wwPDB X-ray Structure Validation Summary Report (i)

Aug 30, 2023 – 06:49 AM EDT

PDB ID	:	3N8Y
Title	:	Structure of Aspirin Acetylated Cyclooxygenase-1 in Complex with Diclofenac
Authors	:	Sidhu, R.S.
Deposited on	:	2010-05-28
Resolution	:	2.60 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) were used in the production of this report:

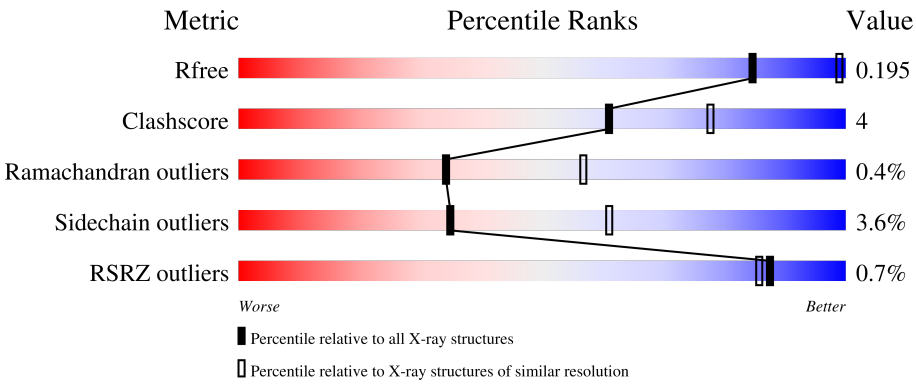
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	553	% 89% 10% •
2	B	553	% 89% 10%
3	C	4	50% 50%
3	F	4	100%
4	D	5	100%

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Mol	Chain	Length	Quality of chain
5	E	2	100%
6	G	4	25% 75%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BMA	C	4	-	-	-	X
8	DIF	B	585[B]	-	-	X	-

2 Entry composition (i)

There are 11 unique types of molecules in this entry. The entry contains 9465 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 Prostaglandin G/H synthase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	553	4427	2870	742	787	28	0	5	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	92	LEU	MET	conflict	UNP P05979

- Molecule 2 is a protein called Prostaglandin G/H synthase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	553	4443	2887	740	789	27	3	4	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	92	LEU	MET	conflict	UNP P05979
B	530	OAS	SER	microheterogeneity	UNP P05979

- Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-6)-beta-D-mannopyranose -(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucop yranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	4	50	28	2	20	0	0	0

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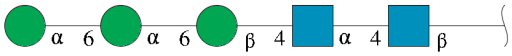
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	F	4	50	28	2	20	0	0	0

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	D	5	61	34	2	25	0	0	0

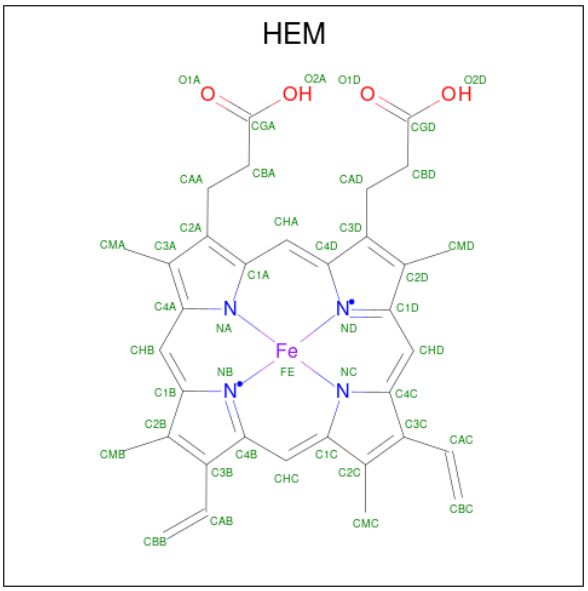
- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

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

- Molecule 6 is an oligosaccharide called beta-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

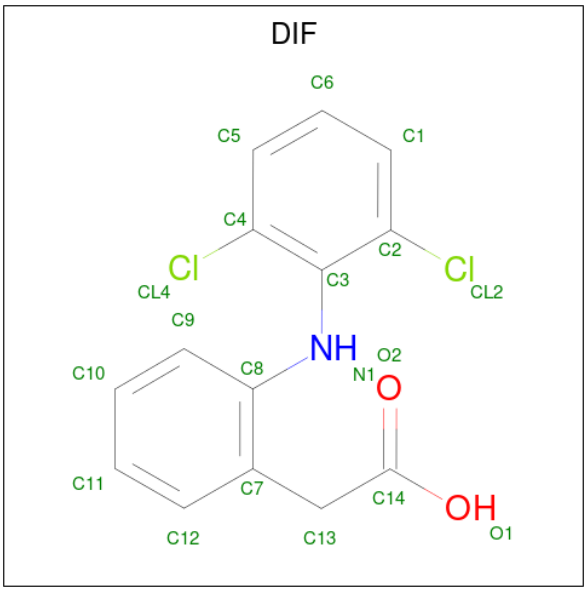
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
6	G	4	50	28	2	20	0	0	0

- Molecule 7 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
7	A	1	43	34	1	4	4	0	0
			Total	C	Fe	N	O		
7	B	1	43	34	1	4	4	0	0

- Molecule 8 is 2-[2,6-DICHLOROPHENYL)AMINO]BENZENEACETIC ACID (three-letter code: DIF) (formula: $C_{14}H_{11}Cl_2NO_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Cl	N	O		
8	A	1	19	14	2	1	2	0	0

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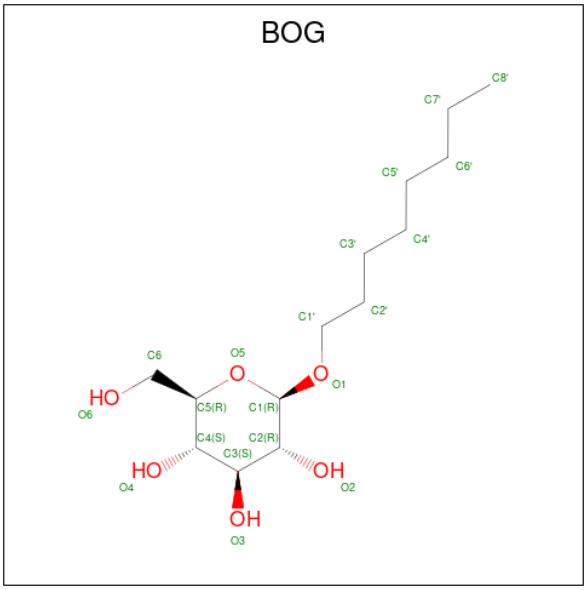
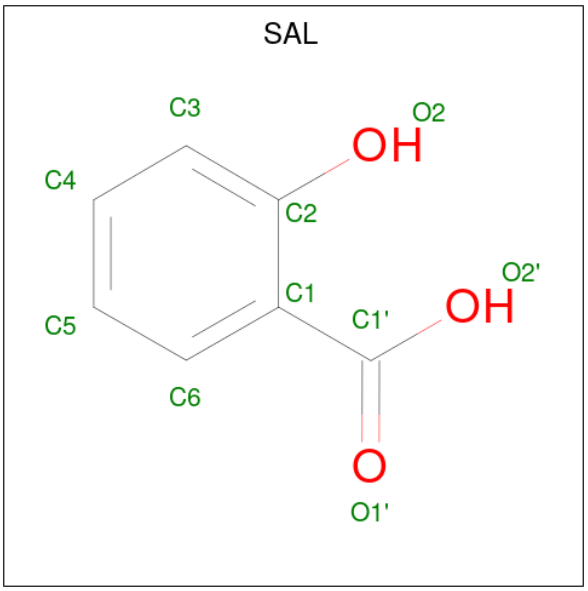
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Cl	N	O		
8	B	1	19	14	2	1	2	0	1

- Molecule 9 is octyl beta-D-glucopyranoside (three-letter code: BOG) (formula: $C_{14}H_{28}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
9	A	1	20	14	6	0	0
			Total	C	O		
9	A	1	20	14	6	0	0
			Total	C	O		
9	B	1	20	14	6	0	0
			Total	C	O		
9	B	1	20	14	6	0	0

- Molecule 10 is 2-HYDROXYBENZOIC ACID (three-letter code: SAL) (formula: $C_7H_6O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	B	1	Total C O 10 7 3	0	1

- Molecule 11 is water.

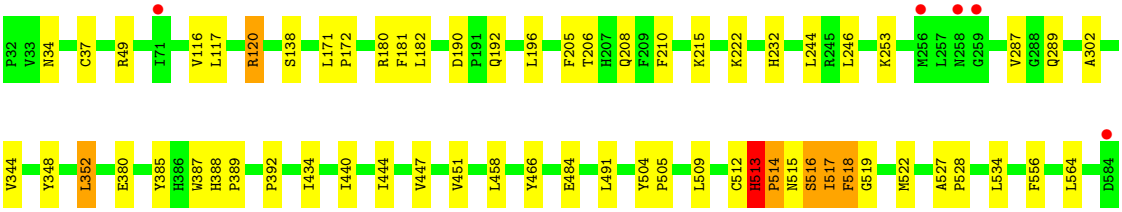
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	74	Total O 74 74	0	0
11	B	68	Total O 68 68	0	0

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.

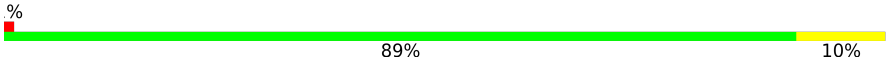
• Molecule 1: Prostaglandin G/H synthase 1

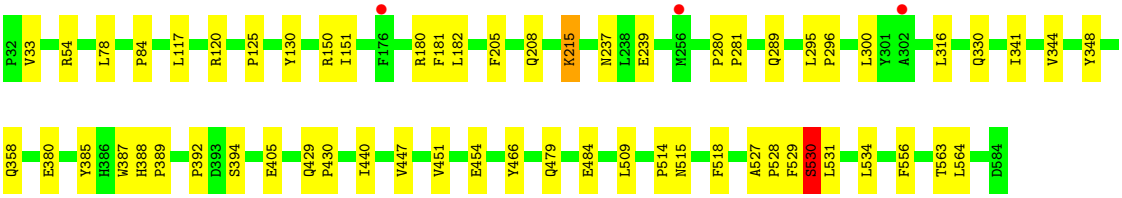
Chain A:



• Molecule 2: Prostaglandin G/H synthase 1

Chain B:



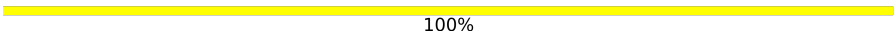


• Molecule 3: beta-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:

• Molecule 3: beta-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:



• Molecule 4: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D: 100%

NAG1
NDG2
BMA3
MAN4
MAN5

• Molecule 5: 2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E: 100%

NAG1
NDG2

• Molecule 6: beta-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G: 25% 75%

NAG1
NDG2
BMA3
BMA4

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α, β, γ	182.33Å 182.33Å 103.11Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.84 – 2.60 29.84 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.8 (29.84-2.60) 99.0 (29.84-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.16 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.184 , 0.200 0.183 , 0.195	Depositor DCC
R_{free} test set	1597 reflections (2.69%)	wwPDB-VP
Wilson B-factor (Å^2)	48.7	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 26.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.057 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.501 for H, K, L 0.499 for -H-K, K, -L	Depositor
Outliers	0 of 59356 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	9465	wwPDB-VP
Average B, all atoms (Å^2)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.37% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: HEM, SAL, BMA, BOG, MAN, DIF, OAS, NDG, NAG

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	0.32	0/4572	0.47	0/6228
2	B	0.33	0/4569	0.47	1/6218 (0.0%)
All	All	0.32	0/9141	0.47	1/12446 (0.0%)

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
2	B	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	479	GLN	CG-CD-OE1	-5.19	111.22	121.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	529[B]	PHE	Mainchain
2	B	530[A]	OAS	Mainchain
2	B	530[B]	OAS	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	4427	0	4236	41	0
2	B	4443	0	4258	33	0
3	C	50	0	43	0	0
3	F	50	0	43	0	0
4	D	61	0	51	0	0
5	E	28	0	24	0	0
6	G	50	0	42	0	0
7	A	43	0	30	2	0
7	B	43	0	30	1	0
8	A	19	0	10	0	0
8	B	19	0	10	14	0
9	A	40	0	56	0	0
9	B	40	0	56	1	0
10	B	10	0	4	0	0
11	A	74	0	0	0	0
11	B	68	0	0	0	0
All	All	9465	0	8893	81	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 4.

The worst 5 of 81 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)
2:B:530[B]:OAS:OG	8:B:585[B]:DIF:O2	1.64	1.15
1:A:513[A]:HIS:HB2	1:A:514[A]:PRO:CD	1.86	1.05
1:A:513[A]:HIS:CB	1:A:514[A]:PRO:HD2	1.88	1.02
1:A:513[A]:HIS:HB2	1:A:514[A]:PRO:HD2	1.01	0.99
8:B:585[B]:DIF:O2	8:B:585[B]:DIF:N1	2.01	0.94

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	556/553 (100%)	534 (96%)	16 (3%)	6 (1%)	14	30
2	B	553/553 (100%)	533 (96%)	19 (3%)	1 (0%)	47	71
All	All	1109/1106 (100%)	1067 (96%)	35 (3%)	7 (1%)	34	47

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	513[A]	HIS
1	A	513[B]	HIS
2	B	33	VAL
1	A	514[A]	PRO
1	A	514[B]	PRO

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	471/488 (96%)	452 (96%)	19 (4%)	31	57
2	B	471/487 (97%)	455 (97%)	16 (3%)	37	63
All	All	942/975 (97%)	907 (96%)	35 (4%)	35	60

5 of 35 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	385	TYR

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Mol	Chain	Res	Type
2	B	454	GLU
2	B	556	PHE
1	A	513[A]	HIS
1	A	484	GLU

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

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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	$\#\|Z\| > 2$
2	OAS	B	530[A]	2	7,8,9	2.51	3 (42%)	5,9,11	2.20	1 (20%)
2	OAS	B	530[B]	2	4,5,9	0.76	0	0,5,11	-	-

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	OAS	B	530[A]	2	-	2/5/7/9	-
2	OAS	B	530[B]	2	-	2/2/4/9	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	530[A]	OAS	C2A-C1A	-4.89	1.32	1.49
2	B	530[A]	OAS	OAC-C1A	3.46	1.33	1.20
2	B	530[A]	OAS	OG-C1A	2.62	1.46	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	530[A]	OAS	OG-CB-CA	4.06	119.96	108.48

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	530[B]	OAS	N-CA-CB-OG
2	B	530[B]	OAS	C-CA-CB-OG
2	B	530[A]	OAS	OAC-C1A-OG-CB
2	B	530[A]	OAS	C2A-C1A-OG-CB

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	530[B]	OAS	3	0

5.5 Carbohydrates (i)

19 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	$\#\|Z\| > 2$
3	NAG	C	1	3,1	14,14,15	0.49	0	17,19,21	0.72	0
3	NAG	C	2	3	14,14,15	0.45	0	17,19,21	0.94	0
3	BMA	C	3	3	11,11,12	0.55	0	15,15,17	1.84	2 (13%)
3	BMA	C	4	3	11,11,12	0.53	0	15,15,17	1.82	3 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	NAG	D	1	4,1	14,14,15	0.57	0	17,19,21	0.90	1 (5%)
4	NDG	D	2	4	14,14,15	0.55	0	17,19,21	1.83	4 (23%)
4	BMA	D	3	4	11,11,12	0.56	0	15,15,17	1.60	1 (6%)
4	MAN	D	4	4	11,11,12	0.58	0	15,15,17	1.33	2 (13%)
4	MAN	D	5	4	11,11,12	0.59	0	15,15,17	0.88	1 (6%)
5	NAG	E	1	5,2	14,14,15	1.48	1 (7%)	17,19,21	1.10	1 (5%)
5	NDG	E	2	5	14,14,15	0.50	0	17,19,21	0.99	1 (5%)
3	NAG	F	1	3,2	14,14,15	0.54	0	17,19,21	0.84	1 (5%)
3	NAG	F	2	3	14,14,15	0.47	0	17,19,21	1.00	2 (11%)
3	BMA	F	3	3	11,11,12	0.59	0	15,15,17	1.27	2 (13%)
3	BMA	F	4	3	11,11,12	0.61	0	15,15,17	1.15	2 (13%)
6	NAG	G	1	6,2	14,14,15	0.55	0	17,19,21	1.04	1 (5%)
6	NDG	G	2	6	14,14,15	0.75	1 (7%)	17,19,21	1.87	4 (23%)
6	BMA	G	3	6	11,11,12	0.61	0	15,15,17	0.87	0
6	BMA	G	4	6	11,11,12	0.53	0	15,15,17	1.81	3 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	C	2	3	-	0/6/23/26	0/1/1/1
3	BMA	C	3	3	-	1/2/19/22	1/1/1/1
3	BMA	C	4	3	-	1/2/19/22	1/1/1/1
4	NAG	D	1	4,1	-	0/6/23/26	0/1/1/1
4	NDG	D	2	4	-	2/6/23/26	0/1/1/1
4	BMA	D	3	4	-	2/2/19/22	0/1/1/1
4	MAN	D	4	4	-	2/2/19/22	0/1/1/1
4	MAN	D	5	4	-	2/2/19/22	0/1/1/1
5	NAG	E	1	5,2	-	2/6/23/26	0/1/1/1
5	NDG	E	2	5	-	1/6/23/26	0/1/1/1
3	NAG	F	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	1/2/19/22	0/1/1/1
3	BMA	F	4	3	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	G	1	6,2	-	2/6/23/26	0/1/1/1
6	NDG	G	2	6	-	3/6/23/26	0/1/1/1
6	BMA	G	3	6	-	2/2/19/22	0/1/1/1
6	BMA	G	4	6	-	2/2/19/22	1/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	1	NAG	O5-C1	-5.31	1.35	1.43
6	G	2	NDG	C1-C2	2.24	1.55	1.52

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	3	BMA	C1-O5-C5	6.15	120.52	112.19
6	G	4	BMA	C1-O5-C5	5.49	119.63	112.19
3	C	4	BMA	C1-O5-C5	5.48	119.62	112.19
6	G	2	NDG	C1-O5-C5	4.99	118.95	112.19
4	D	3	BMA	C1-O5-C5	4.89	118.82	112.19

There are no chirality outliers.

5 of 31 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	2	NDG	C3-C2-N2-C7
6	G	3	BMA	C4-C5-C6-O6
6	G	3	BMA	O5-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
5	E	1	NAG	O5-C5-C6-O6

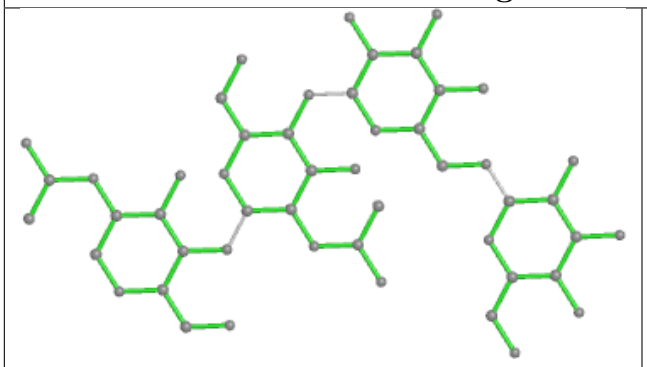
All (3) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	3	BMA	C1-C2-C3-C4-C5-O5
3	C	4	BMA	C1-C2-C3-C4-C5-O5
6	G	4	BMA	C1-C2-C3-C4-C5-O5

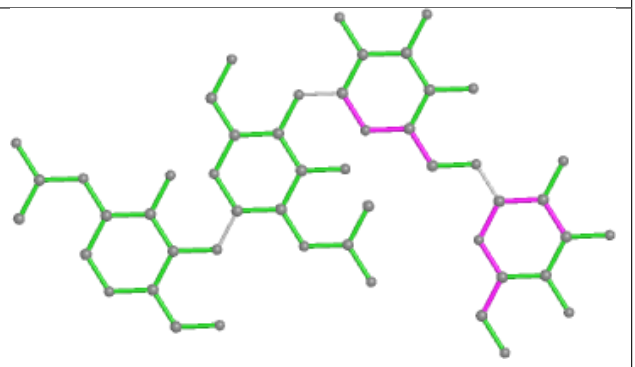
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

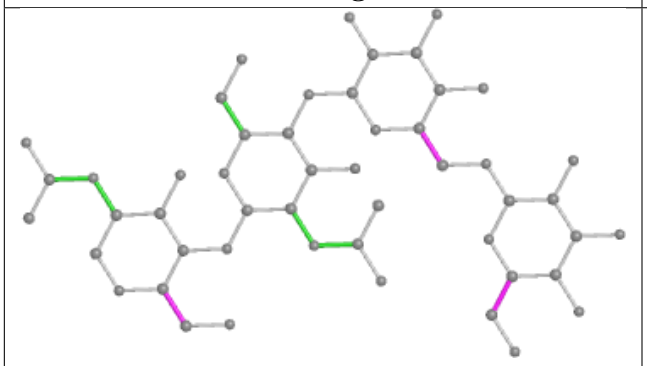
Oligosaccharide Chain C



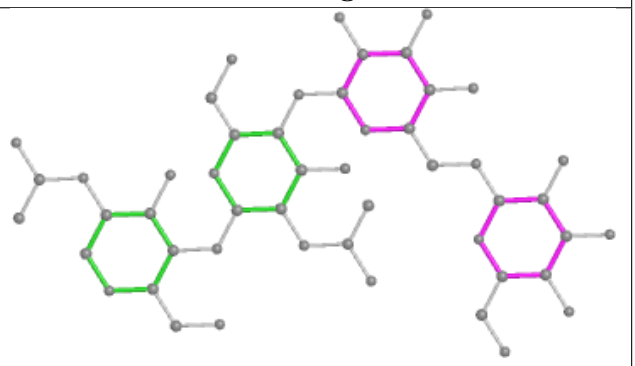
Bond lengths



Bond angles

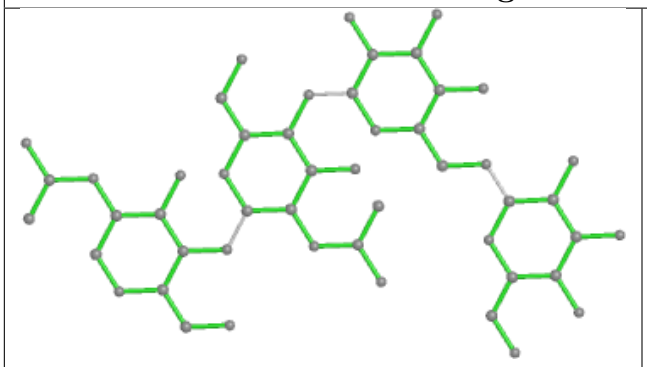


Torsions

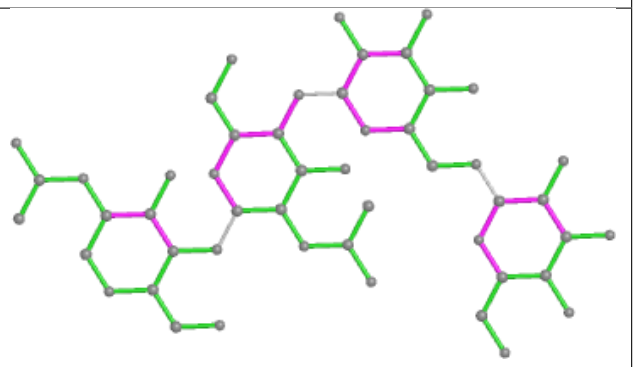


Rings

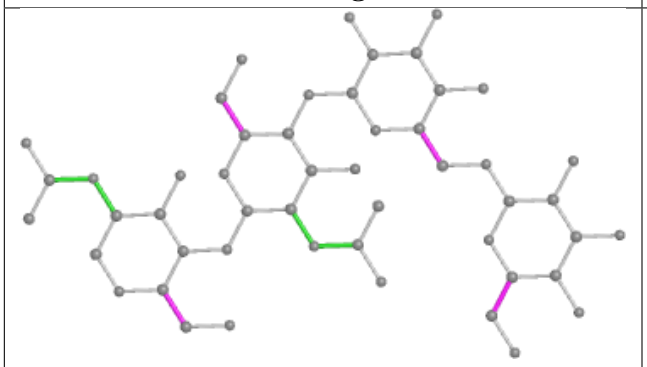
Oligosaccharide Chain F



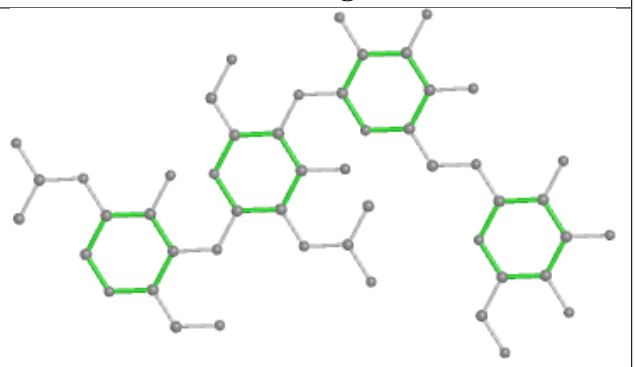
Bond lengths



Bond angles

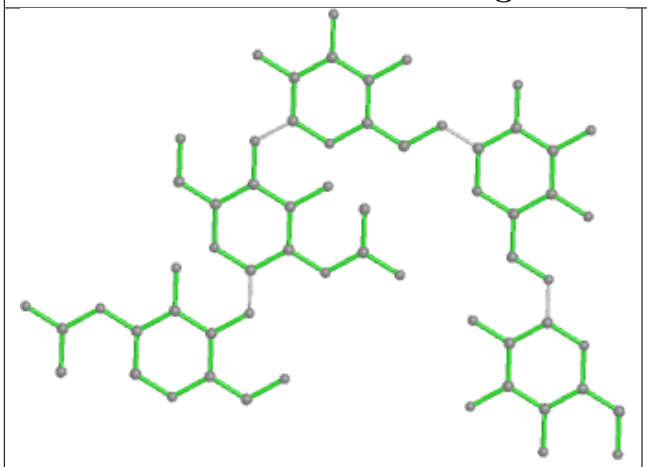


Torsions

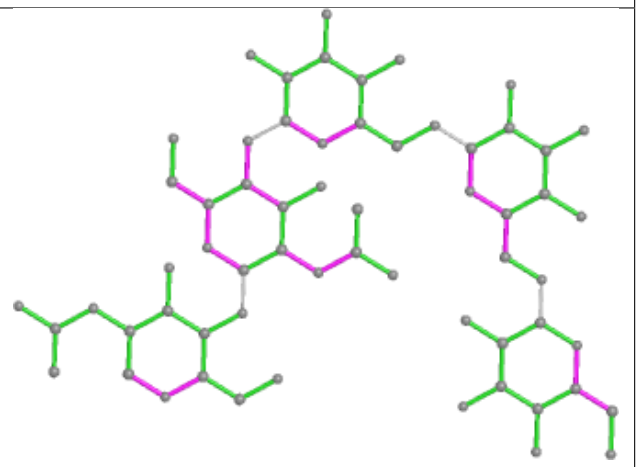


Rings

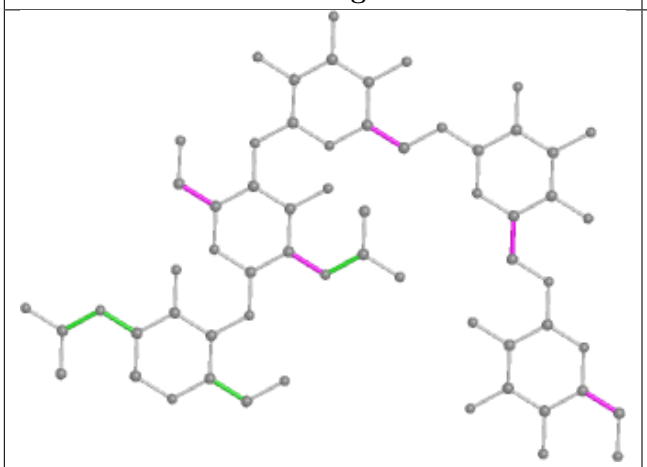
Oligosaccharide Chain D



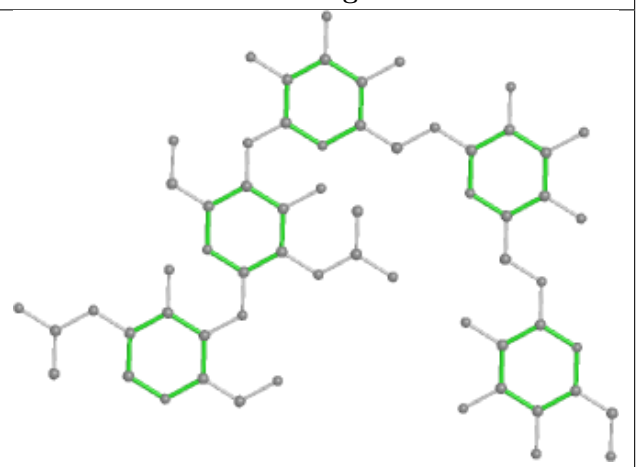
Bond lengths



Bond angles

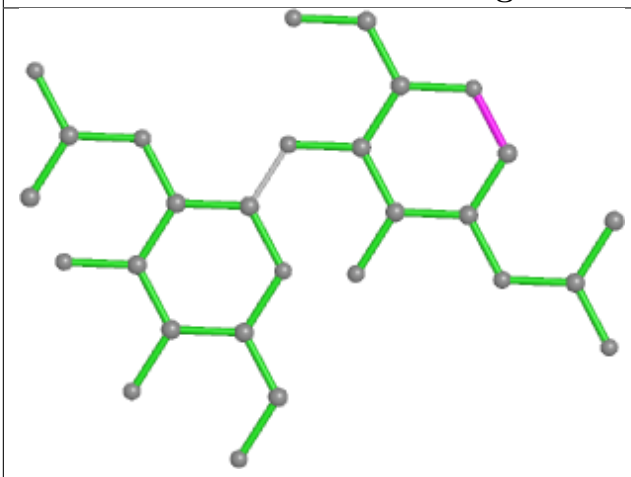


Torsions

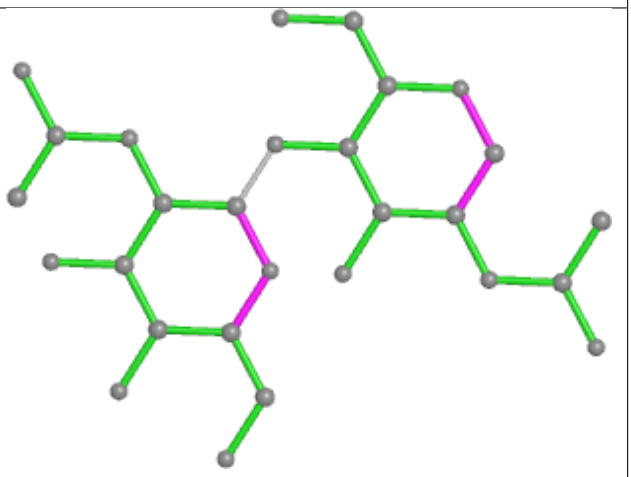


Rings

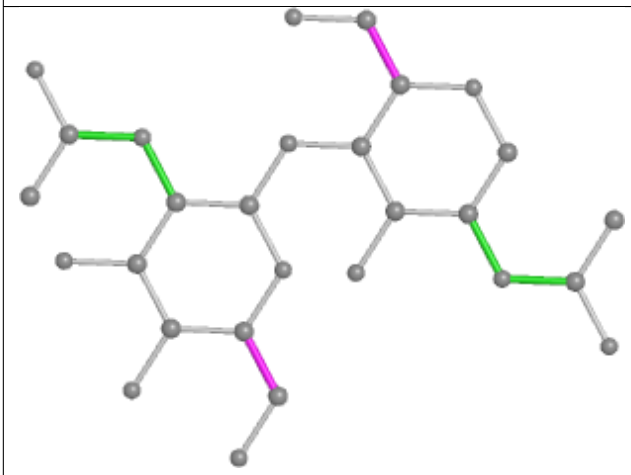
Oligosaccharide Chain E



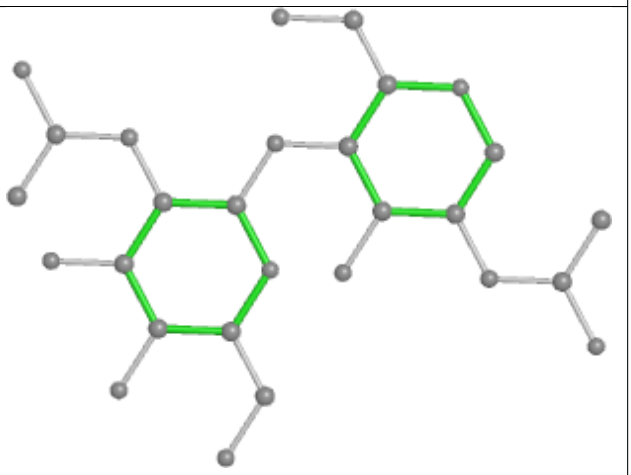
Bond lengths



Bond angles

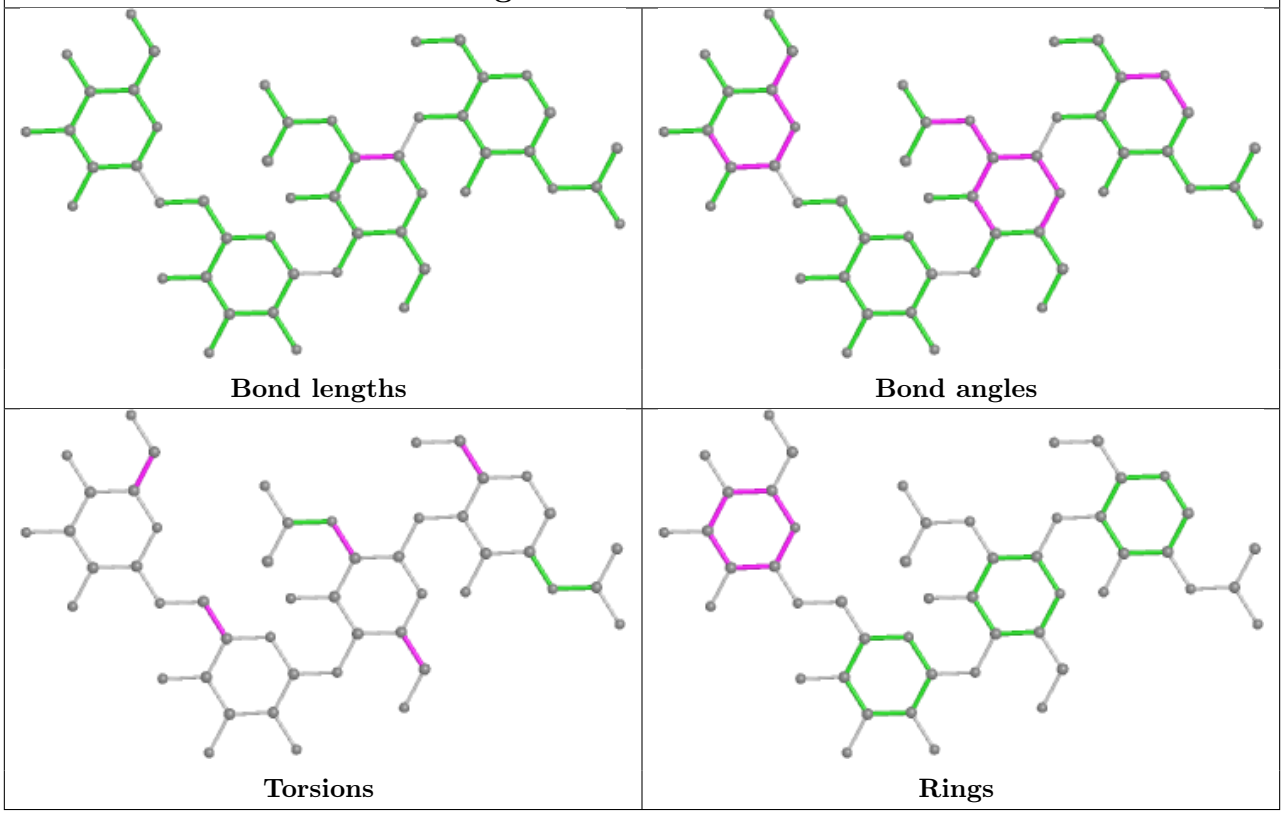


Torsions



Rings

Oligosaccharide Chain G



5.6 Ligand geometry (i)

9 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	$\#\|Z\| > 2$
7	HEM	B	601	2	41,50,50	2.01	9 (21%)	45,82,82	1.61	4 (8%)
9	BOG	B	1750	-	20,20,20	0.53	0	25,25,25	0.58	0
8	DIF	A	701	-	20,20,20	1.06	2 (10%)	27,27,27	1.09	2 (7%)
10	SAL	B	900[A]	-	10,10,10	0.95	1 (10%)	13,13,13	0.72	0
9	BOG	A	754	-	20,20,20	0.54	0	25,25,25	0.58	0
8	DIF	B	585[B]	-	20,20,20	1.13	2 (10%)	27,27,27	1.11	2 (7%)
9	BOG	B	1751	-	20,20,20	0.45	0	25,25,25	0.68	0
7	HEM	A	601	1,11	41,50,50	2.03	8 (19%)	45,82,82	1.74	7 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
9	BOG	A	751	-	20,20,20	0.53	0	25,25,25	1.00	1 (4%)

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
7	HEM	B	601	2	-	4/12/54/54	-
9	BOG	B	1750	-	-	7/11/31/31	0/1/1/1
8	DIF	A	701	-	-	0/8/8/8	0/2/2/2
10	SAL	B	900[A]	-	-	2/4/4/4	0/1/1/1
9	BOG	A	754	-	-	9/11/31/31	0/1/1/1
8	DIF	B	585[B]	-	-	2/8/8/8	0/2/2/2
9	BOG	B	1751	-	-	4/11/31/31	0/1/1/1
7	HEM	A	601	1,11	-	6/12/54/54	-
9	BOG	A	751	-	-	7/11/31/31	0/1/1/1

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	601	HEM	C3D-C2D	8.30	1.54	1.36
7	B	601	HEM	C3D-C2D	8.15	1.54	1.36
7	A	601	HEM	C3C-C2C	-4.28	1.34	1.40
7	B	601	HEM	C3C-C2C	-4.11	1.34	1.40
7	A	601	HEM	C3C-CAC	3.70	1.55	1.47

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	A	601	HEM	C4D-ND-C1D	6.69	111.98	105.07
7	B	601	HEM	C4D-ND-C1D	6.45	111.74	105.07
7	A	601	HEM	C4C-CHD-C1D	3.58	127.28	122.56
7	B	601	HEM	C4C-CHD-C1D	3.19	126.77	122.56
7	A	601	HEM	C1B-NB-C4B	3.05	108.23	105.07

There are no chirality outliers.

5 of 41 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	754	BOG	C2'-C1'-O1-C1
9	A	754	BOG	C4-C5-C6-O6
9	B	1750	BOG	C2-C1-O1-C1'
9	A	754	BOG	O1-C1'-C2'-C3'
9	B	1750	BOG	O5-C1-O1-C1'

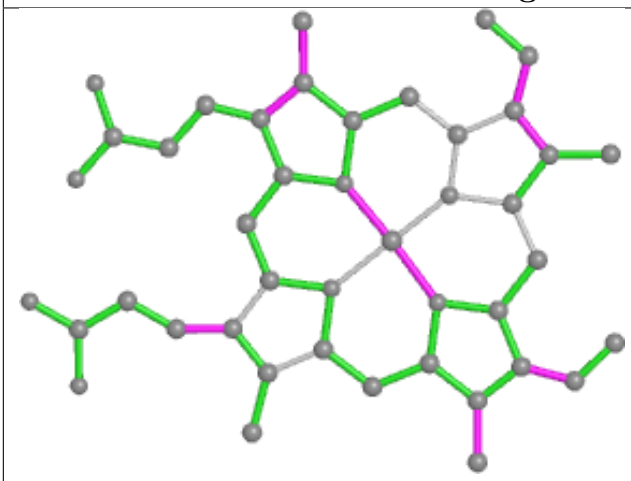
There are no ring outliers.

4 monomers are involved in 18 short contacts:

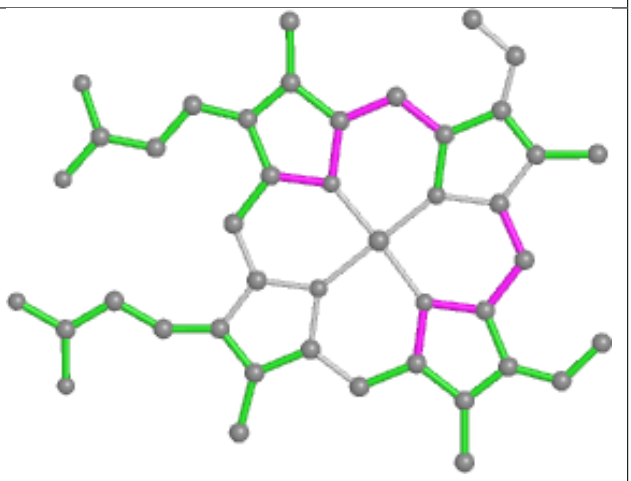
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	601	HEM	1	0
9	B	1750	BOG	1	0
8	B	585[B]	DIF	14	0
7	A	601	HEM	2	0

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 then 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.

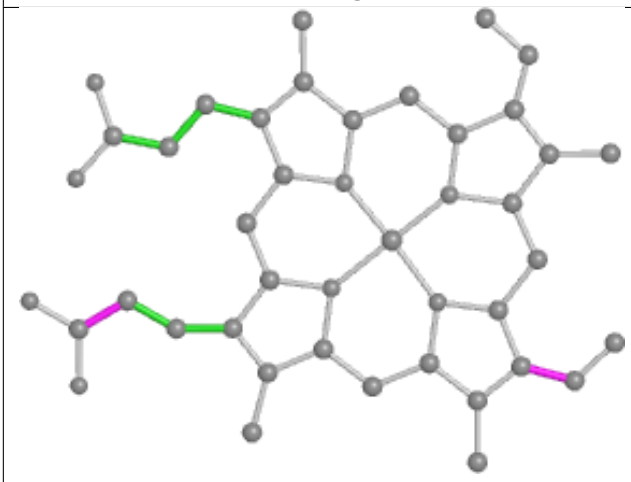
Ligand HEM B 601



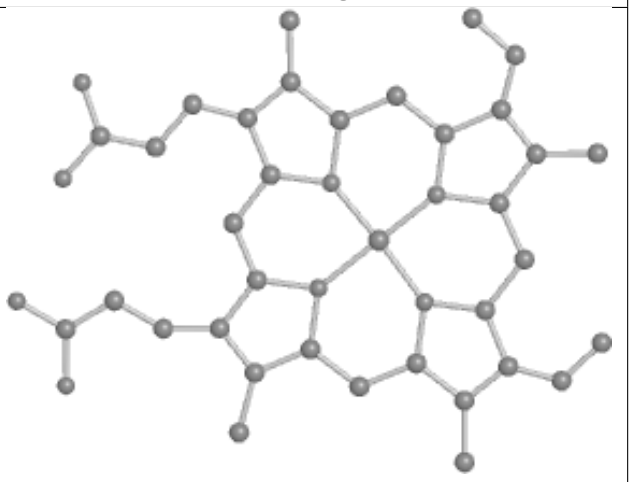
Bond lengths



Bond angles

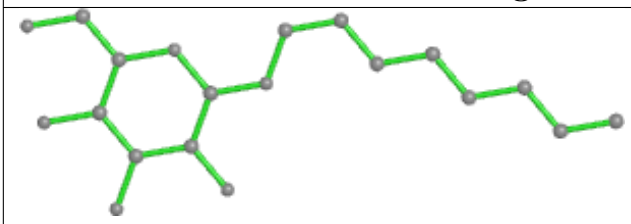


Torsions

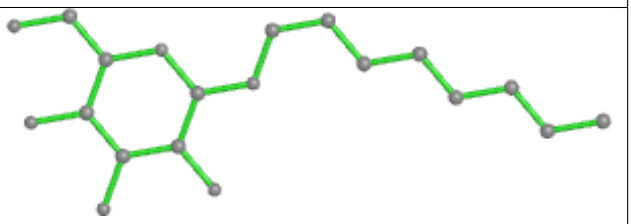


Rings

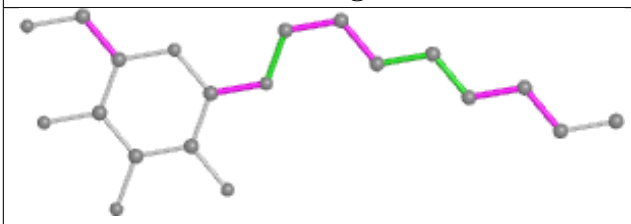
Ligand BOG B 1750



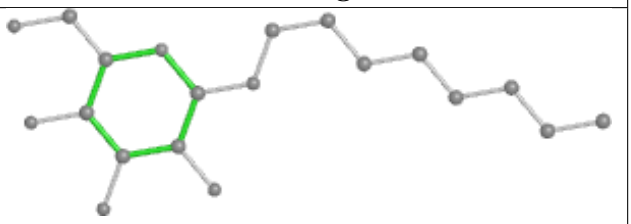
Bond lengths



Bond angles

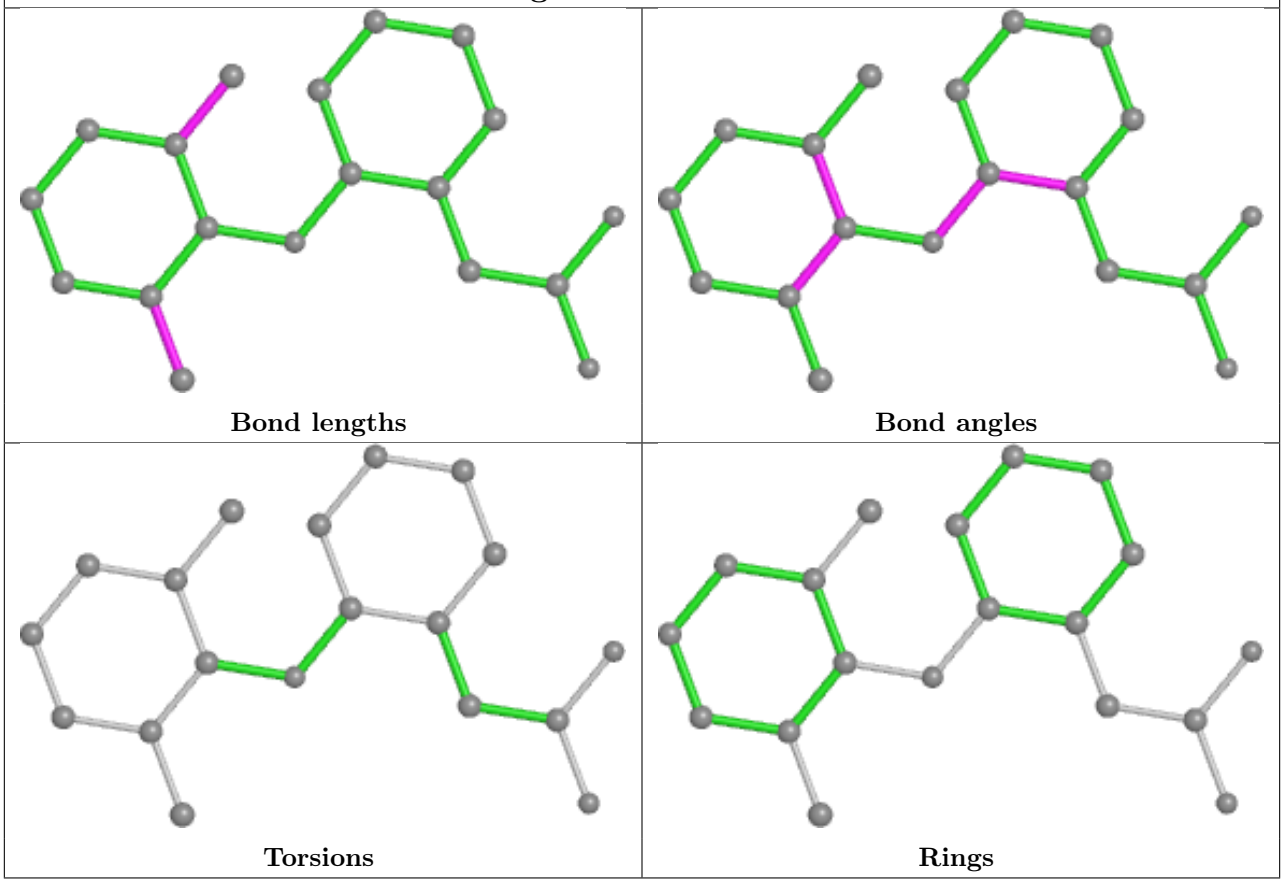


Torsions

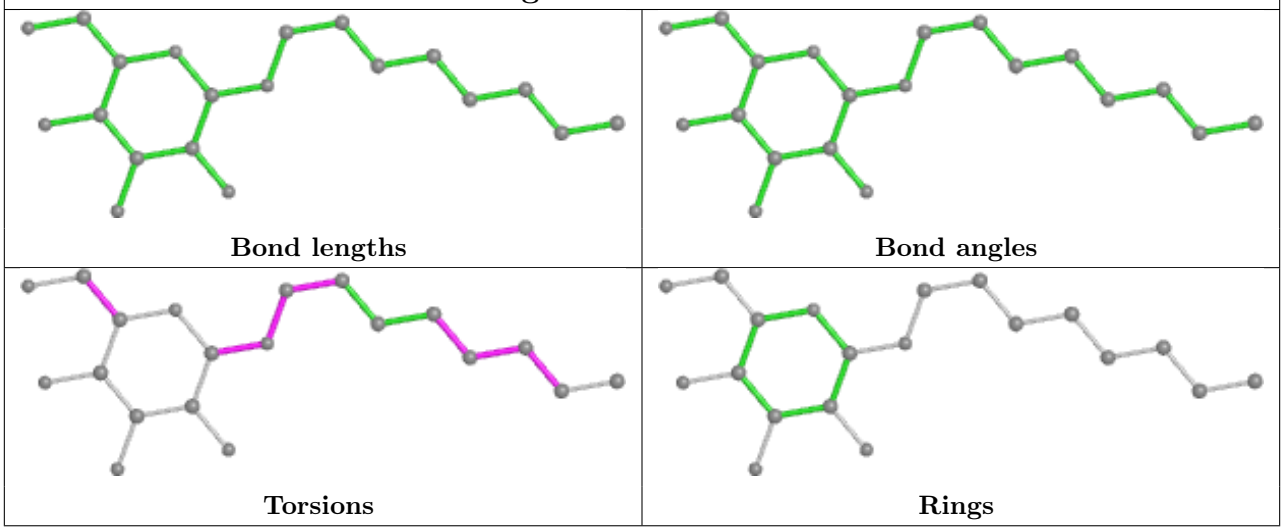


Rings

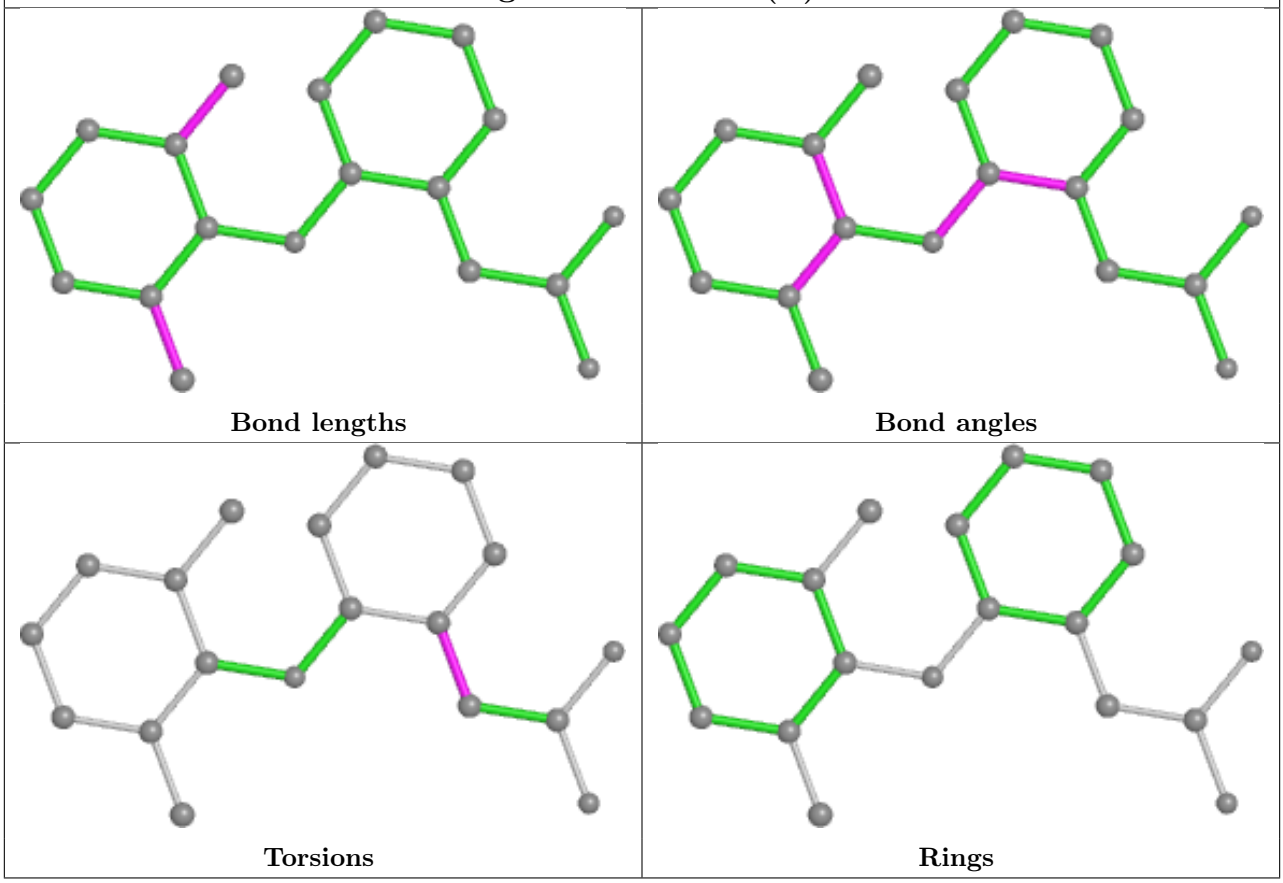
Ligand DIF A 701



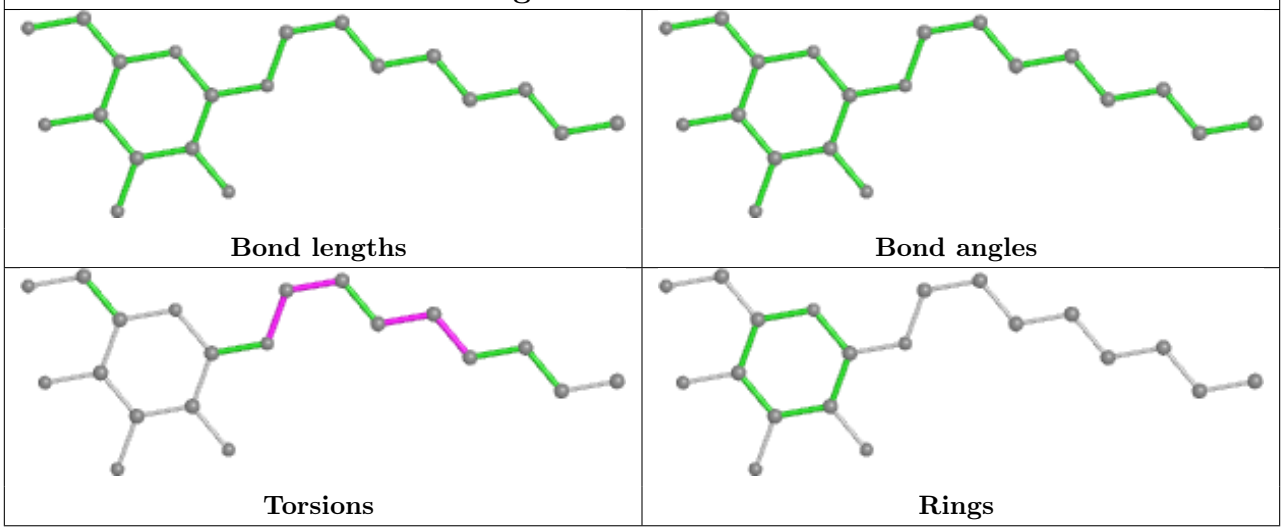
Ligand BOG A 754



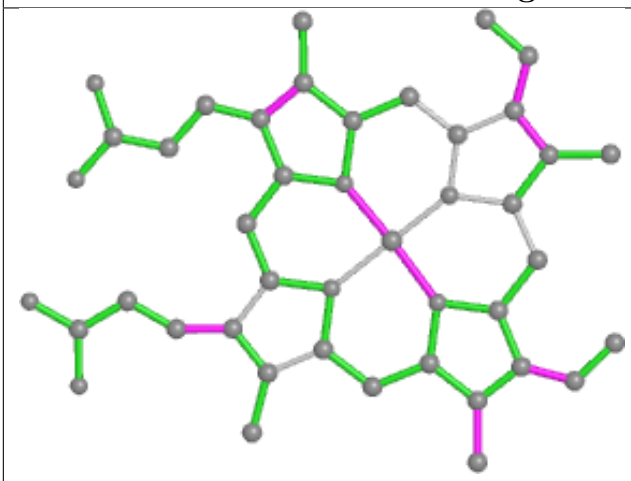
Ligand DIF B 585 (B)



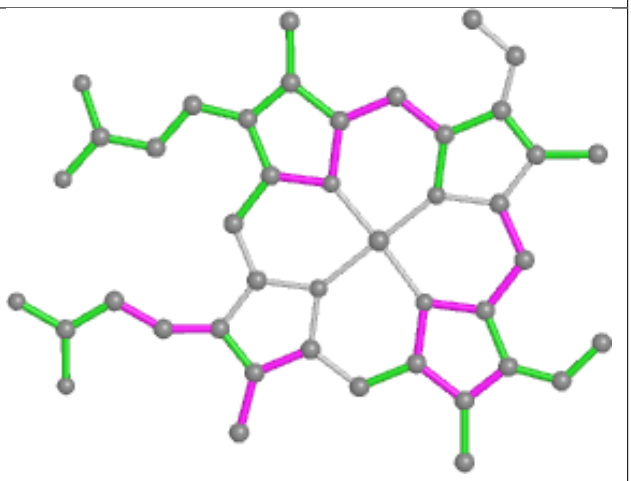
Ligand BOG B 1751



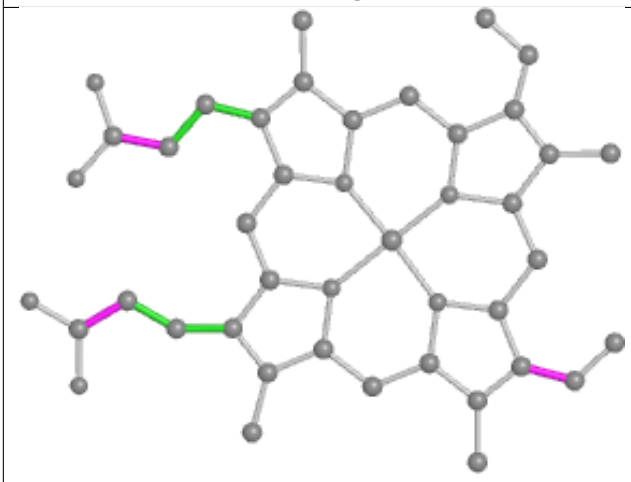
Ligand HEM A 601



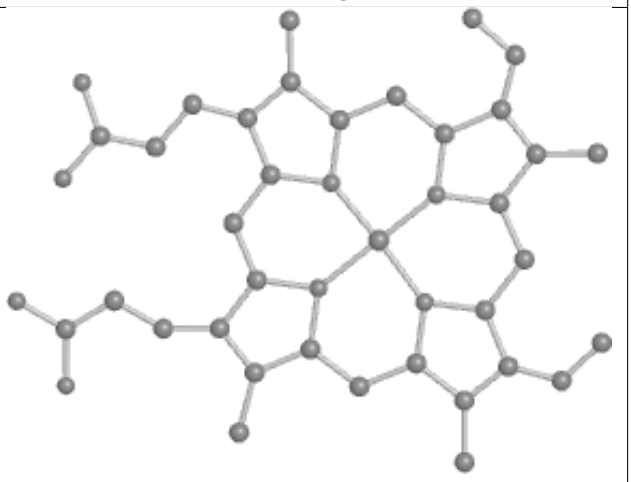
Bond lengths



Bond angles

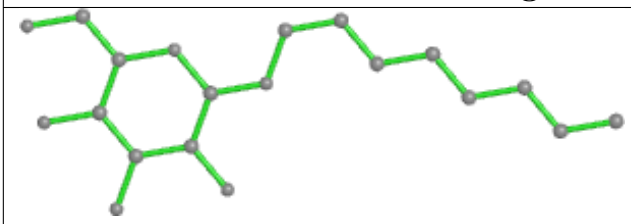


Torsions

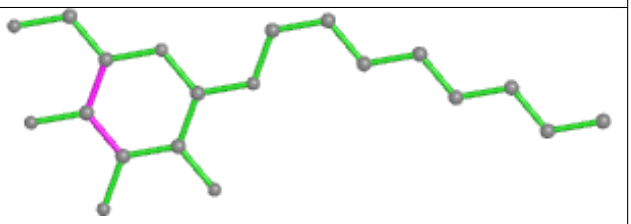


Rings

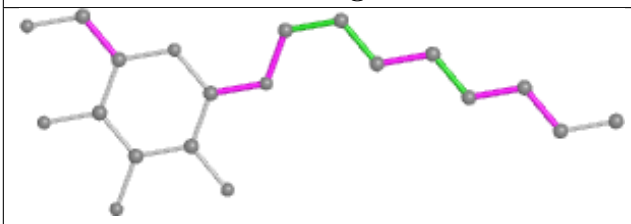
Ligand BOG A 751



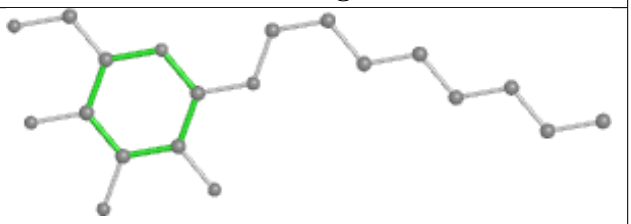
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	553/553 (100%)	-0.47	5 (0%)	84	82	34, 47, 61, 68	1 (0%)
2	B	552/553 (99%)	-0.48	3 (0%)	91	89	34, 48, 61, 69	1 (0%)
All	All	1105/1106 (99%)	-0.48	8 (0%)	87	86	34, 47, 61, 69	2 (0%)

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	256	MET	2.5
2	B	302	ALA	2.4
1	A	584	ASP	2.3
1	A	259	GLY	2.3
1	A	258	ASN	2.2

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, 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	OAS	B	530[A]	9/10	0.99	0.14	37,37,39,39	9
2	OAS	B	530[B]	6/10	0.99	0.14	35,36,36,36	6

6.3 Carbohydrates (i)

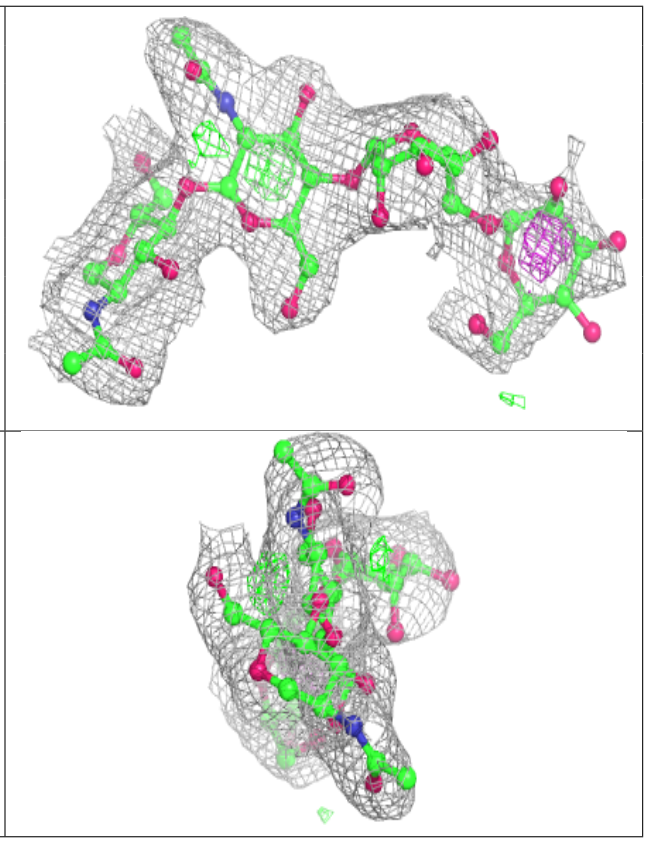
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	BMA	C	4	11/12	0.72	0.41	88,97,116,117	0
4	MAN	D	4	11/12	0.78	0.26	68,69,71,71	0
3	BMA	F	4	11/12	0.82	0.46	91,93,96,97	0
6	BMA	G	4	11/12	0.84	0.37	84,89,98,99	0
3	BMA	F	3	11/12	0.86	0.23	85,86,89,90	0
4	MAN	D	5	11/12	0.87	0.24	69,71,72,73	0
4	NDG	D	2	14/15	0.87	0.20	63,64,65,65	0
3	NAG	F	2	14/15	0.89	0.27	84,85,86,87	0
3	BMA	C	3	11/12	0.89	0.28	67,76,86,87	0
6	NDG	G	2	14/15	0.89	0.25	65,66,68,68	0
4	BMA	D	3	11/12	0.89	0.26	65,66,67,68	0
5	NAG	E	1	14/15	0.90	0.20	58,59,66,73	0
3	NAG	F	1	14/15	0.91	0.14	85,86,87,87	0
5	NDG	E	2	14/15	0.93	0.23	60,66,72,79	0
6	BMA	G	3	11/12	0.93	0.48	71,75,81,83	0
6	NAG	G	1	14/15	0.93	0.15	61,64,67,69	0
3	NAG	C	2	14/15	0.94	0.12	57,59,63,64	0
4	NAG	D	1	14/15	0.94	0.12	58,60,60,62	0
3	NAG	C	1	14/15	0.96	0.08	56,58,61,62	0

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

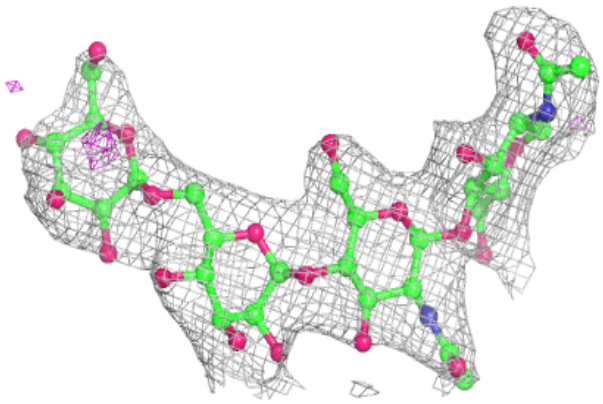
Electron density around Chain C:

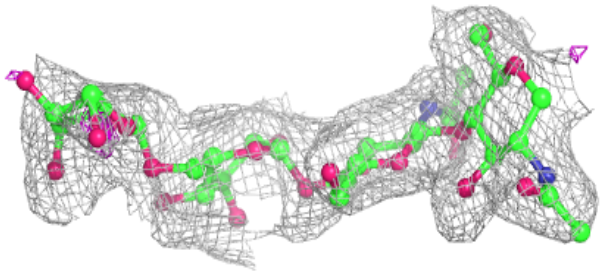
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

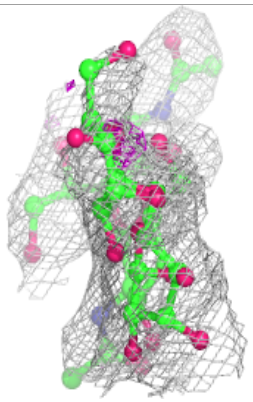


Electron density around Chain F:

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

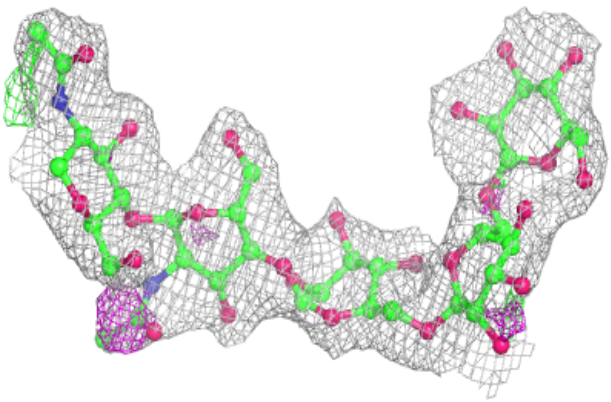


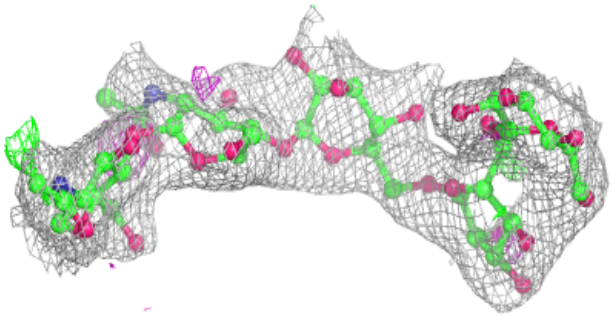


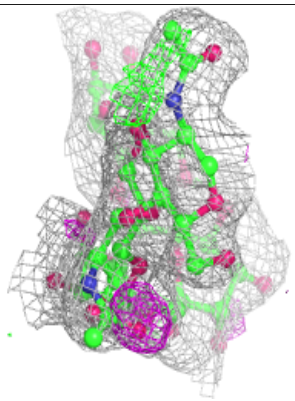


Electron density around Chain D:

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

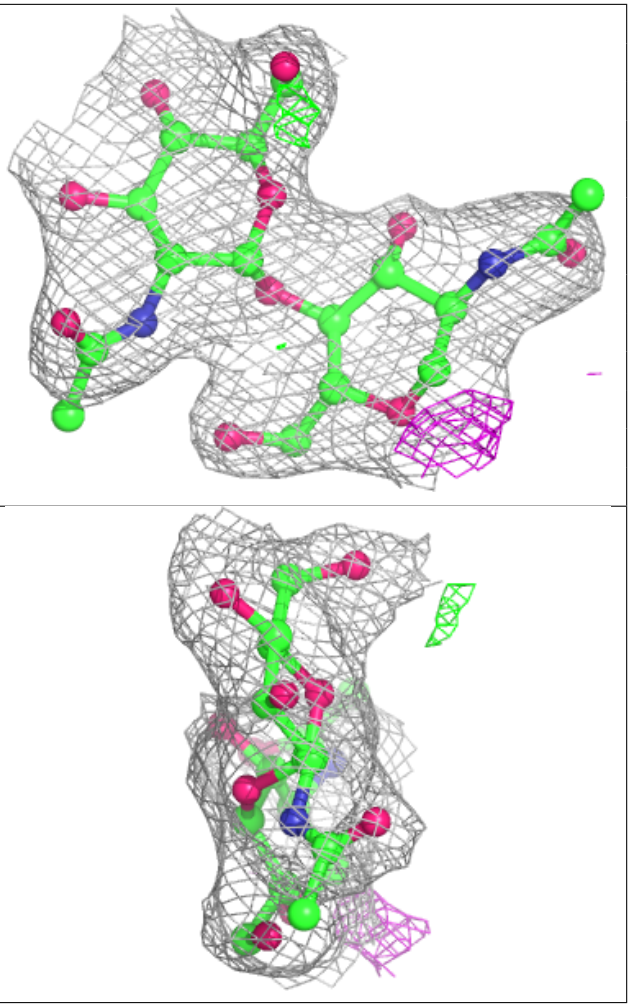






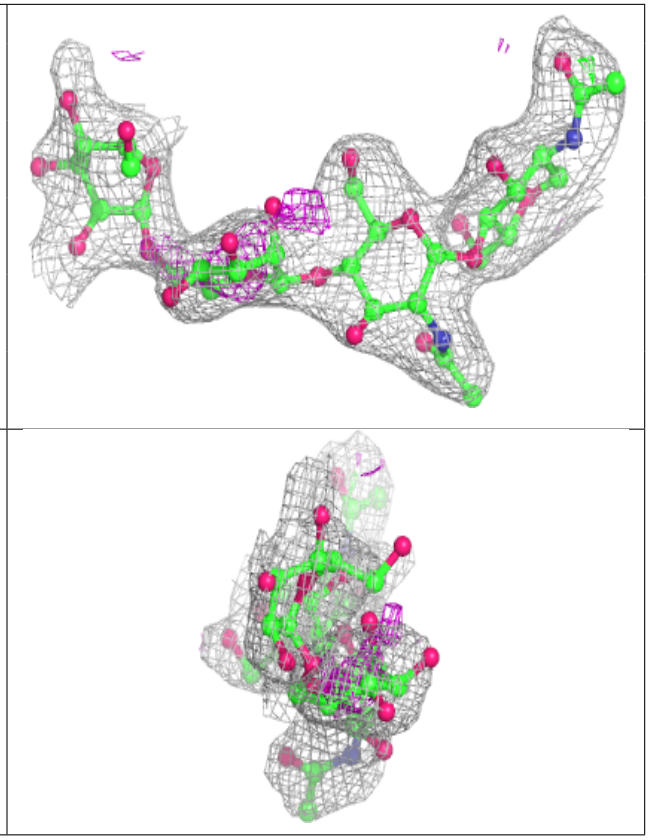
Electron density around Chain E:

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



Electron density around Chain G:

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



6.4 Ligands (i)

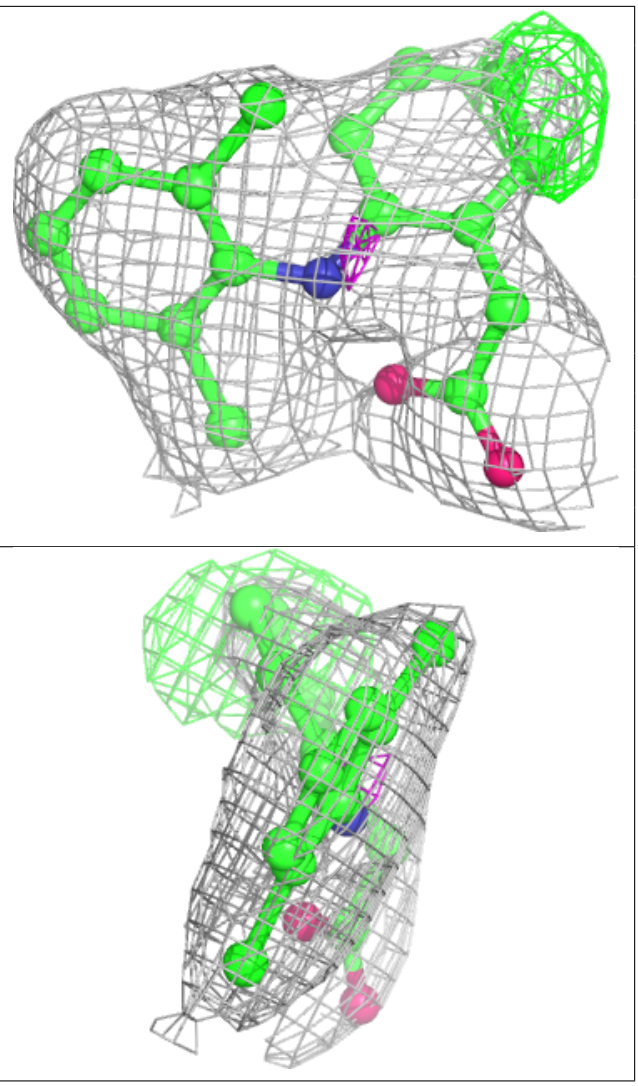
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
8	DIF	B	585[B]	19/19	0.92	0.19	22,23,23,23	19
9	BOG	B	1750	20/20	0.92	0.15	42,43,44,44	0
9	BOG	A	754	20/20	0.93	0.21	46,46,46,46	0
9	BOG	A	751	20/20	0.93	0.16	39,47,64,69	0
7	HEM	A	601	43/43	0.94	0.16	40,45,49,51	0
7	HEM	B	601	43/43	0.95	0.16	41,45,50,51	0
10	SAL	B	900[A]	10/10	0.95	0.12	34,34,35,35	10
8	DIF	A	701	19/19	0.96	0.11	46,47,48,50	0
9	BOG	B	1751	20/20	0.97	0.11	35,37,38,38	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.

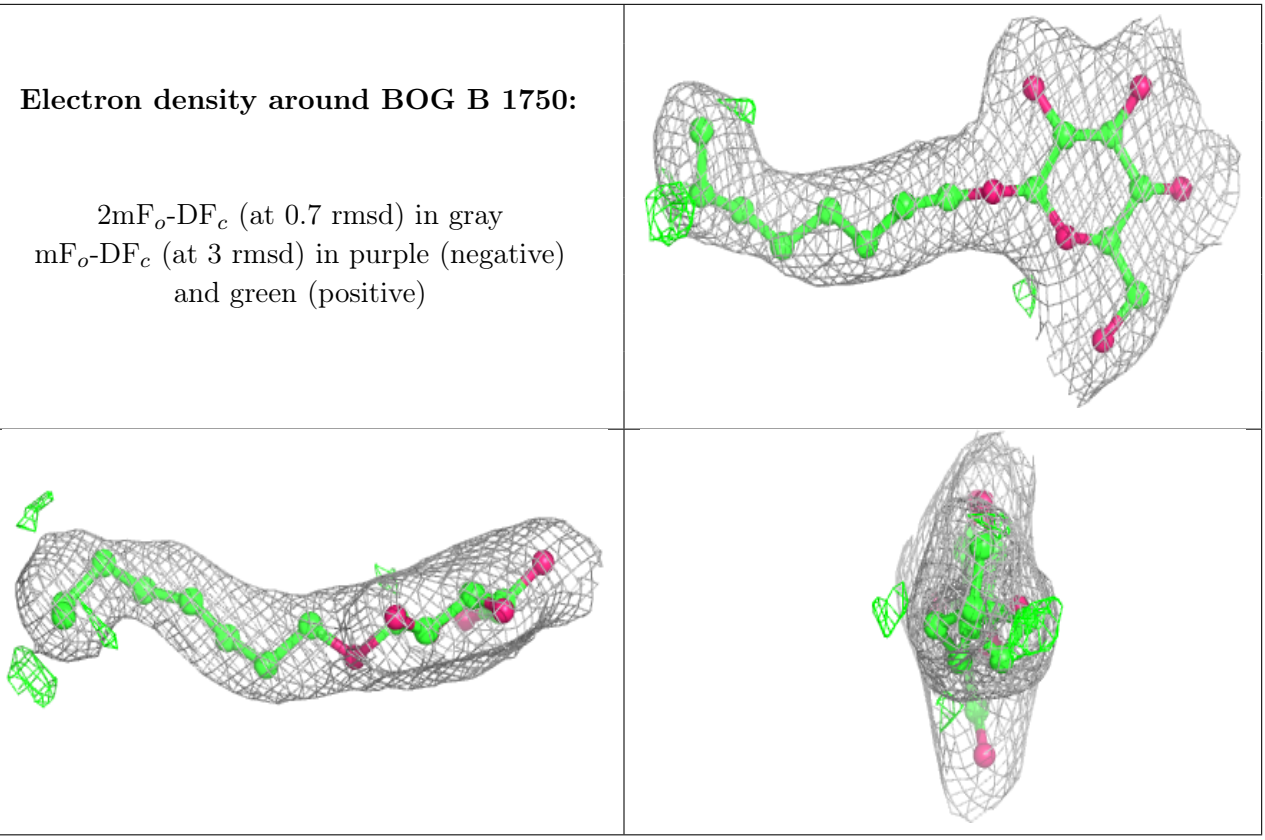
Electron density around DIF B 585 (B):

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



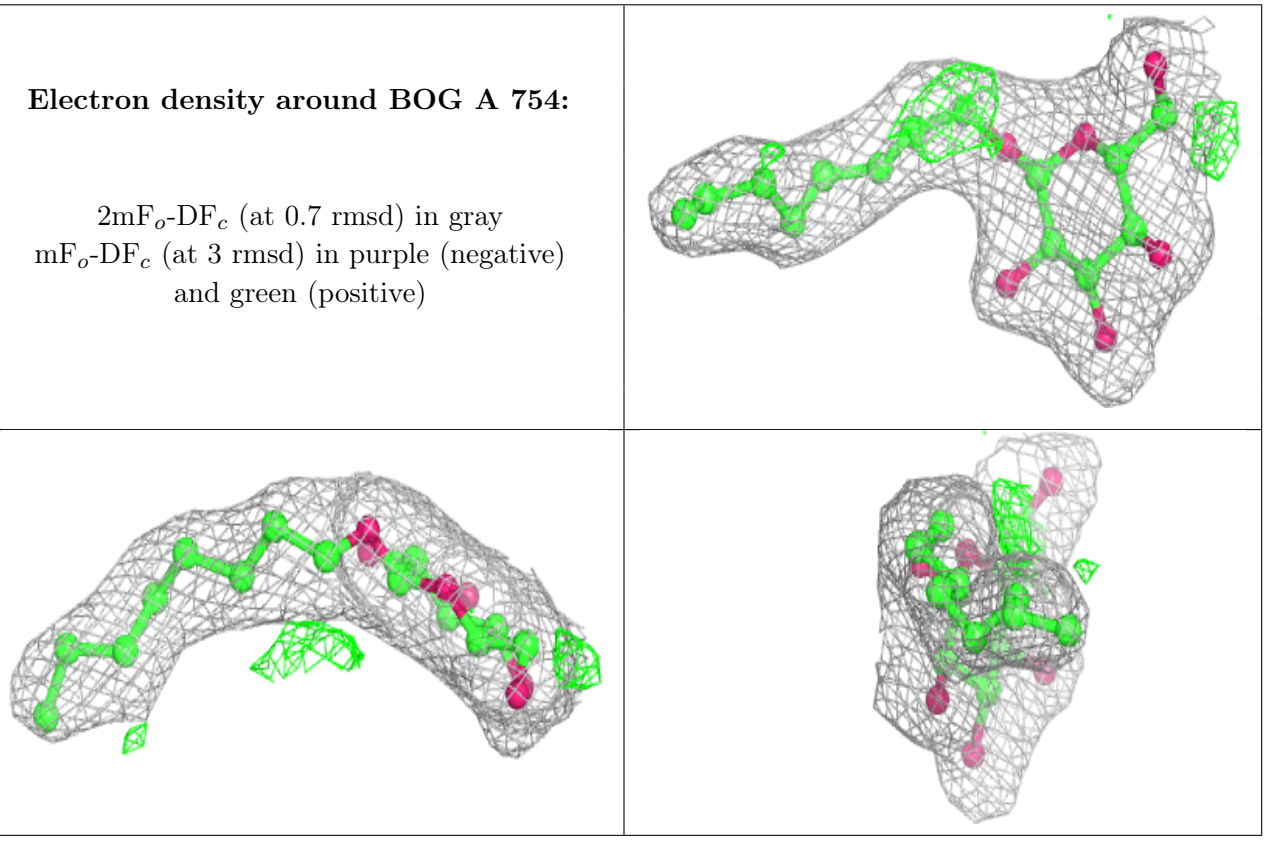
Electron density around BOG B 1750:

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



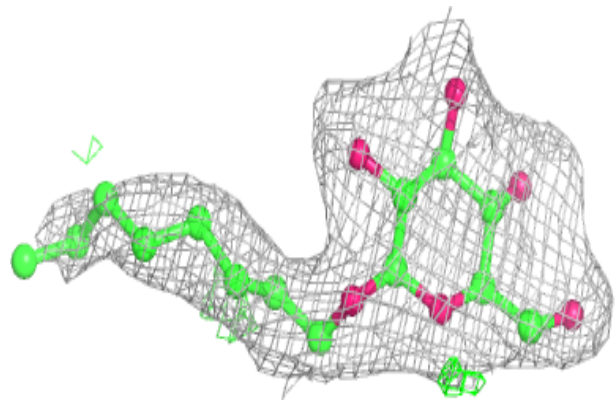
Electron density around BOG A 754:

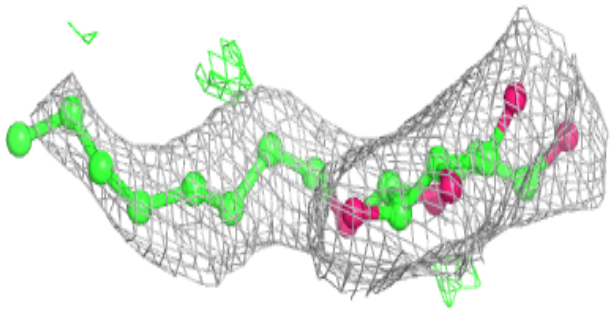
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

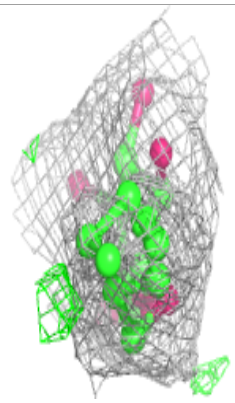


Electron density around BOG A 751:

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

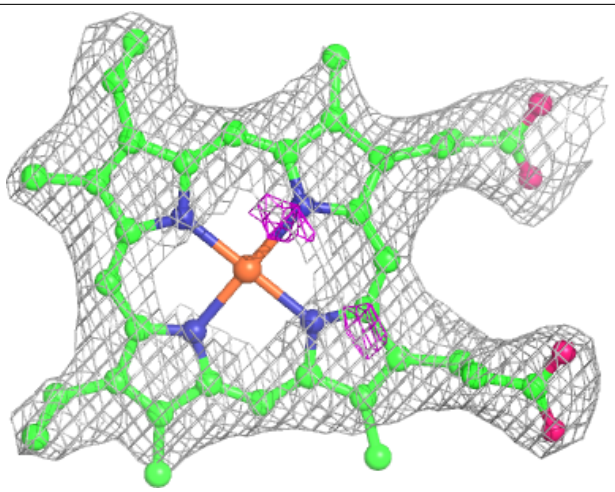


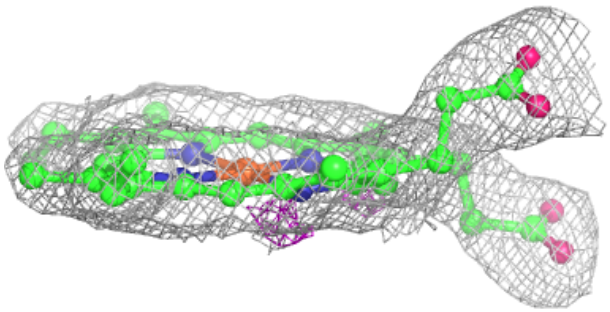


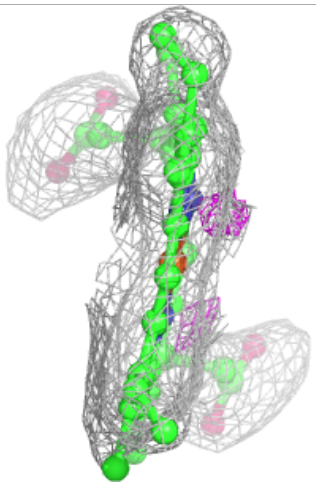


Electron density around HEM A 601:

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

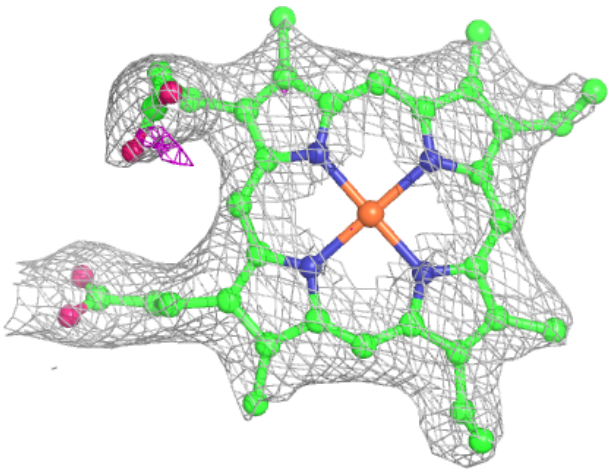


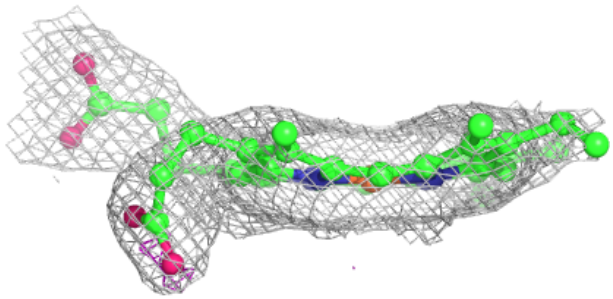


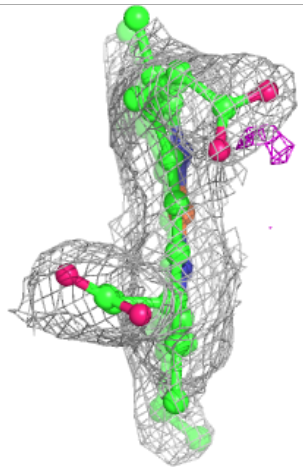


Electron density around HEM B 601:

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

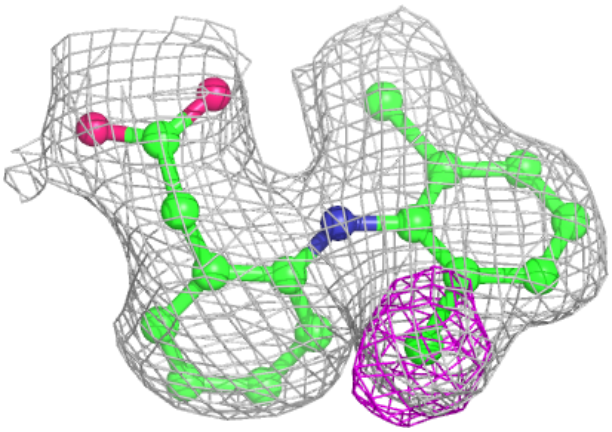


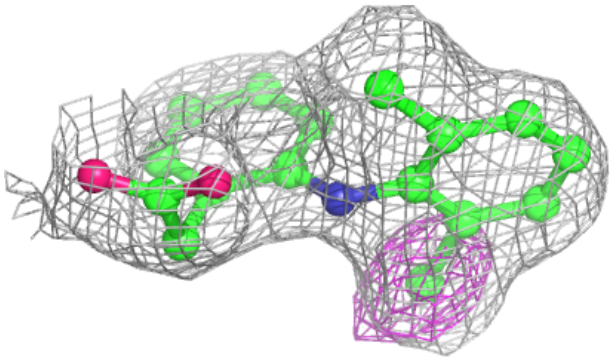


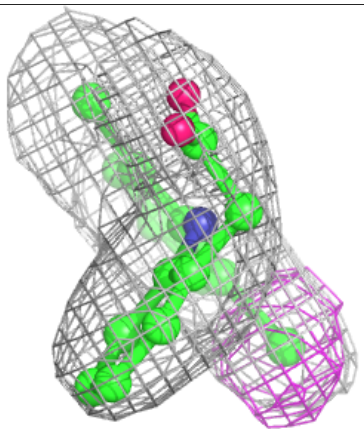


Electron density around DIF A 701:

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

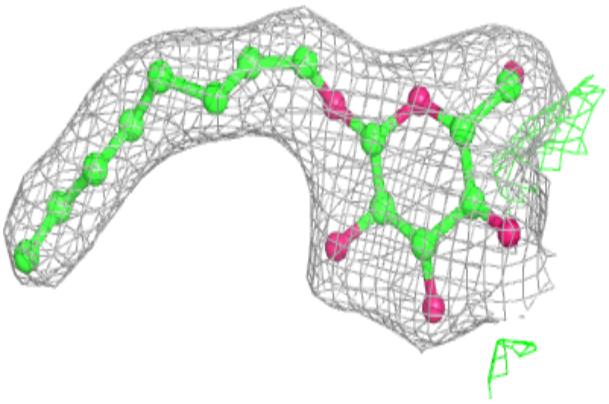


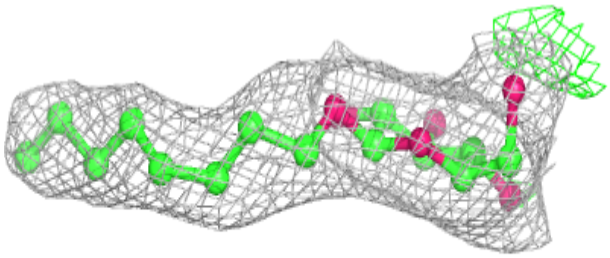


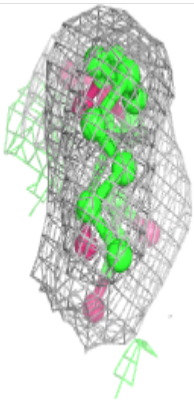


Electron density around BOG B 1751:

$2mF_o$-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.