Full wwPDB X-ray Structure Validation Report (i)

Sep 23, 2023 – 03:34 PM EDT

PDB ID	:	5T6S
Title	:	Crystal structure of the A/Shanghai/2/2013 (H7N9) influenza virus hemagglutinin in complex with the antiviral drug arbidol
Authors	:	Kadam, R.U.; Wilson, I.A.
Deposited on	:	2016-09-01
Resolution	:	2.36 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (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:

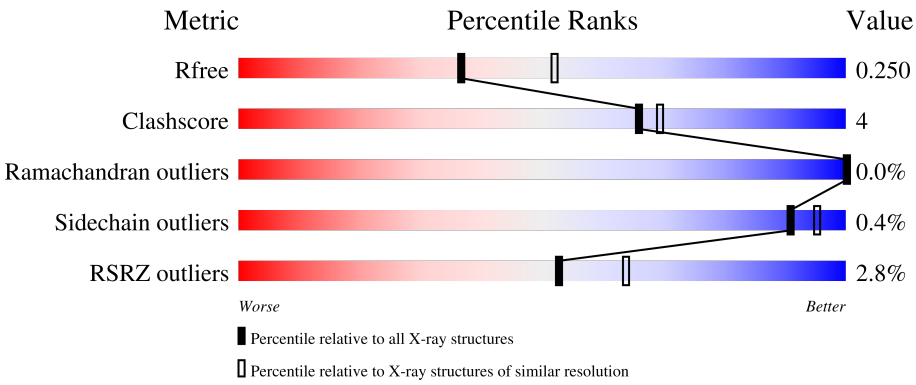
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.36 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	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	321	90% 9% •
1	C	321	89% 9% •
1	E	321	% 88% 10% •
1	G	321	8% 83% 15% •
1	I	321	2% 84% 14% •

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Mol	Chain	Length	Quality of chain
1	K	321	5% 88% 10%
2	B	183	2% 89% 5% 7%
2	D	183	3% 89% 5% 7%
2	F	183	% 88% 5% 7%
2	H	183	3% 89% 5% 7%
2	J	183	6% 85% 8% 7%
2	L	183	% 83% 11% 7%
3	M	3	67% 33%
3	N	3	33% 67%
3	Q	3	33% 67%
3	R	3	100%
3	S	3	33% 33% 33%
4	O	2	100%
4	P	2	100%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BMA	M	3	-	-	-	X
3	BMA	N	3	-	-	-	X
3	BMA	R	3	-	-	-	X
3	BMA	S	3	-	-	-	X
4	NAG	O	2	-	-	-	X
4	NAG	P	2	-	-	-	X
5	NAG	J	203	-	-	-	X

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 24445 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 Hemagglutinin HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	316	2411	1498	436	463	14	0	0	0
1	C	316	2411	1498	436	463	14	0	0	0
1	E	316	2411	1498	436	463	14	0	0	0
1	G	316	2411	1498	436	463	14	0	0	0
1	I	316	2411	1498	436	463	14	0	0	0
1	K	316	2411	1498	436	463	14	0	0	0

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	171	1396	862	243	284	7	0	1	0
2	D	171	1387	857	242	281	7	0	0	0
2	F	171	1393	861	242	283	7	0	1	0
2	H	171	1396	862	243	284	7	0	1	0
2	J	171	1387	857	242	281	7	0	0	0
2	L	171	1396	862	243	284	7	0	1	0

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	177	SER	-	expression tag	UNP R4NN21

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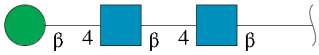
Chain	Residue	Modelled	Actual	Comment	Reference
B	178	GLY	-	expression tag	UNP R4NN21
B	179	ARG	-	expression tag	UNP R4NN21
B	180	LEU	-	expression tag	UNP R4NN21
B	181	VAL	-	expression tag	UNP R4NN21
B	182	PRO	-	expression tag	UNP R4NN21
B	183	ARG	-	expression tag	UNP R4NN21
D	177	SER	-	expression tag	UNP R4NN21
D	178	GLY	-	expression tag	UNP R4NN21
D	179	ARG	-	expression tag	UNP R4NN21
D	180	LEU	-	expression tag	UNP R4NN21
D	181	VAL	-	expression tag	UNP R4NN21
D	182	PRO	-	expression tag	UNP R4NN21
D	183	ARG	-	expression tag	UNP R4NN21
F	177	SER	-	expression tag	UNP R4NN21
F	178	GLY	-	expression tag	UNP R4NN21
F	179	ARG	-	expression tag	UNP R4NN21
F	180	LEU	-	expression tag	UNP R4NN21
F	181	VAL	-	expression tag	UNP R4NN21
F	182	PRO	-	expression tag	UNP R4NN21
F	183	ARG	-	expression tag	UNP R4NN21
H	177	SER	-	expression tag	UNP R4NN21
H	178	GLY	-	expression tag	UNP R4NN21
H	179	ARG	-	expression tag	UNP R4NN21
H	180	LEU	-	expression tag	UNP R4NN21
H	181	VAL	-	expression tag	UNP R4NN21
H	182	PRO	-	expression tag	UNP R4NN21
H	183	ARG	-	expression tag	UNP R4NN21
J	177	SER	-	expression tag	UNP R4NN21
J	178	GLY	-	expression tag	UNP R4NN21
J	179	ARG	-	expression tag	UNP R4NN21
J	180	LEU	-	expression tag	UNP R4NN21
J	181	VAL	-	expression tag	UNP R4NN21
J	182	PRO	-	expression tag	UNP R4NN21
J	183	ARG	-	expression tag	UNP R4NN21
L	177	SER	-	expression tag	UNP R4NN21
L	178	GLY	-	expression tag	UNP R4NN21
L	179	ARG	-	expression tag	UNP R4NN21
L	180	LEU	-	expression tag	UNP R4NN21
L	181	VAL	-	expression tag	UNP R4NN21
L	182	PRO	-	expression tag	UNP R4NN21
L	183	ARG	-	expression tag	UNP R4NN21

- Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-b

eta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

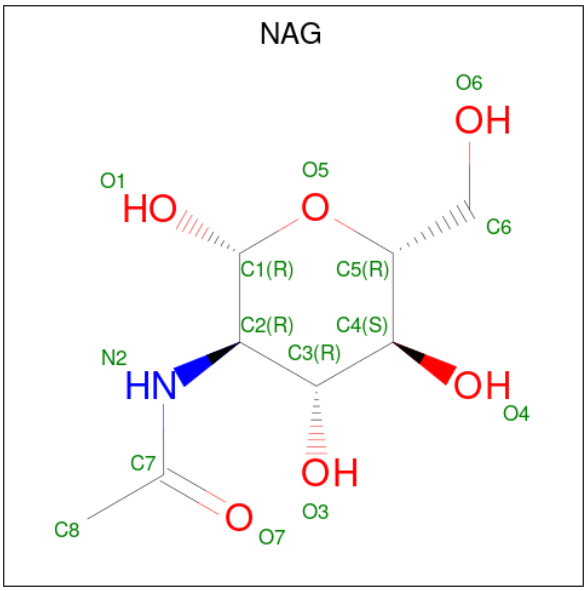


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	M	3	39	22	2	15	0	0	0
3	N	3	39	22	2	15	0	0	0
3	Q	3	39	22	2	15	0	0	0
3	R	3	39	22	2	15	0	0	0
3	S	3	39	22	2	15	0	0	0

- 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
			Total	C	N	O			
4	O	2	28	16	2	10	0	0	0
4	P	2	28	16	2	10	0	0	0

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	A	1	14	8	1	5	0	0
			Total	C	N	O		
5	A	1	14	8	1	5	0	0
			Total	C	N	O		
5	B	1	14	8	1	5	0	0
			Total	C	N	O		
5	B	1	14	8	1	5	0	0
			Total	C	N	O		
5	C	1	14	8	1	5	0	0
			Total	C	N	O		
5	D	1	14	8	1	5	0	0
			Total	C	N	O		
5	D	1	14	8	1	5	0	0
			Total	C	N	O		
5	E	1	14	8	1	5	0	0
			Total	C	N	O		
5	F	1	14	8	1	5	0	0
			Total	C	N	O		
5	H	1	14	8	1	5	0	0
			Total	C	N	O		
5	I	1	14	8	1	5	0	0
			Total	C	N	O		
5	J	1	14	8	1	5	0	0
			Total	C	N	O		
5	J	1	14	8	1	5	0	0
			Total	C	N	O		
5	K	1	14	8	1	5	0	0

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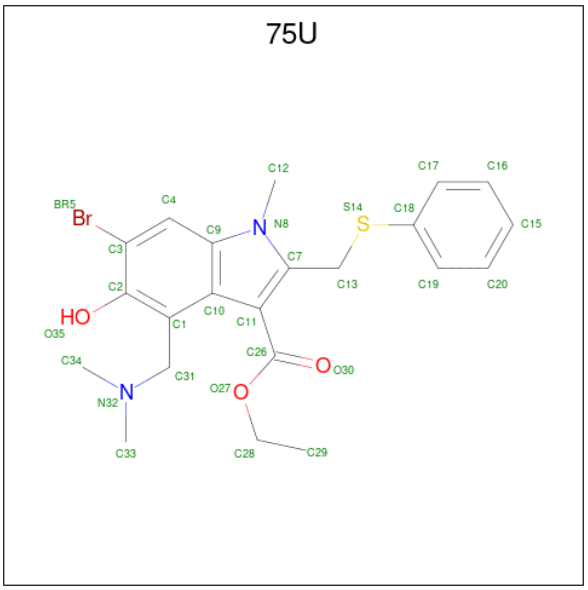
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	L	1	14	8	1	5	0	0

- Molecule 6 is ethyl 6-bromo-4-[(dimethylamino)methyl]-5-hydroxy-1-methyl-2-[(phenylsulfanyl)methyl]-1H-indole-3-carboxylate (three-letter code: 75U) (formula: $C_{22}H_{25}BrN_2O_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	Br	C	N	O	S		
6	B	1	29	1	22	2	3	1	0	0
			Total	Br	C	N	O	S		
6	B	1	29	1	22	2	3	1	0	0
			Total	Br	C	N	O	S		
6	F	1	29	1	22	2	3	1	0	0
			Total	Br	C	N	O	S		
6	H	1	29	1	22	2	3	1	0	0
			Total	Br	C	N	O	S		
6	J	1	29	1	22	2	3	1	0	0
			Total	Br	C	N	O	S		
6	L	1	29	1	22	2	3	1	0	0

- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
7	J	1	1	1	0	0

- Molecule 8 is water.

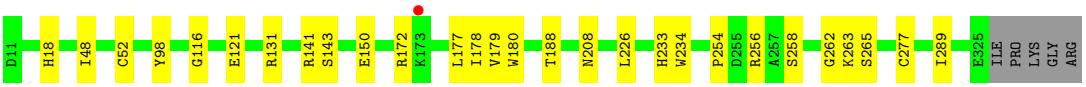
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	116	116	116	0	0
			Total	O		
8	B	86	86	86	0	0
			Total	O		
8	C	142	142	142	0	0
			Total	O		
8	D	81	81	81	0	0
			Total	O		
8	E	143	143	143	0	0
			Total	O		
8	F	80	80	80	0	0
			Total	O		
8	G	33	33	33	0	0
			Total	O		
8	H	34	34	34	0	0
			Total	O		
8	I	90	90	90	0	0
			Total	O		
8	J	53	53	53	0	0
			Total	O		
8	K	59	59	59	0	0
			Total	O		
8	L	71	71	71	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.

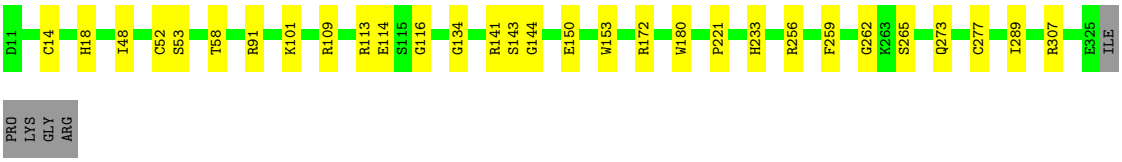
• Molecule 1: Hemagglutinin HA1

Chain A:



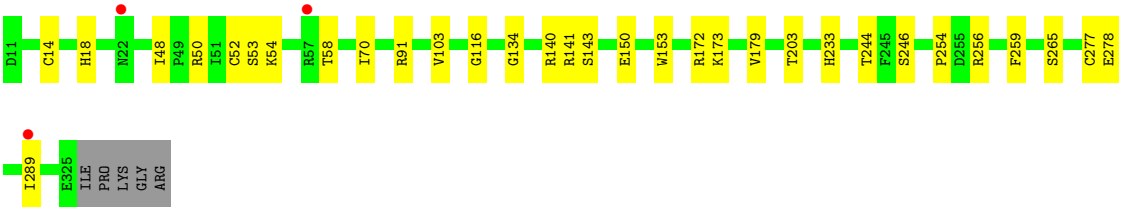
• Molecule 1: Hemagglutinin HA1

Chain C:



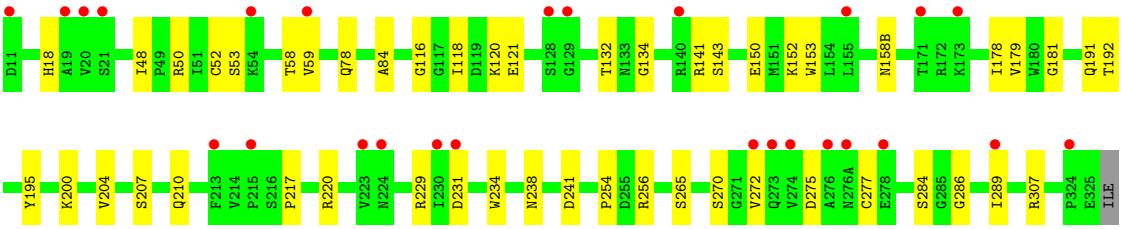
• Molecule 1: Hemagglutinin HA1

Chain E:



• Molecule 1: Hemagglutinin HA1

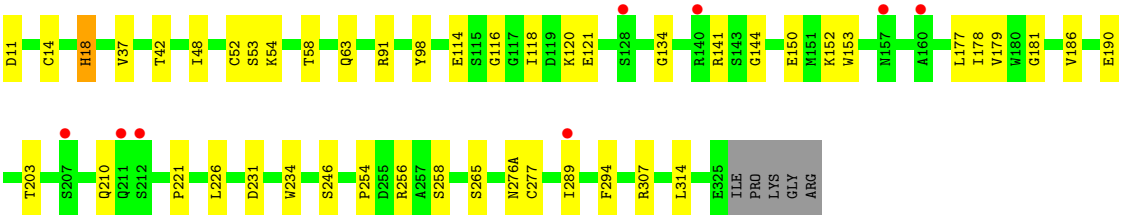
Chain G:



PRO
LYS
GLY
ARG

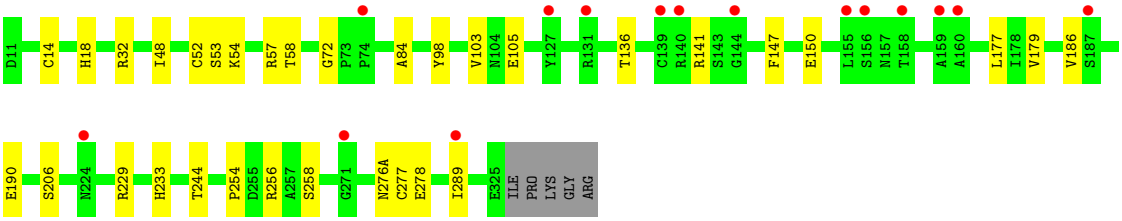
• Molecule 1: Hemagglutinin HA1

Chain I:



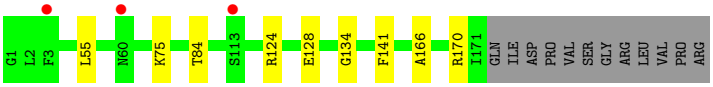
• Molecule 1: Hemagglutinin HA1

Chain K:



• Molecule 2: Hemagglutinin HA2

Chain B:

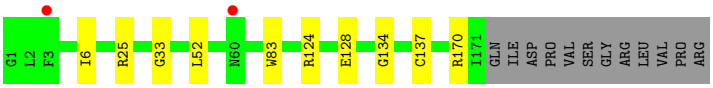


• Molecule 2: Hemagglutinin HA2

Chain D:

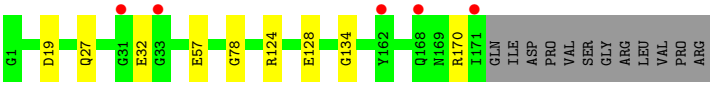
• Molecule 2: Hemagglutinin HA2

Chain F:



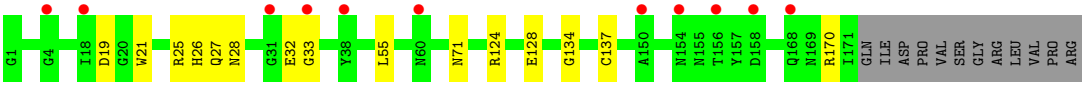
• Molecule 2: Hemagglutinin HA2

Chain H:



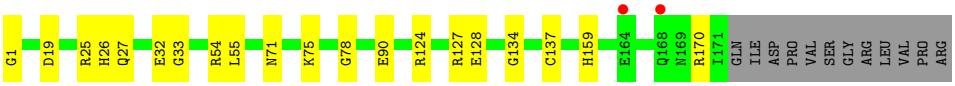
• Molecule 2: Hemagglutinin HA2

Chain J:



• Molecule 2: Hemagglutinin HA2

Chain L:

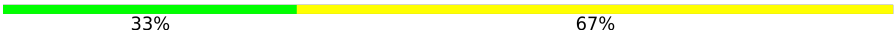


• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:

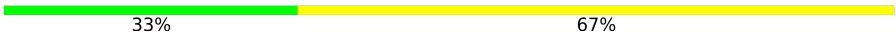
• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:



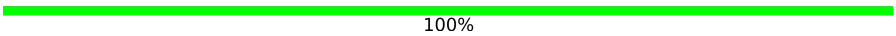
• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:

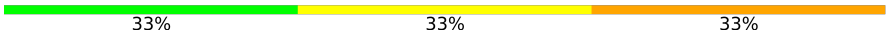


• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

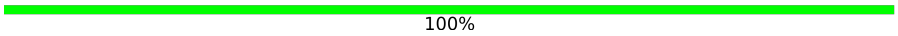
Chain R:



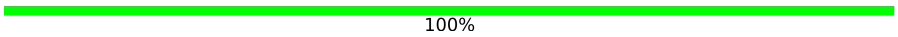
• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S: 

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

Chain O: 

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

Chain P: 

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	67.42Å 231.49Å 127.85Å 90.00° 96.88° 90.00°	Depositor
Resolution (Å)	48.50 – 2.36 49.02 – 2.36	Depositor EDS
% Data completeness (in resolution range)	97.0 (48.50-2.36) 91.2 (49.02-2.36)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.83 (at 2.37Å)	Xtriage
Refinement program	PHENIX (1.10	Depositor
R, R_{free}	0.204 , 0.250 0.204 , 0.250	Depositor DCC
R_{free} test set	7634 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å^2)	32.7	Xtriage
Anisotropy	0.687	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 33.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	24445	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.85% 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: NAG, NA, BMA, 75U

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	0.30	0/2457	0.53	0/3321
1	C	0.31	0/2457	0.54	0/3321
1	E	0.31	0/2457	0.53	0/3321
1	G	0.28	0/2457	0.50	0/3321
1	I	0.29	0/2457	0.51	0/3321
1	K	0.28	0/2457	0.49	0/3321
2	B	0.29	0/1420	0.49	0/1913
2	D	0.36	0/1411	0.50	0/1901
2	F	0.31	0/1420	0.49	0/1913
2	H	0.28	0/1420	0.46	0/1913
2	J	0.33	0/1411	0.48	0/1901
2	L	0.29	0/1420	0.48	0/1913
All	All	0.30	0/23244	0.51	0/31380

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	2411	0	2368	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2411	0	2368	19	0
1	E	2411	0	2368	22	0
1	G	2411	0	2369	29	0
1	I	2411	0	2368	33	0
1	K	2411	0	2368	20	0
2	B	1396	0	1295	8	0
2	D	1387	0	1290	9	0
2	F	1393	0	1296	9	0
2	H	1396	0	1296	8	0
2	J	1387	0	1290	13	0
2	L	1396	0	1295	18	0
3	M	39	0	34	0	0
3	N	39	0	34	1	0
3	Q	39	0	34	1	0
3	R	39	0	34	0	0
3	S	39	0	34	1	0
4	O	28	0	25	0	0
4	P	28	0	25	0	0
5	A	28	0	26	0	0
5	B	28	0	26	1	0
5	C	14	0	13	0	0
5	D	28	0	26	2	0
5	E	14	0	13	0	0
5	F	14	0	13	0	0
5	H	14	0	13	1	0
5	I	14	0	13	0	0
5	J	28	0	26	0	0
5	K	14	0	13	0	0
5	L	14	0	13	2	0
6	B	58	0	0	2	0
6	F	29	0	0	0	0
6	H	29	0	0	1	0
6	J	29	0	0	2	0
6	L	29	0	0	0	0
7	J	1	0	0	0	0
8	A	116	0	0	2	0
8	B	86	0	0	0	0
8	C	142	0	0	3	0
8	D	81	0	0	2	0
8	E	143	0	0	3	0
8	F	80	0	0	2	0
8	G	33	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	H	34	0	0	0	0
8	I	90	0	0	4	0
8	J	53	0	0	2	0
8	K	59	0	0	0	0
8	L	71	0	0	2	0
All	All	24445	0	22386	182	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.

All (182) 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:F:6:ILE:O	8:F:301:HOH:O	2.03	0.76
1:G:231:ASP:OD1	1:I:210:GLN:NE2	2.19	0.74
2:D:134:GLY:HA2	2:F:124:ARG:HD3	1.70	0.71
2:D:19:ASP:OD2	8:D:301:HOH:O	2.14	0.65
2:L:128:GLU:HG3	2:L:170:ARG:HH12	1.63	0.62
1:C:150:GLU:OE1	1:C:256:ARG:HD3	1.99	0.62
1:I:210:GLN:NE2	8:I:503:HOH:O	2.32	0.62
2:H:124:ARG:HD3	2:L:134:GLY:HA2	1.81	0.62
1:G:307:ARG:NH2	2:L:90[B]:GLU:OE1	2.33	0.62
1:C:91:ARG:NH1	8:C:502:HOH:O	2.33	0.61
1:E:134:GLY:HA3	1:E:153:TRP:HB3	1.82	0.61
1:C:48:ILE:HD12	1:C:289:ILE:HD12	1.82	0.61
2:H:134:GLY:HA2	2:J:124:ARG:HD3	1.82	0.61
2:J:134:GLY:HA2	2:L:124:ARG:HD3	1.83	0.59
1:I:18:HIS:CD2	2:J:21:TRP:HA	2.38	0.59
1:G:307:ARG:NH2	2:L:90[A]:GLU:OE1	2.33	0.59
1:A:150:GLU:OE1	1:A:256:ARG:HD3	2.02	0.58
3:S:2:NAG:O3	3:S:3:BMA:O5	2.15	0.58
2:B:134:GLY:HA2	2:D:124:ARG:HD3	1.85	0.57
2:B:124:ARG:HD3	2:F:134:GLY:HA2	1.86	0.57
1:K:48:ILE:HD12	1:K:289:ILE:HD12	1.87	0.57
1:K:186:VAL:HG12	1:K:190:GLU:OE1	2.06	0.56
1:G:116:GLY:HA2	1:G:265:SER:HB3	1.88	0.56
1:G:238:ASN:OD1	8:G:501:HOH:O	2.18	0.56
1:E:150:GLU:OE1	1:E:256:ARG:HD3	2.05	0.56
6:J:201:75U:N32	6:J:201:75U:O30	2.38	0.56
1:G:150:GLU:OE1	1:G:256:ARG:HD3	2.06	0.56
1:K:57:ARG:HG3	1:K:84:ALA:HA	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:27:GLN:OE1	8:J:301:HOH:O	2.18	0.55
2:L:128:GLU:HG3	2:L:170:ARG:NH1	2.22	0.55
2:D:78:GLY:HA3	5:D:201:NAG:H82	1.89	0.54
1:A:48:ILE:HD12	1:A:289:ILE:HD12	1.87	0.54
1:I:152:LYS:NZ	8:I:507:HOH:O	2.40	0.54
1:E:91:ARG:NH1	8:E:512:HOH:O	2.40	0.53
2:B:75:LYS:HD2	5:B:203:NAG:H81	1.91	0.53
2:H:78:GLY:HA3	5:H:202:NAG:H82	1.91	0.53
2:B:128:GLU:HG3	2:B:170:ARG:HH12	1.74	0.53
2:L:78:GLY:HA3	5:L:202:NAG:H82	1.91	0.53
1:G:48:ILE:HD12	1:G:289:ILE:HD12	1.90	0.52
6:J:201:75U:S14	2:L:55:LEU:HD23	2.50	0.52
2:L:25:ARG:HA	2:L:33:GLY:O	2.09	0.52
2:F:128:GLU:HG3	2:F:170:ARG:HH12	1.75	0.52
1:A:98:TYR:CE2	1:A:226:LEU:HD13	2.45	0.52
1:G:132:THR:CG2	1:G:152:LYS:HD3	2.40	0.52
1:G:53:SER:HB2	1:G:58:THR:OG1	2.11	0.51
1:A:131:ARG:NH2	8:A:504:HOH:O	2.40	0.51
1:I:186:VAL:HG12	1:I:190:GLU:OE2	2.10	0.51
1:K:98:TYR:HD2	1:K:136:THR:HG21	1.76	0.51
1:A:179:VAL:O	1:A:254:PRO:HB3	2.11	0.51
2:H:128:GLU:HG3	2:H:170:ARG:HH12	1.76	0.50
1:I:150:GLU:OE1	1:I:256:ARG:HD3	2.11	0.50
1:C:52:CYS:HB3	1:C:277:CYS:O	2.12	0.50
1:A:52:CYS:SG	1:A:277:CYS:CB	2.99	0.50
2:J:128:GLU:HG3	2:J:170:ARG:NH1	2.27	0.50
1:I:91:ARG:NH1	8:I:508:HOH:O	2.41	0.50
1:C:116:GLY:HA2	1:C:265:SER:HB3	1.93	0.49
2:B:128:GLU:HG3	2:B:170:ARG:NH1	2.27	0.49
1:G:220:ARG:HD2	1:G:229:ARG:HG2	1.94	0.49
2:L:75:LYS:HD2	5:L:202:NAG:H81	1.94	0.49
1:E:50:ARG:NH1	8:E:511:HOH:O	2.40	0.49
1:E:172:ARG:HD3	1:E:259:PHE:CZ	2.46	0.49
1:I:179:VAL:O	1:I:254:PRO:HB3	2.12	0.49
2:H:57:GLU:OE1	1:K:32:ARG:HD2	2.13	0.49
2:F:128:GLU:HG3	2:F:170:ARG:NH1	2.28	0.49
1:G:118:ILE:HD12	1:G:120:LYS:HE3	1.95	0.49
1:G:241:ASP:OD2	8:G:502:HOH:O	2.20	0.49
6:B:202:75U:S14	2:D:55:LEU:HD23	2.53	0.49
1:G:78:GLN:O	1:G:78:GLN:HG3	2.12	0.48
3:N:1:NAG:H61	3:N:2:NAG:H82	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:143:SER:HB2	1:I:121:GLU:HA	1.95	0.48
2:L:27:GLN:NE2	2:L:32:GLU:OE2	2.42	0.48
1:I:307:ARG:NH2	8:I:501:HOH:O	2.21	0.48
1:K:52:CYS:SG	1:K:277:CYS:CB	3.02	0.48
1:C:134:GLY:HA3	1:C:153:TRP:HB3	1.96	0.48
2:L:1:GLY:N	8:L:305:HOH:O	2.47	0.47
1:G:204:VAL:O	1:G:210:GLN:HA	2.13	0.47
1:I:54:LYS:HD2	1:I:276(A):ASN:C	2.35	0.47
1:K:72:GLY:O	1:K:141:ARG:NH2	2.47	0.47
1:C:143:SER:N	8:C:505:HOH:O	2.43	0.47
1:C:172:ARG:HD3	1:C:259:PHE:CZ	2.50	0.47
1:K:150:GLU:OE1	1:K:256:ARG:HD3	2.15	0.47
1:G:50:ARG:NE	1:G:275:ASP:OD2	2.47	0.47
2:L:127:ARG:HG3	2:L:159:HIS:CG	2.50	0.47
1:K:141:ARG:NH1	1:K:147:PHE:O	2.48	0.46
1:A:177:LEU:HB3	1:A:258:SER:HB2	1.96	0.46
1:G:179:VAL:O	1:G:254:PRO:HB3	2.16	0.46
1:I:116:GLY:HA2	1:I:265:SER:HB3	1.97	0.46
2:L:54:ARG:NH2	8:L:307:HOH:O	2.48	0.46
1:E:70:ILE:HG21	1:E:179:VAL:HG21	1.98	0.46
1:A:116:GLY:HA2	1:A:265:SER:HB3	1.97	0.46
1:E:116:GLY:HA2	1:E:265:SER:HB3	1.97	0.46
1:I:14:CYS:HA	2:J:137:CYS:HA	1.97	0.45
1:C:273:GLN:HG3	8:C:550:HOH:O	2.16	0.45
1:G:134:GLY:HA3	1:G:153:TRP:HB3	1.99	0.45
1:A:98:TYR:CD2	1:A:226:LEU:HD13	2.51	0.45
1:A:178:ILE:O	1:A:234:TRP:HA	2.16	0.45
1:G:181:GLY:HA2	1:G:231:ASP:O	2.17	0.45
1:E:173:LYS:NZ	8:E:513:HOH:O	2.40	0.45
1:K:14:CYS:HA	2:L:137:CYS:HA	1.99	0.45
1:I:134:GLY:HA3	1:I:153:TRP:HB3	1.98	0.45
1:A:121:GLU:OE1	1:A:172:ARG:HD2	2.17	0.45
1:G:270:SER:HA	1:G:284:SER:HA	1.99	0.45
2:J:19:ASP:OD1	2:J:19:ASP:N	2.47	0.45
2:J:128:GLU:HG3	2:J:170:ARG:HH12	1.81	0.45
1:E:179:VAL:O	1:E:254:PRO:HB3	2.17	0.45
2:L:19:ASP:OD1	2:L:19:ASP:N	2.49	0.44
2:L:26:HIS:CE1	2:L:33:GLY:HA3	2.52	0.44
1:I:118:ILE:HD12	1:I:120:LYS:HE3	1.99	0.44
2:H:19:ASP:OD1	2:H:19:ASP:N	2.47	0.44
1:K:103:VAL:HG21	1:K:233:HIS:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:203:THR:OG1	1:I:246:SER:HB3	2.17	0.44
1:E:140:ARG:NH2	1:I:121:GLU:OE1	2.43	0.44
1:E:48:ILE:HD12	1:E:289:ILE:HD12	2.00	0.44
1:E:53:SER:HB2	1:E:58:THR:OG1	2.18	0.44
1:E:203:THR:OG1	1:E:246:SER:HB3	2.18	0.44
2:D:19:ASP:OD1	2:D:19:ASP:N	2.50	0.44
2:H:128:GLU:HG3	2:H:170:ARG:NH1	2.33	0.43
1:I:18:HIS:CE1	1:I:37:VAL:HG11	2.53	0.43
1:A:141:ARG:O	1:A:143:SER:OG	2.33	0.43
2:F:25:ARG:HA	2:F:33:GLY:O	2.18	0.43
2:F:124:ARG:NH1	8:F:303:HOH:O	2.29	0.43
6:H:201:75U:S14	2:J:55:LEU:HD23	2.59	0.43
1:I:42:THR:HG22	1:I:314:LEU:O	2.19	0.43
1:C:109:ARG:O	1:C:113:ARG:HG3	2.19	0.43
2:D:54:ARG:NH2	8:D:310:HOH:O	2.51	0.43
1:G:158(B):ASN:HA	1:G:192:THR:O	2.18	0.43
2:H:27:GLN:NE2	2:H:32:GLU:HG2	2.33	0.43
1:I:114:GLU:HA	1:I:265:SER:O	2.19	0.43
1:I:178:ILE:O	1:I:234:TRP:HA	2.18	0.43
1:C:114:GLU:HA	1:C:265:SER:O	2.19	0.43
1:E:14:CYS:HA	2:F:137:CYS:HA	2.00	0.43
1:I:221:PRO:HG2	1:K:206:SER:HA	2.01	0.43
2:D:75:LYS:HD2	5:D:201:NAG:H81	2.01	0.43
1:E:54:LYS:HB2	1:E:278:GLU:HB2	2.00	0.43
1:I:181:GLY:HA2	1:I:231:ASP:O	2.19	0.43
1:K:54:LYS:HB2	1:K:278:GLU:HB2	2.01	0.43
1:K:105:GLU:OE1	2:L:71:ASN:ND2	2.52	0.43
1:G:178:ILE:O	1:G:234:TRP:HA	2.19	0.42
2:B:141:PHE:O	2:B:166:ALA:HA	2.20	0.42
1:E:103:VAL:HG21	1:E:233:HIS:CE1	2.54	0.42
1:C:141:ARG:C	1:C:144:GLY:H	2.22	0.42
1:G:52:CYS:SG	1:G:277:CYS:CB	3.07	0.42
1:I:221:PRO:HD3	1:K:244:THR:HB	2.00	0.42
1:E:52:CYS:SG	1:E:277:CYS:CB	3.08	0.42
1:G:195:TYR:HB2	1:G:200:LYS:HE3	2.02	0.42
2:B:55:LEU:HD23	6:B:201:75U:S14	2.60	0.42
1:C:52:CYS:SG	1:C:277:CYS:CB	3.08	0.42
1:C:221:PRO:HD3	1:E:244:THR:HB	2.00	0.42
1:I:48:ILE:HD12	1:I:289:ILE:HD12	2.01	0.42
1:K:53:SER:HB2	1:K:58:THR:OG1	2.20	0.42
1:K:179:VAL:O	1:K:254:PRO:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:141:ARG:O	8:G:503:HOH:O	2.22	0.42
1:I:52:CYS:SG	1:I:277:CYS:CB	3.08	0.42
2:B:84:THR:HG22	2:F:83:TRP:HZ3	1.85	0.42
1:I:42:THR:HG23	1:I:294:PHE:HD2	1.84	0.42
1:K:54:LYS:HD2	1:K:276(A):ASN:C	2.40	0.42
1:E:141:ARG:O	1:E:143:SER:OG	2.31	0.41
1:I:98:TYR:CE2	1:I:226:LEU:HD13	2.55	0.41
1:I:141:ARG:C	1:I:144:GLY:H	2.22	0.41
1:K:177:LEU:HB3	1:K:258:SER:HB2	2.02	0.41
1:A:180:TRP:CE2	1:A:233:HIS:HB2	2.55	0.41
1:E:143:SER:HA	1:I:121:GLU:HG3	2.01	0.41
1:C:307:ARG:HD3	1:C:307:ARG:HA	1.79	0.41
2:J:25:ARG:HA	2:J:33:GLY:O	2.20	0.41
1:A:208:ASN:HA	8:A:585:HOH:O	2.21	0.41
1:I:53:SER:HB2	1:I:58:THR:OG1	2.20	0.41
1:C:101:LYS:HB2	1:C:101:LYS:HE3	1.84	0.41
2:J:26:HIS:O	2:J:32:GLU:HA	2.21	0.41
1:A:262:GLY:O	1:A:263:LYS:HD2	2.21	0.41
1:C:180:TRP:CE2	1:C:233:HIS:HB2	2.56	0.41
1:E:52:CYS:HB3	1:E:277:CYS:O	2.21	0.41
1:G:207:SER:HA	1:K:229:ARG:HH21	1.86	0.41
3:Q:1:NAG:H61	3:Q:2:NAG:N2	2.36	0.41
1:C:53:SER:HB2	1:C:58:THR:OG1	2.20	0.41
1:G:59:VAL:HG23	1:G:84:ALA:HB2	2.02	0.41
2:J:71:ASN:ND2	8:J:309:HOH:O	2.53	0.41
1:C:14:CYS:HA	2:D:137:CYS:HA	2.03	0.40
1:G:141:ARG:O	1:G:143:SER:OG	2.37	0.40
1:I:11:ASP:OD1	2:J:28:ASN:HA	2.21	0.40
1:G:191:GLN:HE22	1:G:217:PRO:HD3	1.86	0.40
1:G:272:VAL:HG21	1:G:286:GLY:HA2	2.03	0.40
1:I:177:LEU:HB3	1:I:258:SER:HB2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	314/321 (98%)	305 (97%)	9 (3%)	0	100	100
1	C	314/321 (98%)	305 (97%)	8 (2%)	1 (0%)	41	47
1	E	314/321 (98%)	305 (97%)	9 (3%)	0	100	100
1	G	314/321 (98%)	303 (96%)	11 (4%)	0	100	100
1	I	314/321 (98%)	305 (97%)	9 (3%)	0	100	100
1	K	314/321 (98%)	306 (98%)	8 (2%)	0	100	100
2	B	170/183 (93%)	164 (96%)	6 (4%)	0	100	100
2	D	169/183 (92%)	165 (98%)	4 (2%)	0	100	100
2	F	170/183 (93%)	165 (97%)	5 (3%)	0	100	100
2	H	170/183 (93%)	166 (98%)	4 (2%)	0	100	100
2	J	169/183 (92%)	165 (98%)	4 (2%)	0	100	100
2	L	170/183 (93%)	165 (97%)	5 (3%)	0	100	100
All	All	2902/3024 (96%)	2819 (97%)	82 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	262	GLY

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	264/269 (98%)	262 (99%)	2 (1%)	81	89
1	C	264/269 (98%)	263 (100%)	1 (0%)	91	95
1	E	264/269 (98%)	263 (100%)	1 (0%)	91	95
1	G	264/269 (98%)	262 (99%)	2 (1%)	81	89
1	I	264/269 (98%)	262 (99%)	2 (1%)	81	89

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	K	264/269 (98%)	263 (100%)	1 (0%)	91	95
2	B	147/157 (94%)	147 (100%)	0	100	100
2	D	146/157 (93%)	145 (99%)	1 (1%)	84	91
2	F	147/157 (94%)	146 (99%)	1 (1%)	84	91
2	H	147/157 (94%)	147 (100%)	0	100	100
2	J	146/157 (93%)	146 (100%)	0	100	100
2	L	147/157 (94%)	147 (100%)	0	100	100
All	All	2464/2556 (96%)	2453 (100%)	11 (0%)	91	95

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	188	THR
1	C	18	HIS
2	D	59	THR
1	E	18	HIS
2	F	52	LEU
1	G	18	HIS
1	G	121	GLU
1	I	18	HIS
1	I	63	GLN
1	K	18	HIS

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

Mol	Chain	Res	Type
1	I	18	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

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	M	1	3,1	14,14,15	0.28	0	17,19,21	0.62	1 (5%)
3	NAG	M	2	3	14,14,15	0.22	0	17,19,21	0.46	0
3	BMA	M	3	3	11,11,12	0.76	0	15,15,17	0.76	0
3	NAG	N	1	3,1	14,14,15	0.26	0	17,19,21	0.43	0
3	NAG	N	2	3	14,14,15	0.25	0	17,19,21	0.52	0
3	BMA	N	3	3	11,11,12	0.83	0	15,15,17	0.76	0
4	NAG	O	1	4,2	14,14,15	0.21	0	17,19,21	0.37	0
4	NAG	O	2	4	14,14,15	0.17	0	17,19,21	0.39	0
4	NAG	P	1	4,1	14,14,15	0.28	0	17,19,21	0.45	0
4	NAG	P	2	4	14,14,15	0.27	0	17,19,21	0.49	0
3	NAG	Q	1	3,1	14,14,15	0.49	0	17,19,21	0.45	0
3	NAG	Q	2	3	14,14,15	0.28	0	17,19,21	0.51	0
3	BMA	Q	3	3	11,11,12	0.71	0	15,15,17	0.76	0
3	NAG	R	1	3,1	14,14,15	0.18	0	17,19,21	0.57	0
3	NAG	R	2	3	14,14,15	0.26	0	17,19,21	0.54	0
3	BMA	R	3	3	11,11,12	0.84	0	15,15,17	0.75	0
3	NAG	S	1	3,2	14,14,15	0.24	0	17,19,21	0.49	0
3	NAG	S	2	3	14,14,15	0.60	0	17,19,21	0.84	0
3	BMA	S	3	3	11,11,12	0.55	0	15,15,17	1.34	1 (6%)

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	M	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	M	2	3	-	2/6/23/26	0/1/1/1
3	BMA	M	3	3	-	2/2/19/22	0/1/1/1
3	NAG	N	1	3,1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	N	2	3	-	2/6/23/26	0/1/1/1
3	BMA	N	3	3	-	1/2/19/22	0/1/1/1
4	NAG	O	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	O	2	4	-	2/6/23/26	0/1/1/1
4	NAG	P	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	P	2	4	-	2/6/23/26	0/1/1/1
3	NAG	Q	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	Q	2	3	-	2/6/23/26	0/1/1/1
3	BMA	Q	3	3	-	2/2/19/22	0/1/1/1
3	NAG	R	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	R	2	3	-	2/6/23/26	0/1/1/1
3	BMA	R	3	3	-	1/2/19/22	0/1/1/1
3	NAG	S	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	S	2	3	-	2/6/23/26	0/1/1/1
3	BMA	S	3	3	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	S	3	BMA	C1-O5-C5	4.37	118.11	112.19
3	M	1	NAG	C1-O5-C5	2.18	115.14	112.19

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	R	2	NAG	O5-C5-C6-O6
3	N	2	NAG	O5-C5-C6-O6
3	Q	2	NAG	O5-C5-C6-O6
3	Q	3	BMA	O5-C5-C6-O6
3	M	2	NAG	O5-C5-C6-O6
3	M	3	BMA	O5-C5-C6-O6
3	R	2	NAG	C4-C5-C6-O6
3	N	2	NAG	C4-C5-C6-O6
3	Q	2	NAG	C4-C5-C6-O6
3	S	2	NAG	O5-C5-C6-O6
4	P	2	NAG	O5-C5-C6-O6
3	M	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	P	2	NAG	C4-C5-C6-O6
3	M	3	BMA	C4-C5-C6-O6
3	Q	3	BMA	C4-C5-C6-O6
4	O	2	NAG	O5-C5-C6-O6
4	O	2	NAG	C4-C5-C6-O6
3	N	3	BMA	O5-C5-C6-O6
3	R	3	BMA	O5-C5-C6-O6
3	S	2	NAG	C4-C5-C6-O6
3	S	3	BMA	O5-C5-C6-O6

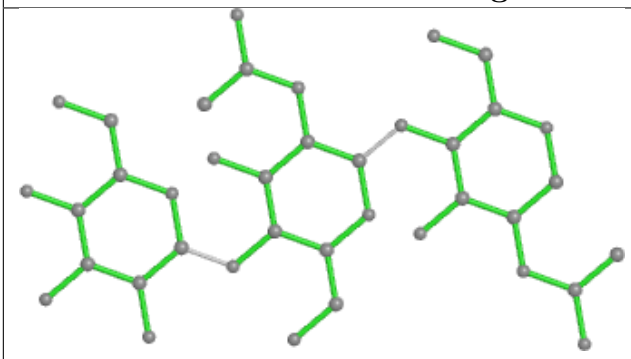
There are no ring outliers.

6 monomers are involved in 3 short contacts:

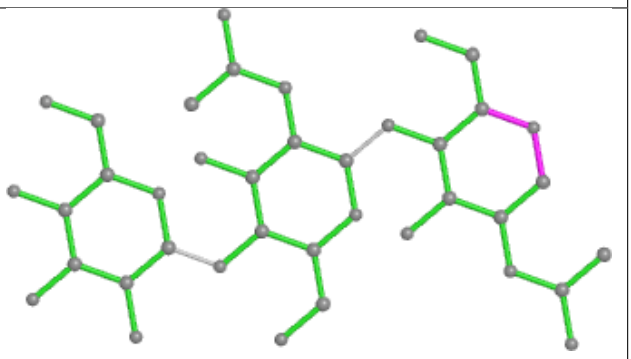
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	N	2	NAG	1	0
3	Q	2	NAG	1	0
3	S	3	BMA	1	0
3	N	1	NAG	1	0
3	Q	1	NAG	1	0
3	S	2	NAG	1	0

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

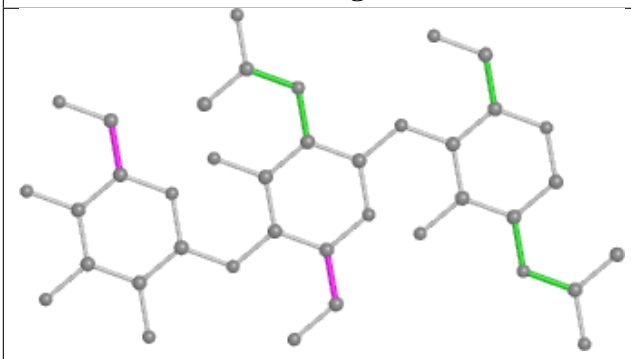
Oligosaccharide Chain M



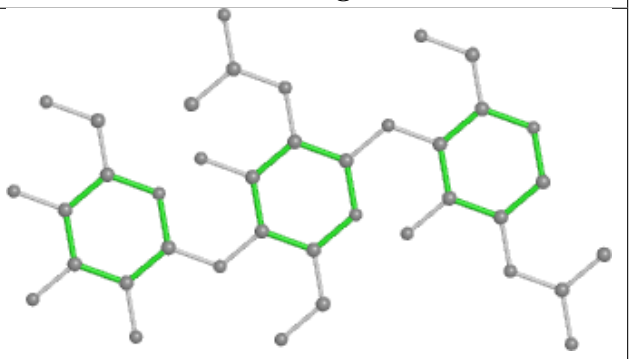
Bond lengths



Bond angles

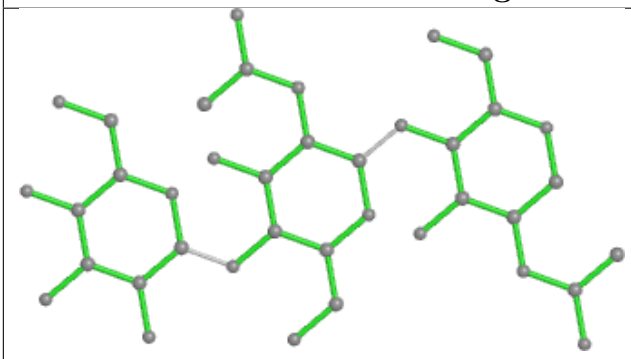


Torsions

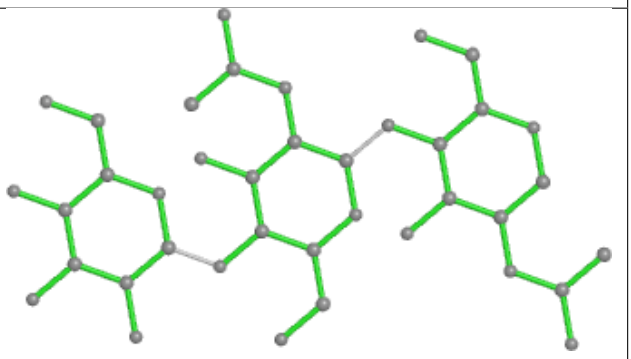


Rings

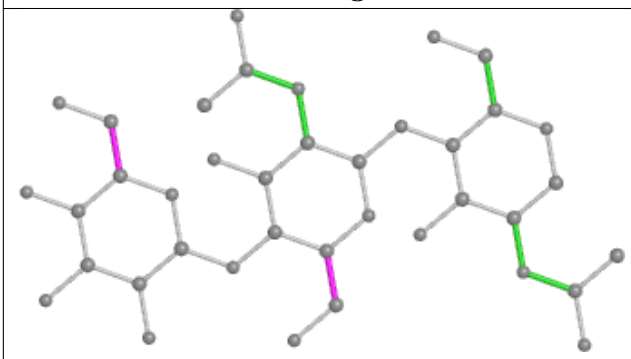
Oligosaccharide Chain N



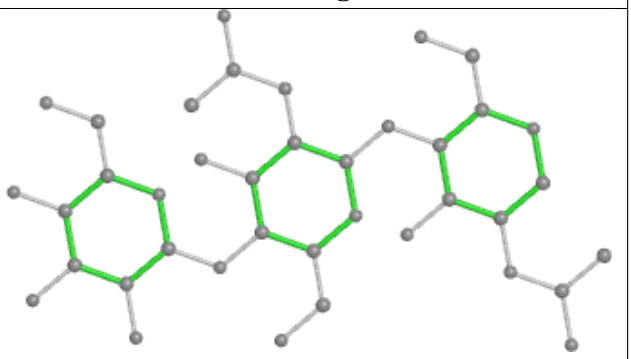
Bond lengths



Bond angles

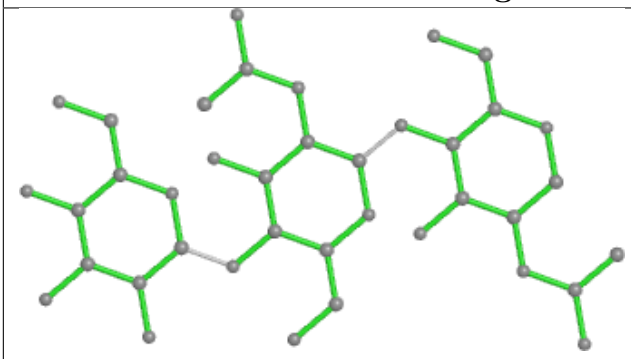


Torsions

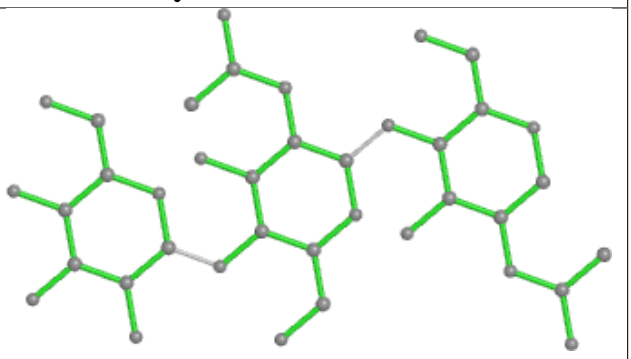


Rings

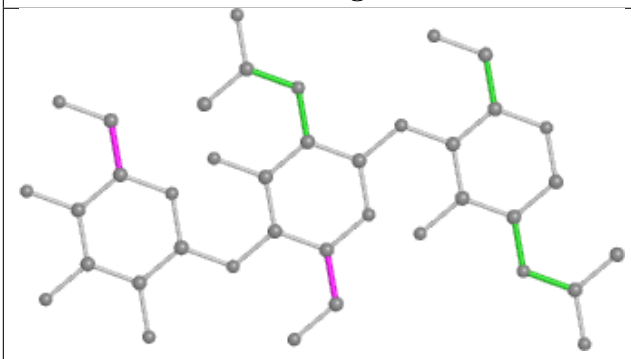
Oligosaccharide Chain Q



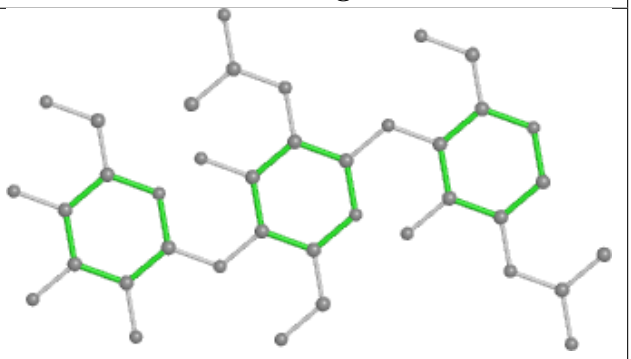
Bond lengths



Bond angles

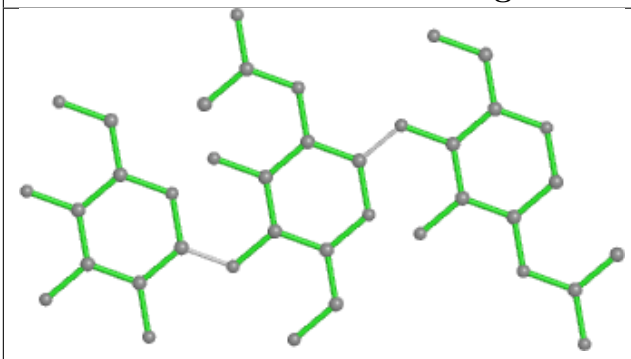


Torsions

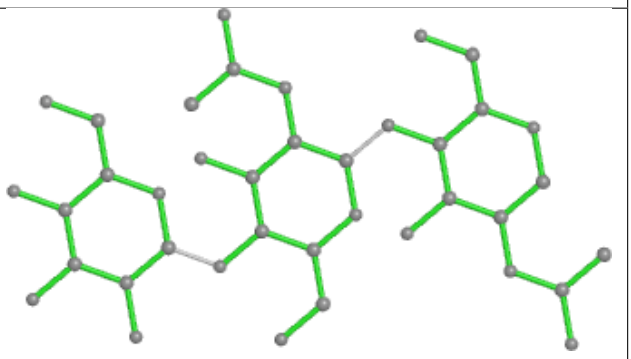


Rings

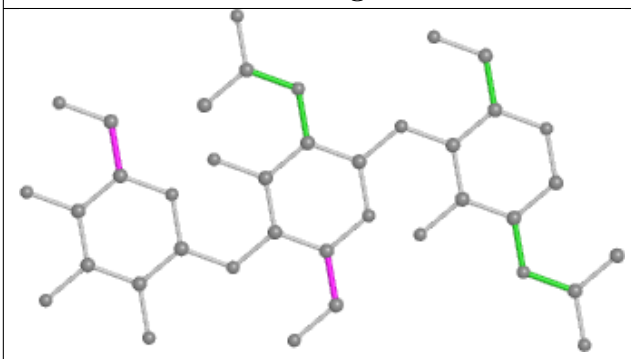
Oligosaccharide Chain R



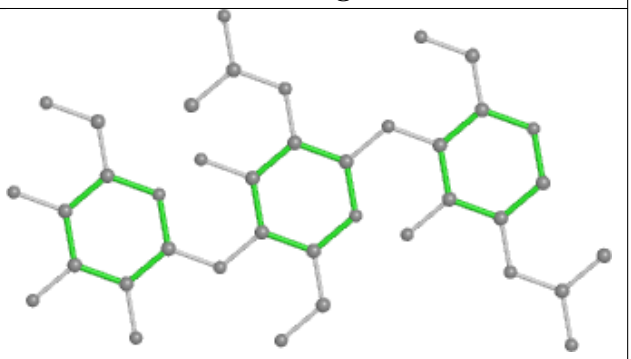
Bond lengths



Bond angles

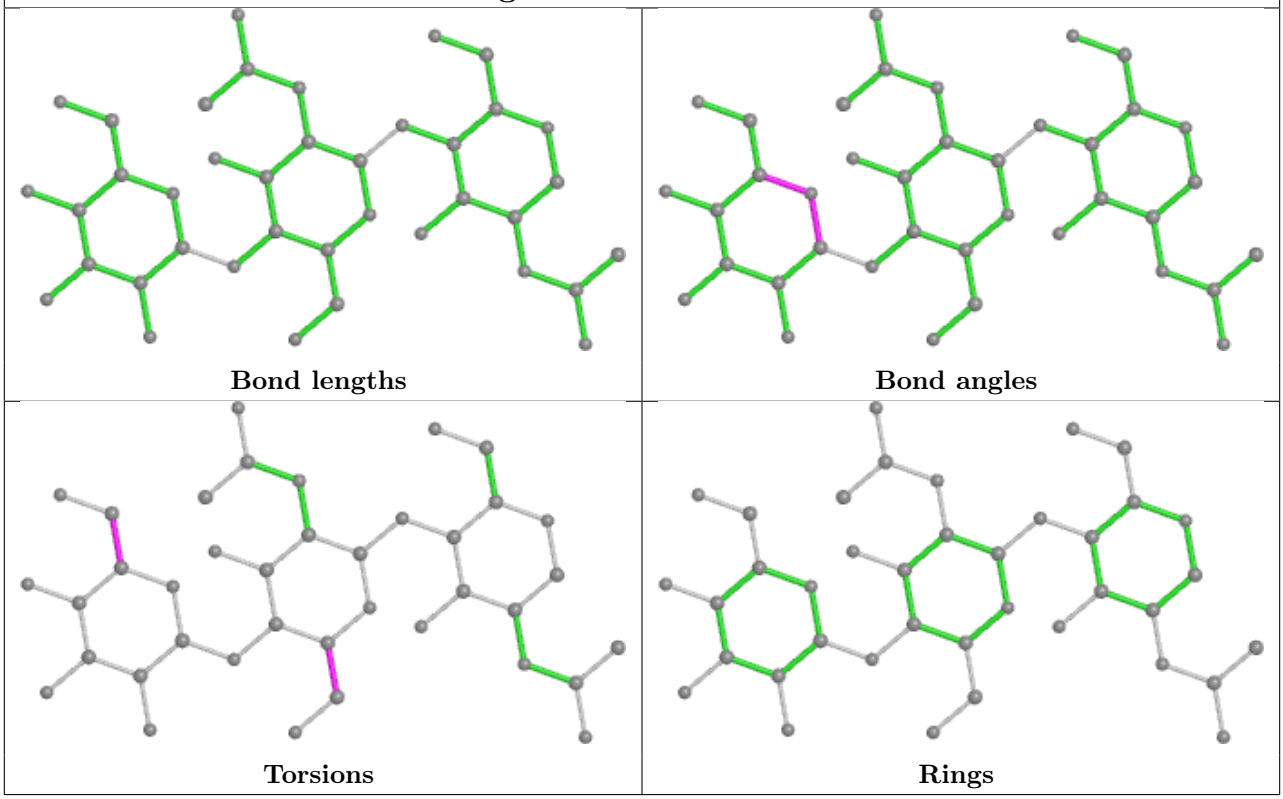


Torsions

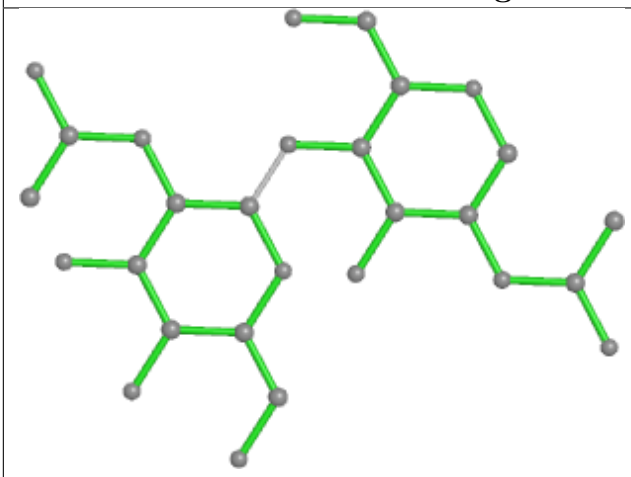


Rings

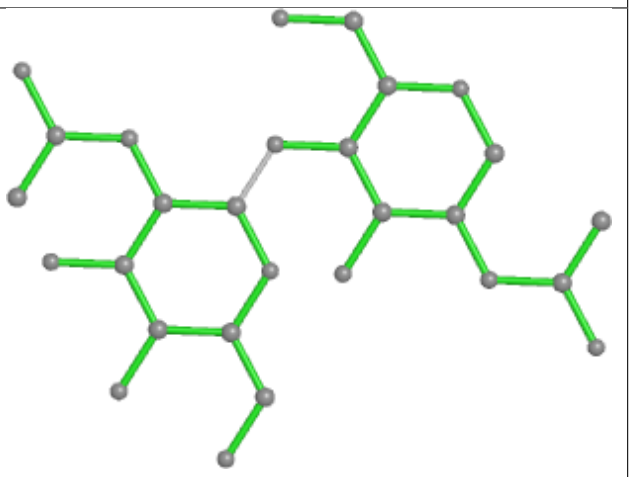
Oligosaccharide Chain S



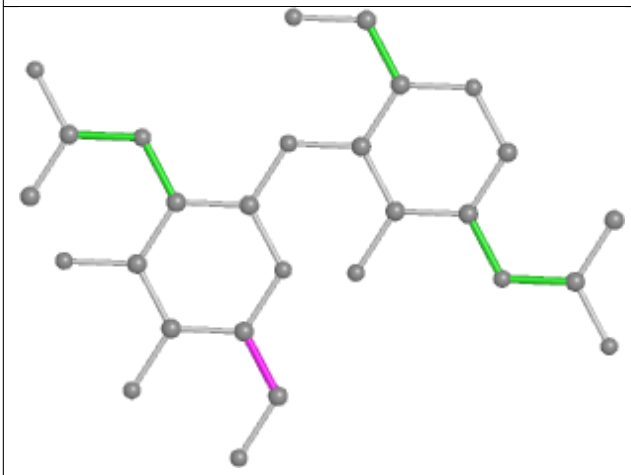
Oligosaccharide Chain O



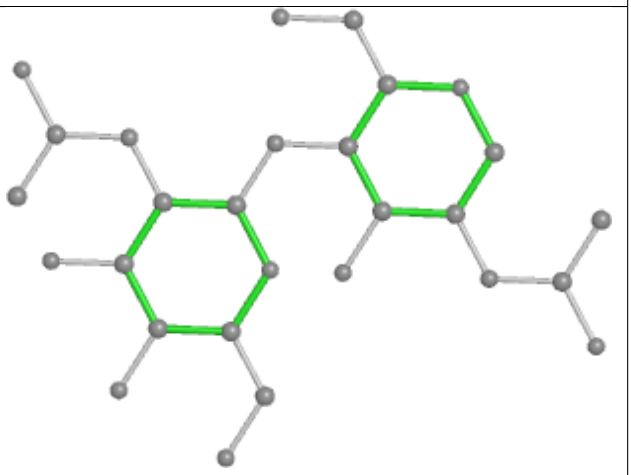
Bond lengths



Bond angles

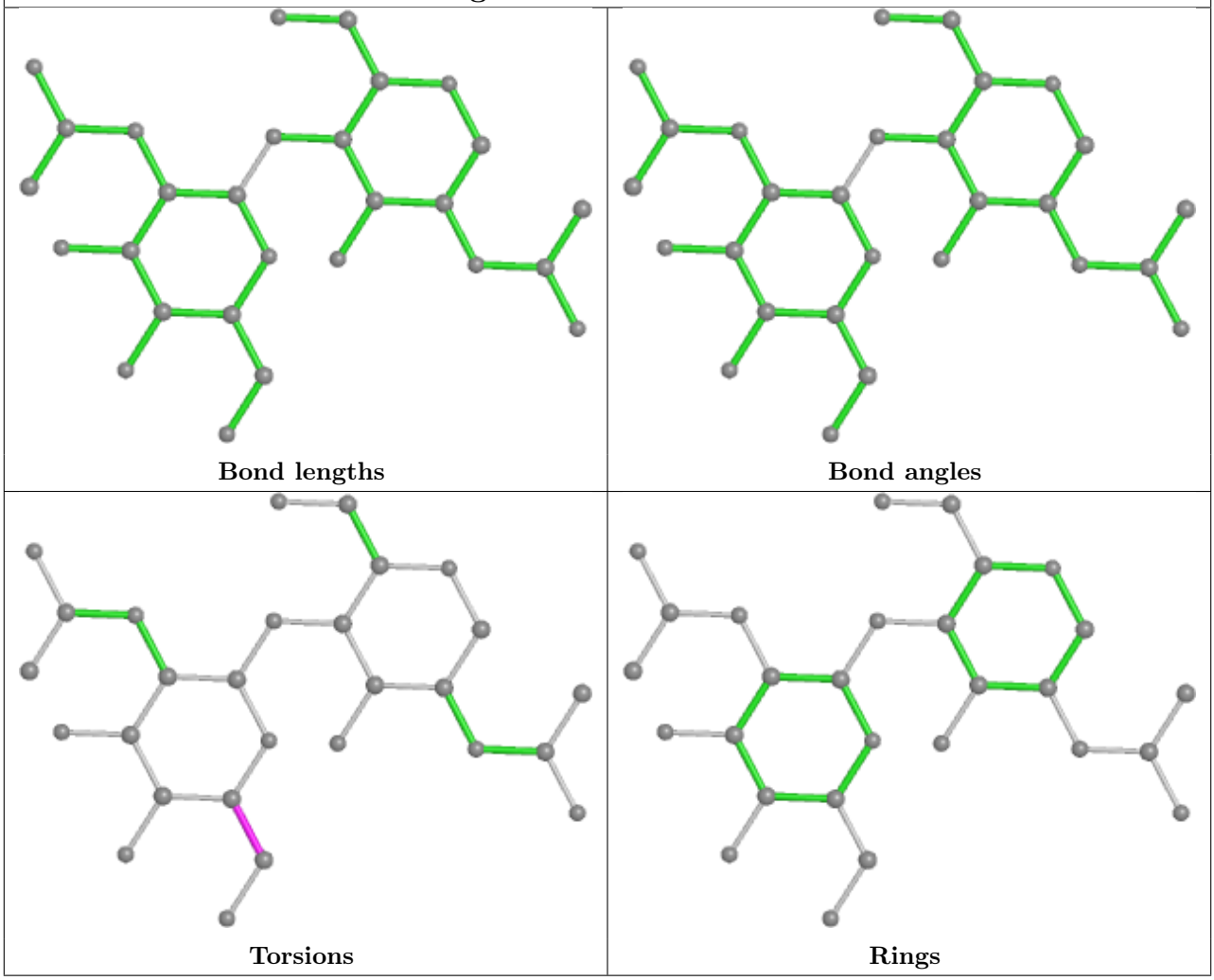


Torsions



Rings

Oligosaccharide Chain P



5.6 Ligand geometry (i)

Of 22 ligands modelled in this entry, 1 is monoatomic - leaving 21 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	$\#\|Z\| > 2$
6	75U	F	201	-	28,31,31	3.82	7 (25%)	35,44,44	1.91	7 (20%)
5	NAG	I	404	1	14,14,15	0.15	0	17,19,21	0.53	0
5	NAG	C	404	1	14,14,15	0.23	0	17,19,21	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	NAG	F	202	2	14,14,15	0.44	0	17,19,21	0.67	1 (5%)
6	75U	B	201	-	28,31,31	4.00	7 (25%)	35,44,44	2.22	8 (22%)
5	NAG	H	202	2	14,14,15	0.49	0	17,19,21	0.64	1 (5%)
5	NAG	B	203	2	14,14,15	0.55	0	17,19,21	0.52	0
5	NAG	A	402	1	14,14,15	0.35	0	17,19,21	0.43	0
6	75U	J	201	-	28,31,31	3.56	7 (25%)	35,44,44	1.97	7 (20%)
5	NAG	D	202	2	14,14,15	0.17	0	17,19,21	0.39	0
6	75U	H	201	-	28,31,31	3.82	7 (25%)	35,44,44	2.16	8 (22%)
6	75U	L	201	-	28,31,31	3.86	7 (25%)	35,44,44	2.07	7 (20%)
5	NAG	J	203	2	14,14,15	0.19	0	17,19,21	0.40	0
5	NAG	L	202	2	14,14,15	0.51	0	17,19,21	0.73	1 (5%)
5	NAG	J	202	2	14,14,15	0.37	0	17,19,21	0.59	1 (5%)
5	NAG	B	204	2	14,14,15	0.31	0	17,19,21	0.42	0
5	NAG	K	404	1	14,14,15	0.25	0	17,19,21	0.48	0
5	NAG	E	404	1	14,14,15	0.29	0	17,19,21	0.43	0
6	75U	B	202	-	28,31,31	3.63	7 (25%)	35,44,44	2.24	8 (22%)
5	NAG	A	401	1	14,14,15	0.21	0	17,19,21	0.46	0
5	NAG	D	201	2	14,14,15	0.62	1 (7%)	17,19,21	0.61	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
6	75U	F	201	-	-	5/16/16/16	0/3/3/3
5	NAG	I	404	1	-	0/6/23/26	0/1/1/1
5	NAG	C	404	1	-	2/6/23/26	0/1/1/1
5	NAG	F	202	2	-	0/6/23/26	0/1/1/1
6	75U	B	201	-	-	5/16/16/16	0/3/3/3
5	NAG	H	202	2	-	0/6/23/26	0/1/1/1
5	NAG	B	203	2	-	2/6/23/26	0/1/1/1
5	NAG	A	402	1	-	1/6/23/26	0/1/1/1
6	75U	J	201	-	-	6/16/16/16	0/3/3/3
5	NAG	D	202	2	-	0/6/23/26	0/1/1/1
6	75U	H	201	-	-	4/16/16/16	0/3/3/3
6	75U	L	201	-	-	5/16/16/16	0/3/3/3
5	NAG	J	203	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	L	202	2	-	2/6/23/26	0/1/1/1
5	NAG	J	202	2	-	0/6/23/26	0/1/1/1
5	NAG	B	204	2	-	0/6/23/26	0/1/1/1
5	NAG	K	404	1	-	2/6/23/26	0/1/1/1
5	NAG	E	404	1	-	2/6/23/26	0/1/1/1
6	75U	B	202	-	-	4/16/16/16	0/3/3/3
5	NAG	A	401	1	-	0/6/23/26	0/1/1/1
5	NAG	D	201	2	-	0/6/23/26	0/1/1/1

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	201	75U	C31-C1	-18.22	1.39	1.51
6	L	201	75U	C31-C1	-17.22	1.39	1.51
6	H	201	75U	C31-C1	-17.02	1.40	1.51
6	F	201	75U	C31-C1	-16.89	1.40	1.51
6	B	202	75U	C31-C1	-15.93	1.40	1.51
6	J	201	75U	C31-C1	-14.98	1.41	1.51
6	J	201	75U	C11-C26	-6.02	1.34	1.50
6	L	201	75U	C11-C26	-5.72	1.35	1.50
6	F	201	75U	C11-C26	-5.72	1.35	1.50
6	B	201	75U	C10-C9	-5.72	1.35	1.42
6	J	201	75U	C10-C9	-5.62	1.35	1.42
6	B	202	75U	C11-C26	-5.61	1.35	1.50
6	H	201	75U	C10-C9	-5.60	1.35	1.42
6	H	201	75U	C11-C26	-5.56	1.35	1.50
6	L	201	75U	C10-C9	-5.53	1.35	1.42
6	B	201	75U	C11-C26	-5.47	1.35	1.50
6	B	202	75U	C10-C9	-5.43	1.35	1.42
6	F	201	75U	C10-C9	-5.41	1.35	1.42
6	F	201	75U	C1-C10	-4.49	1.36	1.44
6	J	201	75U	C1-C10	-4.41	1.36	1.44
6	H	201	75U	C1-C10	-4.26	1.36	1.44
6	L	201	75U	C1-C10	-4.25	1.36	1.44
6	B	201	75U	C1-C10	-4.19	1.36	1.44
6	B	202	75U	C4-C9	-4.03	1.32	1.40
6	B	202	75U	C1-C10	-4.02	1.36	1.44
6	J	201	75U	C4-C9	-3.97	1.32	1.40
6	F	201	75U	C4-C9	-3.89	1.32	1.40
6	H	201	75U	C4-C9	-3.86	1.33	1.40
6	L	201	75U	C4-C9	-3.73	1.33	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	201	75U	C4-C9	-3.71	1.33	1.40
6	J	201	75U	C11-C10	-3.16	1.35	1.41
6	J	201	75U	C11-C7	-3.08	1.34	1.40
6	L	201	75U	C11-C7	-3.04	1.34	1.40
6	F	201	75U	C11-C7	-3.00	1.34	1.40
6	H	201	75U	C11-C7	-2.84	1.34	1.40
6	F	201	75U	C11-C10	-2.77	1.36	1.41
6	L	201	75U	C11-C10	-2.77	1.36	1.41
6	B	202	75U	C11-C7	-2.63	1.34	1.40
6	B	201	75U	C11-C7	-2.61	1.35	1.40
6	H	201	75U	C11-C10	-2.60	1.36	1.41
6	B	202	75U	C11-C10	-2.55	1.36	1.41
6	B	201	75U	C11-C10	-2.42	1.36	1.41
5	D	201	NAG	C1-C2	2.03	1.55	1.52

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	B	201	75U	C4-C3-C2	-6.55	116.03	121.58
6	B	202	75U	BR5-C3-C2	6.16	125.72	118.80
6	L	201	75U	C4-C3-C2	-6.13	116.39	121.58
6	H	201	75U	C4-C3-C2	-6.03	116.47	121.58
6	B	202	75U	C4-C3-C2	-6.02	116.48	121.58
6	F	201	75U	C4-C3-C2	-5.76	116.70	121.58
6	J	201	75U	C4-C3-C2	-5.65	116.80	121.58
6	J	201	75U	BR5-C3-C2	5.20	124.63	118.80
6	H	201	75U	C28-O27-C26	-4.96	104.91	116.46
6	H	201	75U	BR5-C3-C2	4.87	124.26	118.80
6	B	202	75U	C31-C1-C10	4.71	126.64	120.96
6	B	201	75U	BR5-C3-C2	4.54	123.89	118.80
6	F	201	75U	BR5-C3-C2	4.47	123.81	118.80
6	B	201	75U	C31-C1-C10	4.36	126.22	120.96
6	L	201	75U	C28-O27-C26	-4.28	106.50	116.46
6	B	202	75U	C28-O27-C26	-4.27	106.52	116.46
6	F	201	75U	C28-O27-C26	-4.22	106.64	116.46
6	H	201	75U	C31-C1-C10	4.22	126.05	120.96
6	H	201	75U	C12-N8-C7	4.16	128.17	124.09
6	L	201	75U	BR5-C3-C2	4.11	123.42	118.80
6	L	201	75U	C12-N8-C7	4.09	128.10	124.09
6	J	201	75U	C31-C1-C10	4.08	125.88	120.96
6	B	201	75U	C28-O27-C26	-4.04	107.06	116.46
6	B	201	75U	C31-C1-C2	-4.02	113.96	120.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	L	201	75U	C31-C1-C10	3.96	125.74	120.96
6	B	201	75U	C12-N8-C7	3.92	127.93	124.09
6	B	201	75U	C11-C10-C1	3.92	137.20	132.03
6	B	202	75U	C11-C10-C1	3.91	137.19	132.03
6	B	202	75U	C31-C1-C2	-3.76	114.37	120.34
6	L	201	75U	C11-C10-C1	3.72	136.93	132.03
6	J	201	75U	C12-N8-C7	3.69	127.71	124.09
6	F	201	75U	C31-C1-C10	3.67	125.39	120.96
6	H	201	75U	C11-C10-C1	3.59	136.77	132.03
6	H	201	75U	C31-C1-C2	-3.58	114.64	120.34
6	J	201	75U	C28-O27-C26	-3.53	108.25	116.46
6	L	201	75U	C31-C1-C2	-3.41	114.93	120.34
6	F	201	75U	C11-C10-C1	3.30	136.39	132.03
6	B	202	75U	C12-N8-C7	3.30	127.32	124.09
6	F	201	75U	C31-C1-C2	-2.95	115.65	120.34
6	J	201	75U	C11-C10-C1	2.92	135.88	132.03
6	J	201	75U	C31-C1-C2	-2.91	115.72	120.34
6	F	201	75U	C12-N8-C7	2.67	126.71	124.09
5	L	202	NAG	C1-O5-C5	2.60	115.71	112.19
5	H	202	NAG	C1-O5-C5	2.24	115.23	112.19
5	F	202	NAG	C1-O5-C5	2.14	115.09	112.19
6	B	201	75U	O27-C26-O30	-2.11	119.38	123.67
6	H	201	75U	O27-C26-C11	2.03	117.96	112.10
5	J	202	NAG	C1-O5-C5	2.02	114.92	112.19
6	B	202	75U	C4-C9-C10	2.01	123.96	120.92

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	201	75U	C11-C26-O27-C28
6	B	202	75U	C11-C26-O27-C28
6	H	201	75U	C11-C26-O27-C28
5	B	203	NAG	O5-C5-C6-O6
5	L	202	NAG	O5-C5-C6-O6
5	L	202	NAG	C4-C5-C6-O6
6	B	201	75U	O30-C26-O27-C28
5	E	404	NAG	C4-C5-C6-O6
6	B	202	75U	O30-C26-O27-C28
6	L	201	75U	C11-C26-O27-C28
5	C	404	NAG	O5-C5-C6-O6
5	K	404	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
6	F	201	75U	C11-C26-O27-C28
5	B	203	NAG	C4-C5-C6-O6
5	E	404	NAG	O5-C5-C6-O6
5	J	203	NAG	C4-C5-C6-O6
6	H	201	75U	O30-C26-O27-C28
5	K	404	NAG	C4-C5-C6-O6
6	L	201	75U	O30-C26-O27-C28
5	C	404	NAG	C4-C5-C6-O6
6	F	201	75U	O30-C26-O27-C28
5	J	203	NAG	O5-C5-C6-O6
5	A	402	NAG	O5-C5-C6-O6
6	J	201	75U	C7-C11-C26-O27
6	J	201	75U	C7-C11-C26-O30
6	F	201	75U	C7-C13-S14-C18
6	J	201	75U	C17-C18-S14-C13
6	B	201	75U	C19-C18-S14-C13
6	L	201	75U	C19-C18-S14-C13
6	B	202	75U	C17-C18-S14-C13
6	F	201	75U	C19-C18-S14-C13
6	H	201	75U	C17-C18-S14-C13
6	H	201	75U	C19-C18-S14-C13
6	B	202	75U	C19-C18-S14-C13
6	B	201	75U	C17-C18-S14-C13
6	J	201	75U	C19-C18-S14-C13
6	F	201	75U	C17-C18-S14-C13
6	L	201	75U	C17-C18-S14-C13
6	B	201	75U	C7-C13-S14-C18
6	L	201	75U	C7-C13-S14-C18
6	J	201	75U	C29-C28-O27-C26
6	J	201	75U	C11-C26-O27-C28

There are no ring outliers.

8 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	201	75U	1	0
5	H	202	NAG	1	0
5	B	203	NAG	1	0
6	J	201	75U	2	0
6	H	201	75U	1	0
5	L	202	NAG	2	0
6	B	202	75U	1	0

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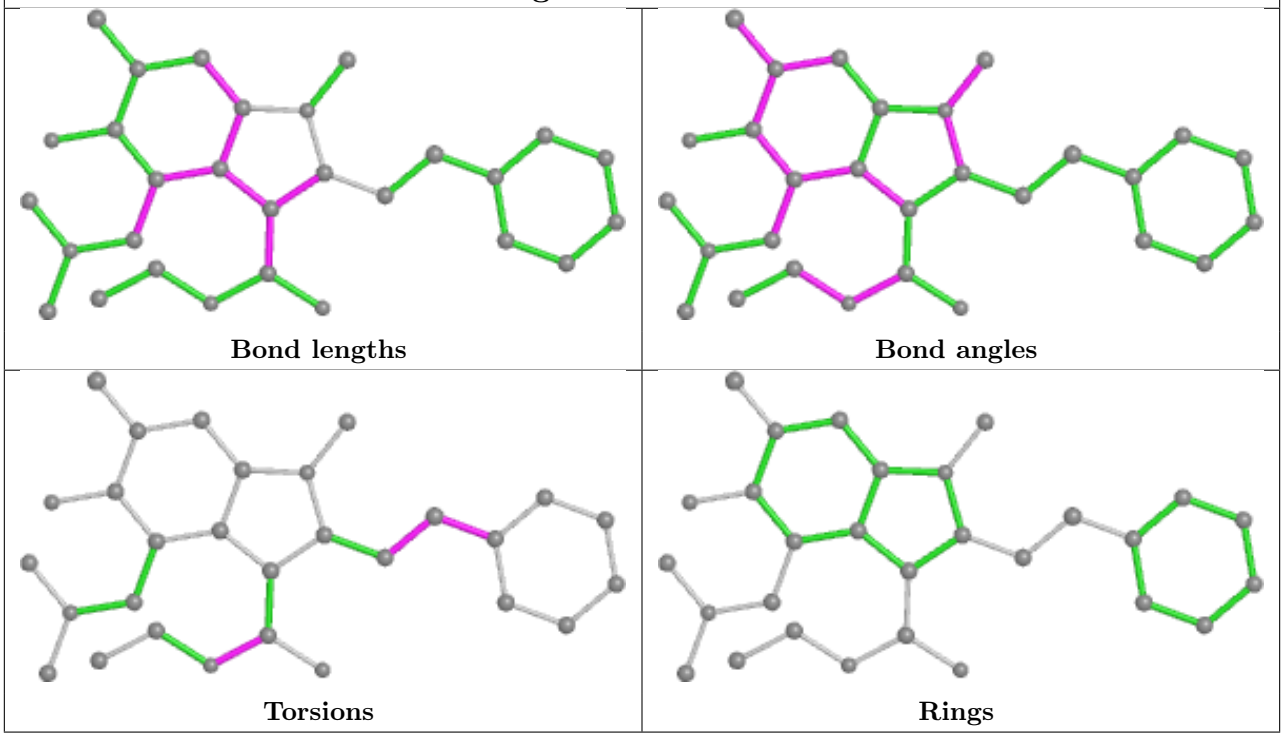
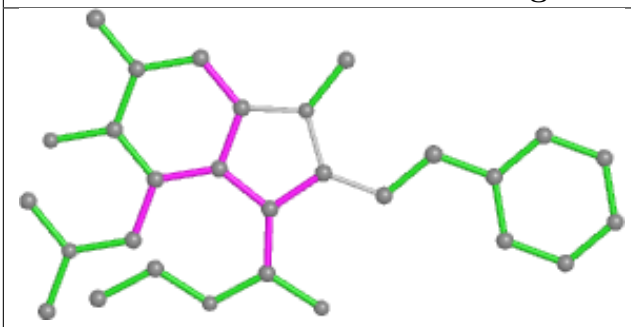
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	201	NAG	2	0

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.

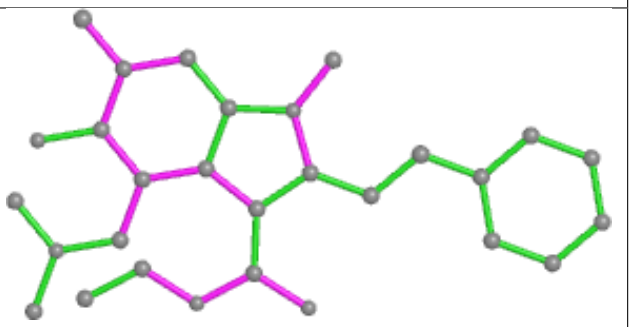
Ligand 75U F 201



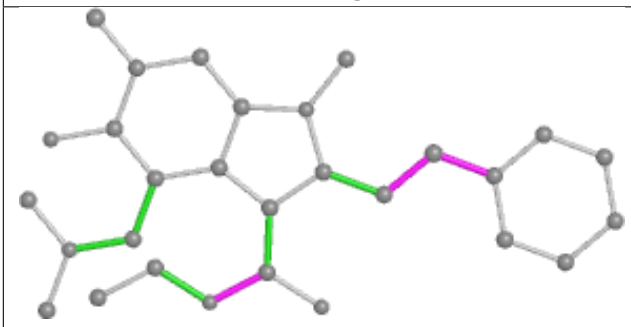
Ligand 75U B 201



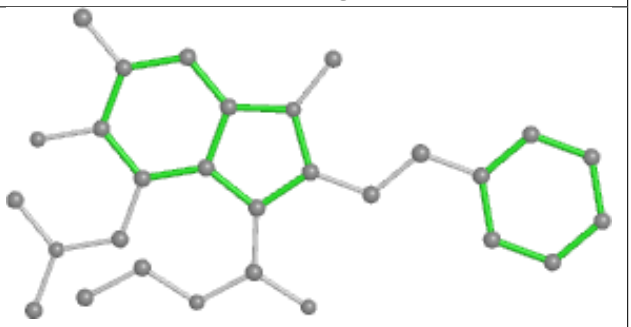
Bond lengths



Bond angles

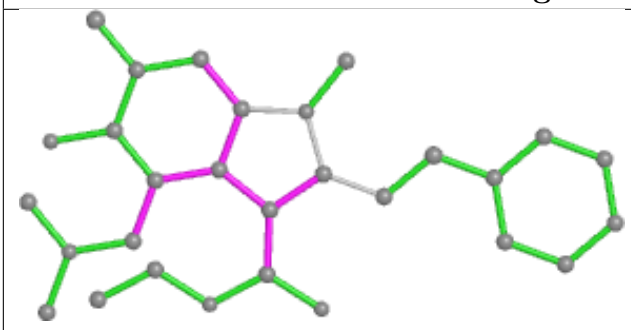


Torsions

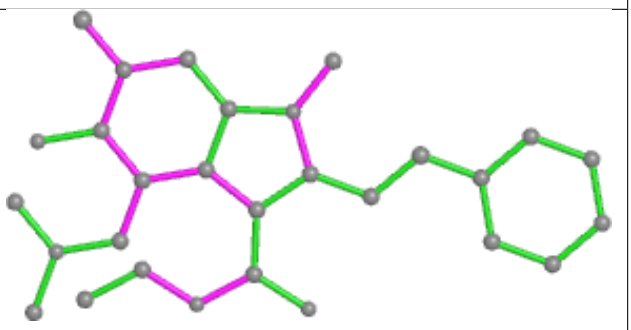


Rings

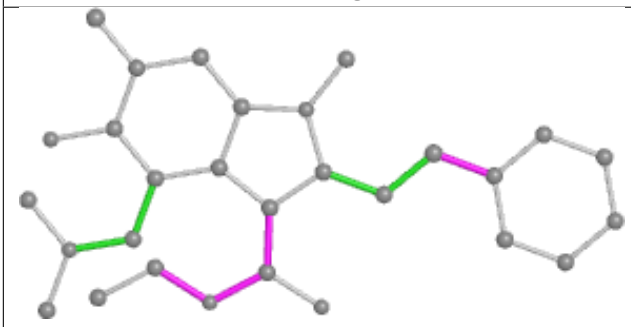
Ligand 75U J 201



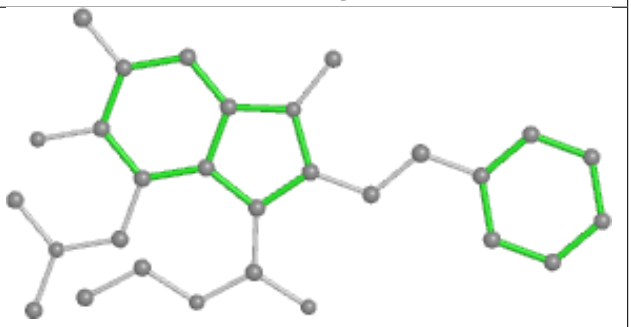
Bond lengths



Bond angles

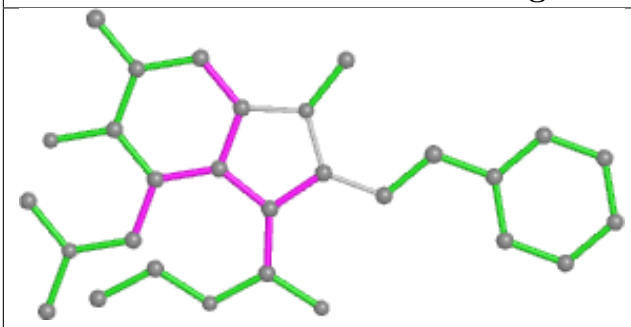


Torsions

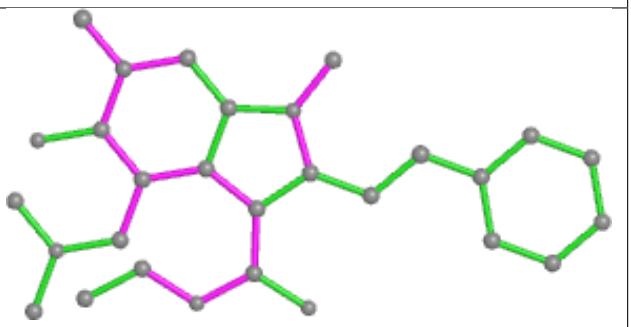


Rings

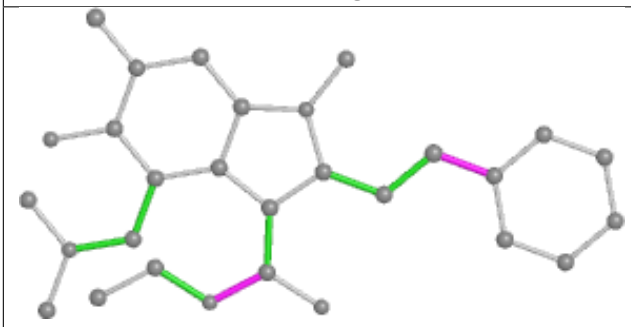
Ligand 75U H 201



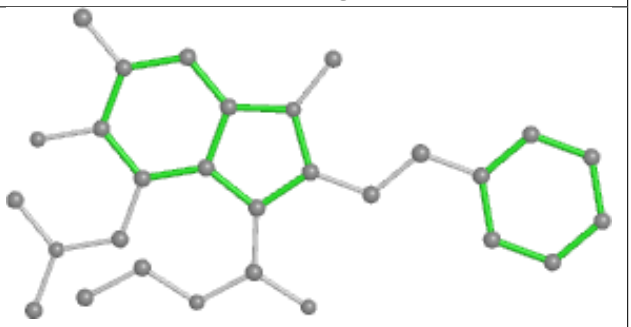
Bond lengths



Bond angles

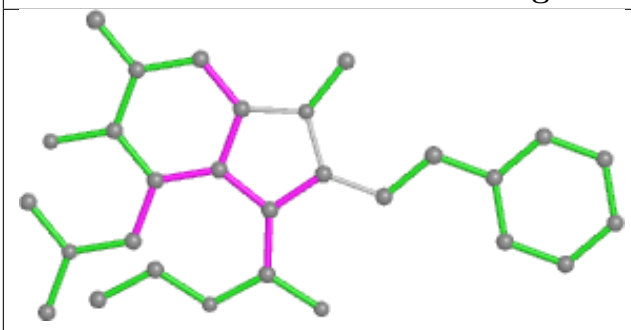


Torsions

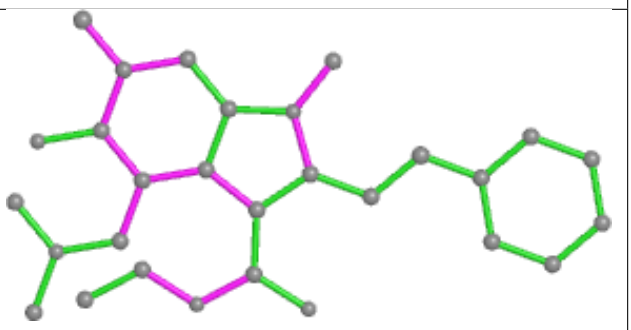


Rings

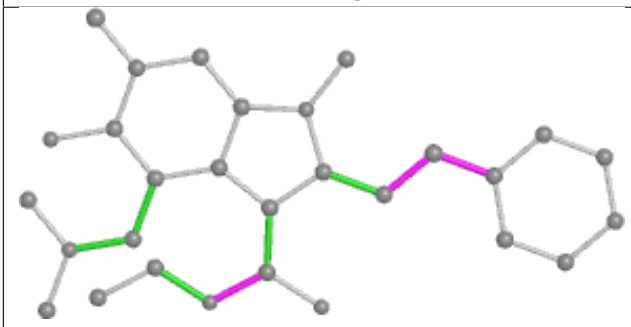
Ligand 75U L 201



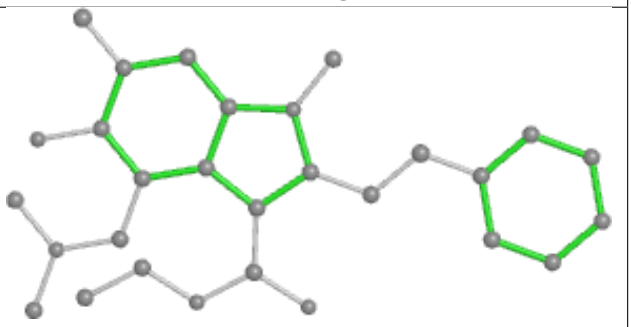
Bond lengths



Bond angles

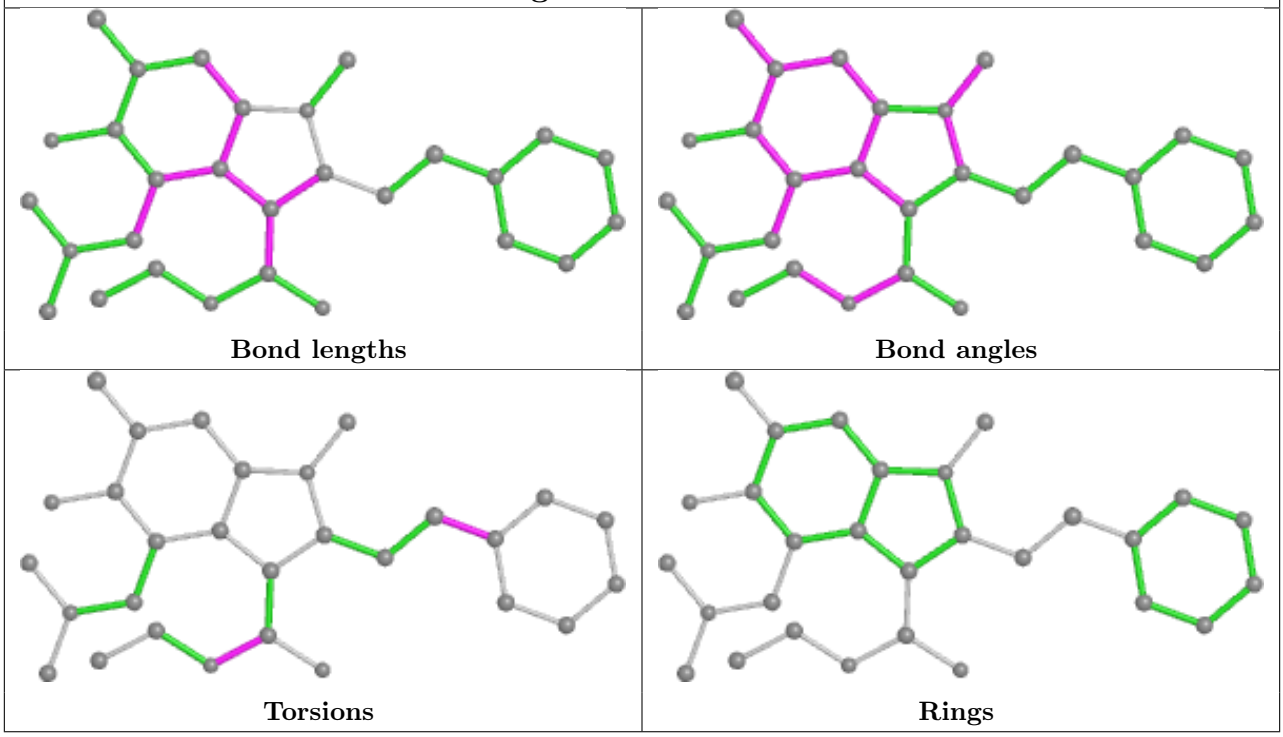


Torsions



Rings

Ligand 75U B 202



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	316/321 (98%)	-0.17	1 (0%) 94 97	24, 39, 53, 73	0
1	C	316/321 (98%)	-0.16	0 100 100	24, 36, 57, 74	1 (0%)
1	E	316/321 (98%)	-0.03	3 (0%) 84 90	24, 38, 58, 73	1 (0%)
1	G	316/321 (98%)	0.56	26 (8%) 11 17	34, 58, 83, 103	0
1	I	316/321 (98%)	0.16	8 (2%) 57 67	29, 46, 75, 107	0
1	K	316/321 (98%)	0.21	15 (4%) 31 44	31, 51, 74, 118	0
2	B	171/183 (93%)	0.14	3 (1%) 68 77	25, 40, 58, 117	0
2	D	171/183 (93%)	0.27	5 (2%) 51 62	23, 39, 57, 73	0
2	F	171/183 (93%)	0.23	2 (1%) 79 86	25, 36, 58, 93	0
2	H	171/183 (93%)	0.40	5 (2%) 51 62	31, 53, 79, 104	0
2	J	171/183 (93%)	0.38	11 (6%) 19 28	28, 49, 74, 103	0
2	L	171/183 (93%)	0.18	2 (1%) 79 86	29, 40, 62, 87	0
All	All	2922/3024 (96%)	0.15	81 (2%) 53 64	23, 43, 72, 118	2 (0%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	224	ASN	4.3
1	K	289	ILE	4.1
1	G	140	ARG	4.0
1	G	155	LEU	3.9
1	G	289	ILE	3.7
2	D	4	GLY	3.7
2	B	60	ASN	3.6
2	H	31	GLY	3.4
1	G	215	PRO	3.3
1	K	160	ALA	3.3
2	J	38	TYR	3.2

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Mol	Chain	Res	Type	RSRZ
2	J	150	ALA	3.1
2	B	3	PHE	3.1
1	I	212	SER	3.1
1	E	289	ILE	3.1
2	H	168	GLN	3.0
1	I	211	GLN	2.9
2	J	4	GLY	2.9
1	K	140	ARG	2.9
1	K	131	ARG	2.9
1	G	324	PRO	2.9
1	K	127	TYR	2.9
1	I	289	ILE	2.8
1	I	207	SER	2.8
1	G	213	PHE	2.8
2	F	60	ASN	2.8
1	G	278	GLU	2.8
1	I	140	ARG	2.8
2	D	33	GLY	2.7
1	K	271	GLY	2.7
2	J	60	ASN	2.6
1	G	128	SER	2.6
1	G	20	VAL	2.6
1	K	139	CYS	2.6
2	J	168	GLN	2.6
1	G	231	ASP	2.6
2	F	3	PHE	2.5
1	K	144	GLY	2.5
1	G	11	ASP	2.5
1	I	157	ASN	2.5
1	G	276	ALA	2.4
1	G	173	LYS	2.4
2	H	162	TYR	2.4
2	D	171	ILE	2.4
1	G	129	GLY	2.4
1	G	59	VAL	2.4
2	D	3	PHE	2.4
1	I	128	SER	2.4
1	E	22	ASN	2.4
1	G	272	VAL	2.4
1	G	273	GLN	2.3
1	G	274	VAL	2.3
1	K	159	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
2	J	154	ASN	2.3
2	L	164	GLU	2.3
1	K	158	THR	2.3
1	K	74	PRO	2.3
2	H	33	GLY	2.3
1	G	230	ILE	2.2
2	J	33	GLY	2.2
1	A	173	LYS	2.2
2	J	18	ILE	2.2
1	I	160	ALA	2.2
2	L	168	GLN	2.2
1	K	155	LEU	2.2
2	J	158	ASP	2.2
1	G	19	ALA	2.1
2	J	31	GLY	2.1
2	J	156	THR	2.1
2	H	171	ILE	2.1
1	G	171	THR	2.1
1	G	223	VAL	2.1
1	G	54	LYS	2.1
2	D	150	ALA	2.1
2	B	113	SER	2.1
1	G	276(A)	ASN	2.0
1	K	156	SER	2.0
1	E	57	ARG	2.0
1	K	224	ASN	2.0
1	G	21	SER	2.0
1	K	187	SER	2.0

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

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

6.3 Carbohydrates (i)

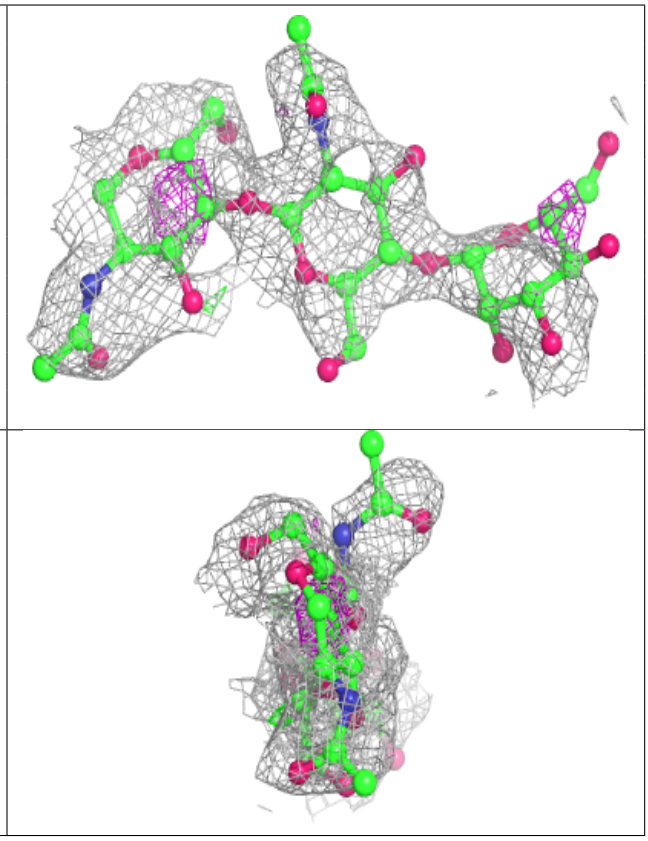
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	BMA	R	3	11/12	0.40	0.55	141,146,148,148	0
3	NAG	Q	2	14/15	0.57	0.39	102,112,117,125	0
3	BMA	Q	3	11/12	0.58	0.36	130,133,134,134	0
3	BMA	N	3	11/12	0.62	0.43	119,126,128,128	0
3	NAG	M	1	14/15	0.63	0.27	79,85,93,98	0
3	BMA	M	3	11/12	0.63	0.46	121,122,125,125	0
4	NAG	O	2	14/15	0.67	0.49	114,122,126,126	0
4	NAG	P	2	14/15	0.67	0.58	105,113,116,117	0
3	BMA	S	3	11/12	0.68	0.45	108,114,116,118	0
4	NAG	O	1	14/15	0.73	0.32	73,78,90,102	0
3	NAG	S	1	14/15	0.74	0.32	72,79,85,94	0
3	NAG	S	2	14/15	0.75	0.35	103,107,112,112	0
3	NAG	R	1	14/15	0.78	0.32	67,76,86,87	0
3	NAG	R	2	14/15	0.80	0.40	74,93,106,122	0
4	NAG	P	1	14/15	0.80	0.26	100,106,110,113	0
3	NAG	N	2	14/15	0.80	0.30	71,84,92,106	0
3	NAG	M	2	14/15	0.81	0.36	95,105,111,116	0
3	NAG	Q	1	14/15	0.81	0.26	77,81,92,99	0
3	NAG	N	1	14/15	0.84	0.24	61,67,77,77	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.

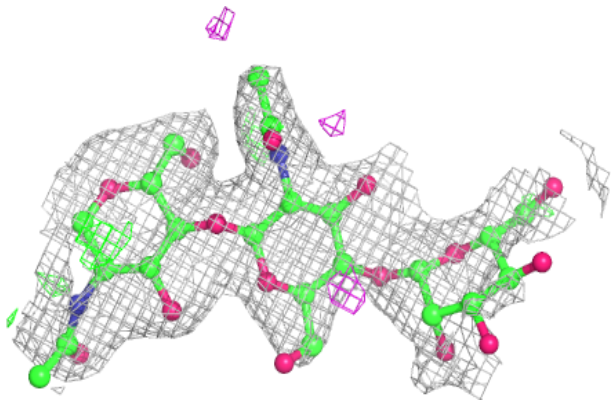
Electron density around Chain M:

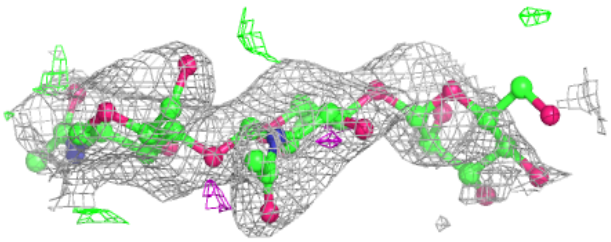
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

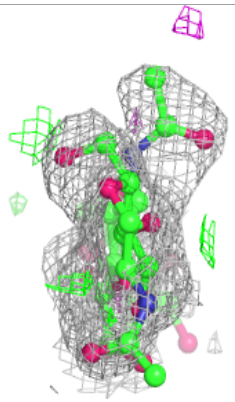


Electron density around Chain N:

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

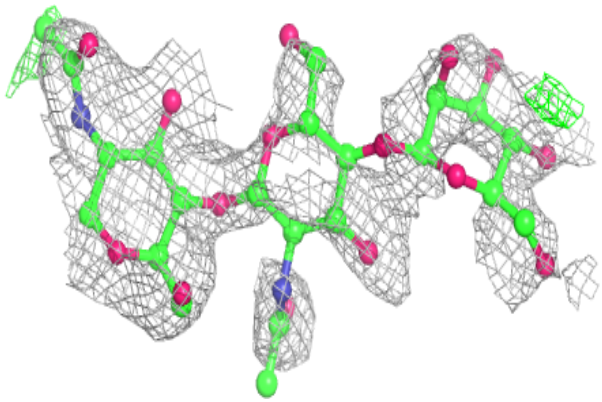


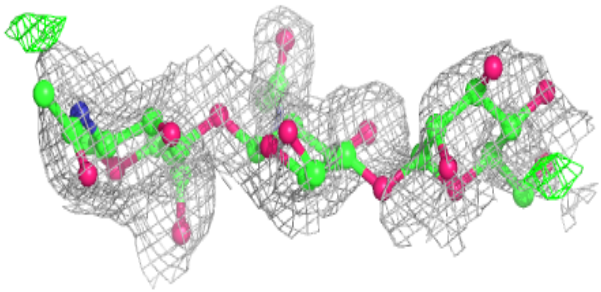


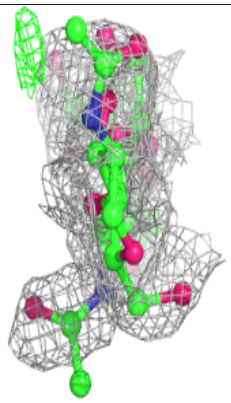


Electron density around Chain Q:

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

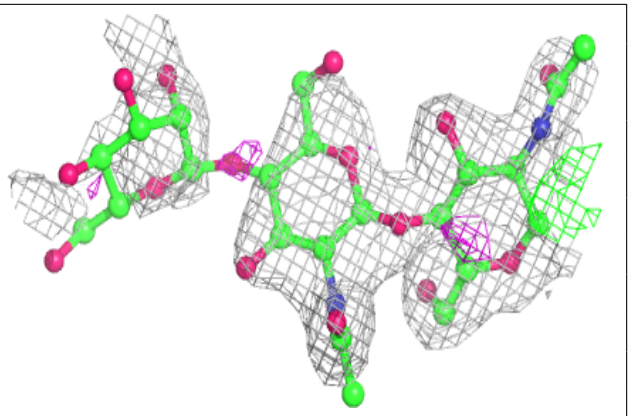


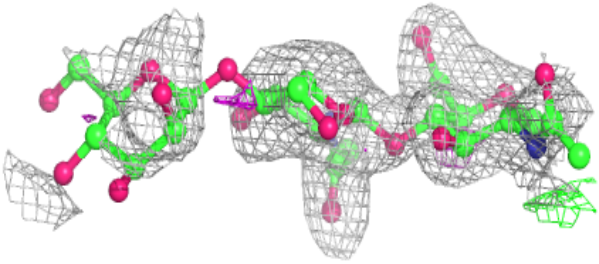


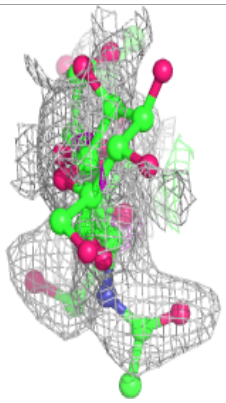


Electron density around Chain R:

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

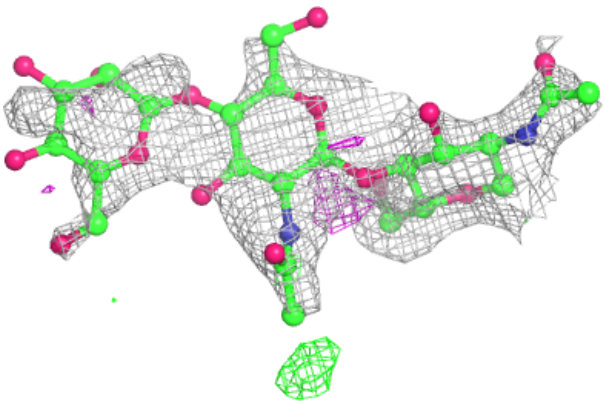


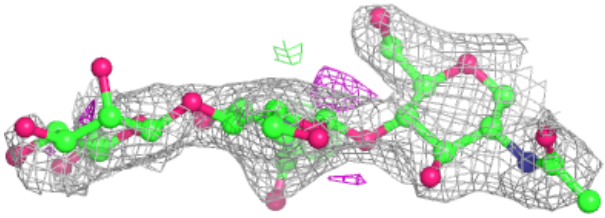


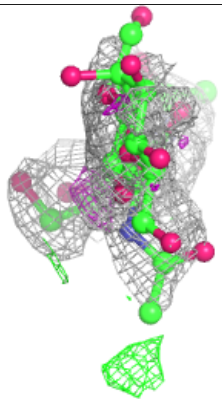


Electron density around Chain S:

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

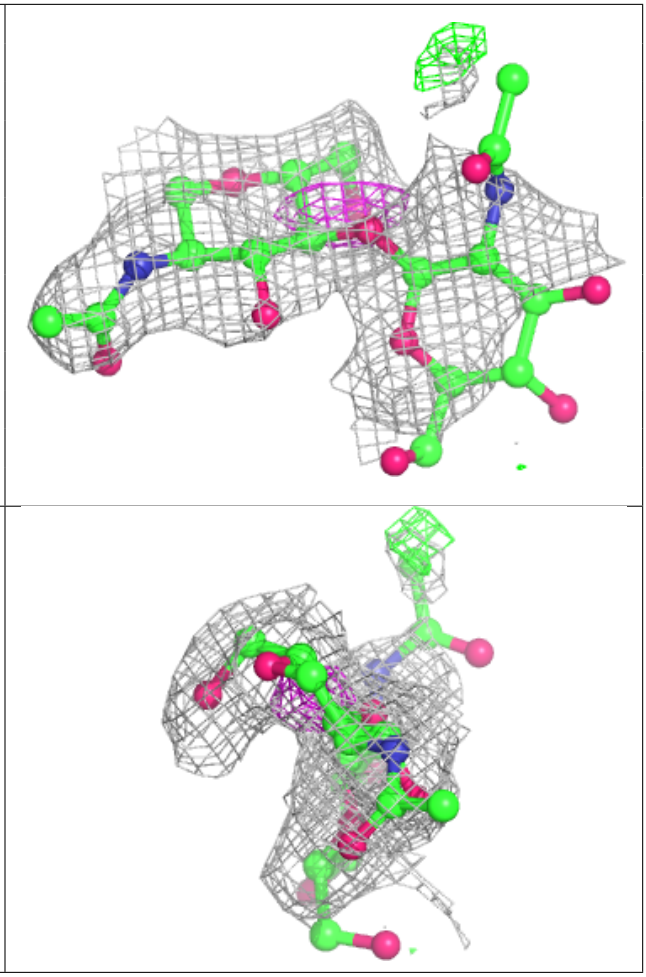


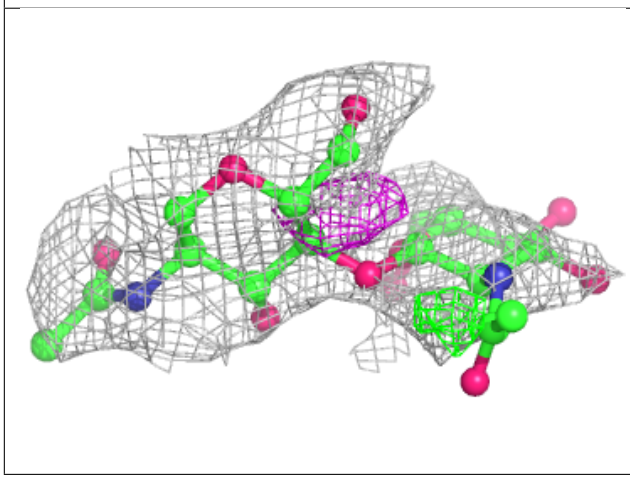




Electron density around Chain O:

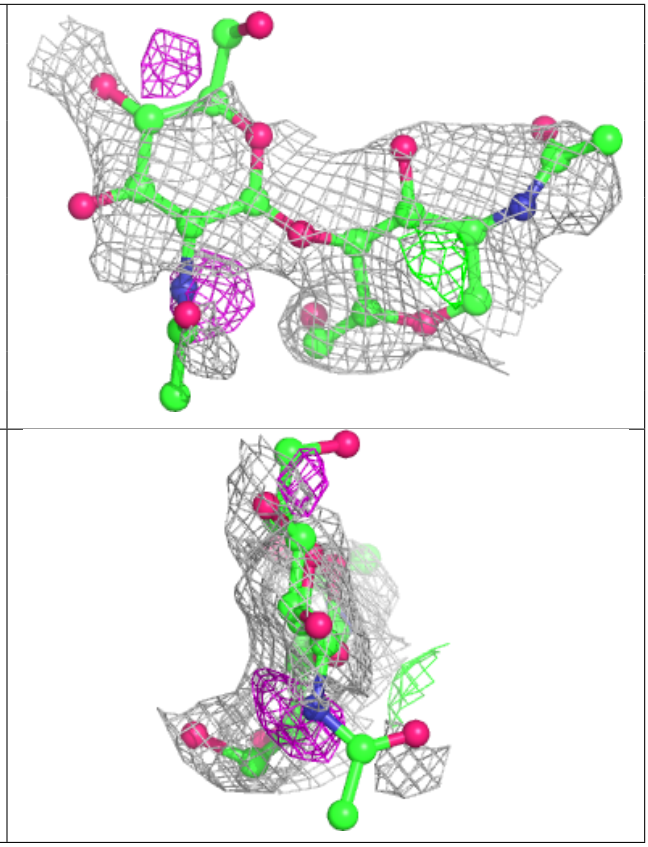
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





Electron density around Chain P:

$2mF_o$-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
5	NAG	A	401	14/15	0.45	0.40	100,112,118,118	0
5	NAG	D	202	14/15	0.55	0.40	101,106,111,114	0
5	NAG	I	404	14/15	0.57	0.35	97,100,102,105	0
5	NAG	J	203	14/15	0.67	0.46	96,108,113,116	0
5	NAG	K	404	14/15	0.68	0.33	100,105,108,108	0
5	NAG	E	404	14/15	0.72	0.27	94,98,104,105	0
5	NAG	B	204	14/15	0.73	0.28	90,98,101,102	0
5	NAG	L	202	14/15	0.75	0.32	57,67,69,71	0
6	75U	H	201	29/29	0.78	0.26	58,74,84,109	0
5	NAG	D	201	14/15	0.79	0.24	50,57,63,65	0
5	NAG	B	203	14/15	0.80	0.21	58,62,65,68	0
6	75U	J	201	29/29	0.80	0.24	54,73,88,105	0
6	75U	F	201	29/29	0.81	0.23	44,60,77,99	0
5	NAG	J	202	14/15	0.81	0.20	59,66,69,70	0

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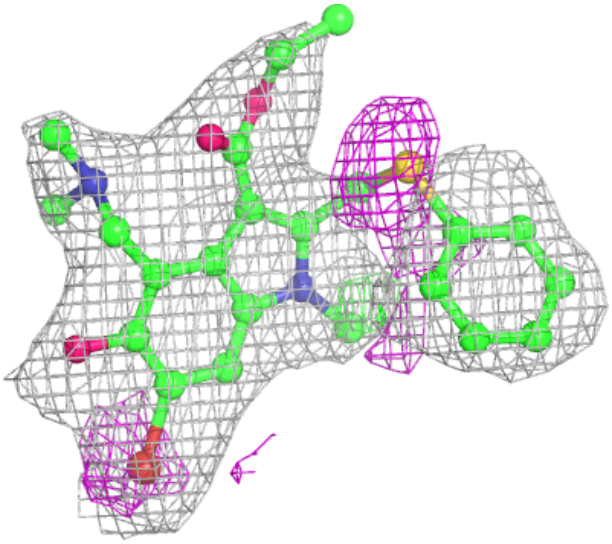
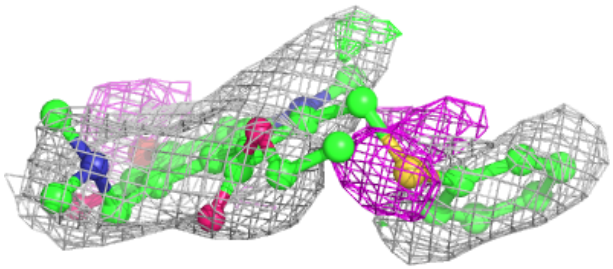
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	NAG	H	202	14/15	0.81	0.24	46,63,73,77	0
6	75U	B	202	29/29	0.82	0.25	51,61,75,98	0
6	75U	L	201	29/29	0.84	0.23	48,61,75,94	0
6	75U	B	201	29/29	0.86	0.19	48,58,73,90	0
5	NAG	C	404	14/15	0.86	0.41	94,97,102,105	0
5	NAG	A	402	14/15	0.87	0.38	66,74,81,84	0
5	NAG	F	202	14/15	0.87	0.17	46,53,65,68	0
7	NA	J	204	1/1	0.94	0.56	56,56,56,56	0

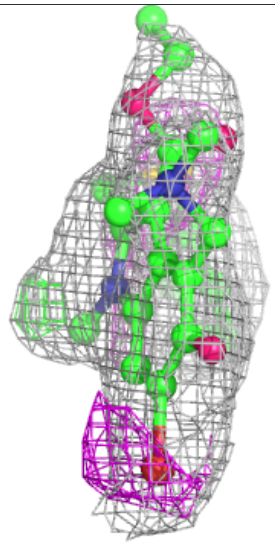
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 75U H 201:

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

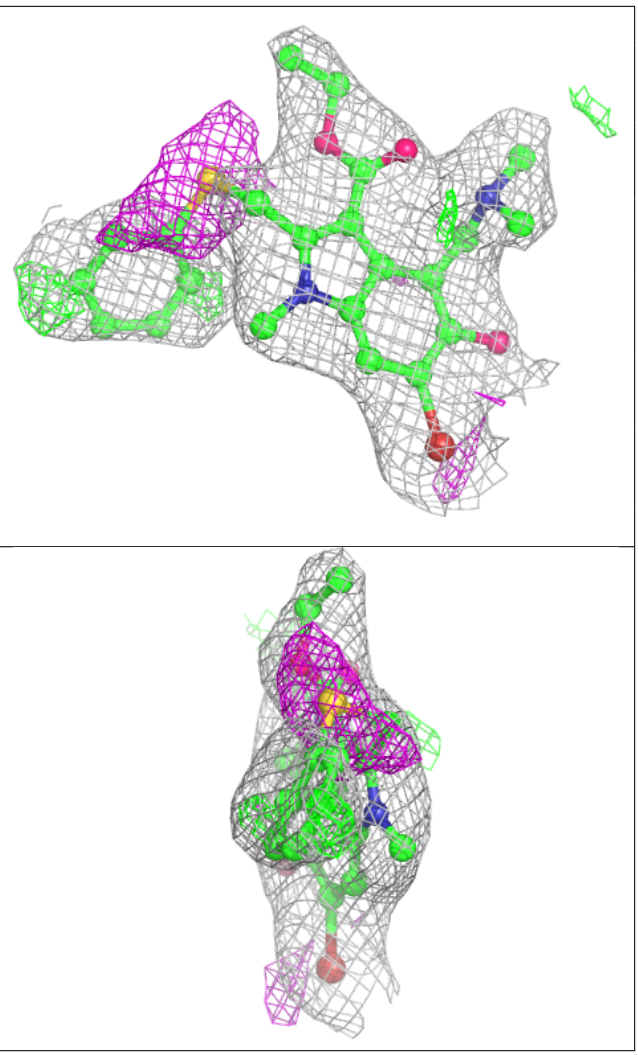






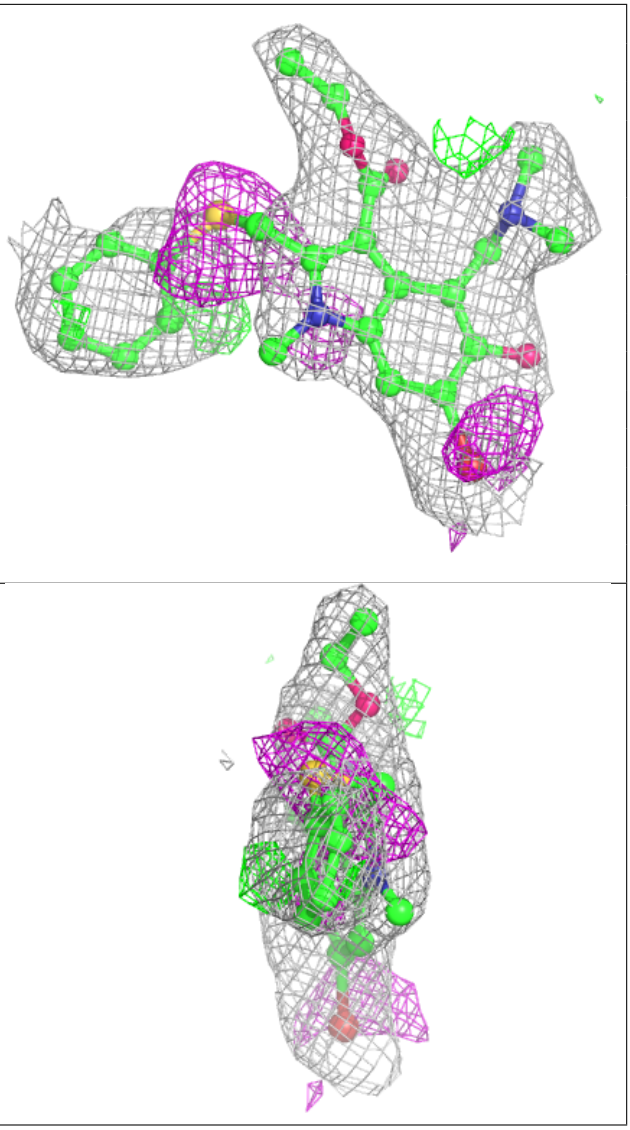
Electron density around 75U J 201:

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



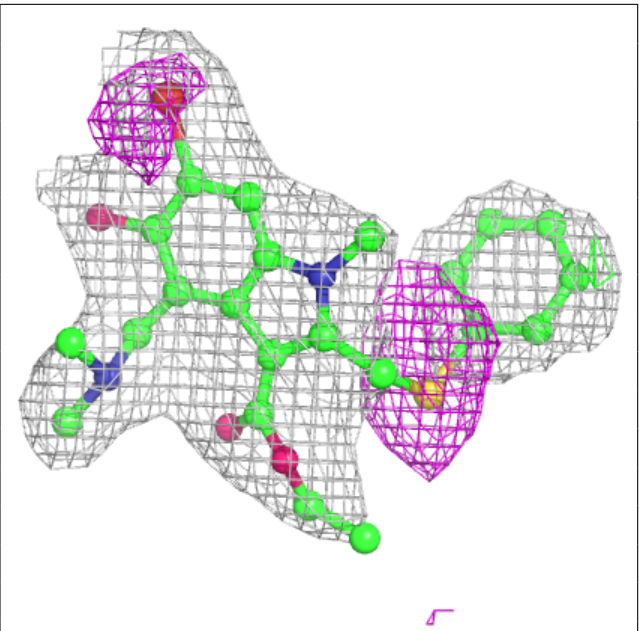
Electron density around 75U F 201:

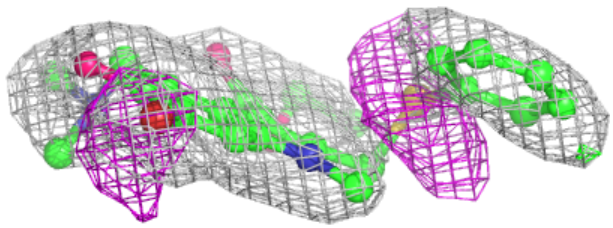
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

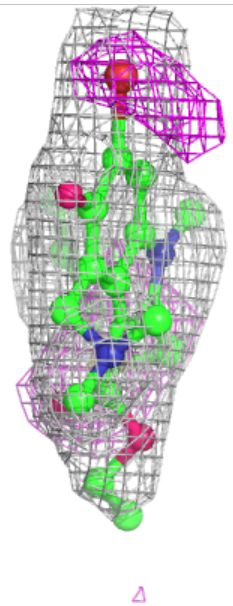


Electron density around 75U B 202:

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

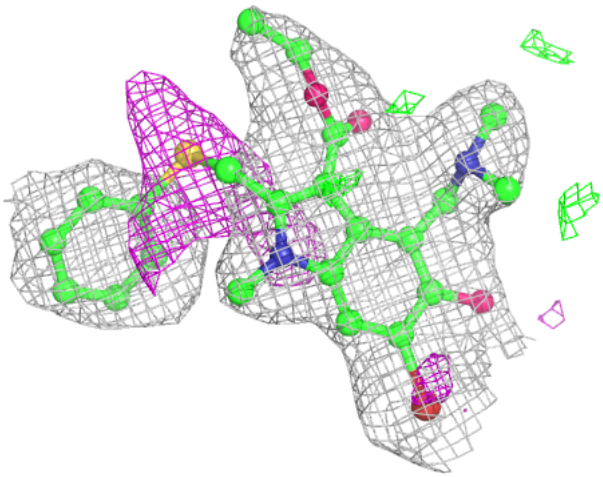


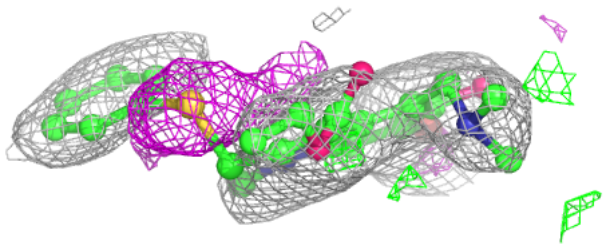


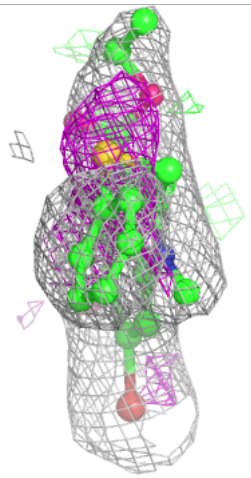


Electron density around 75U L 201:

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

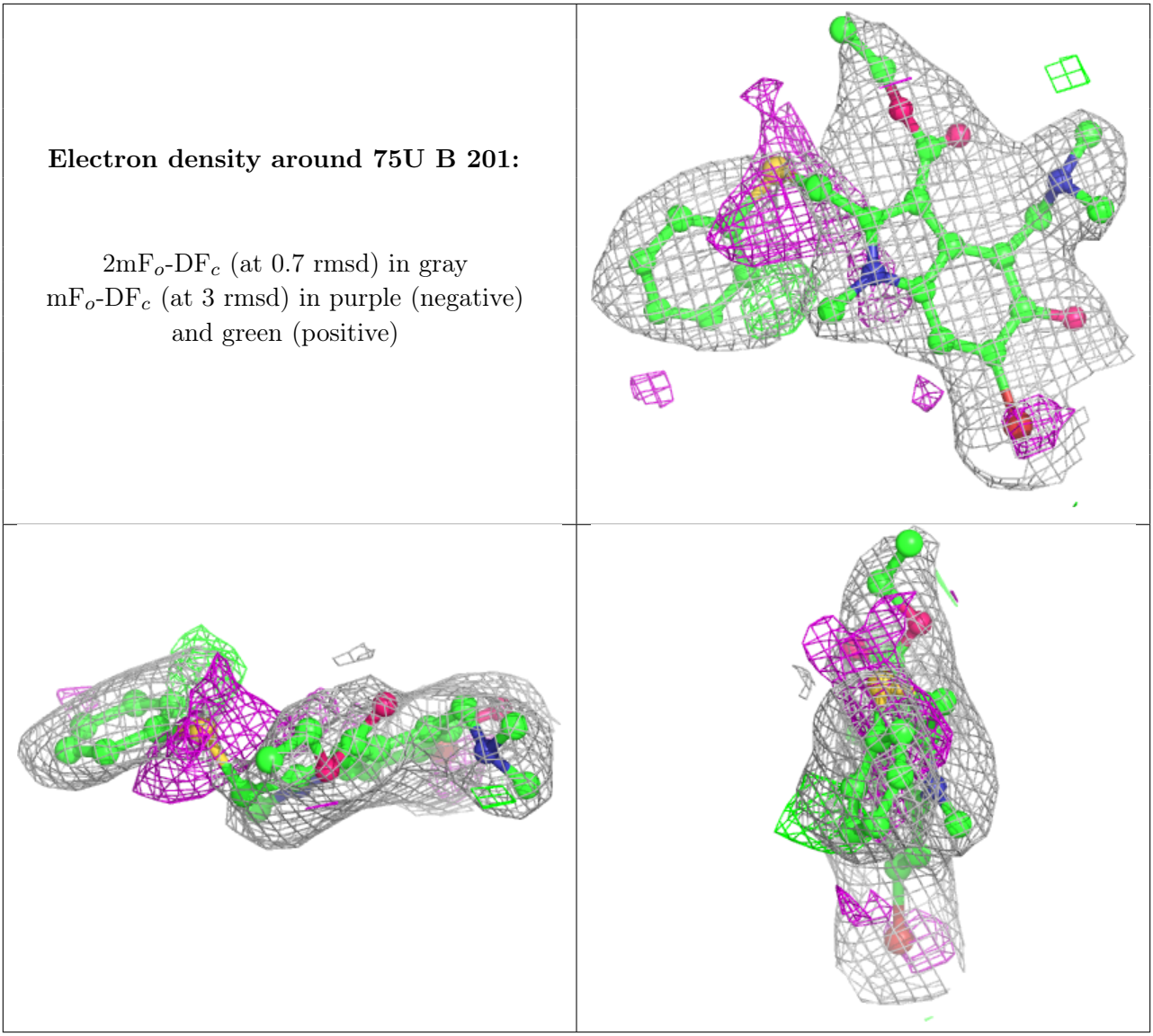






Electron density around 75U B 201:

$2mF_o$-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.