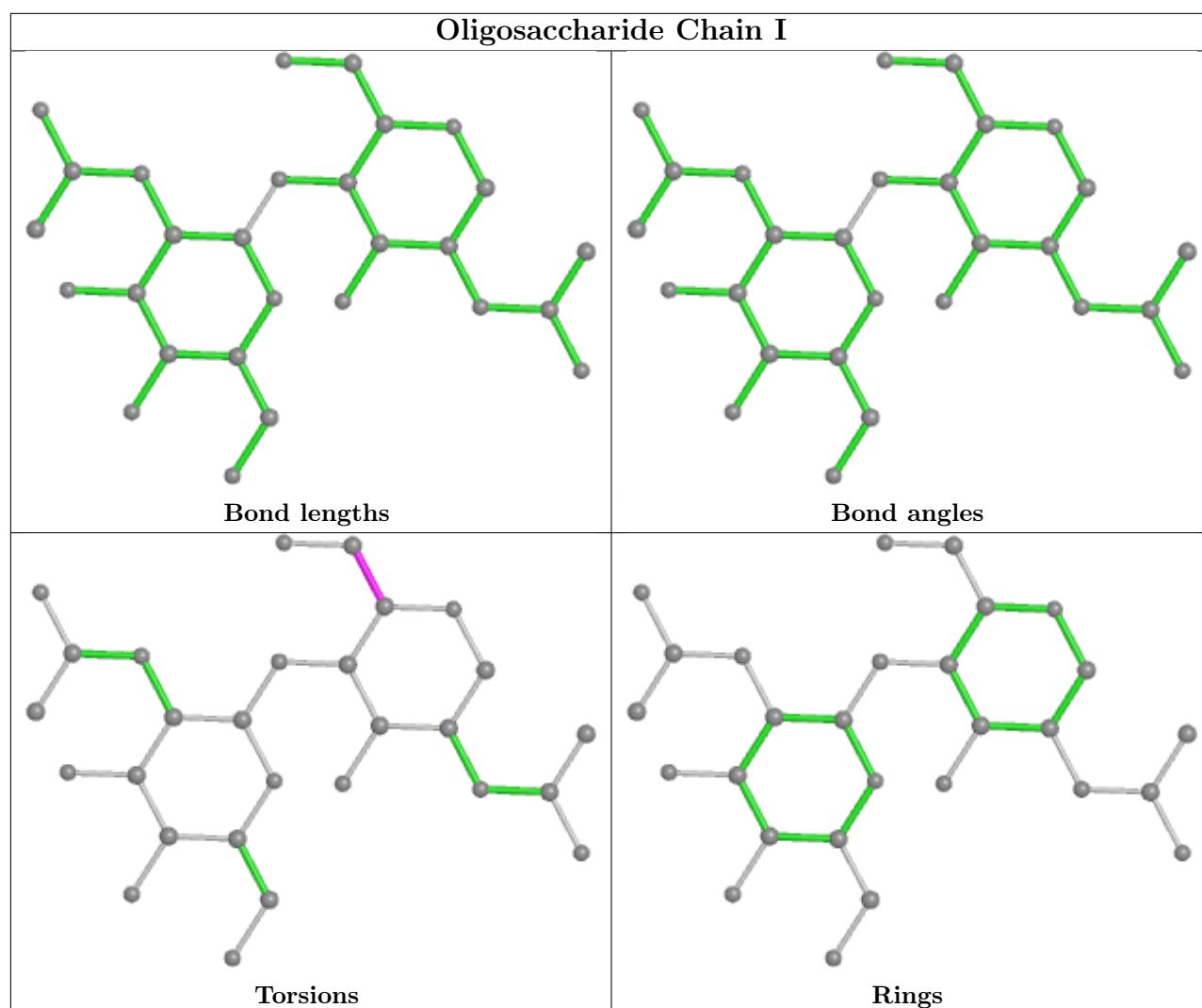
There are no ring outliers.

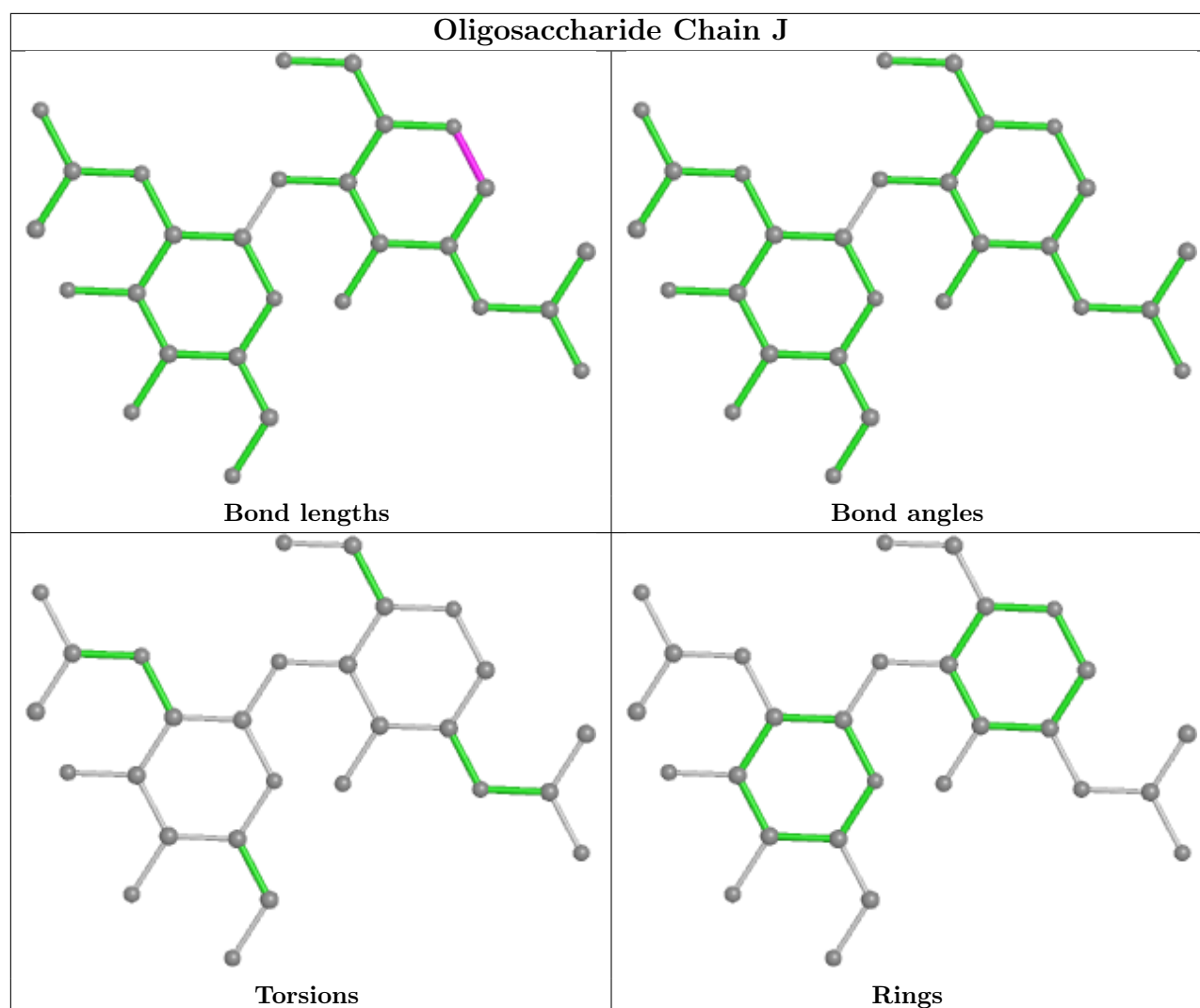
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	I	1	NAG	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](#)

Of 26 ligands modelled in this entry, 6 are monoatomic - leaving 20 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	$\# Z  > 2$
8	PCW	C	1106	-	21,21,53	1.68	6 (28%)	27,29,61	1.23	1 (3%)
9	DGX	A	1110	-	62,62,62	1.21	3 (4%)	89,98,98	1.29	7 (7%)
8	PCW	C	1107	-	21,21,53	1.69	5 (23%)	27,29,61	1.45	3 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
7	CLR	C	1104	-	31,31,31	1.92	9 (29%)	48,48,48	1.61	16 (33%)
8	PCW	A	1107	-	21,21,53	1.68	6 (28%)	27,29,61	1.16	1 (3%)
7	CLR	B	402	-	31,31,31	1.96	12 (38%)	48,48,48	1.63	13 (27%)
8	PCW	B	403	-	21,21,53	1.69	5 (23%)	27,29,61	1.23	1 (3%)
9	DGX	C	1121	-	62,62,62	1.18	2 (3%)	89,98,98	1.24	8 (8%)
7	CLR	D	402	-	31,31,31	1.99	10 (32%)	48,48,48	1.55	12 (25%)
8	PCW	A	1105	-	21,21,53	1.71	5 (23%)	27,29,61	1.33	1 (3%)
10	NAG	D	401	2	14,14,15	0.31	0	17,19,21	0.53	0
7	CLR	A	1104	-	31,31,31	1.91	11 (35%)	48,48,48	1.57	10 (20%)
7	CLR	E	101	-	31,31,31	1.74	9 (29%)	48,48,48	1.57	11 (22%)
8	PCW	C	1105	-	21,21,53	1.72	4 (19%)	27,29,61	1.15	1 (3%)
8	PCW	C	1108	-	21,21,53	1.68	6 (28%)	27,29,61	1.20	1 (3%)
8	PCW	A	1109	-	21,21,53	1.69	5 (23%)	27,29,61	1.35	2 (7%)
8	PCW	A	1108	-	21,21,53	1.68	5 (23%)	27,29,61	1.20	1 (3%)
10	NAG	B	401	2	14,14,15	0.33	0	17,19,21	0.43	0
8	PCW	A	1106	-	21,21,53	1.67	4 (19%)	27,29,61	1.17	1 (3%)
7	CLR	G	101	-	31,31,31	1.77	8 (25%)	48,48,48	1.62	12 (25%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PCW	C	1106	-	-	9/23/23/57	-
9	DGX	A	1110	-	-	4/16/141/141	0/8/8/8
8	PCW	C	1107	-	-	10/23/23/57	-
7	CLR	C	1104	-	-	6/10/68/68	0/4/4/4
8	PCW	A	1107	-	-	11/23/23/57	-
7	CLR	B	402	-	-	1/10/68/68	0/4/4/4
8	PCW	B	403	-	-	11/23/23/57	-
9	DGX	C	1121	-	-	4/16/141/141	0/8/8/8
7	CLR	D	402	-	-	1/10/68/68	0/4/4/4
8	PCW	A	1105	-	-	12/23/23/57	-
10	NAG	D	401	2	-	4/6/23/26	0/1/1/1
7	CLR	A	1104	-	-	6/10/68/68	0/4/4/4
7	CLR	E	101	-	-	3/10/68/68	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PCW	C	1105	-	-	11/23/23/57	-
8	PCW	C	1108	-	-	13/23/23/57	-
8	PCW	A	1109	-	-	14/23/23/57	-
8	PCW	A	1108	-	-	10/23/23/57	-
10	NAG	B	401	2	-	1/6/23/26	0/1/1/1
8	PCW	A	1106	-	-	16/23/23/57	-
7	CLR	G	101	-	-	3/10/68/68	0/4/4/4

The worst 5 of 115 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	402	CLR	C10-C9	5.14	1.64	1.56
7	A	1104	CLR	C10-C9	4.90	1.64	1.56
7	D	402	CLR	C10-C9	4.77	1.64	1.56
7	C	1104	CLR	C10-C9	4.73	1.64	1.56
7	G	101	CLR	C10-C9	4.17	1.63	1.56

The worst 5 of 102 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	1110	DGX	O21-C23-C22	5.45	115.79	108.67
9	C	1121	DGX	O21-C23-C22	5.32	115.63	108.67
8	A	1105	PCW	O2-C31-C32	5.21	120.67	111.09
8	C	1107	PCW	O2-C31-C32	5.12	120.50	111.09
8	A	1109	PCW	O2-C31-C32	4.98	120.24	111.09

There are no chirality outliers.

5 of 150 torsion outliers are listed below:

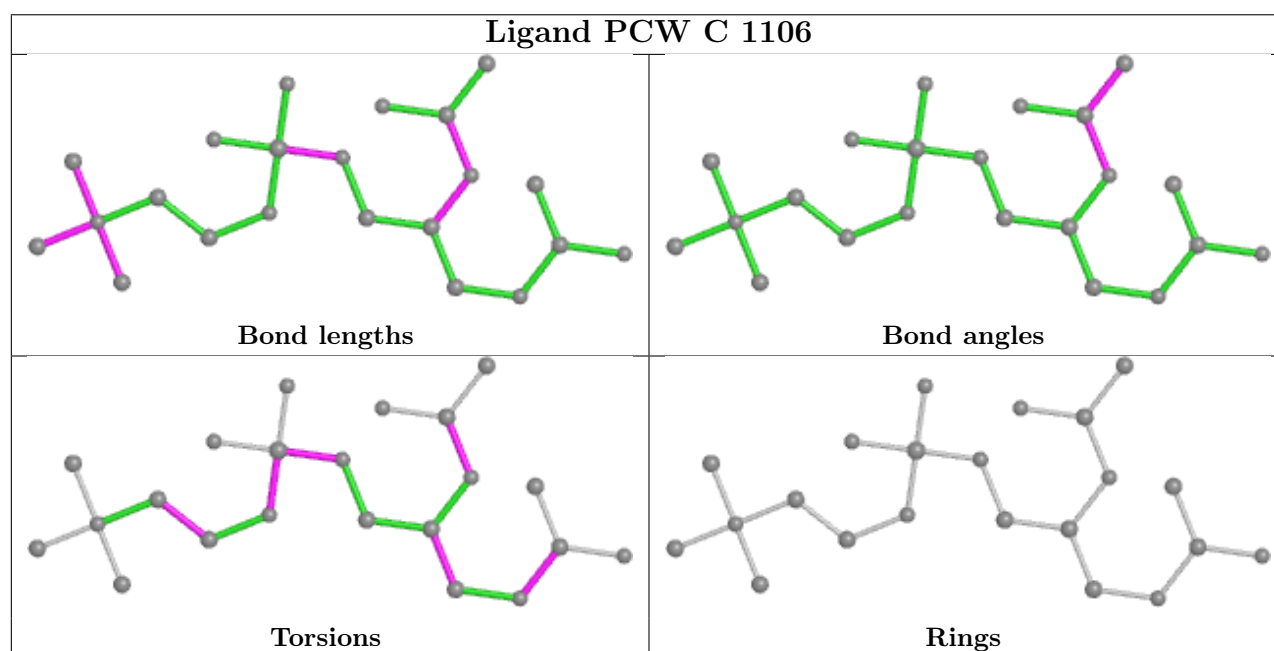
Mol	Chain	Res	Type	Atoms
8	C	1107	PCW	C4-O4P-P-O2P
8	A	1107	PCW	C1-O3P-P-O2P
8	A	1105	PCW	C4-O4P-P-O1P
8	C	1105	PCW	C1-O3P-P-O1P
8	C	1105	PCW	C1-O3P-P-O2P

There are no ring outliers.

14 monomers are involved in 31 short contacts:

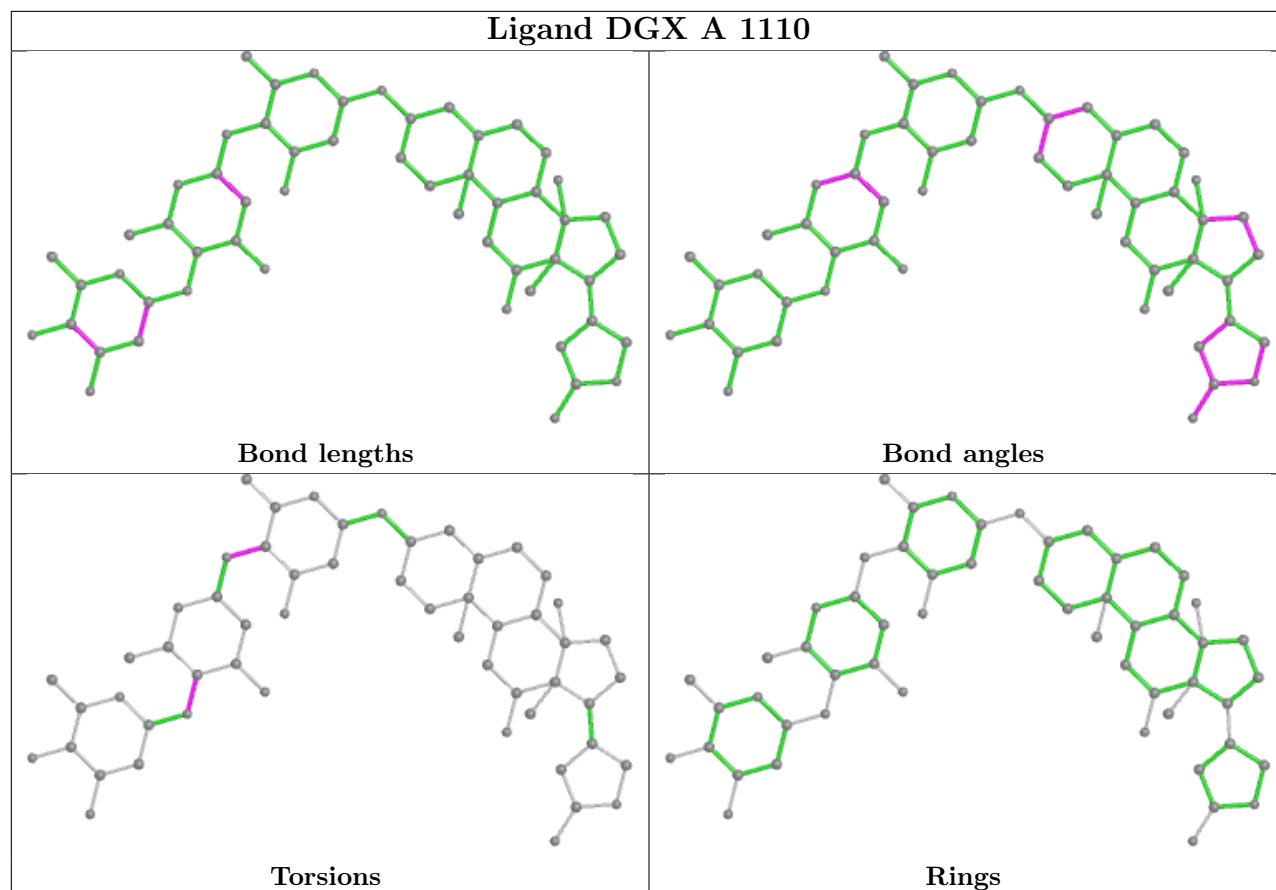
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	1110	DGX	1	0
8	C	1107	PCW	2	0
7	C	1104	CLR	3	0
8	B	403	PCW	2	0
9	C	1121	DGX	2	0
7	D	402	CLR	1	0
8	A	1105	PCW	2	0
7	A	1104	CLR	2	0
7	E	101	CLR	6	0
8	C	1105	PCW	2	0
8	A	1109	PCW	1	0
8	A	1108	PCW	2	0
8	A	1106	PCW	4	0
7	G	101	CLR	1	0

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.

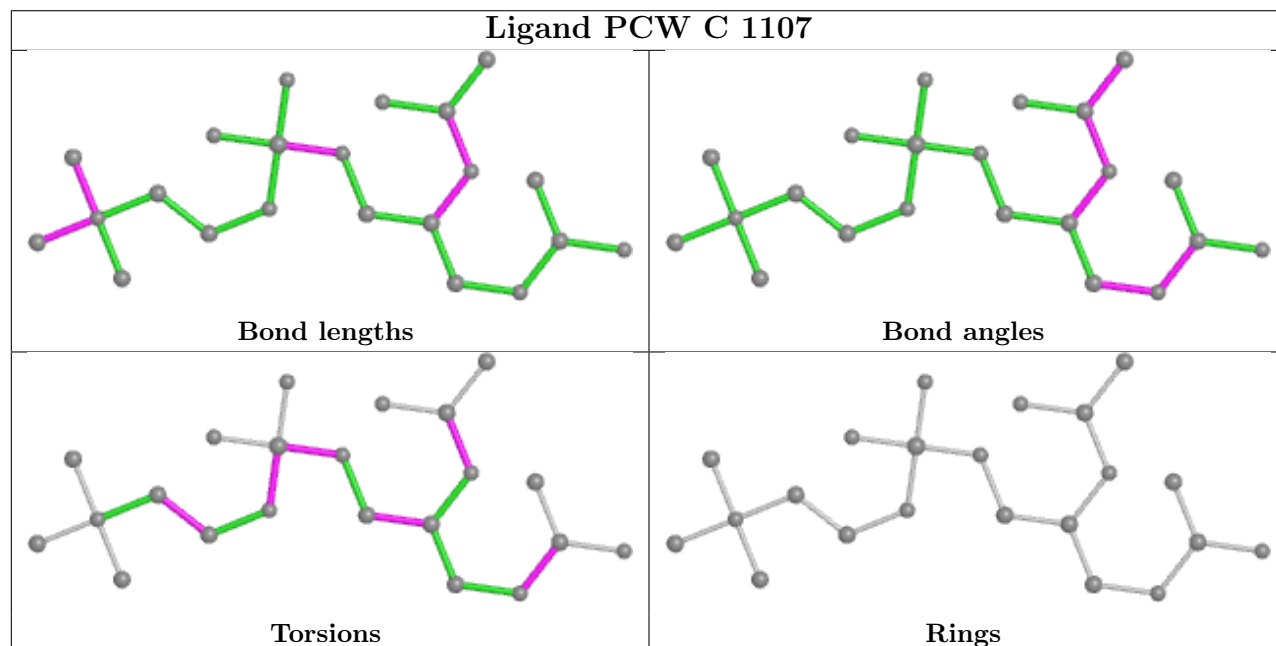




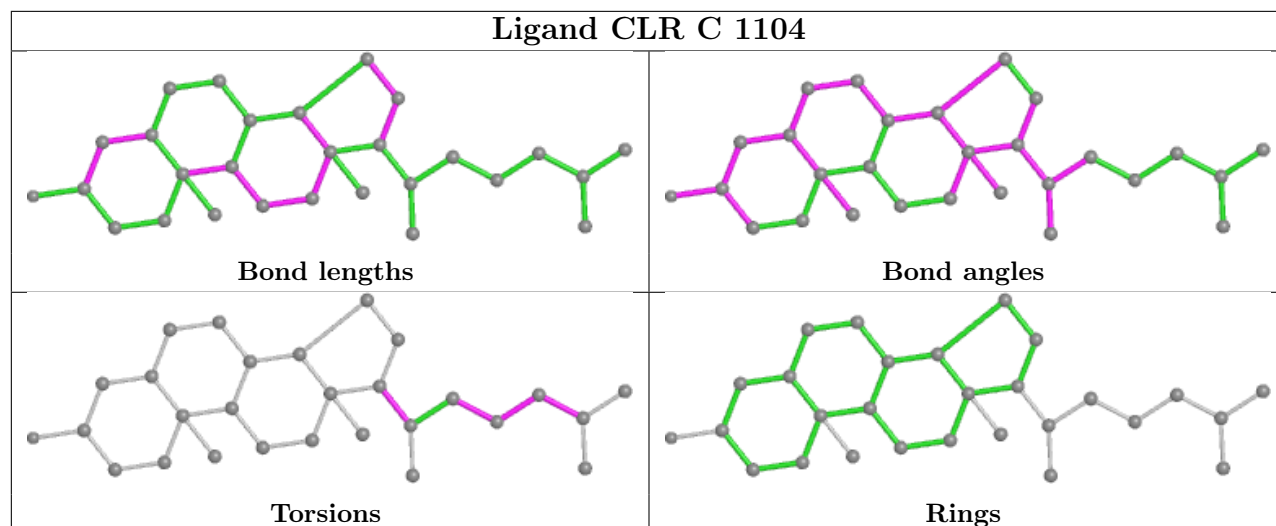
## Ligand DGX A 1110



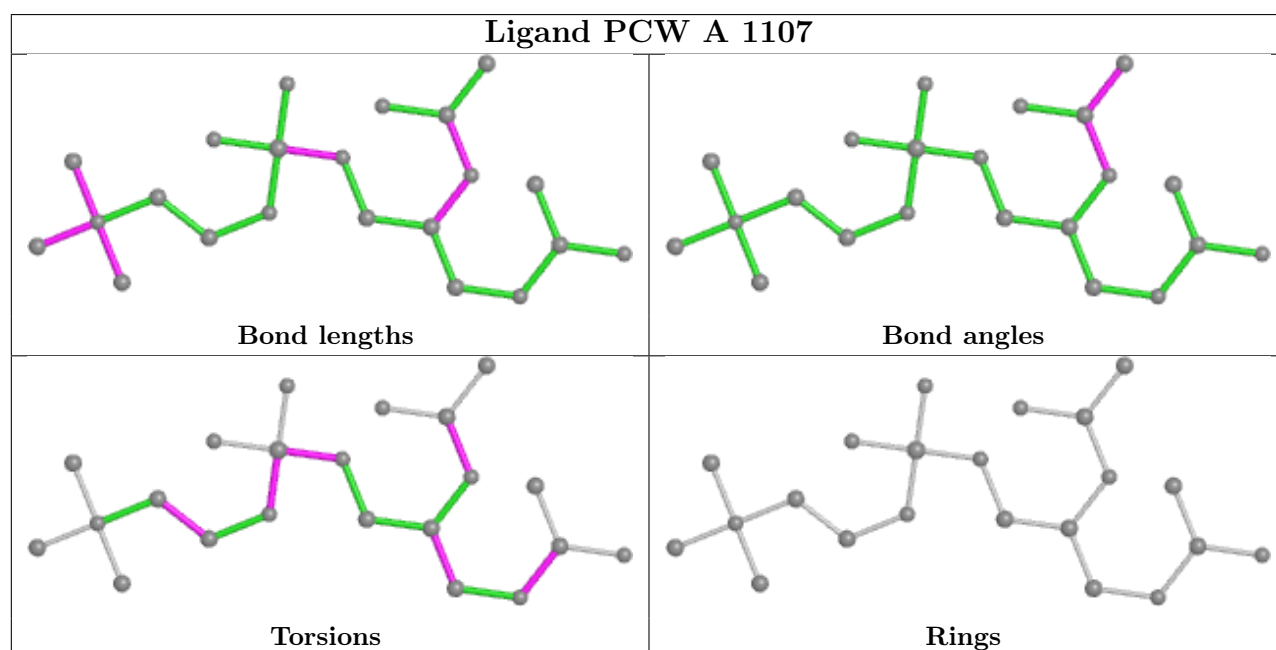
## Ligand PCW C 1107



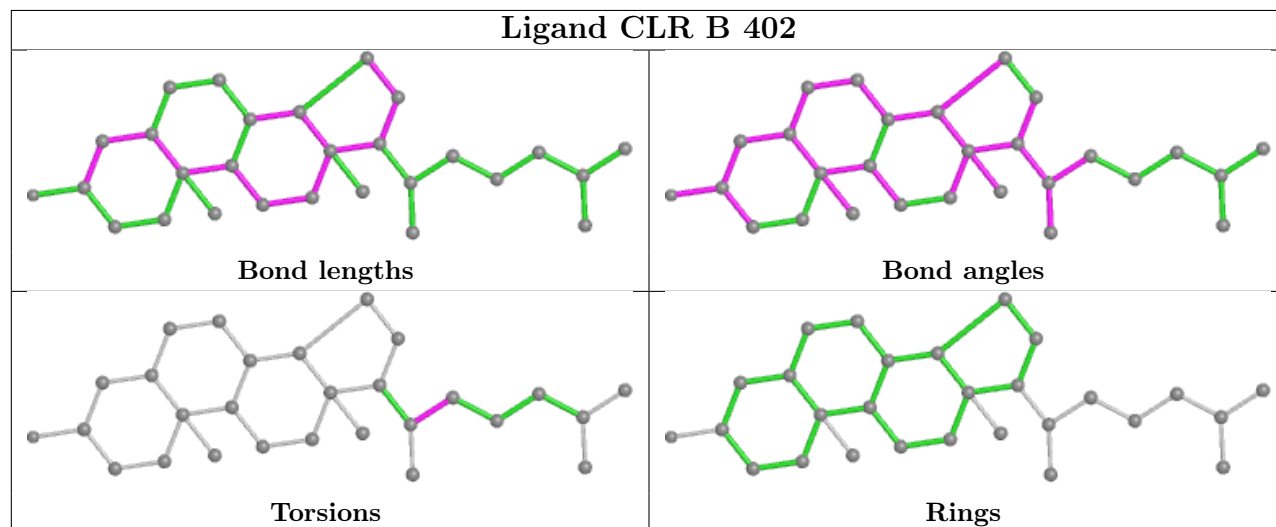
## Ligand CLR C 1104

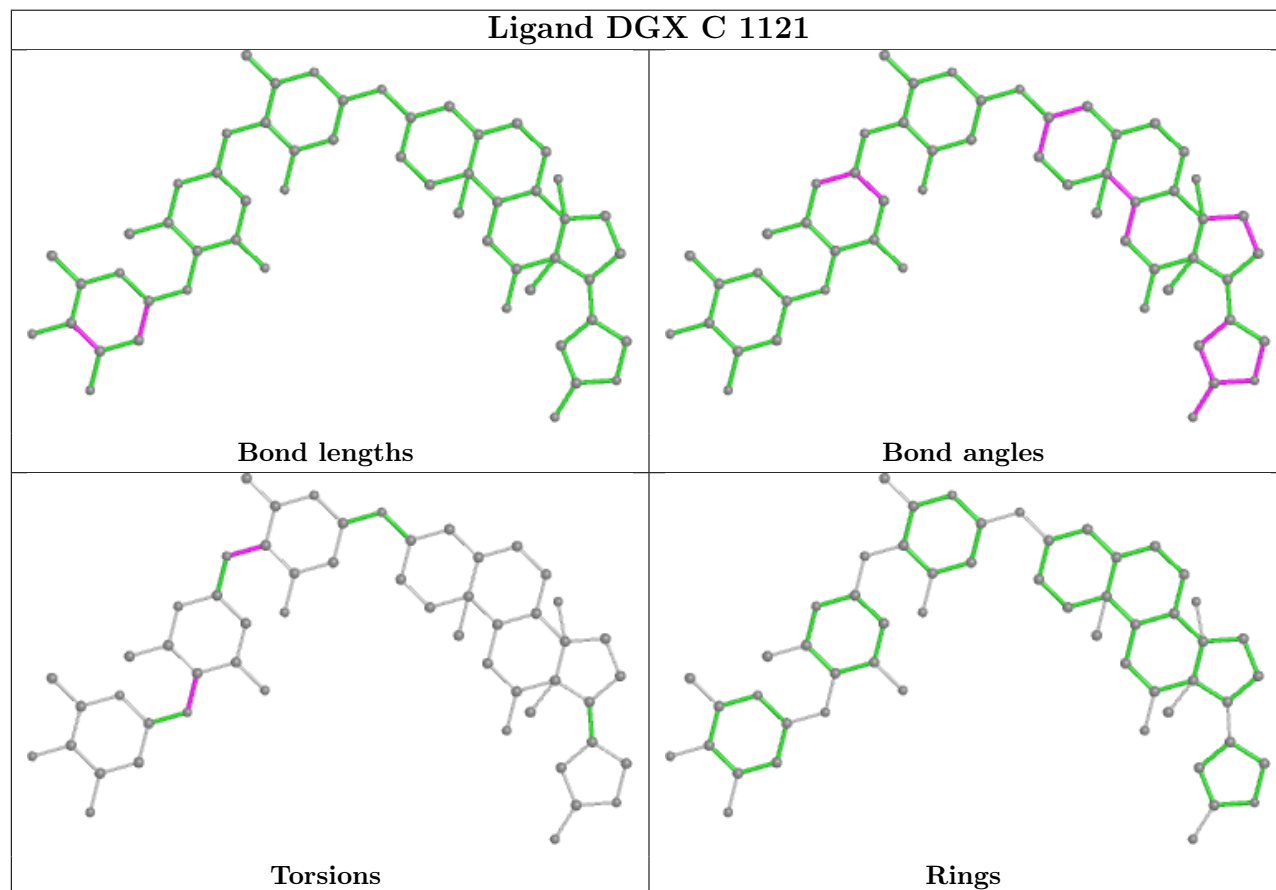
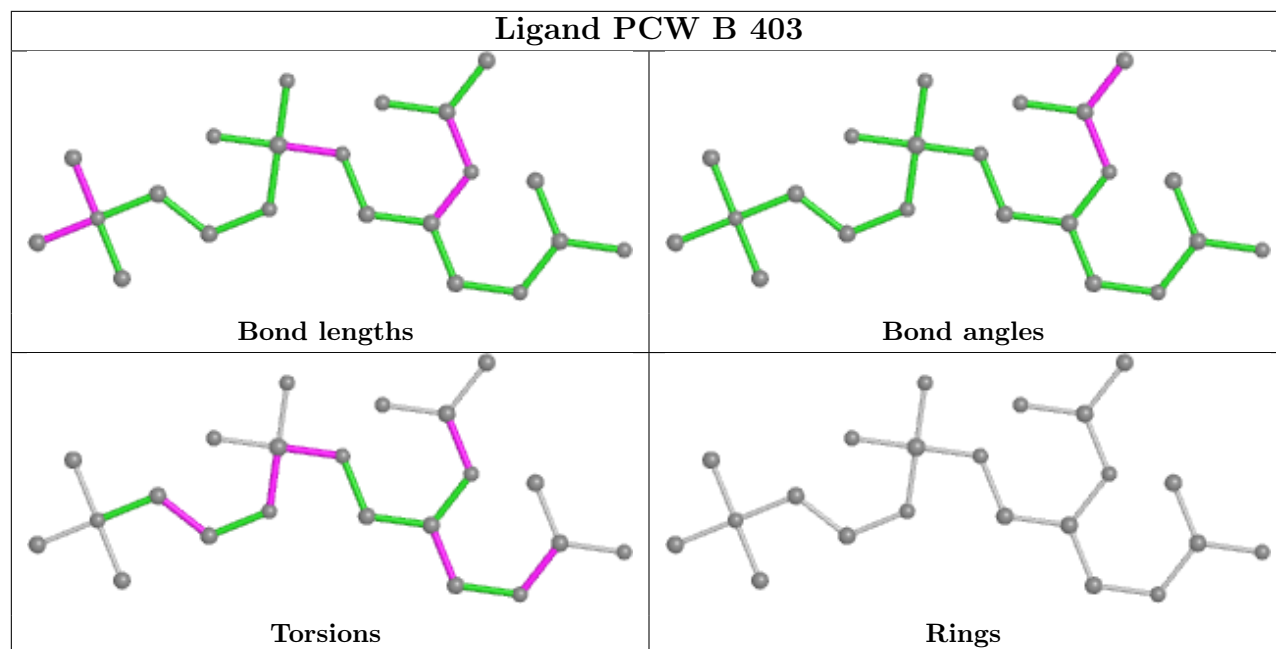


## Ligand PCW A 1107

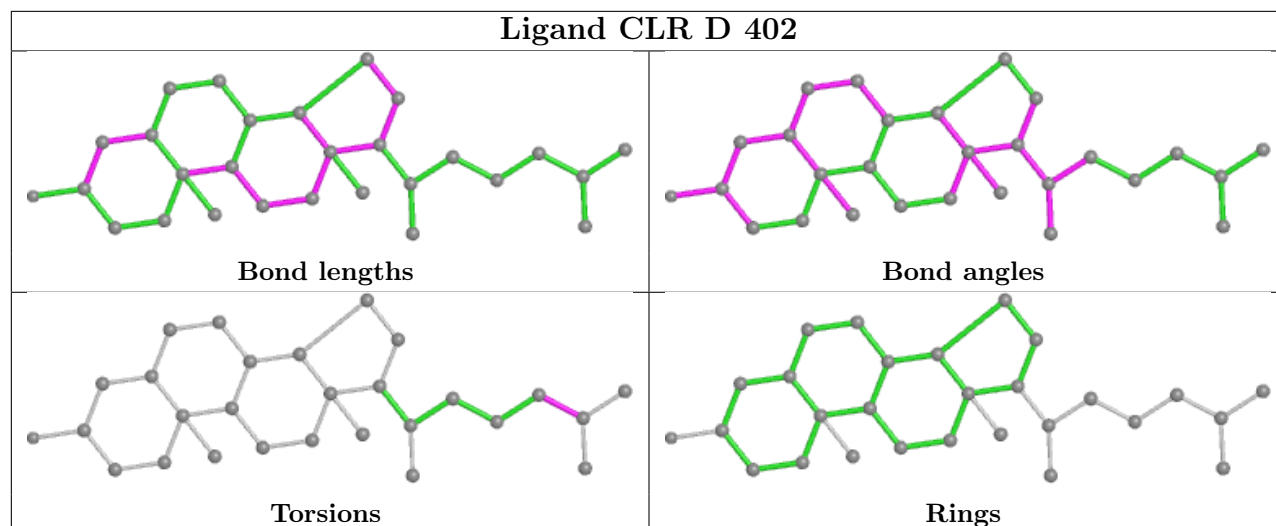


## Ligand CLR B 402

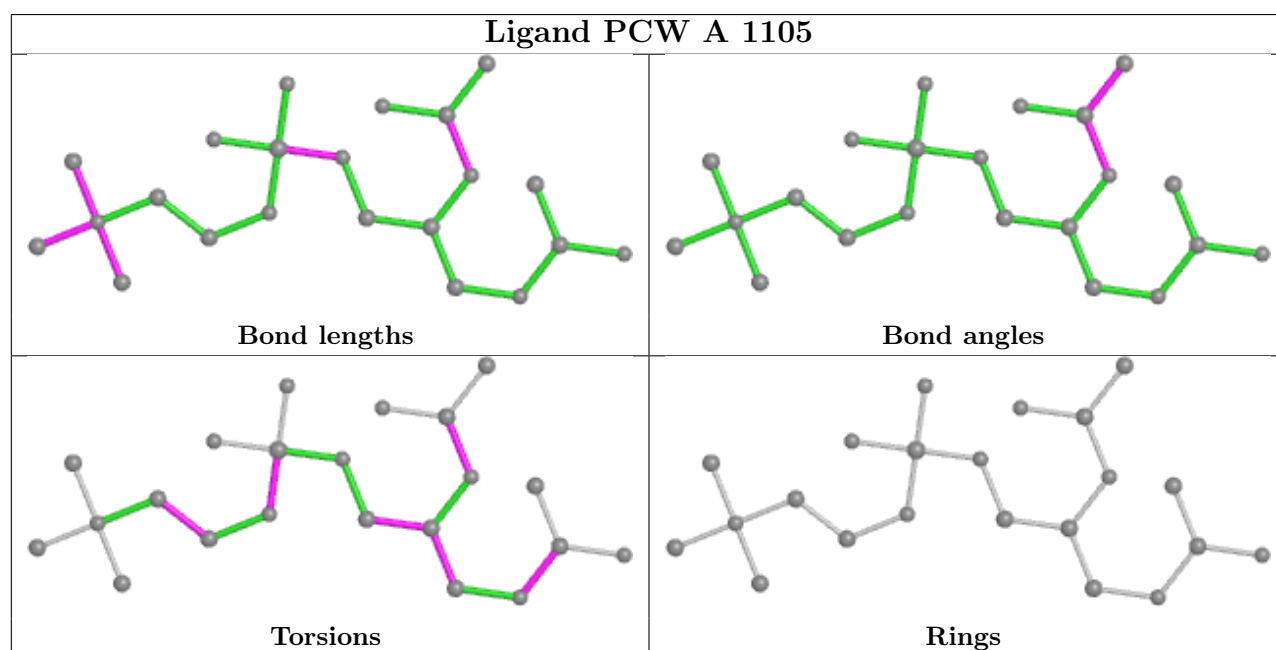




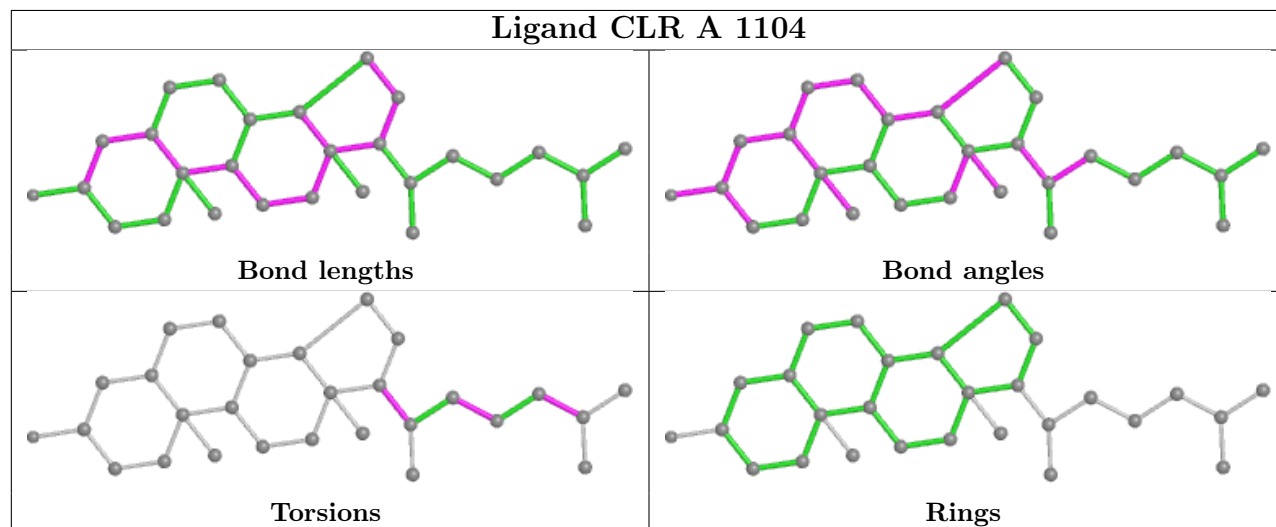
## Ligand CLR D 402



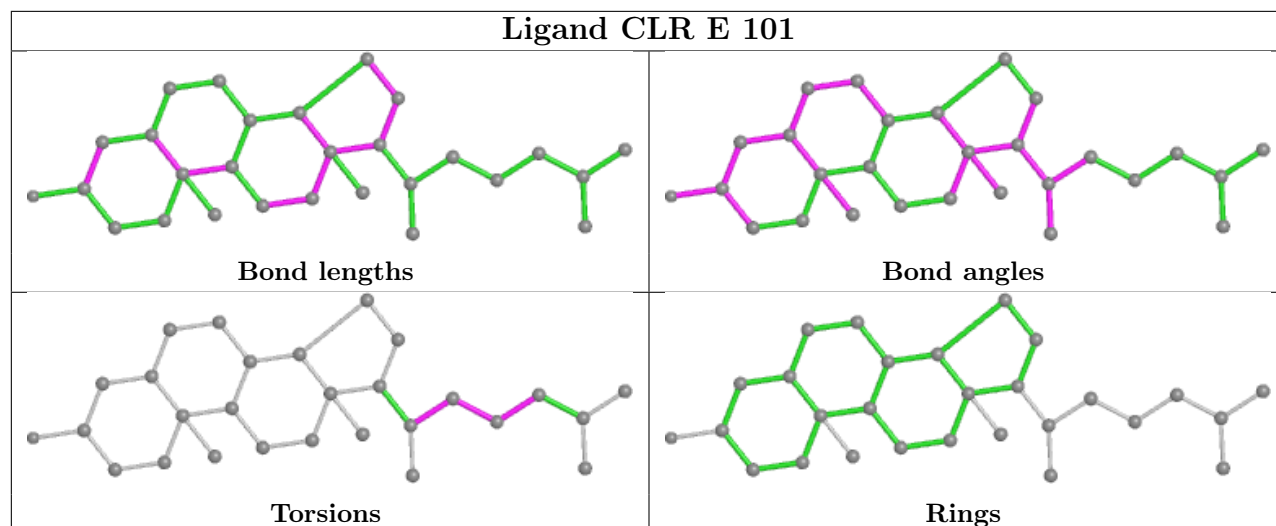
## Ligand PCW A 1105



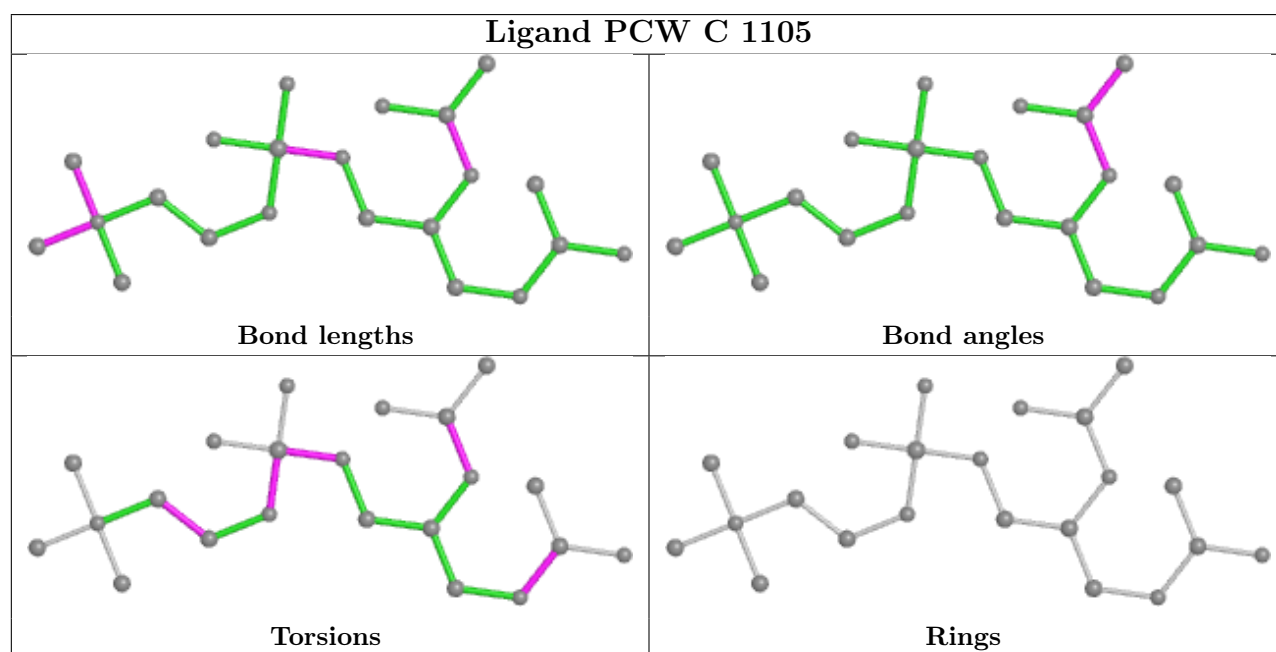
## Ligand CLR A 1104



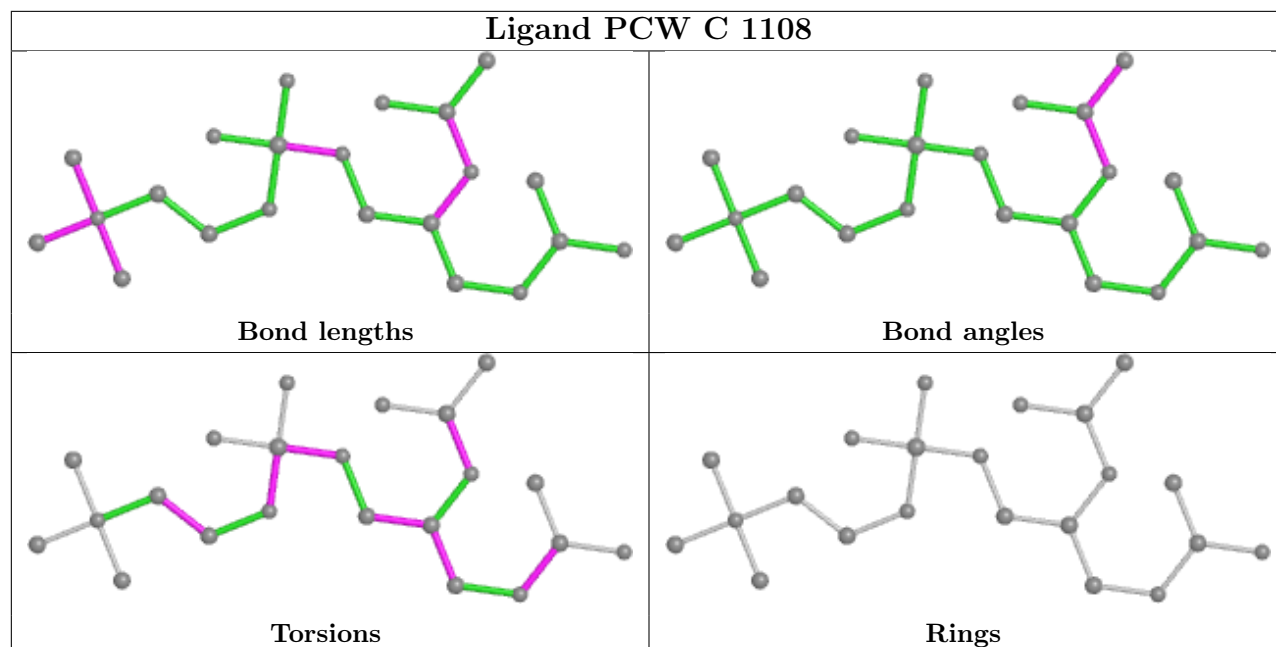
## Ligand CLR E 101



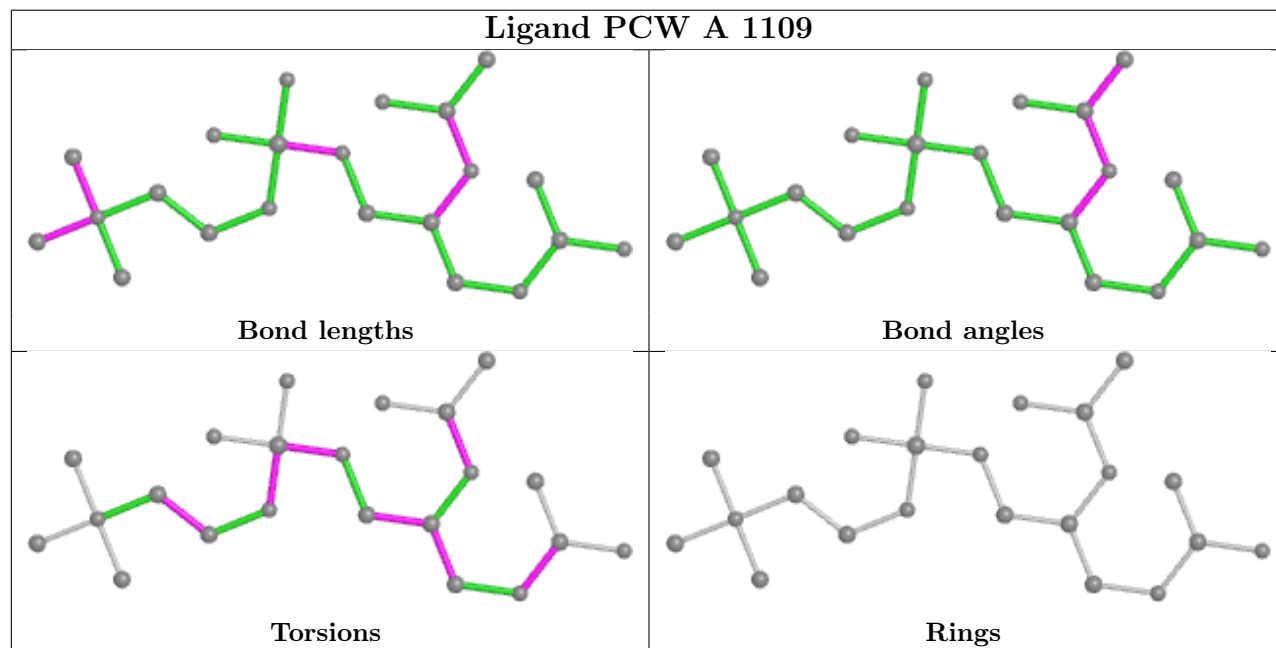
## Ligand PCW C 1105



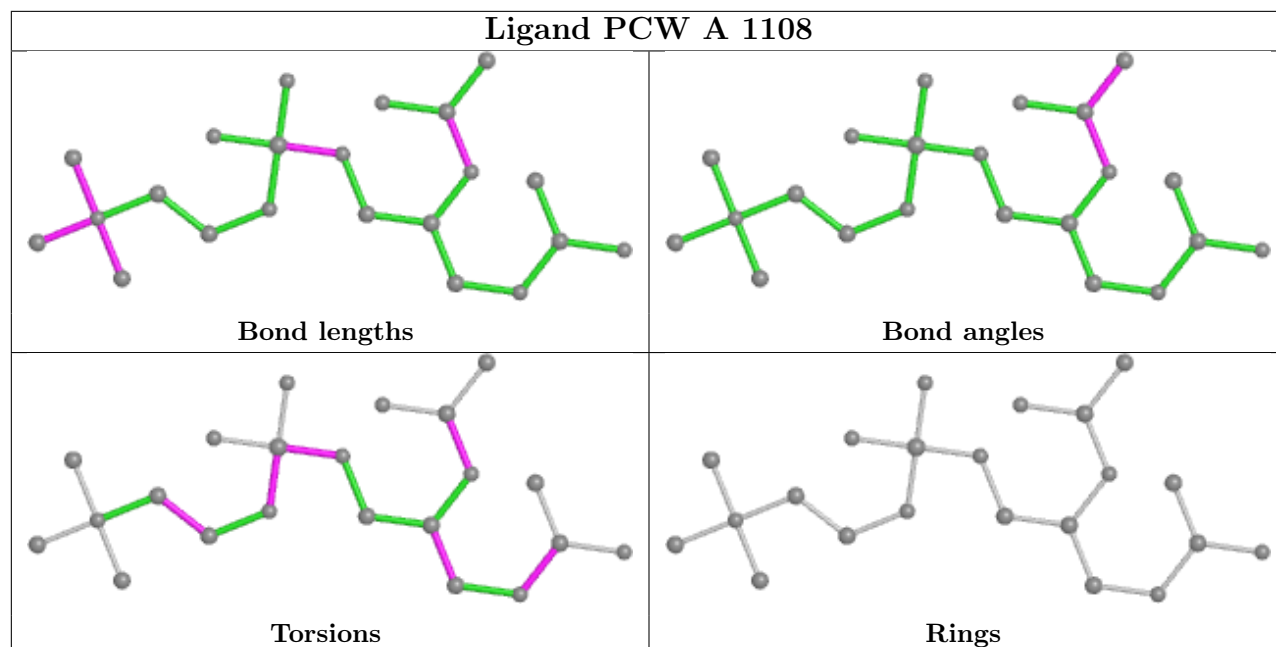
## Ligand PCW C 1108



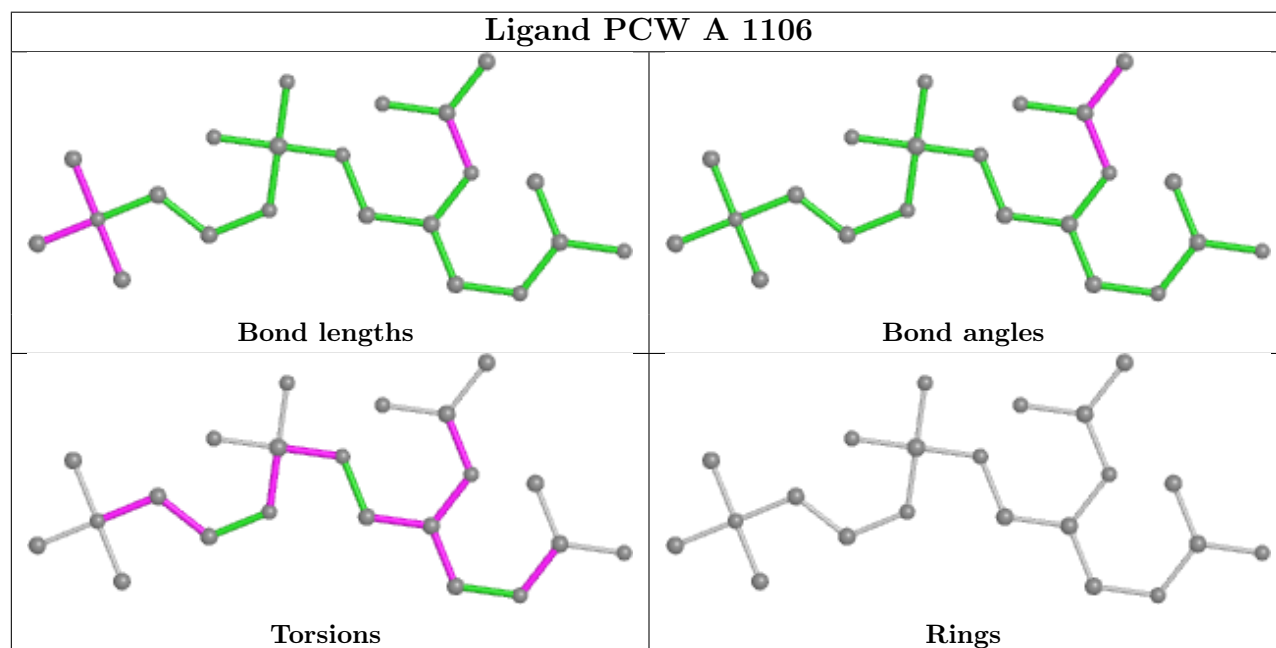
## Ligand PCW A 1109

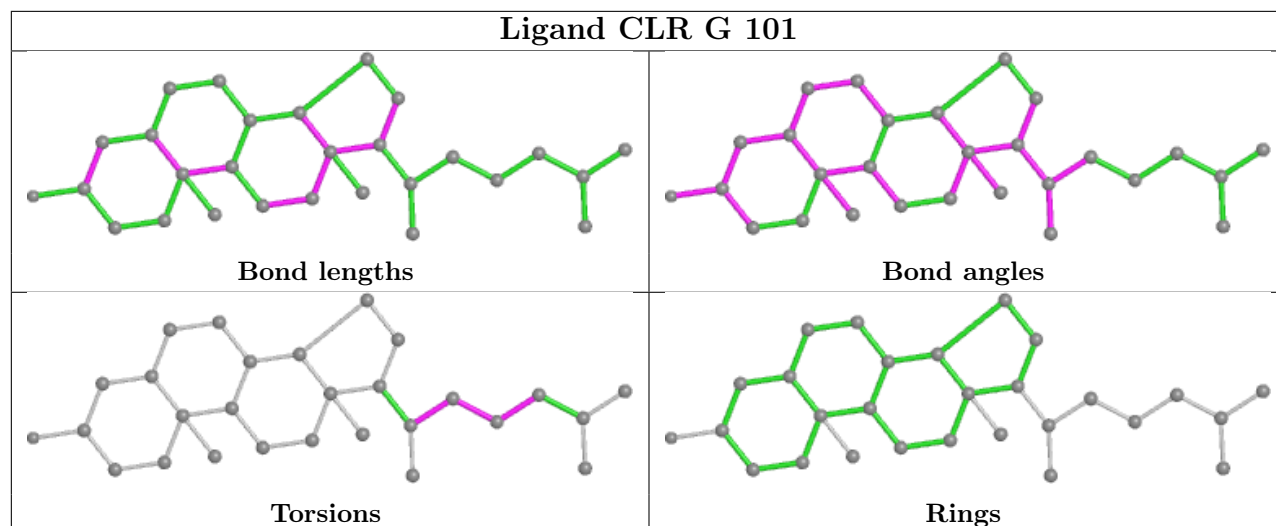


## Ligand PCW A 1108



## Ligand PCW A 1106





## 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	995/1016 (97%)	0.16	43 (4%)	35 34	50, 103, 209, 236	0
1	C	995/1016 (97%)	0.14	32 (3%)	47 46	51, 107, 185, 211	0
2	B	291/303 (96%)	0.10	21 (7%)	15 17	66, 134, 178, 209	0
2	D	285/303 (94%)	0.00	11 (3%)	39 38	58, 124, 176, 202	0
3	E	32/65 (49%)	-0.01	0	100 100	60, 84, 119, 143	0
3	G	32/65 (49%)	0.08	0	100 100	57, 85, 120, 130	0
All	All	2630/2768 (95%)	0.12	107 (4%)	37 36	50, 110, 196, 236	0

The worst 5 of 107 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	571	PHE	8.7
2	B	165	GLU	8.0
1	C	579	VAL	6.1
1	A	550	HIS	6.1
1	A	489	PRO	6.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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.

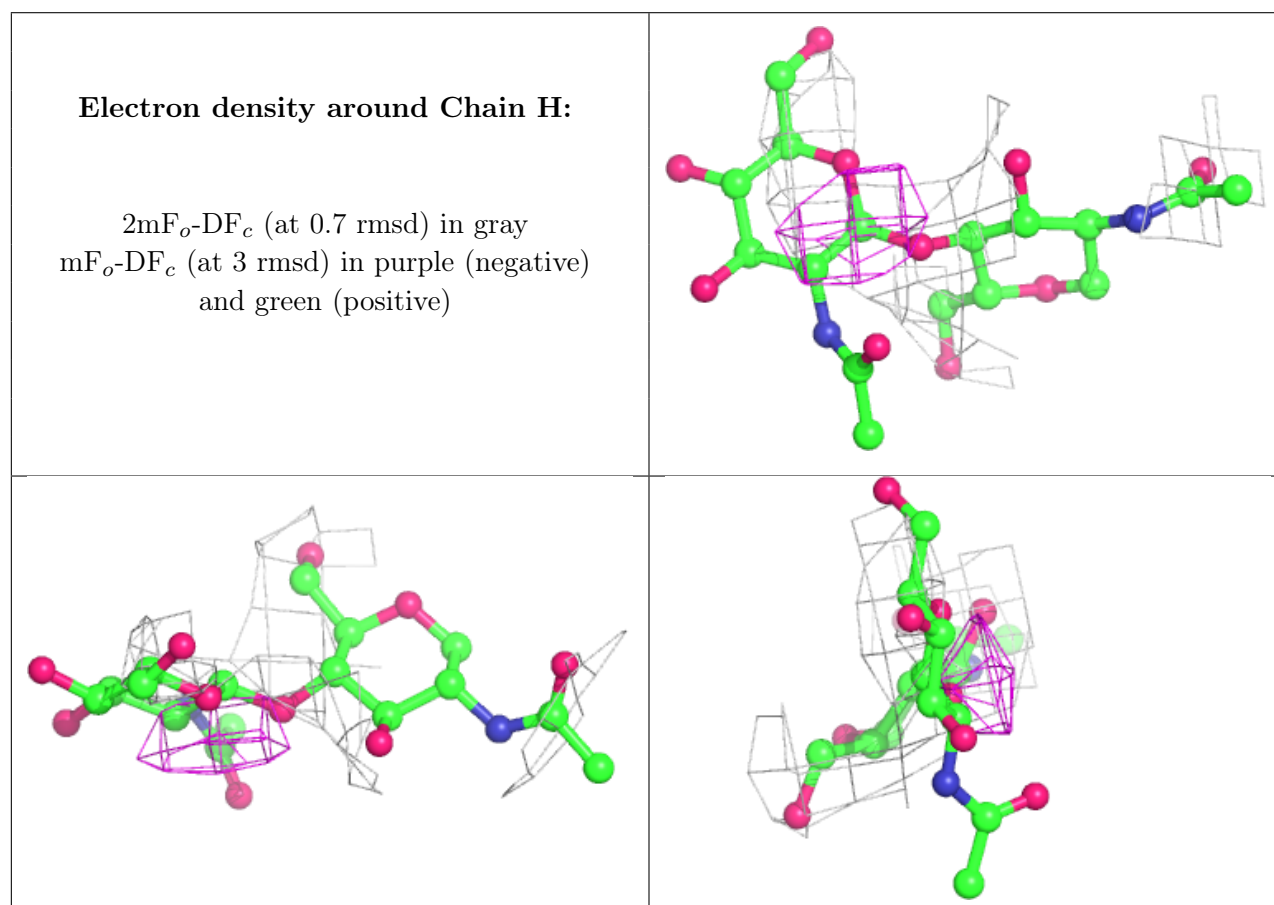
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	PHD	A	369	12/13	0.98	0.22	73,73,97,101	0
1	PHD	C	369	12/13	0.99	0.21	86,86,118,125	0

### 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, 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.

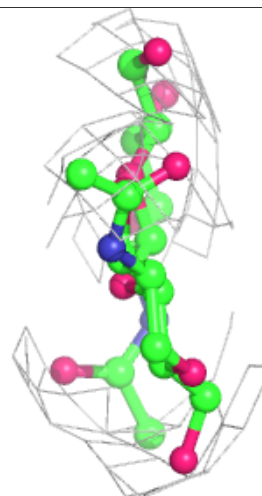
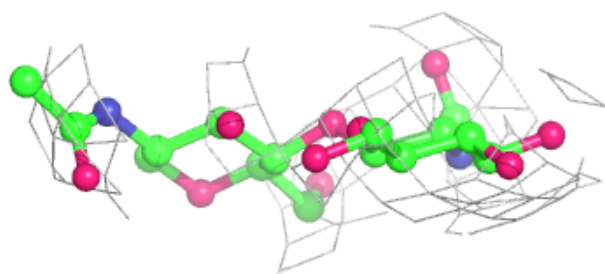
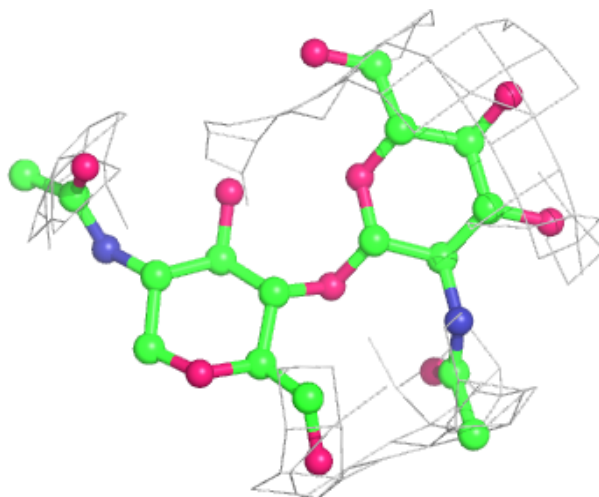
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	NAG	H	2	14/15	0.53	0.69	193,205,230,231	0
4	NAG	H	1	14/15	0.74	0.20	179,187,213,237	0
4	NAG	J	2	14/15	0.80	0.57	184,196,213,213	0
4	NAG	F	1	14/15	0.80	0.13	155,159,180,196	0
4	NAG	J	1	14/15	0.84	0.29	176,187,213,219	0
4	NAG	F	2	14/15	0.87	0.38	168,187,196,201	0
4	NAG	I	1	14/15	0.88	0.14	143,151,166,188	0
4	NAG	I	2	14/15	0.93	0.11	150,162,174,178	0

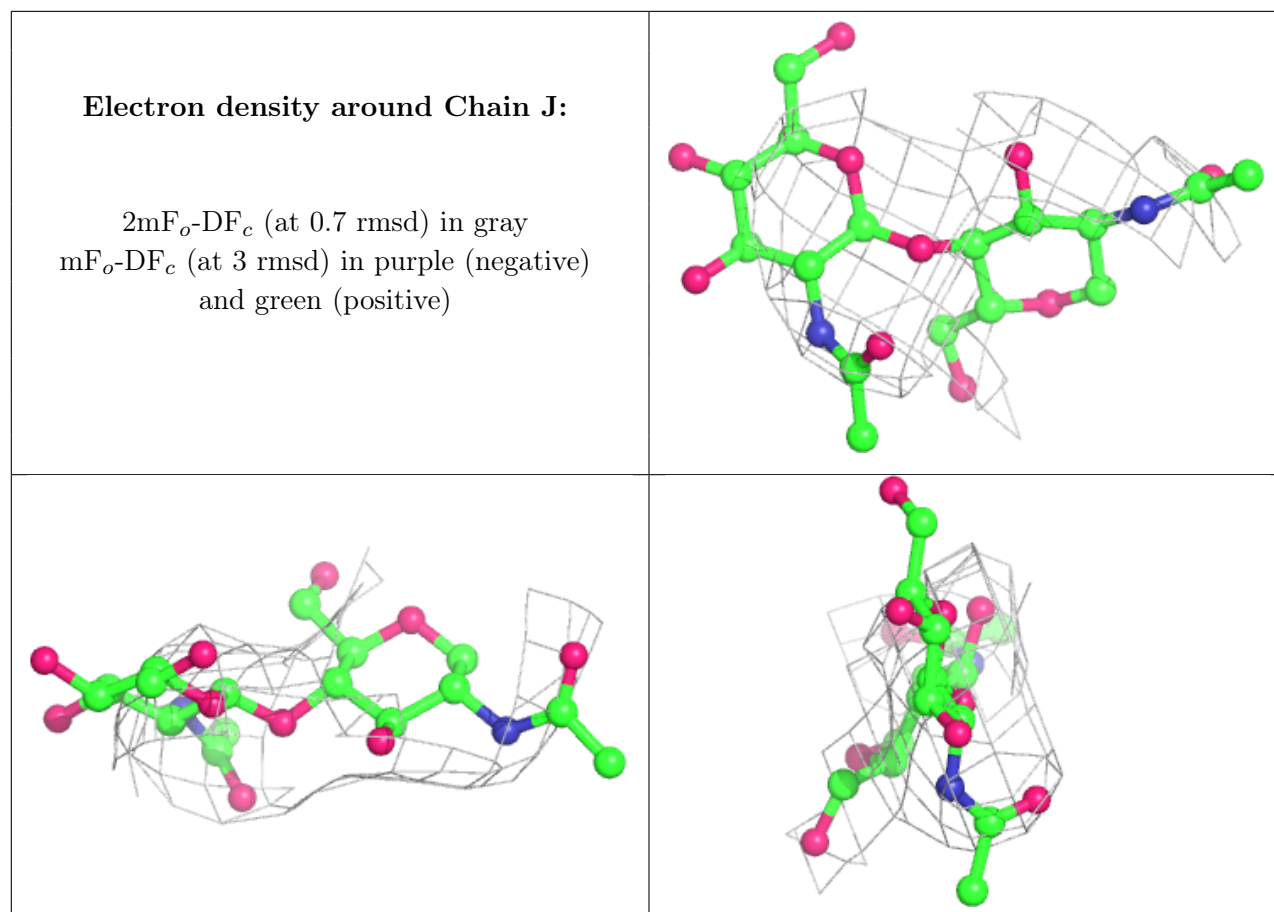
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



**Electron density around Chain I:**

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





## 6.4 Ligands [i](#)

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

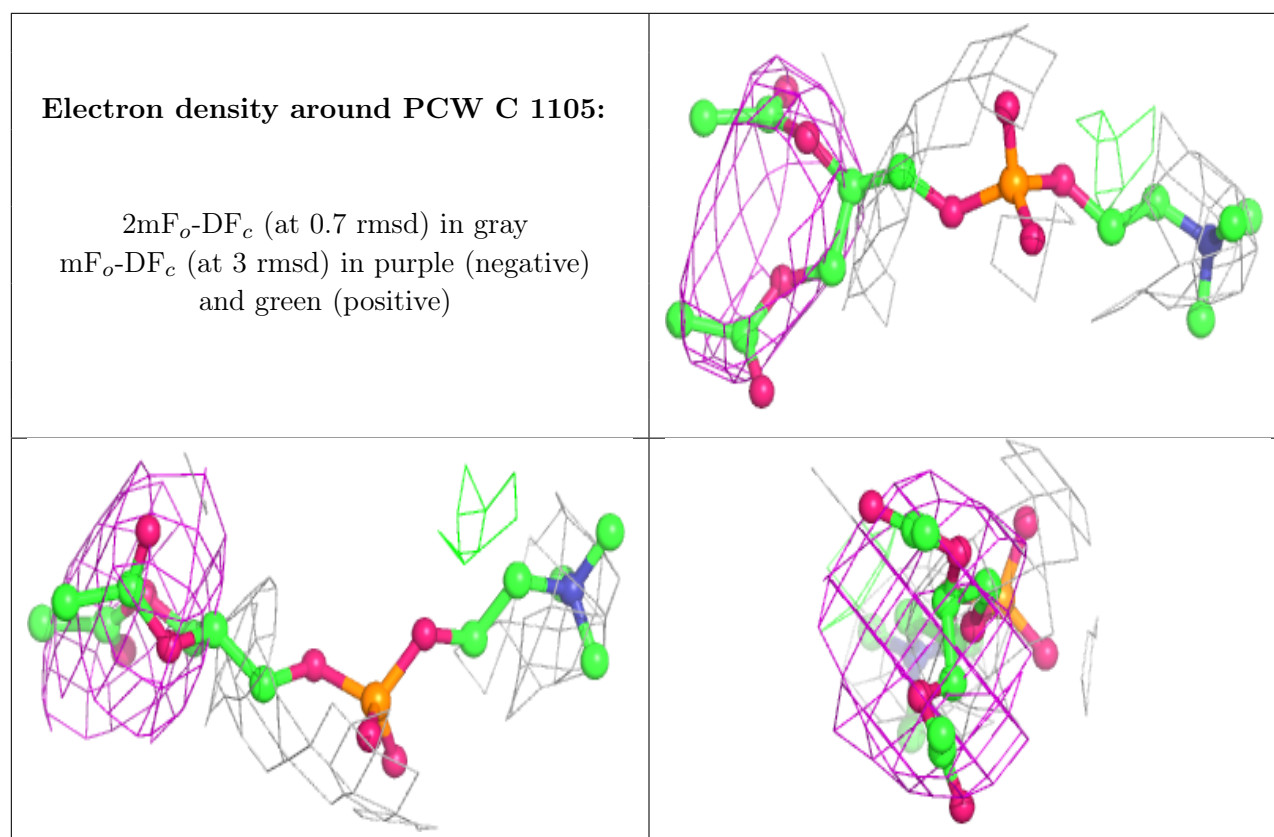
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
8	PCW	C	1105	22/54	0.55	0.65	119,157,191,204	0
8	PCW	A	1107	22/54	0.67	0.42	154,180,205,211	0
10	NAG	B	401	14/15	0.74	0.20	161,177,190,195	0
8	PCW	A	1108	22/54	0.75	0.69	130,163,204,216	0
8	PCW	A	1106	22/54	0.75	0.42	118,158,195,209	0
10	NAG	D	401	14/15	0.77	0.18	164,178,198,201	0
8	PCW	C	1106	22/54	0.78	0.36	122,158,180,196	0
7	CLR	D	402	28/28	0.85	0.71	97,120,149,166	0
7	CLR	A	1104	28/28	0.86	0.91	83,111,135,145	0
8	PCW	A	1105	22/54	0.86	0.22	93,140,181,190	0
7	CLR	C	1104	28/28	0.88	0.70	67,104,133,144	0
9	DGX	A	1110	55/55	0.88	0.26	102,119,149,165	0

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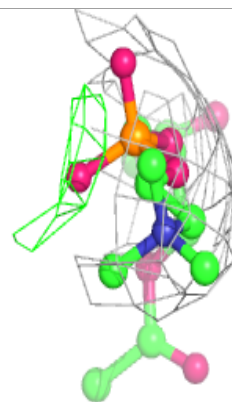
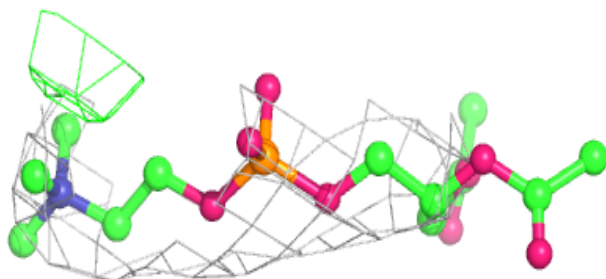
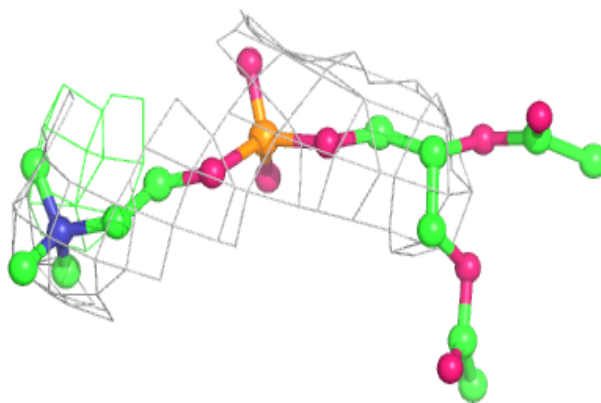
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
8	PCW	C	1108	22/54	0.90	0.21	106,125,142,161	0
8	PCW	B	403	22/54	0.90	0.45	118,165,189,206	0
8	PCW	A	1109	22/54	0.90	0.39	83,118,149,157	0
7	CLR	B	402	28/28	0.92	0.83	107,128,142,152	0
9	DGX	C	1121	55/55	0.92	0.30	80,90,126,132	0
7	CLR	E	101	28/28	0.93	0.46	54,55,75,88	0
8	PCW	C	1107	22/54	0.93	0.32	82,112,165,171	0
7	CLR	G	101	28/28	0.95	0.47	61,63,95,121	0
6	NA	C	1102	1/1	0.97	0.38	41,41,41,41	0
6	NA	A	1102	1/1	0.97	0.29	50,50,50,50	0
5	MG	C	1103	1/1	0.97	0.20	55,55,55,55	0
5	MG	A	1101	1/1	0.98	0.20	87,87,87,87	0
5	MG	A	1103	1/1	0.98	0.28	67,67,67,67	0
5	MG	C	1101	1/1	0.99	0.22	87,87,87,87	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.

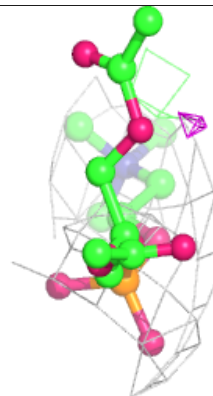
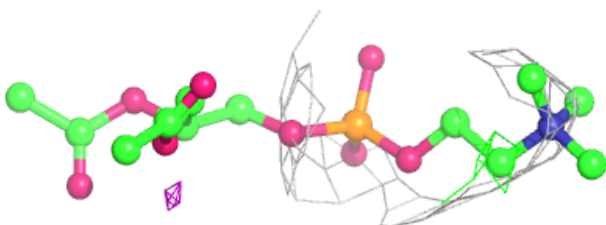
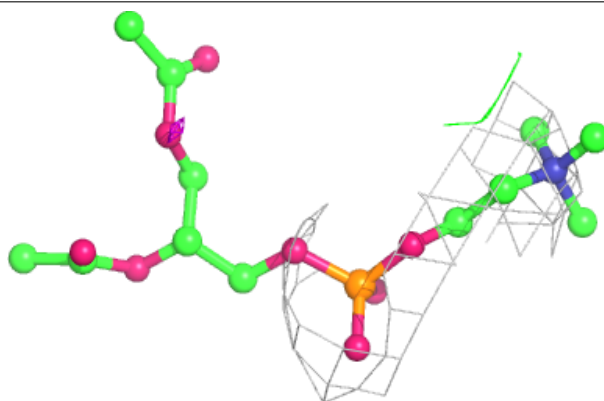


**Electron density around PCW A 1107:**

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

**Electron density around PCW A 1108:**

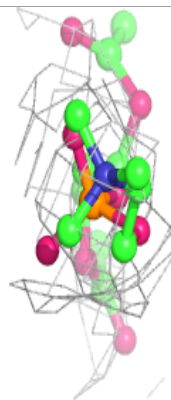
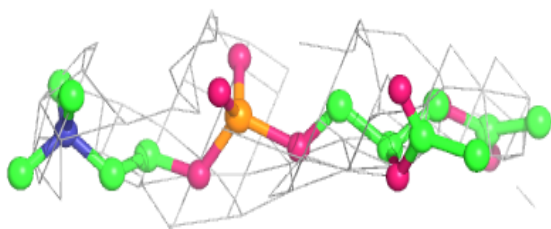
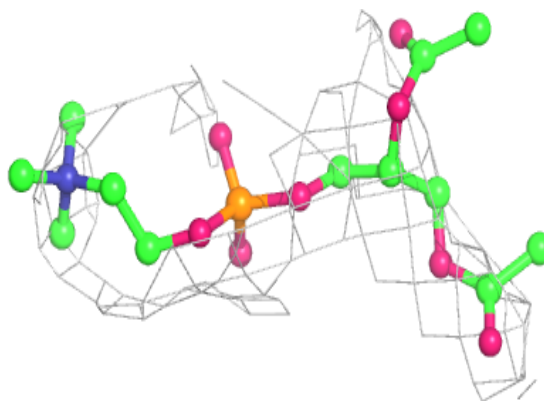
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



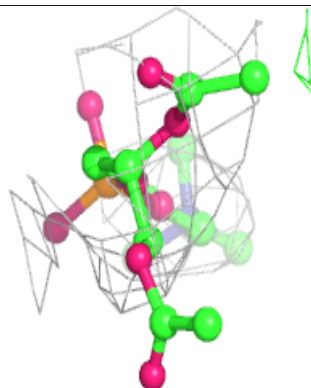
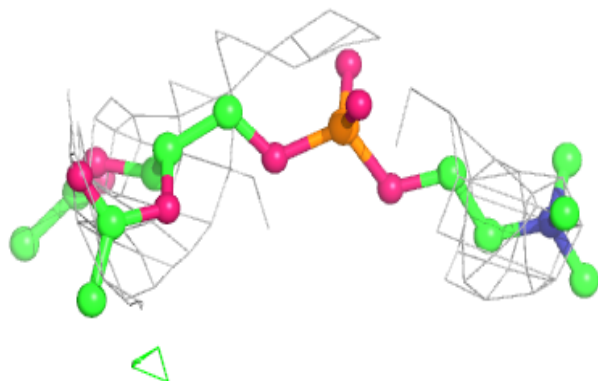
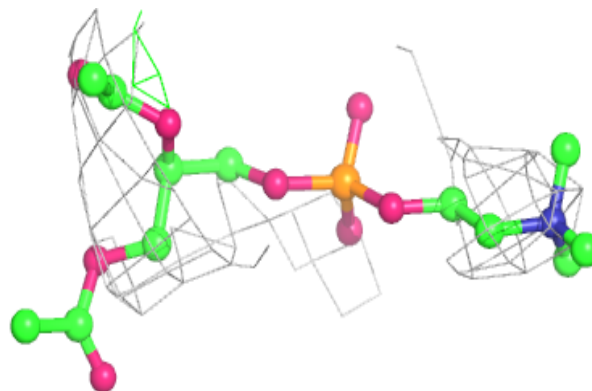


**Electron density around PCW A 1106:**

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

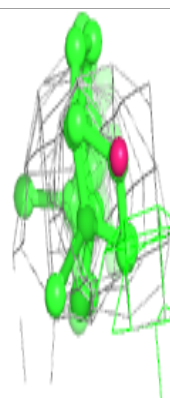
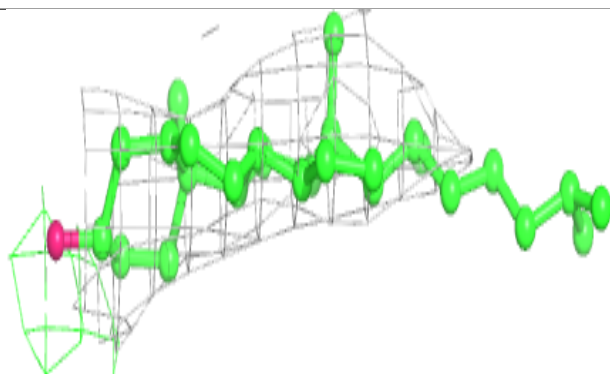
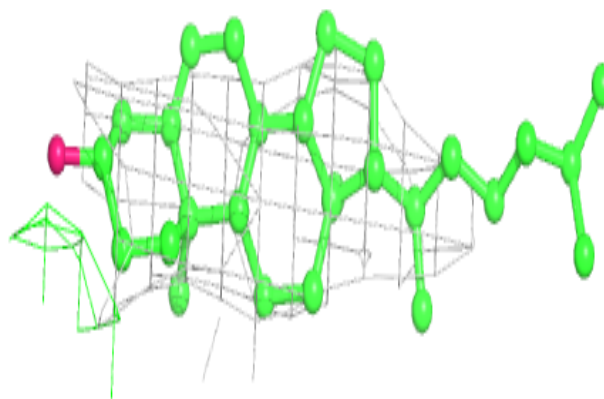
**Electron density around PCW C 1106:**

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

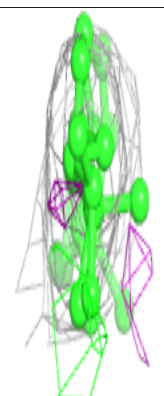
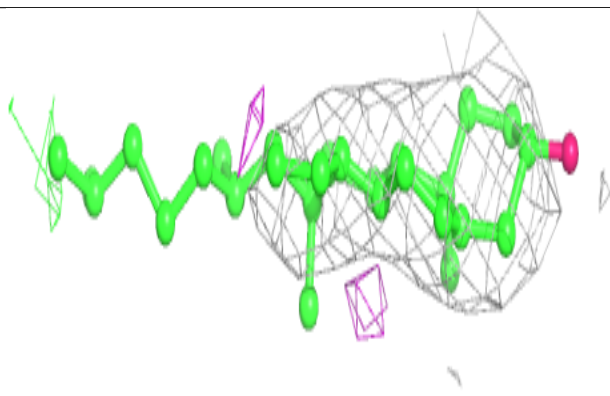
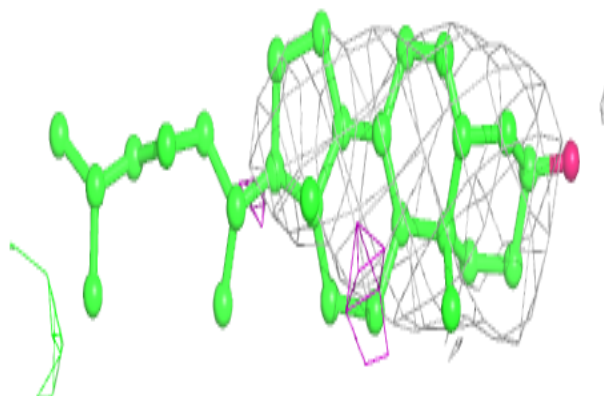


**Electron density around CLR D 402:**

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

**Electron density around CLR A 1104:**

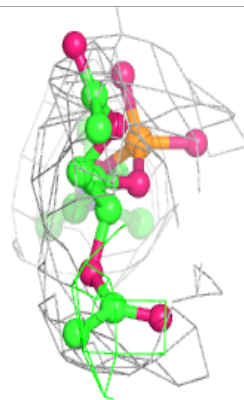
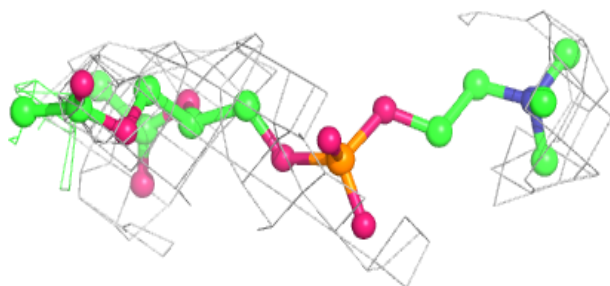
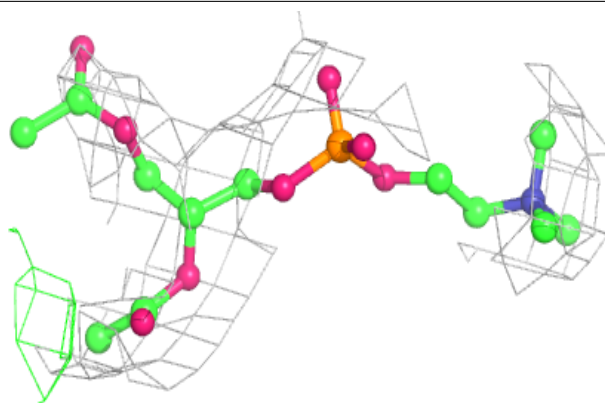
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



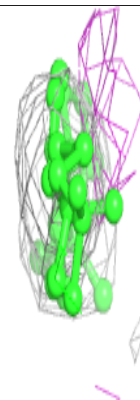
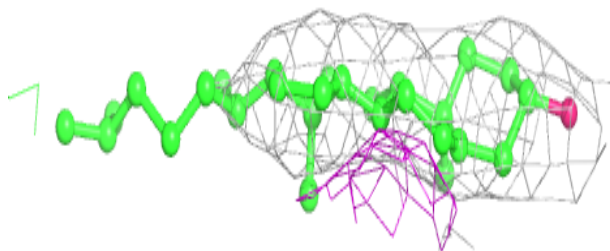
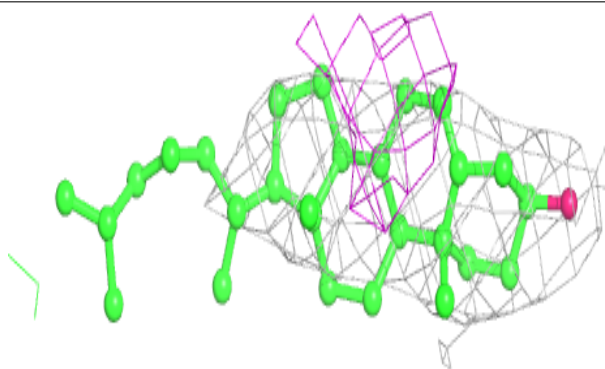


**Electron density around PCW A 1105:**

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

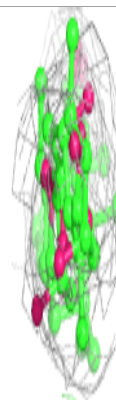
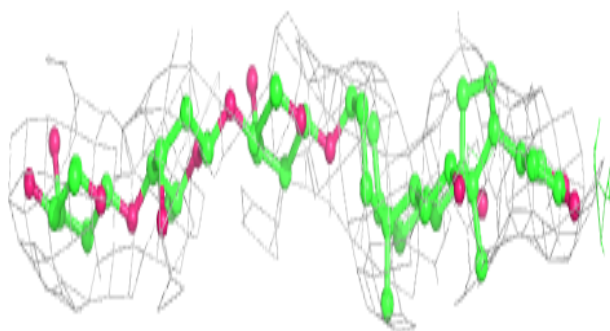
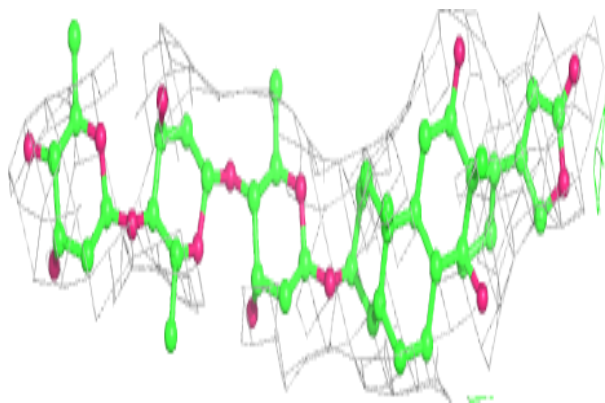
**Electron density around CLR C 1104:**

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

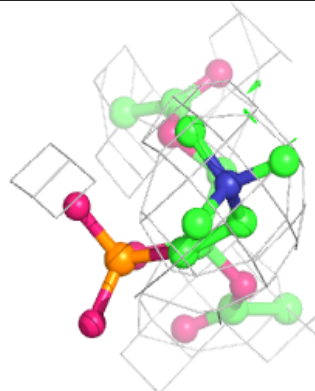
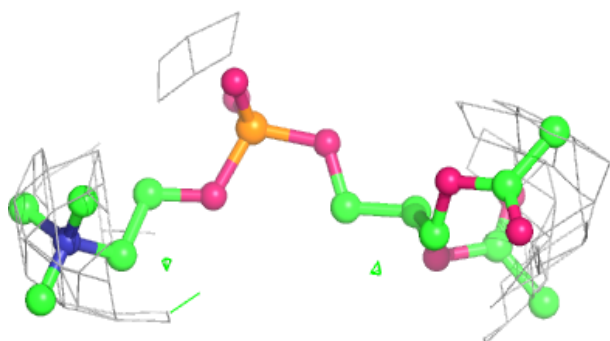
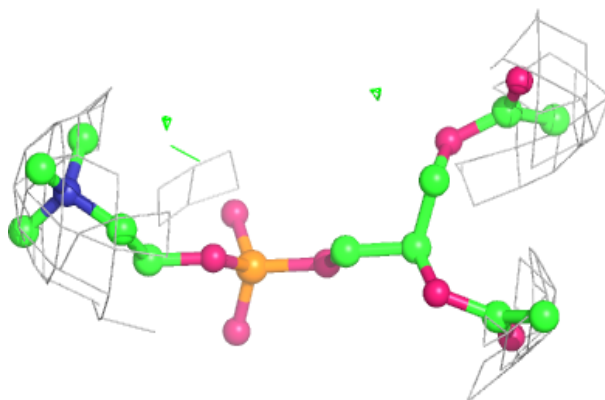


**Electron density around DGX A 1110:**

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

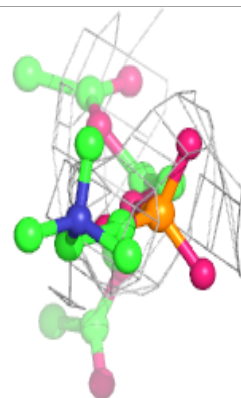
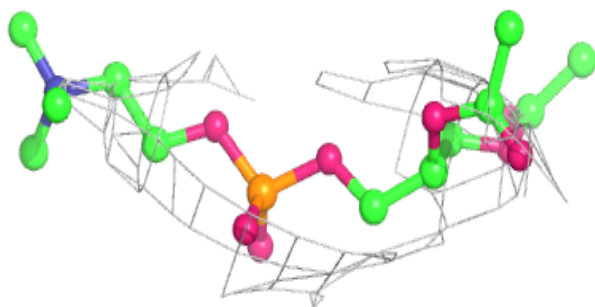
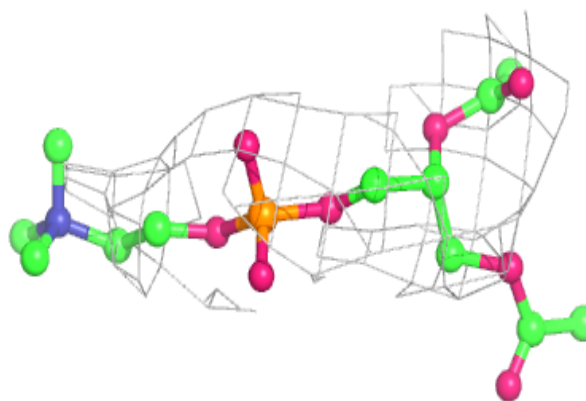
**Electron density around PCW C 1108:**

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

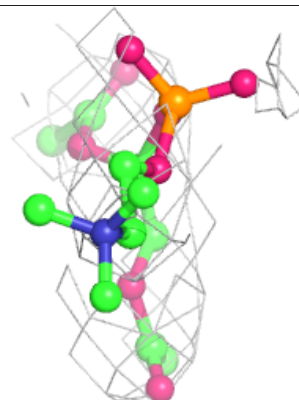
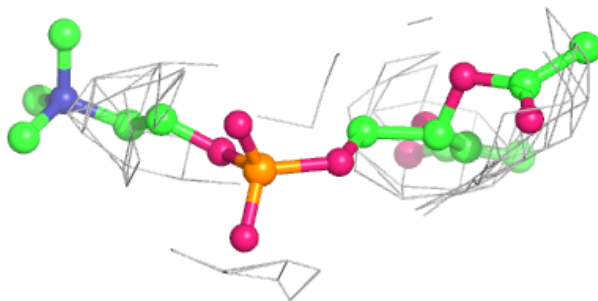
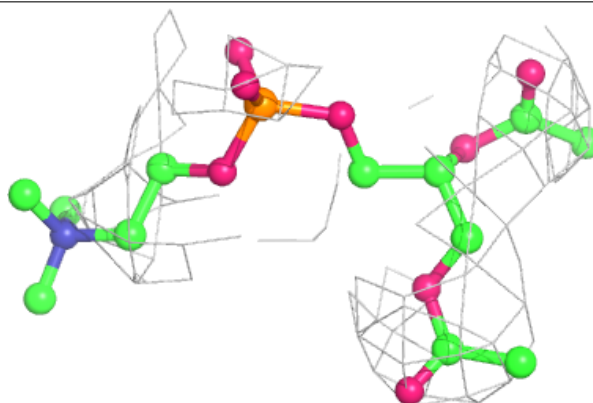


**Electron density around PCW B 403:**

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

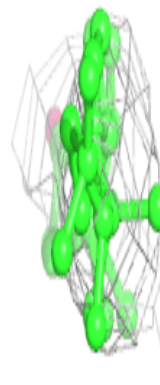
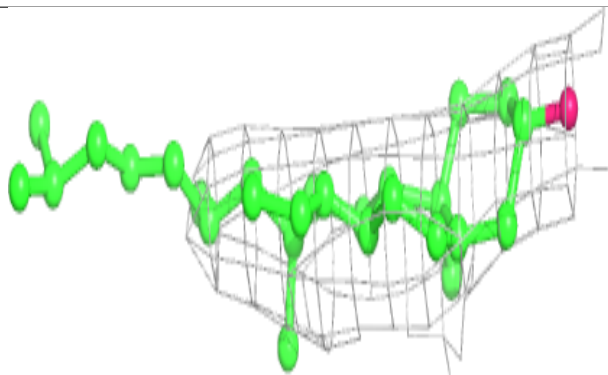
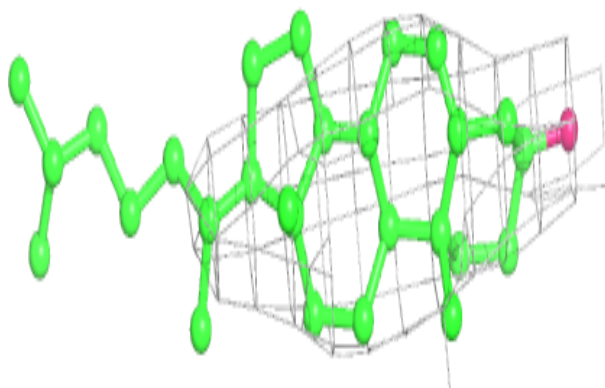
**Electron density around PCW A 1109:**

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

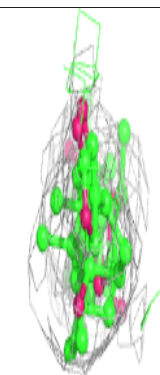
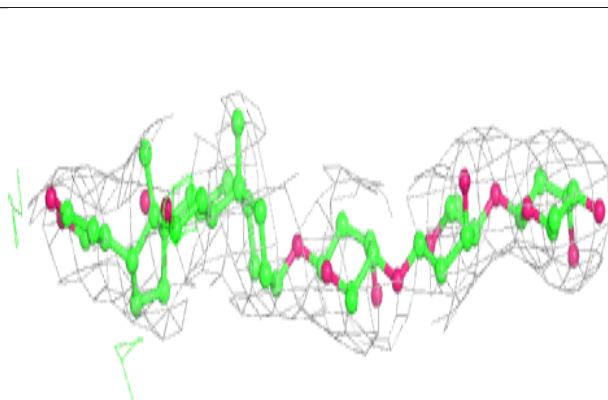
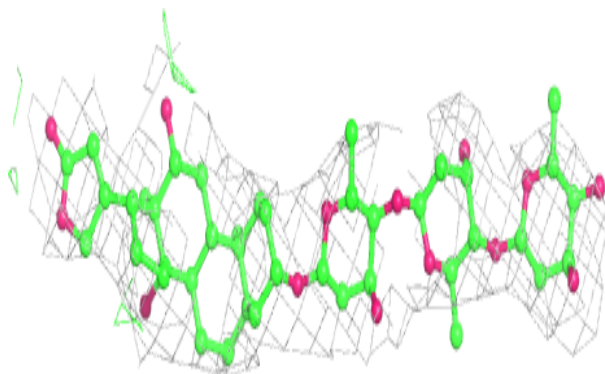


**Electron density around CLR B 402:**

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

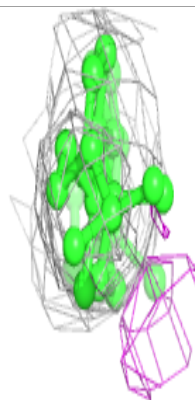
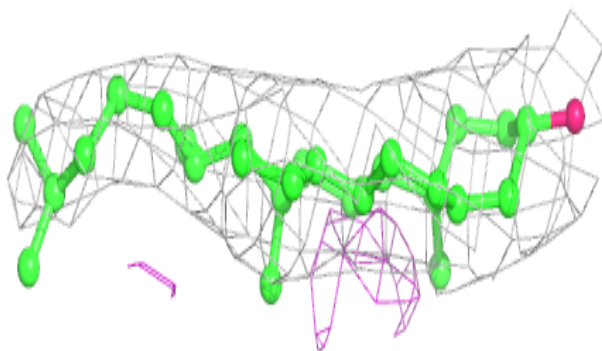
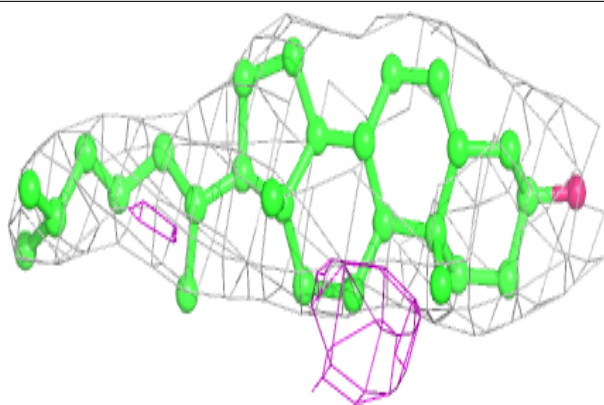
**Electron density around DGX C 1121:**

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

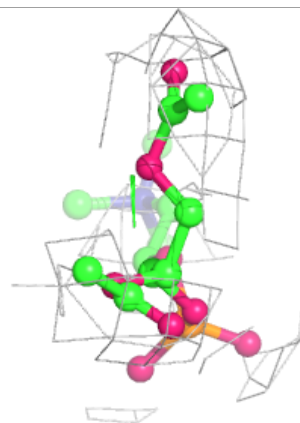
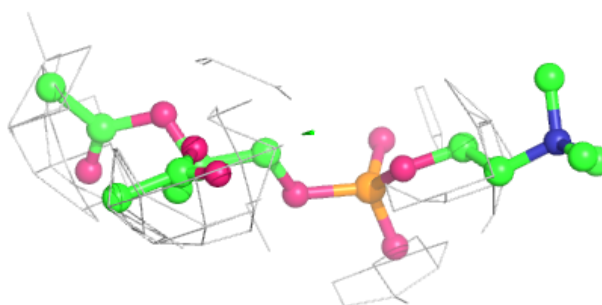
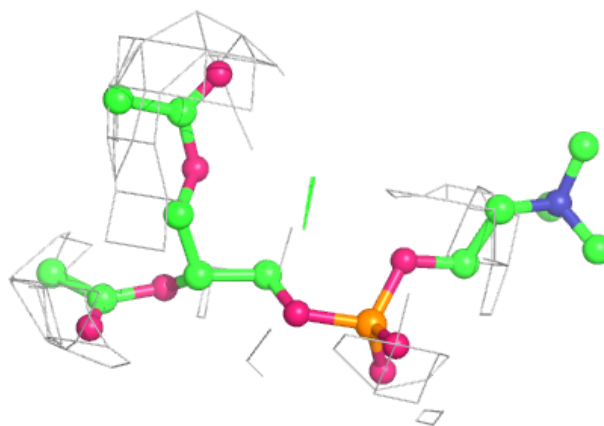


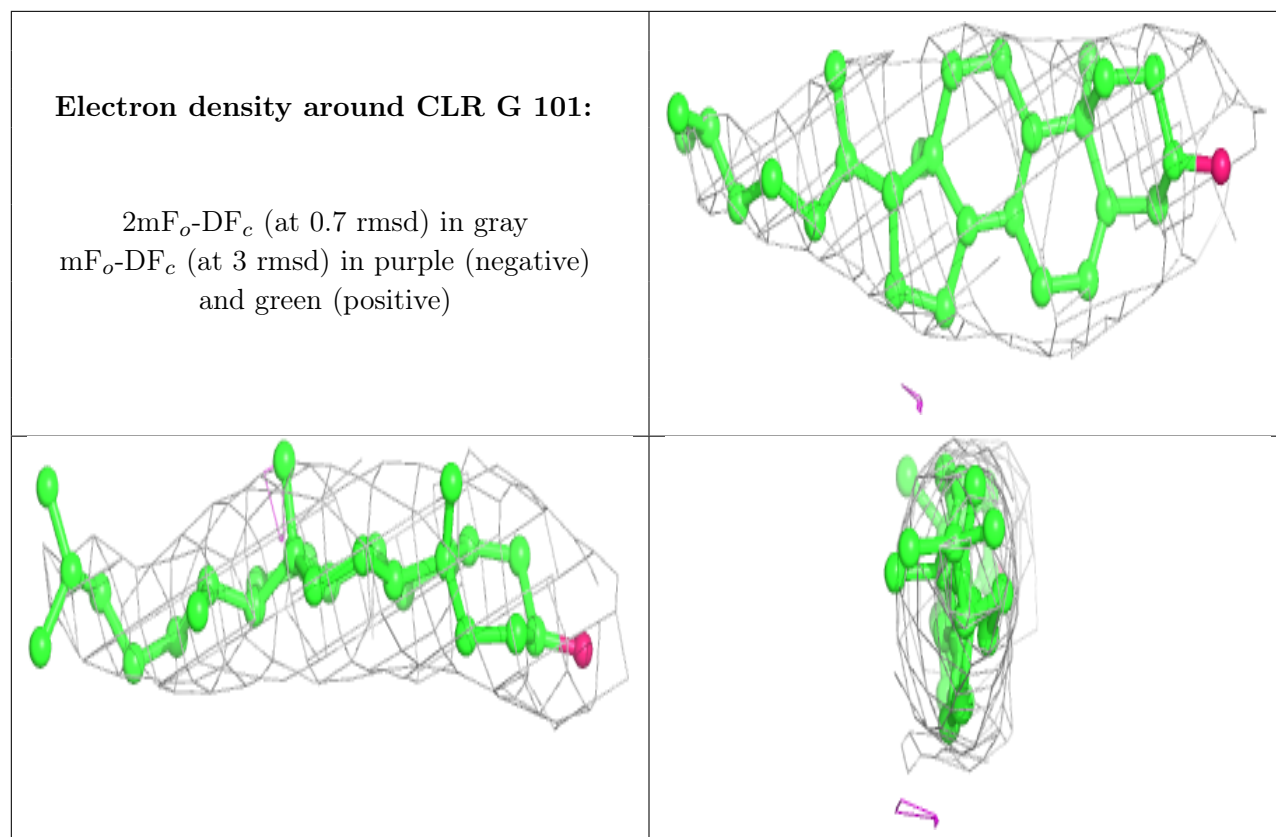
**Electron density around CLR E 101:**

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

**Electron density around PCW C 1107:**

$2mF_o-DF_c$  (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.