Full wwPDB X-ray Structure Validation Report ⓘ

Dec 15, 2020 – 11:10 AM GMT

PDB ID	:	6YGG
Title	:	NADase from Aspergillus fumigatus complexed with a substrate anologue
Authors	:	Stromland, O.; Ziegler, M.; Kallio, J.P.
Deposited on	:	2020-03-27
Resolution	:	1.85 Å(reported)

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

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

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

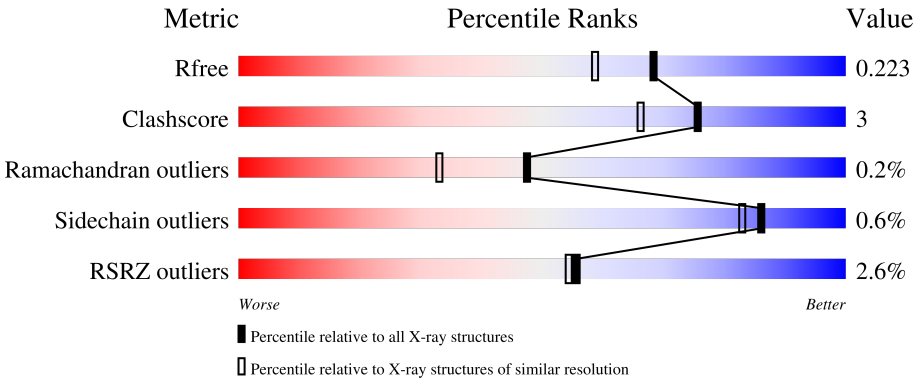
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.15.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.15.1

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 1.85 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	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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	248	2% 78% 6% 16%
1	B	248	3% 79% 6% 16%
2	C	3	33% 67%
2	E	3	67% 33%
2	F	3	100%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	G	3	
3	D	2	

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 3998 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 AfNADase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	0	0
			1651	1069	262	313	7			
1	B	209	Total	C	N	O	S	0	0	0
			1651	1069	262	313	7			

There are 28 discrepancies between the modelled and reference sequences:

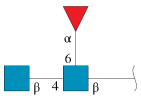
Chain	Residue	Modelled	Actual	Comment	Reference
A	235	ASP	-	expression tag	UNP Q4WL81
A	236	VAL	-	expression tag	UNP Q4WL81
A	237	LEU	-	expression tag	UNP Q4WL81
A	238	PHE	-	expression tag	UNP Q4WL81
A	239	GLN	-	expression tag	UNP Q4WL81
A	240	GLY	-	expression tag	UNP Q4WL81
A	241	PRO	-	expression tag	UNP Q4WL81
A	242	GLY	-	expression tag	UNP Q4WL81
A	243	HIS	-	expression tag	UNP Q4WL81
A	244	HIS	-	expression tag	UNP Q4WL81
A	245	HIS	-	expression tag	UNP Q4WL81
A	246	HIS	-	expression tag	UNP Q4WL81
A	247	HIS	-	expression tag	UNP Q4WL81
A	248	HIS	-	expression tag	UNP Q4WL81
B	235	ASP	-	expression tag	UNP Q4WL81
B	236	VAL	-	expression tag	UNP Q4WL81
B	237	LEU	-	expression tag	UNP Q4WL81
B	238	PHE	-	expression tag	UNP Q4WL81
B	239	GLN	-	expression tag	UNP Q4WL81
B	240	GLY	-	expression tag	UNP Q4WL81
B	241	PRO	-	expression tag	UNP Q4WL81
B	242	GLY	-	expression tag	UNP Q4WL81
B	243	HIS	-	expression tag	UNP Q4WL81
B	244	HIS	-	expression tag	UNP Q4WL81
B	245	HIS	-	expression tag	UNP Q4WL81

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	246	HIS	-	expression tag	UNP Q4WL81
B	247	HIS	-	expression tag	UNP Q4WL81
B	248	HIS	-	expression tag	UNP Q4WL81

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.

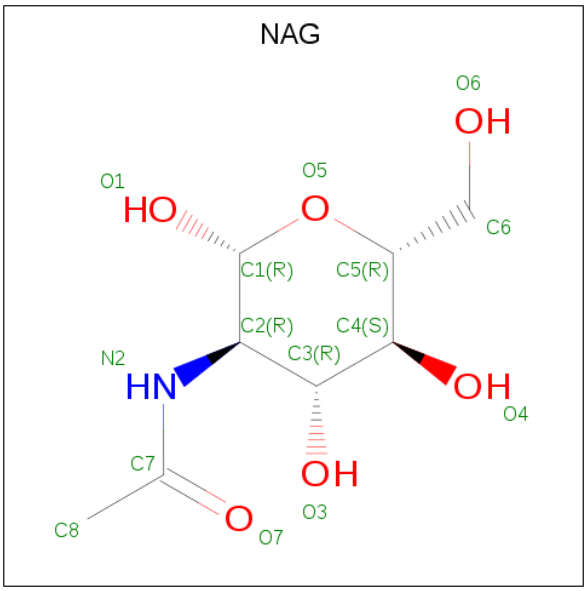


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	E	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	F	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	G	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 3 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	2	Total	C	N	O	0	0	0
			24	14	1	9			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

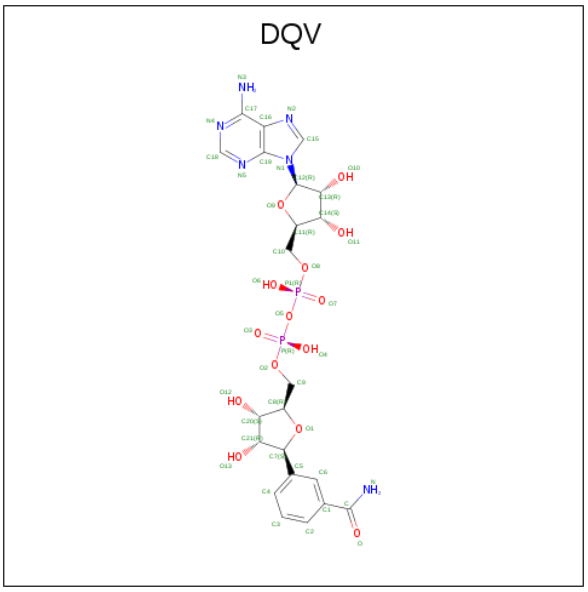
- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Ca 1 1	0	0
5	A	1	Total Ca 1 1	0	0

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

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

- Molecule 7 is [(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-3,4-dihydroxytetrahydrofuran-2-yl]methyl [(2R,3S,4R,5S)-5-(3-carbamoylphenyl)-3,4-dihydroxytetrahydrofuran-2-yl]methyl dihydrogen diphosphate (non-preferred name) (three-letter code: DQV) (formula: $C_{22}H_{28}N_6O_{14}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	0	0
			44	22	6	14	2		

- Molecule 8 is water.

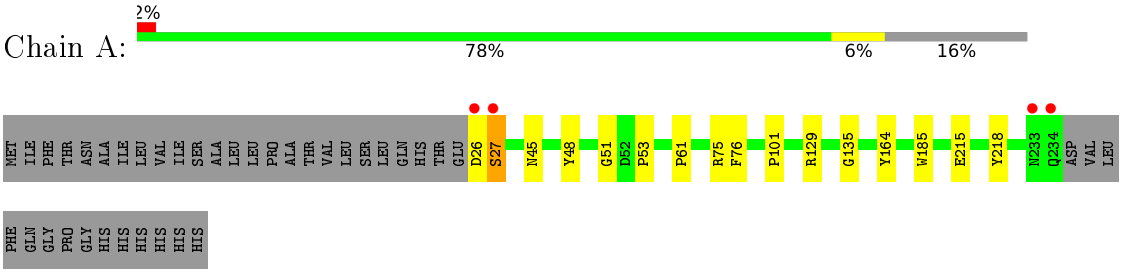
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	240	Total	O	0	0
			240	240		
8	B	218	Total	O	0	1
			219	219		

3 Residue-property plots

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

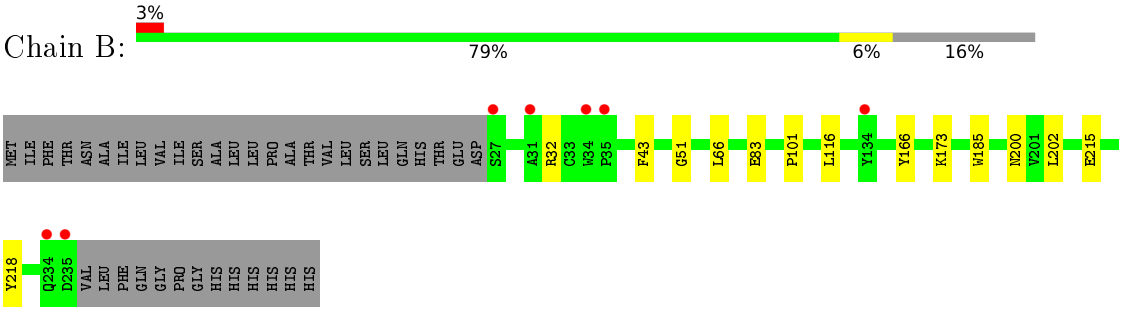
• Molecule 1: AfNADase

Chain A:



• Molecule 1: AfNADase

Chain B:



• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:

• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:

• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:

• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:

• Molecule 3: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	51.05Å 64.32Å 161.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.66 – 1.85 48.66 – 1.85	Depositor EDS
% Data completeness (in resolution range)	98.5 (48.66-1.85) 98.6 (48.66-1.85)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.44 (at 1.86Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.185 , 0.226 0.181 , 0.223	Depositor DCC
R_{free} test set	2313 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	25.0	Xtriage
Anisotropy	0.840	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 50.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3998	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, DQV, CA, NAG, FUC

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	0.39	0/1708	0.57	0/2343
1	B	0.38	0/1708	0.56	0/2343
All	All	0.39	0/3416	0.56	0/4686

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1651	0	1552	9	0
1	B	1651	0	1552	9	0
2	C	38	0	34	0	0
2	E	38	0	34	0	0
2	F	38	0	34	0	0
2	G	38	0	34	2	0
3	D	24	0	22	0	0
4	A	14	0	13	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	44	0	0	1	0
8	A	240	0	0	0	0
8	B	219	0	0	1	0
All	All	3998	0	3275	21	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 3.

All (21) 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)
7:B:510:DQV:C12	7:B:510:DQV:O9	1.67	1.22
1:A:26:ASP:CG	1:A:27:SER:H	2.00	0.65
1:B:32:ARG:NH2	1:B:83:GLU:OE2	2.36	0.55
1:A:129:ARG:NH2	1:A:135:GLY:HA3	2.24	0.52
1:B:215:GLU:HA	1:B:218:TYR:CE1	2.45	0.52
1:B:200:ASN:ND2	8:B:605:HOH:O	2.39	0.50
1:A:26:ASP:O	1:A:27:SER:OG	2.23	0.49
2:G:2:NAG:H82	2:G:2:NAG:O3	2.13	0.47
1:B:66:LEU:HD21	1:B:166:TYR:CG	2.49	0.47
1:A:45:ASN:O	1:A:51:GLY:HA2	2.15	0.47
2:G:1:NAG:O3	2:G:1:NAG:O7	2.20	0.46
1:A:75:ARG:HD2	1:A:76:PHE:CZ	2.50	0.45
1:B:43:PHE:CD1	1:B:51:GLY:HA3	2.52	0.45
1:B:101:PRO:HD3	1:B:185:TRP:CD1	2.53	0.44
1:A:61:PRO:HG3	1:A:164:TYR:CD2	2.52	0.43
1:B:202:LEU:HD12	1:B:202:LEU:HA	1.80	0.43
1:B:215:GLU:HA	1:B:218:TYR:CZ	2.54	0.43
1:A:101:PRO:HB3	1:A:185:TRP:CE3	2.54	0.42
1:B:66:LEU:HA	1:B:66:LEU:HD23	1.74	0.42
1:A:48:TYR:CG	1:A:53:PRO:HA	2.56	0.41
1:A:215:GLU:HA	1:A:218:TYR:CE1	2.56	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	207/248 (84%)	204 (99%)	2 (1%)	1 (0%)	29 15
1	B	207/248 (84%)	204 (99%)	3 (1%)	0	100 100
All	All	414/496 (84%)	408 (99%)	5 (1%)	1 (0%)	47 33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	27	SER

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	177/211 (84%)	177 (100%)	0	100 100
1	B	177/211 (84%)	175 (99%)	2 (1%)	73 65
All	All	354/422 (84%)	352 (99%)	2 (1%)	86 83

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	116	LEU
1	B	173	LYS

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

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

14 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	NAG	C	1	1,2	14,14,15	0.20	0	17,19,21	0.58	0
2	NAG	C	2	2	14,14,15	1.11	1 (7%)	17,19,21	0.77	0
2	FUC	C	3	2	10,10,11	1.04	1 (10%)	14,14,16	1.04	1 (7%)
3	NAG	D	1	1,3	14,14,15	0.21	0	17,19,21	0.59	0
3	FUC	D	2	3	10,10,11	1.15	1 (10%)	14,14,16	0.94	1 (7%)
2	NAG	E	1	1,2	14,14,15	0.36	0	17,19,21	0.44	0
2	NAG	E	2	2	14,14,15	1.17	1 (7%)	17,19,21	0.81	1 (5%)
2	FUC	E	3	2	10,10,11	0.62	0	14,14,16	0.86	0
2	NAG	F	1	1,2	14,14,15	0.57	0	17,19,21	0.53	0
2	NAG	F	2	2	14,14,15	0.26	0	17,19,21	0.52	0
2	FUC	F	3	2	10,10,11	1.07	0	14,14,16	0.79	0
2	NAG	G	1	1,2	14,14,15	0.72	1 (7%)	17,19,21	0.89	1 (5%)
2	NAG	G	2	2	14,14,15	2.39	3 (21%)	17,19,21	2.06	5 (29%)
2	FUC	G	3	2	10,10,11	1.03	0	14,14,16	0.93	1 (7%)

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	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	FUC	C	3	2	-	-	0/1/1/1
3	NAG	D	1	1,3	-	0/6/23/26	0/1/1/1
3	FUC	D	2	3	-	-	0/1/1/1
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	FUC	E	3	2	-	-	0/1/1/1
2	NAG	F	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	FUC	F	3	2	-	-	0/1/1/1
2	NAG	G	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	G	2	2	-	4/6/23/26	0/1/1/1
2	FUC	G	3	2	-	-	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	2	NAG	C1-C2	6.90	1.62	1.52
2	G	2	NAG	O5-C1	4.75	1.51	1.43
2	E	2	NAG	O5-C1	4.25	1.50	1.43
2	C	2	NAG	O5-C1	-3.86	1.37	1.43
2	G	1	NAG	C1-C2	2.38	1.55	1.52
3	D	2	FUC	C1-C2	2.29	1.57	1.52
2	C	3	FUC	C1-C2	2.15	1.57	1.52
2	G	2	NAG	C3-C2	2.08	1.56	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	G	2	NAG	C4-C3-C2	4.67	117.87	111.02
2	G	2	NAG	C2-N2-C7	4.21	128.90	122.90
2	G	1	NAG	C2-N2-C7	2.95	127.10	122.90
2	G	2	NAG	O3-C3-C2	2.87	115.41	109.47
2	G	2	NAG	O4-C4-C3	-2.49	104.59	110.35
3	D	2	FUC	C1-C2-C3	2.26	112.44	109.67
2	G	3	FUC	O2-C2-C1	2.13	113.50	109.15
2	C	3	FUC	C1-O5-C5	2.08	117.49	112.78
2	G	2	NAG	O4-C4-C5	-2.04	104.22	109.30
2	E	2	NAG	C1-O5-C5	2.02	114.93	112.19

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	1	NAG	C3-C2-N2-C7
2	E	2	NAG	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	G	2	NAG	C8-C7-N2-C2
2	G	2	NAG	O7-C7-N2-C2
2	G	2	NAG	O5-C5-C6-O6
2	G	1	NAG	C1-C2-N2-C7
2	G	2	NAG	C1-C2-N2-C7

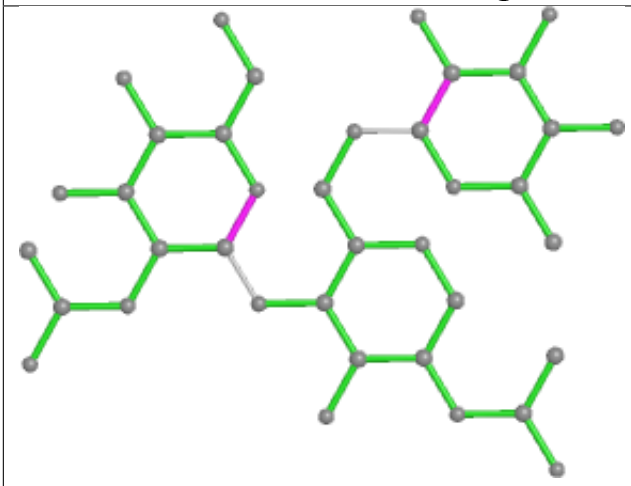
There are no ring outliers.

2 monomers are involved in 2 short contacts:

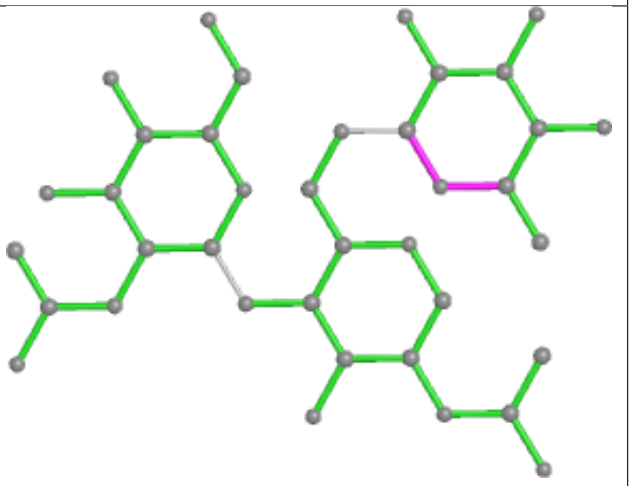
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	1	NAG	1	0
2	G	2	NAG	1	0

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

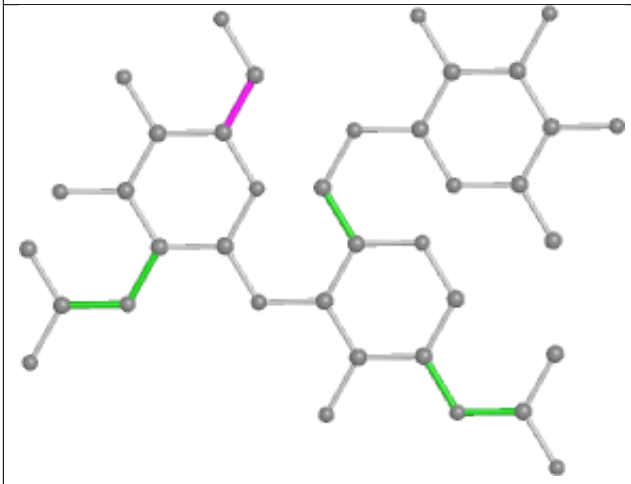
Oligosaccharide Chain C



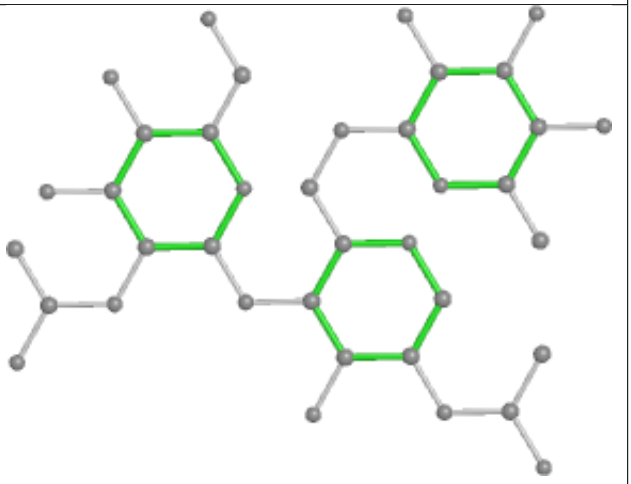
Bond lengths



Bond angles

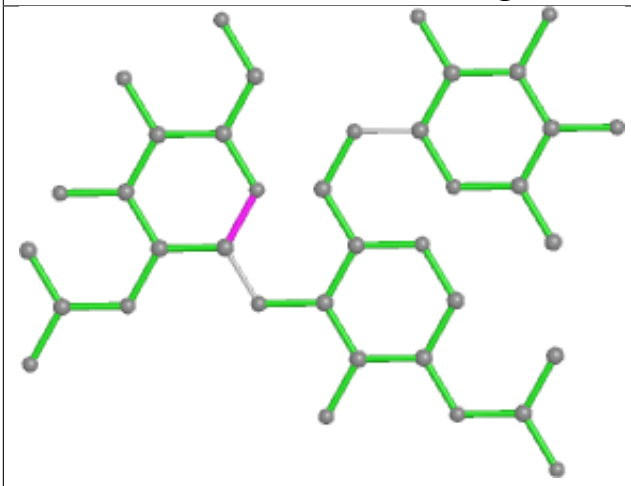


Torsions

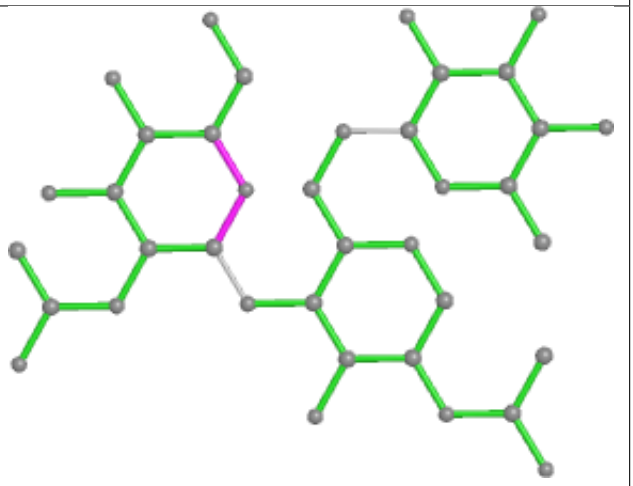


Rings

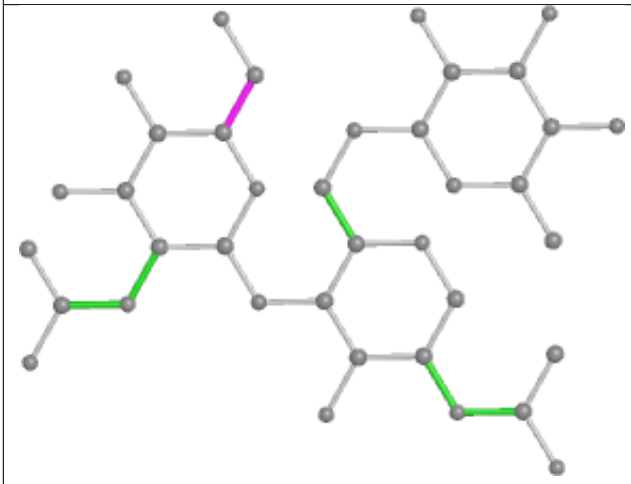
Oligosaccharide Chain E



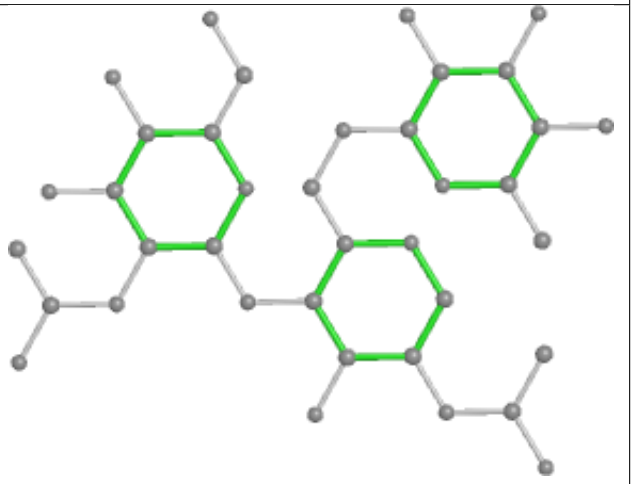
Bond lengths



Bond angles

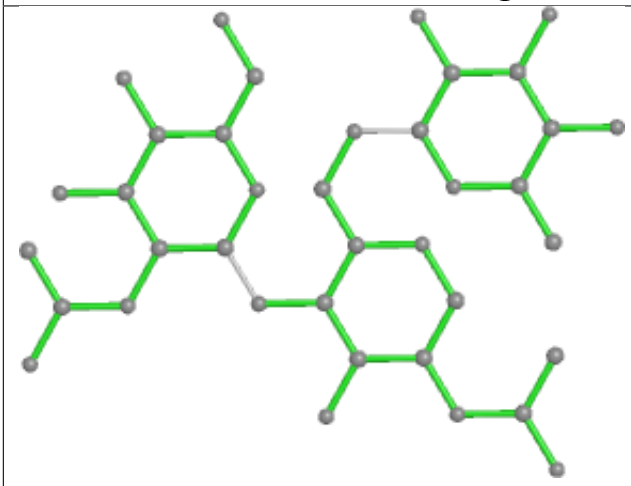


Torsions

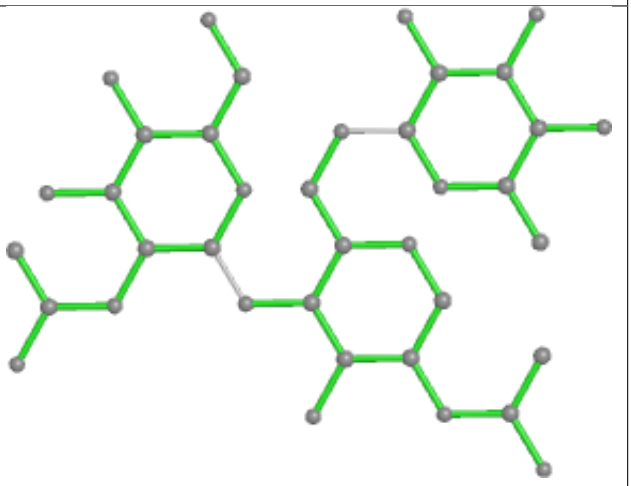


Rings

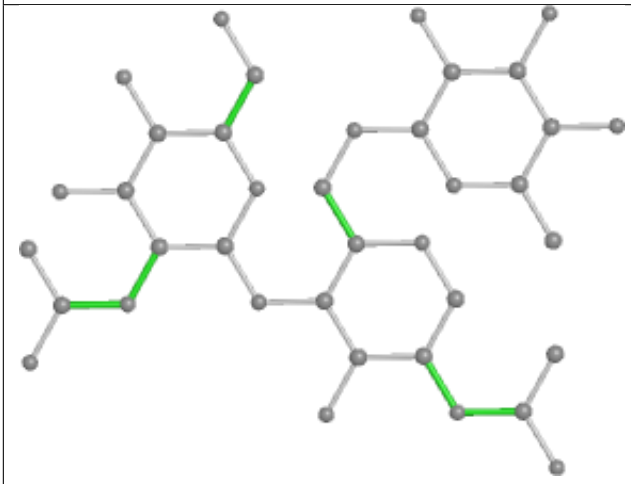
Oligosaccharide Chain F



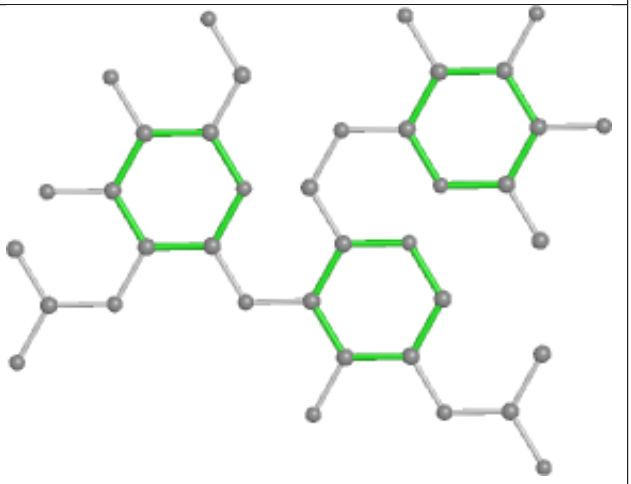
Bond lengths



Bond angles

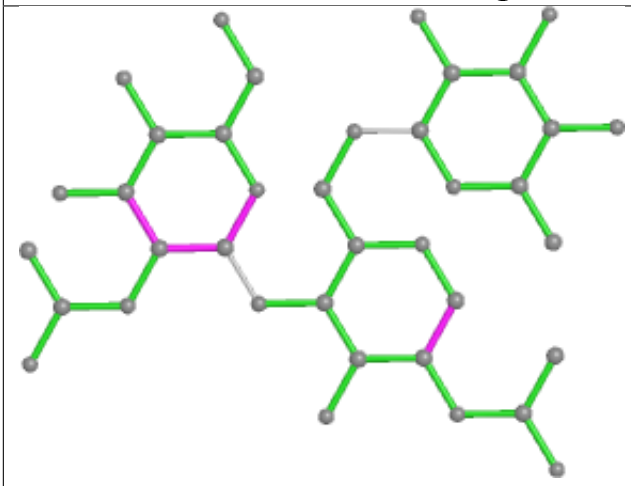


Torsions

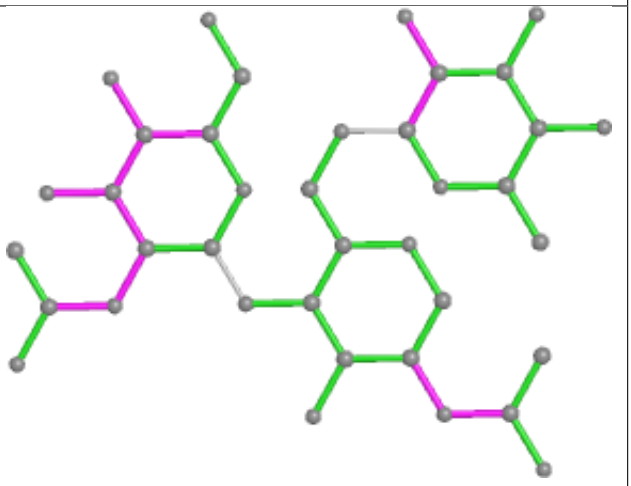


Rings

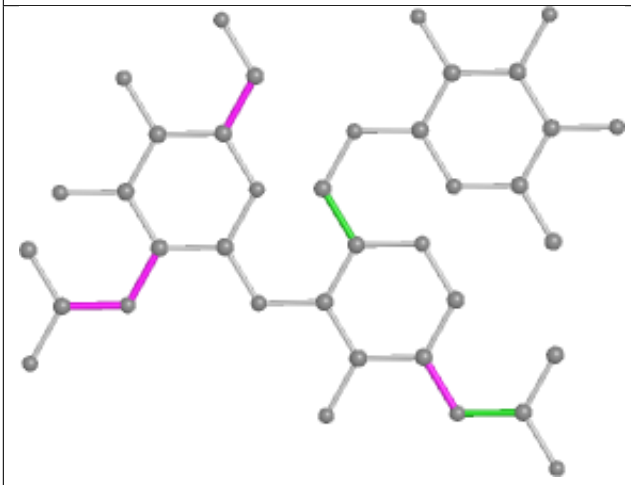
Oligosaccharide Chain G



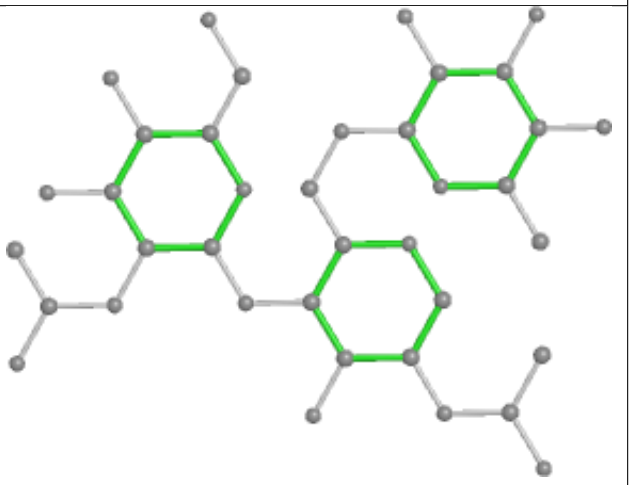
Bond lengths



Bond angles

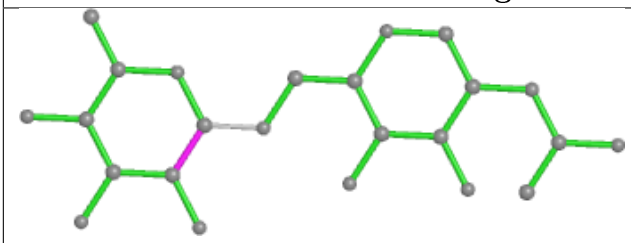


Torsions

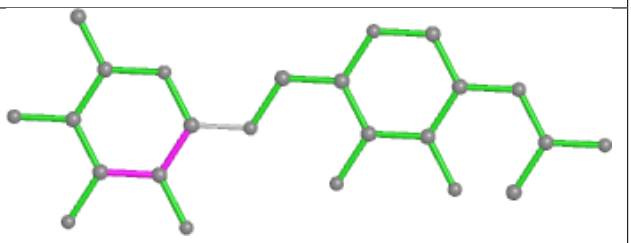


Rings

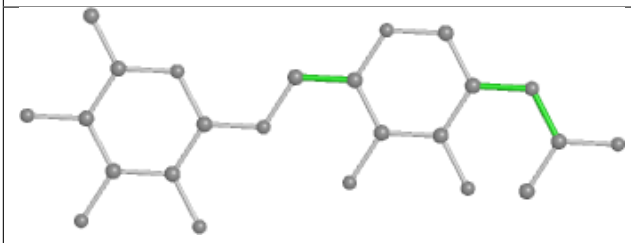
Oligosaccharide Chain D



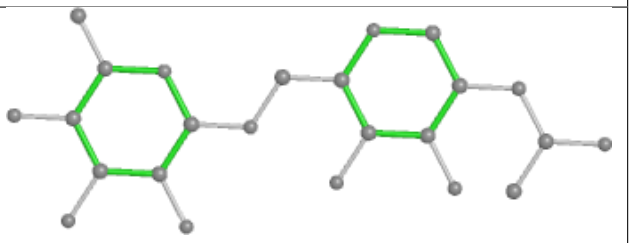
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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	$\#\|Z\| > 2$
7	DQV	B	510	-	43,48,48	4.62	17 (39%)	54,73,73	2.27	11 (20%)
4	NAG	A	306	1	14,14,15	0.36	0	17,19,21	0.44	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
7	DQV	B	510	-	-	3/26/62/62	0/5/5/5
4	NAG	A	306	1	-	2/6/23/26	0/1/1/1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	510	DQV	O9-C12	18.87	1.67	1.41
7	B	510	DQV	C13-C12	-11.69	1.36	1.53
7	B	510	DQV	C21-C20	-8.37	1.30	1.53
7	B	510	DQV	O1-C7	7.12	1.55	1.44
7	B	510	DQV	C-N	7.01	1.46	1.33
7	B	510	DQV	O1-C8	-6.69	1.30	1.45
7	B	510	DQV	O9-C11	-6.16	1.31	1.45
7	B	510	DQV	O11-C14	-5.31	1.30	1.43
7	B	510	DQV	C20-C8	4.95	1.65	1.53
7	B	510	DQV	C5-C7	-4.62	1.43	1.51
7	B	510	DQV	C17-N3	3.71	1.47	1.34
7	B	510	DQV	C21-C7	3.49	1.58	1.54
7	B	510	DQV	C14-C11	3.45	1.61	1.53
7	B	510	DQV	O-C	-3.43	1.17	1.24
7	B	510	DQV	O13-C21	3.15	1.50	1.43
7	B	510	DQV	C18-N5	2.78	1.36	1.32

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	510	DQV	C16-C19	-2.39	1.34	1.40

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	B	510	DQV	C12-N1-C19	8.61	141.77	126.64
7	B	510	DQV	C16-C17-N3	6.07	129.58	120.35
7	B	510	DQV	N5-C18-N4	-5.60	119.92	128.68
7	B	510	DQV	C20-C21-C7	4.56	107.19	101.93
7	B	510	DQV	C8-O1-C7	-4.16	104.29	109.42
7	B	510	DQV	N3-C17-N4	-4.01	110.26	118.57
7	B	510	DQV	C4-C5-C7	3.60	127.34	120.64
7	B	510	DQV	O1-C7-C21	-3.23	99.42	104.66
7	B	510	DQV	C6-C5-C7	-2.56	114.43	119.83
7	B	510	DQV	O10-C13-C14	-2.28	104.45	111.82
7	B	510	DQV	C14-C13-C12	2.05	104.06	100.98

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	306	NAG	O5-C5-C6-O6
4	A	306	NAG	C4-C5-C6-O6
7	B	510	DQV	C4-C5-C7-C21
7	B	510	DQV	C6-C5-C7-C21
7	B	510	DQV	P-O5-P1-O8

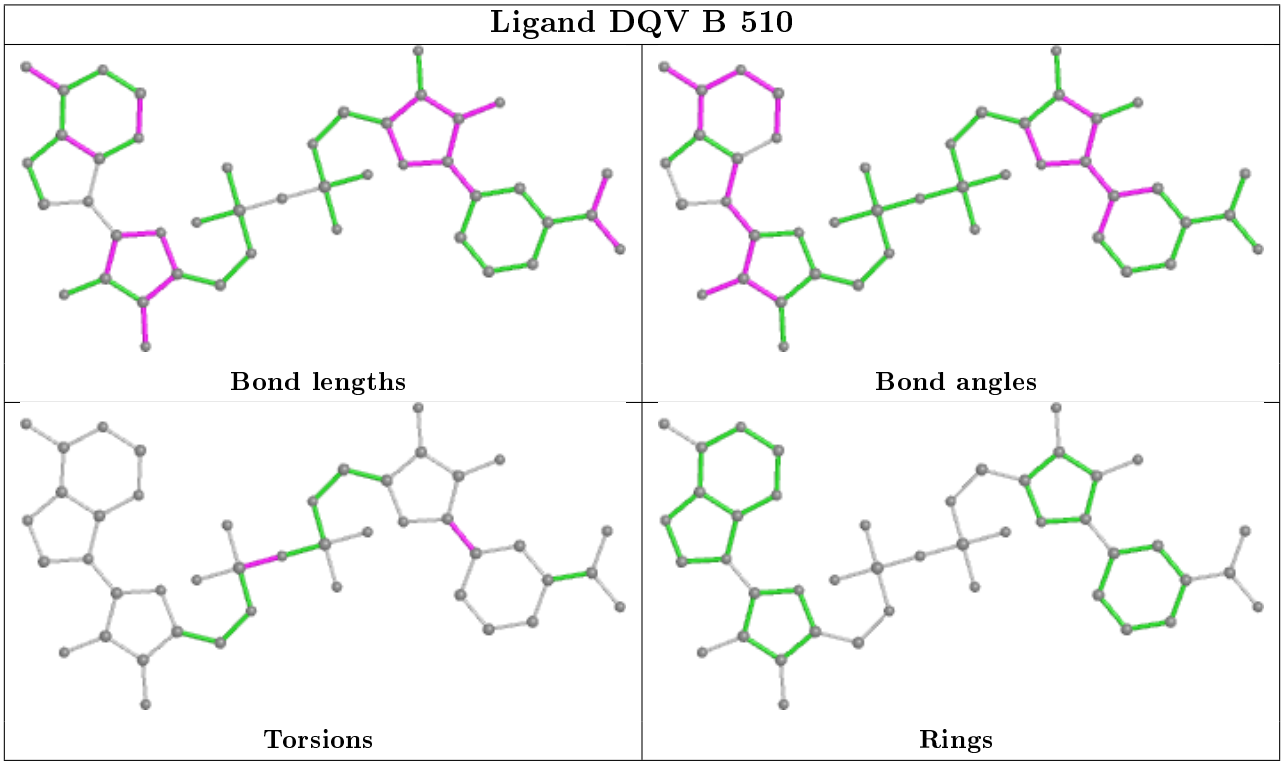
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	510	DQV	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	209/248 (84%)	-0.24	4 (1%) 66 66	19, 28, 44, 68	0
1	B	209/248 (84%)	-0.02	7 (3%) 46 44	19, 27, 52, 74	0
All	All	418/496 (84%)	-0.13	11 (2%) 56 54	19, 28, 51, 74	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	34	TRP	5.4
1	B	235	ASP	3.5
1	B	35	PRO	3.2
1	B	134	TYR	3.1
1	B	27	SER	2.8
1	B	234	GLN	2.7
1	B	31	ALA	2.7
1	A	26	ASP	2.7
1	A	234	GLN	2.7
1	A	233	ASN	2.4
1	A	27	SER	2.3

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

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

6.3 Carbohydrates

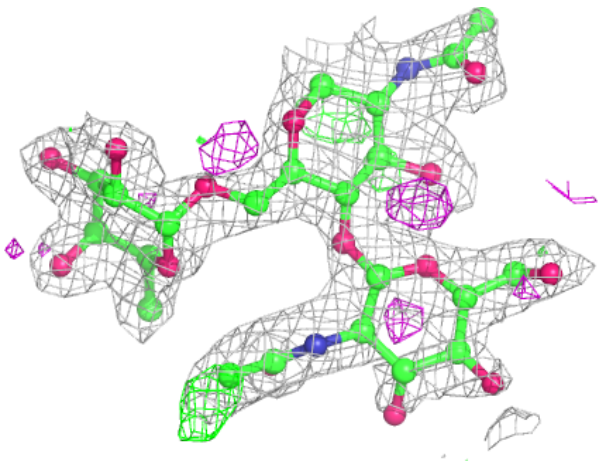
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

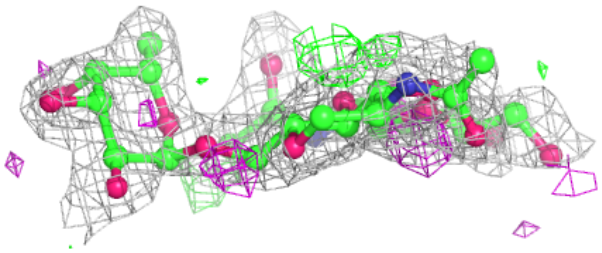
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	NAG	G	2	14/15	0.54	0.37	72,78,86,88	0
2	NAG	E	2	14/15	0.56	0.40	74,83,87,92	0
2	NAG	C	2	14/15	0.69	0.24	52,62,74,75	0
2	FUC	G	3	10/11	0.72	0.20	56,60,63,63	0
3	FUC	D	2	10/11	0.80	0.32	53,65,69,71	0
2	NAG	G	1	14/15	0.80	0.18	51,62,70,75	0
2	FUC	F	3	10/11	0.83	0.15	31,39,40,43	0
2	FUC	C	3	10/11	0.83	0.27	32,48,54,55	0
2	NAG	F	2	14/15	0.84	0.16	55,61,72,80	0
2	NAG	C	1	14/15	0.84	0.19	44,55,61,62	0
2	FUC	E	3	10/11	0.85	0.15	42,50,55,56	0
2	NAG	F	1	14/15	0.87	0.12	35,44,49,51	0
3	NAG	D	1	14/15	0.90	0.13	31,43,57,63	0
2	NAG	E	1	14/15	0.93	0.10	40,44,52,63	0

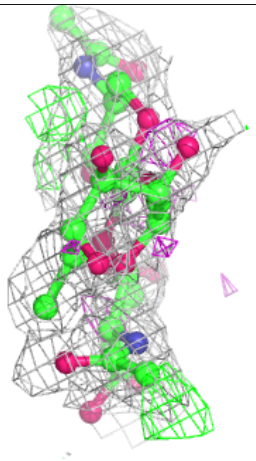
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

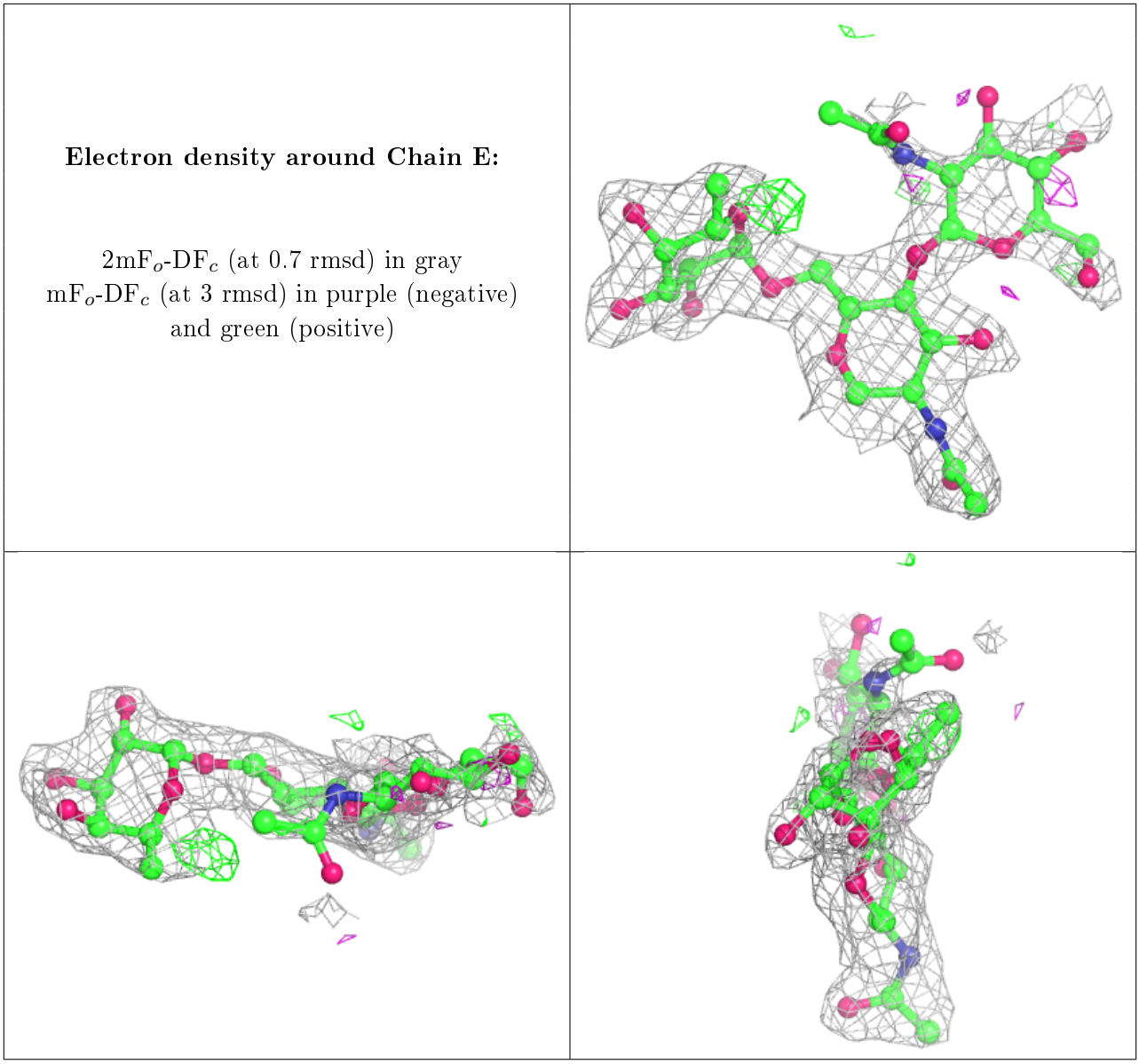






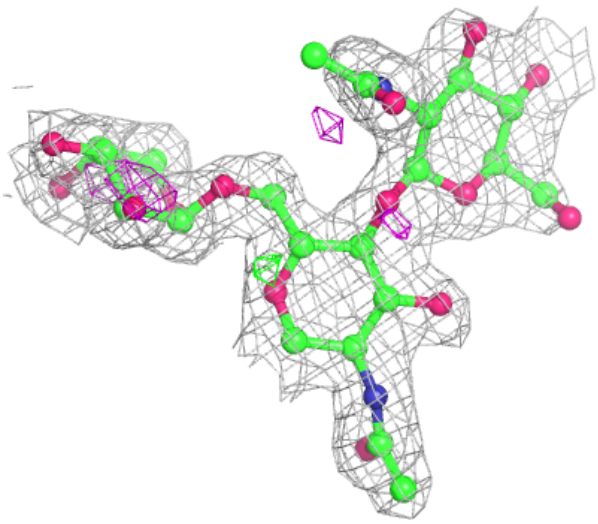
Electron density around Chain E:

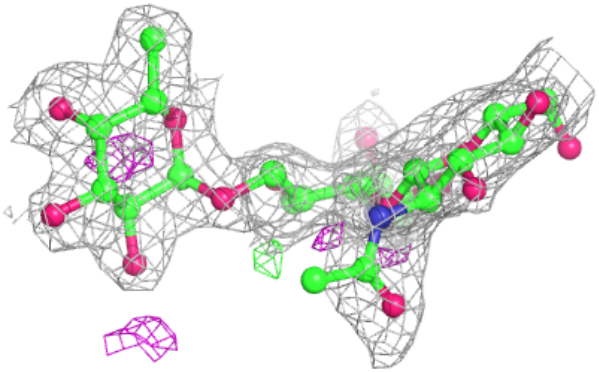
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

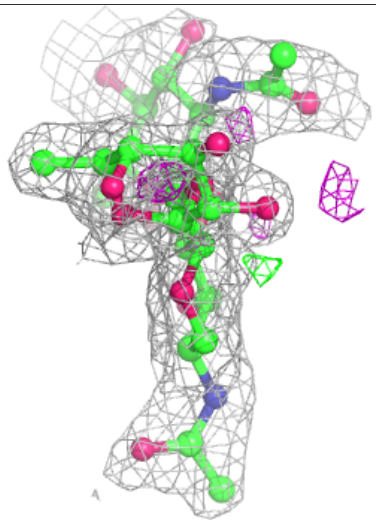


Electron density around Chain F:

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

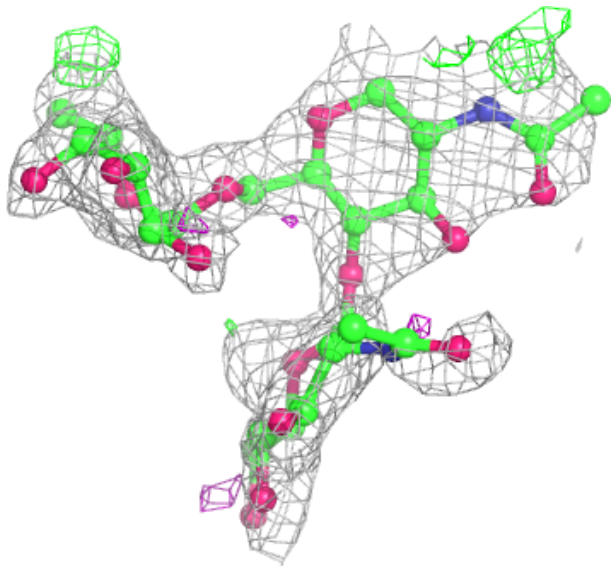


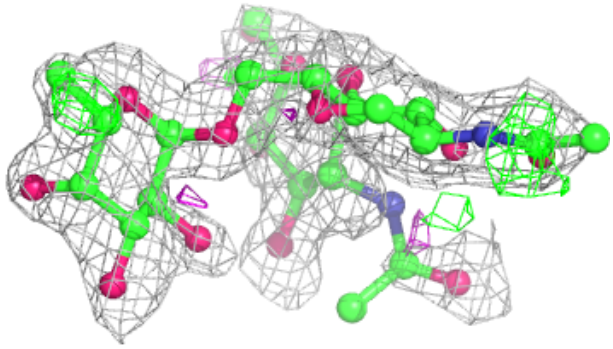


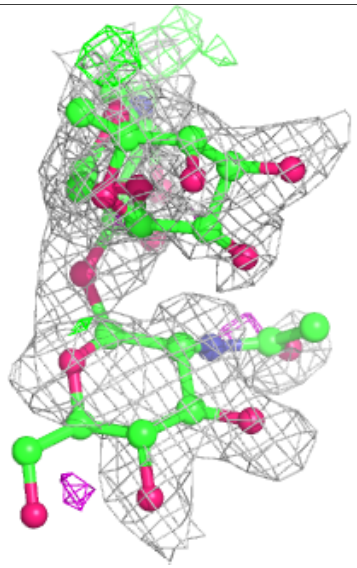


Electron density around Chain G:

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

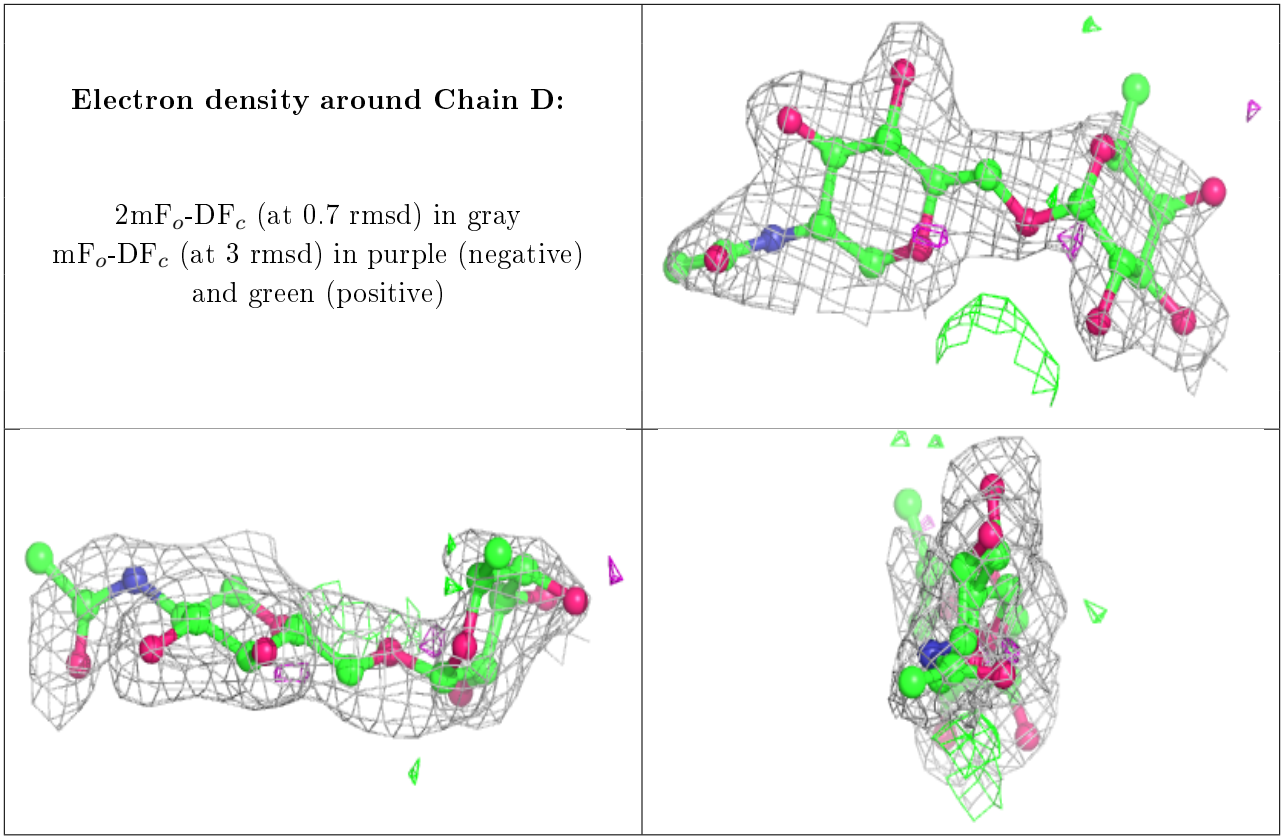






Electron density around Chain D:

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



6.4 Ligands

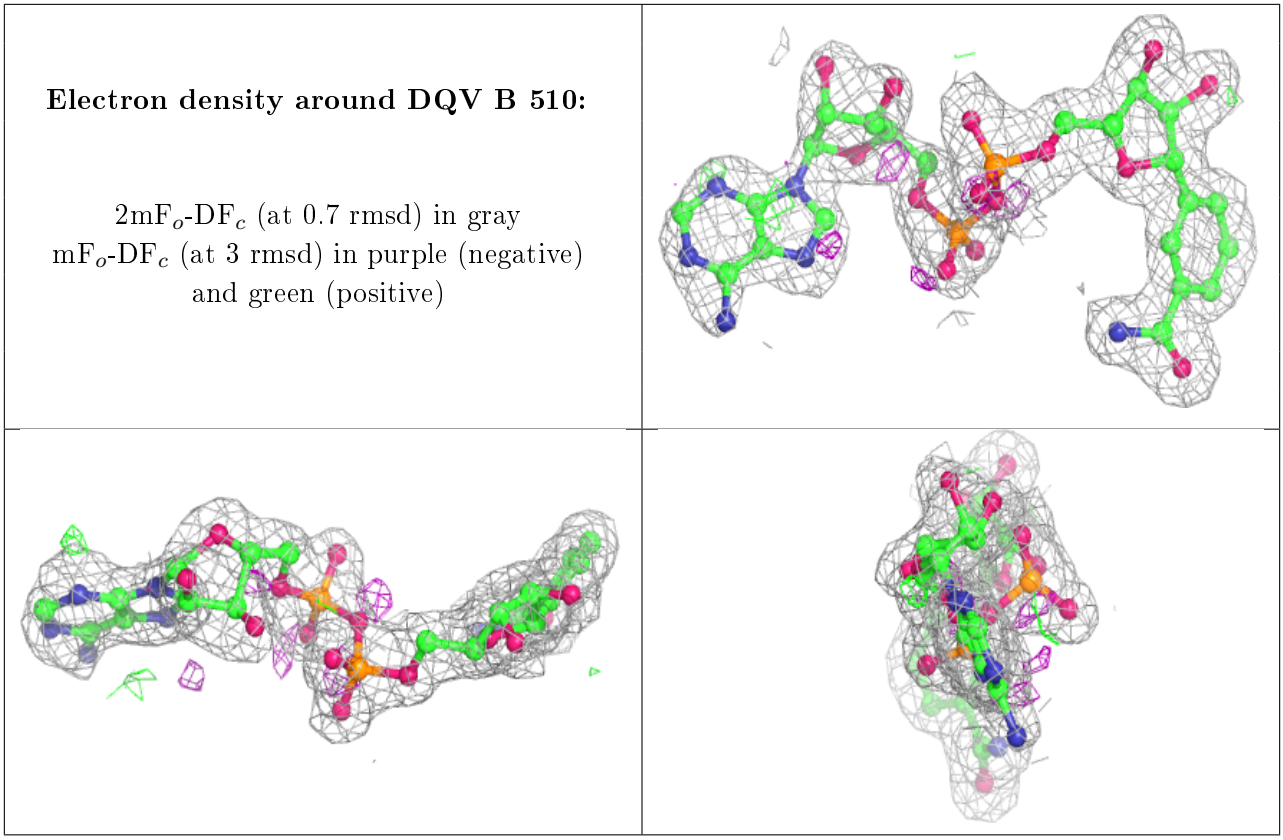
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
4	NAG	A	306	14/15	0.76	0.20	50,63,69,72	0
7	DQV	B	510	44/44	0.94	0.12	19,33,52,54	0
6	NA	A	308	1/1	0.97	0.09	38,38,38,38	0
5	CA	A	307	1/1	0.99	0.08	37,37,37,37	0
5	CA	B	511	1/1	0.99	0.06	27,27,27,27	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.

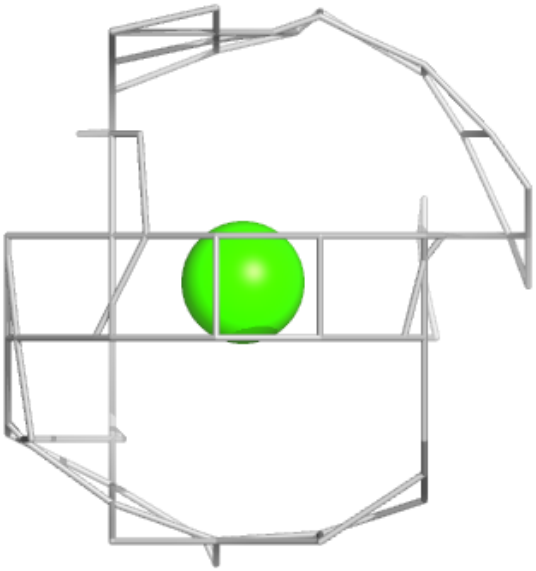
Electron density around DQV B 510:

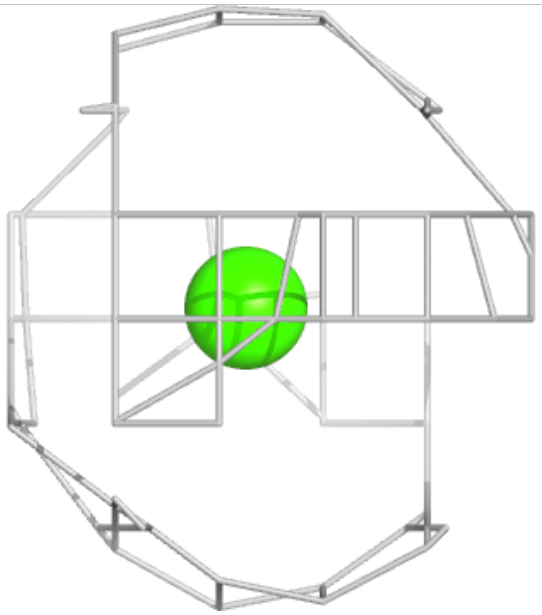
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

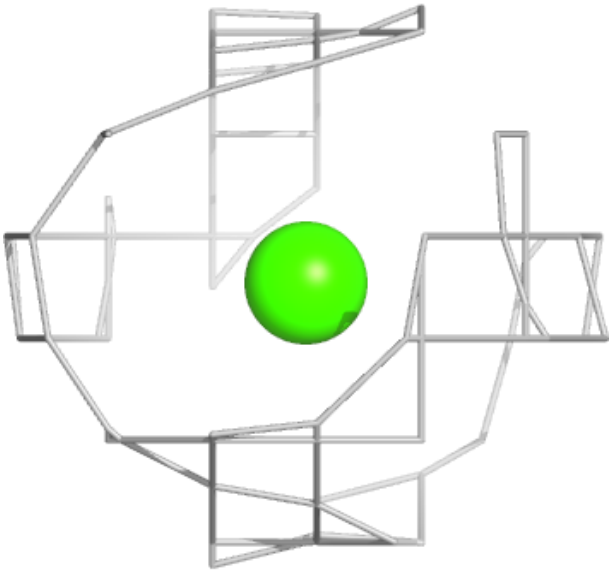


Electron density around CA A 307:

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

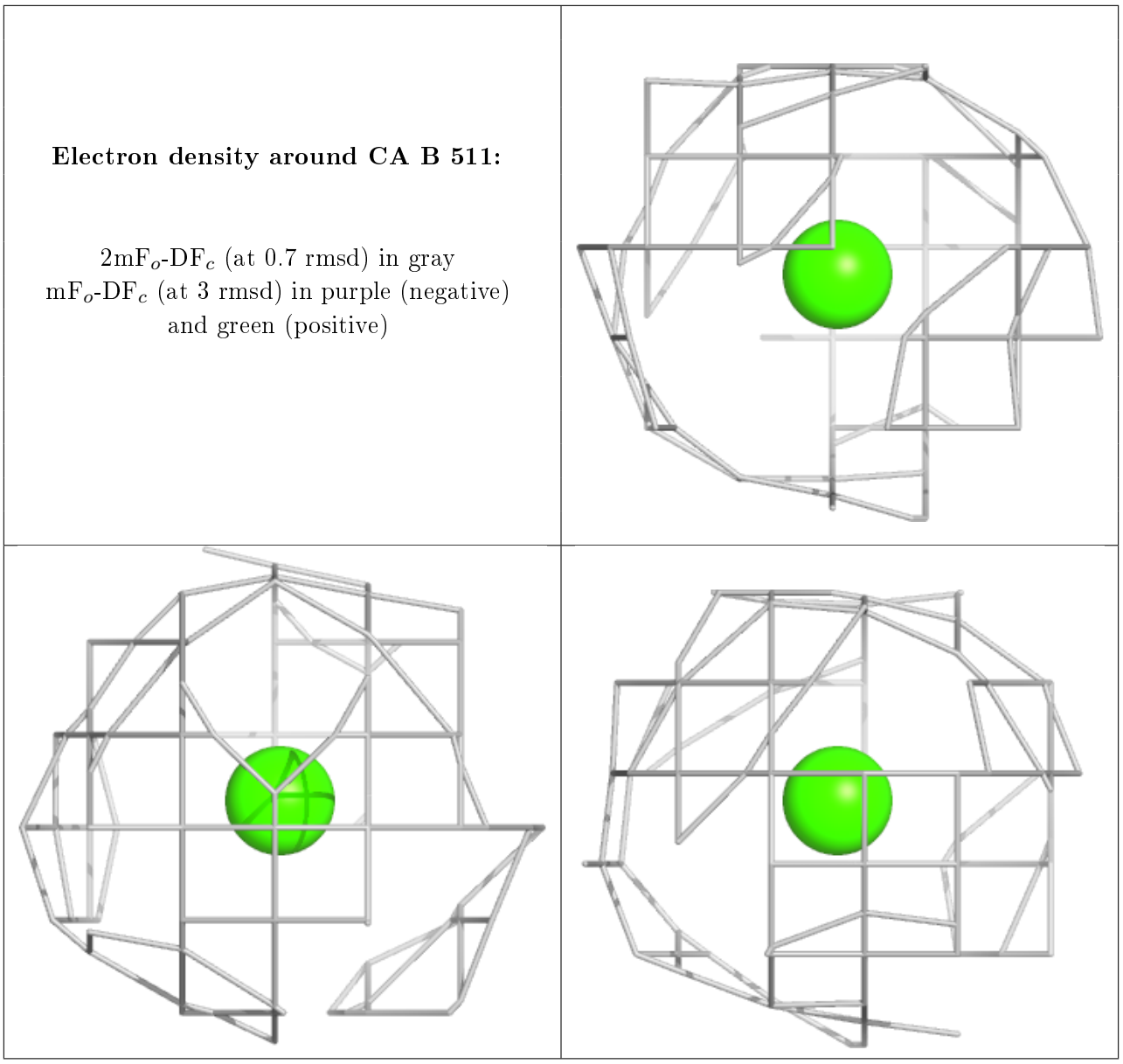






Electron density around CA B 511:

$2mF_o$-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

There are no such residues in this entry.