



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 08:12 AM UTC

PDB ID : 6X5U / pdb_00006x5u
Title : Human Alpha-1,6-fucosyltransferase (FUT8) bound to GDP and NM5M2-Asn
Authors : Kadirvelraj, R.; Wood, Z.A.
Deposited on : 2020-05-26
Resolution : 3.20 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

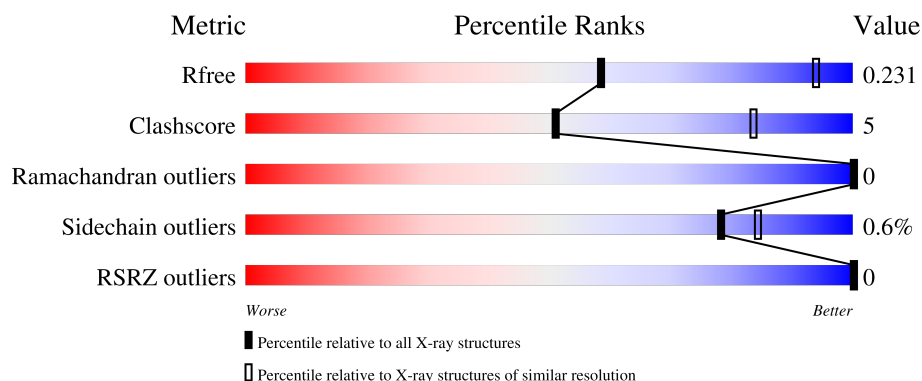
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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.



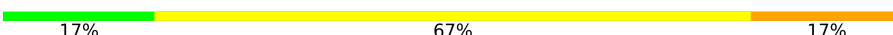
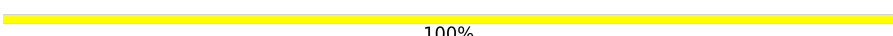
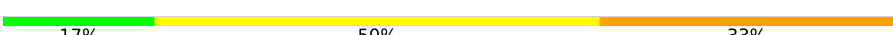
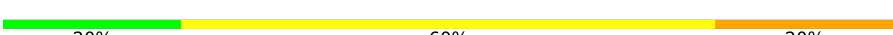
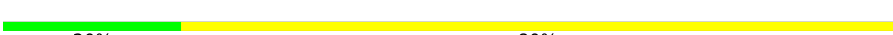
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	535	 79% 8% 13%
1	B	535	 78% 9% 13%
1	C	535	 75% 12% 13%
1	D	535	 78% 9% 13%
1	E	535	 76% 11% 13%

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Mol	Chain	Length	Quality of chain
1	F	535	
1	G	535	
1	H	535	
2	I	6	
2	K	6	
2	P	6	
3	J	5	
3	L	5	
3	N	5	
4	M	4	
5	O	5	

2 Entry composition [i](#)

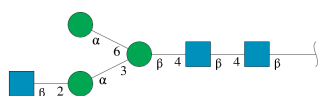
There are 8 unique types of molecules in this entry. The entry contains 31445 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Alpha-(1,6)-fucosyltransferase.

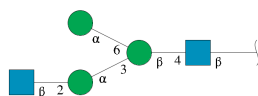
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	1	0
			3809	2424	672	699	14			
1	B	467	Total	C	N	O	S	0	0	0
			3801	2419	669	699	14			
1	C	467	Total	C	N	O	S	0	0	0
			3801	2419	669	699	14			
1	D	467	Total	C	N	O	S	0	0	0
			3801	2419	669	699	14			
1	E	467	Total	C	N	O	S	0	0	0
			3801	2419	669	699	14			
1	F	467	Total	C	N	O	S	0	0	0
			3801	2419	669	699	14			
1	G	467	Total	C	N	O	S	0	0	0
			3801	2419	669	699	14			
1	H	467	Total	C	N	O	S	0	0	0
			3801	2419	669	699	14			

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



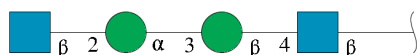
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	6	Total	C	N	O	0	0	0
			75	42	3	30			
2	K	6	Total	C	N	O	0	0	0
			75	42	3	30			
2	P	6	Total	C	N	O	0	0	0
			76	42	3	31			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	J	5	Total	C	N	O	0	0	0
			62	34	2	26			
3	L	5	Total	C	N	O	0	0	0
			62	34	2	26			
3	N	5	Total	C	N	O	0	0	0
			62	34	2	26			

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



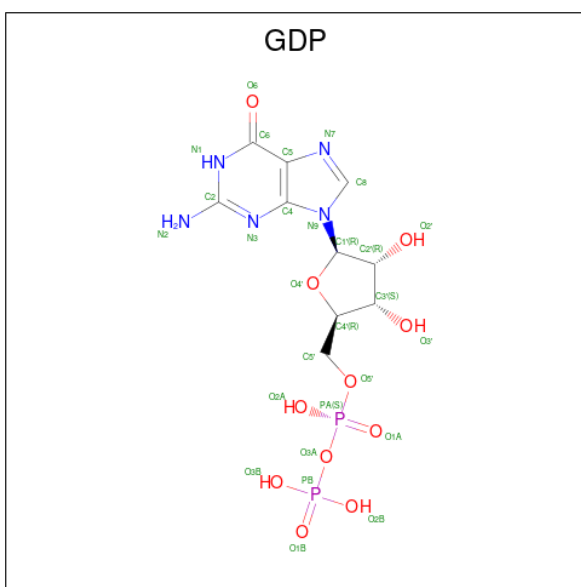
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	M	4	Total	C	N	O	0	0	0
			51	28	2	21			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



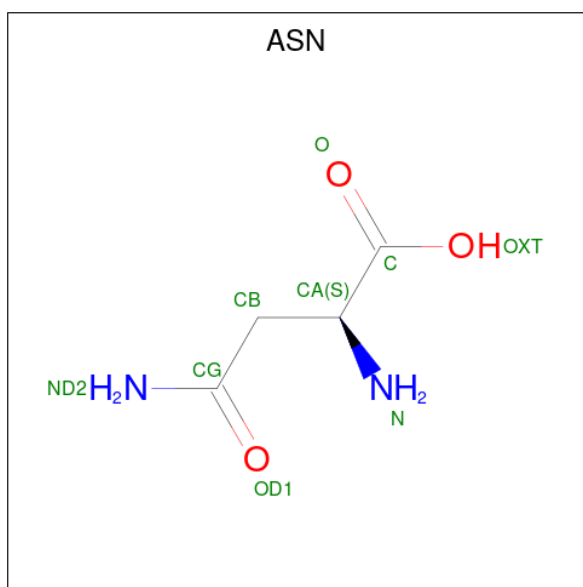
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	O	5	Total	C	N	O	0	0	0
			65	36	3	26			

- Molecule 6 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total 28	C 10	N 5	O 11	P 2	0	0
6	B	1	Total 28	C 10	N 5	O 11	P 2	0	0
6	C	1	Total 28	C 10	N 5	O 11	P 2	0	0
6	D	1	Total 28	C 10	N 5	O 11	P 2	0	0
6	E	1	Total 28	C 10	N 5	O 11	P 2	0	0
6	F	1	Total 28	C 10	N 5	O 11	P 2	0	0
6	G	1	Total 28	C 10	N 5	O 11	P 2	0	0
6	H	1	Total 28	C 10	N 5	O 11	P 2	0	0

- Molecule 7 is ASPARAGINE (CCD ID: ASN) (formula: $C_4H_8N_2O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			8	4	2	2		
7	C	1	Total	C	N	O	0	0
			8	4	2	2		

- Molecule 8 is water.

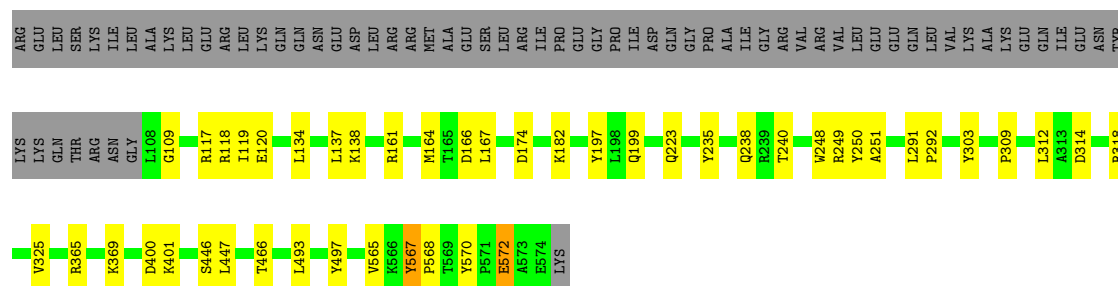
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	34	Total	O	0	0
			34	34		
8	B	34	Total	O	0	0
			34	34		
8	C	33	Total	O	0	0
			33	33		
8	D	31	Total	O	0	0
			31	31		
8	E	28	Total	O	0	0
			28	28		
8	F	40	Total	O	0	0
			40	40		
8	G	25	Total	O	0	0
			25	25		
8	H	36	Total	O	0	0
			36	36		

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.

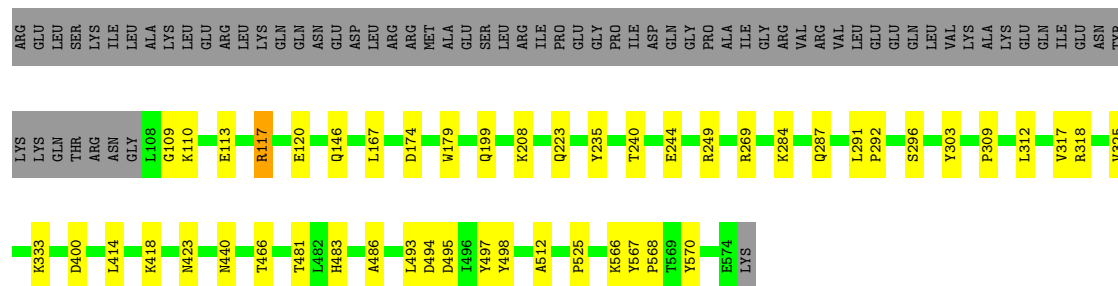
- Molecule 1: Alpha-(1,6)-fucosyltransferase

Chain A: 



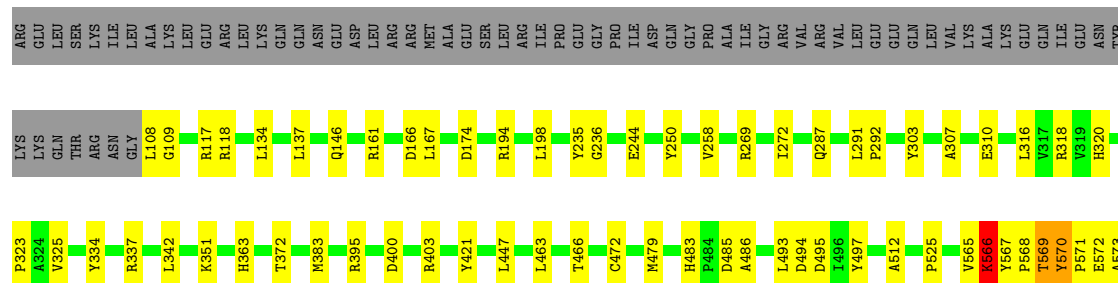
- Molecule 1: Alpha-(1,6)-fucosyltransferase

Chain B: 



- Molecule 1: Alpha-(1,6)-fucosyltransferase

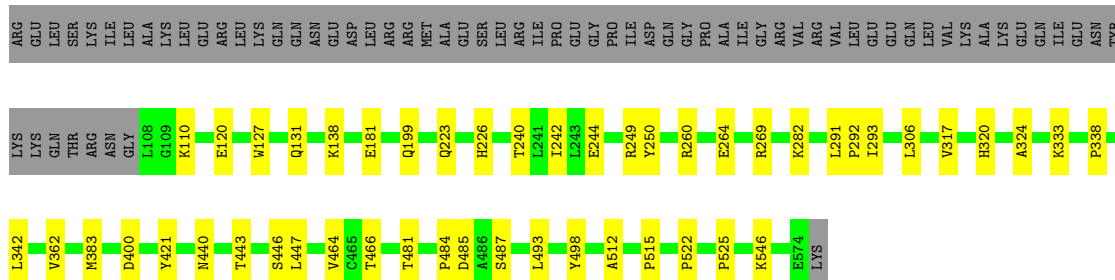
Chain C: 



E574
LYS

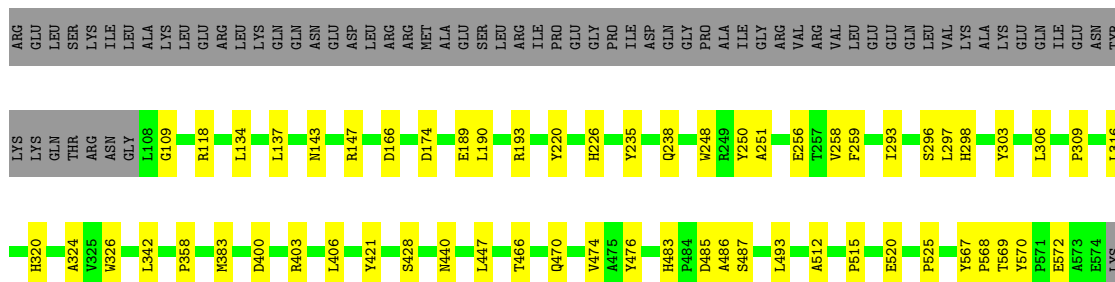
- Molecule 1: Alpha-(1,6)-fucosyltransferase

Chain D: 78% 9% 13%



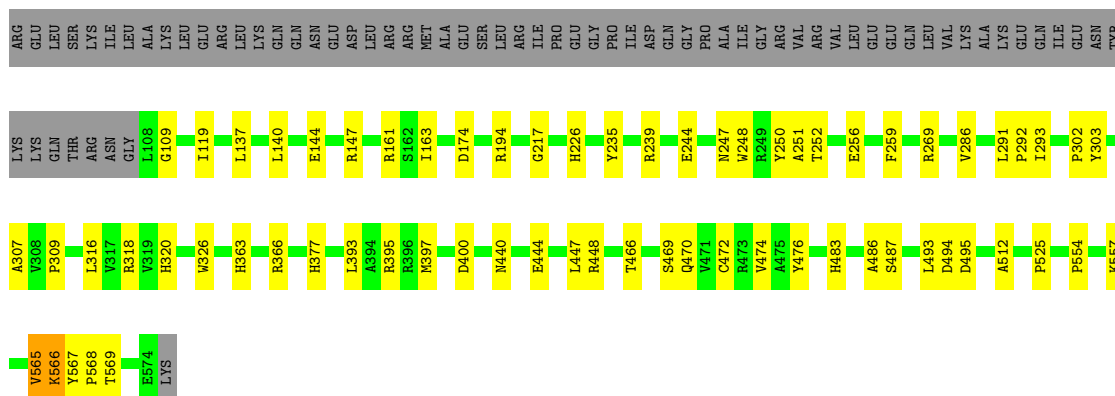
- Molecule 1: Alpha-(1,6)-fucosyltransferase

Chain E: 76% 11% 13%



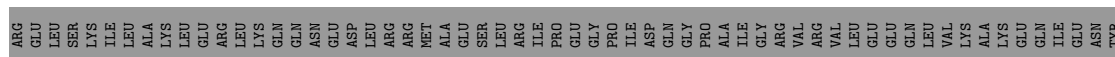
- Molecule 1: Alpha-(1,6)-fucosyltransferase

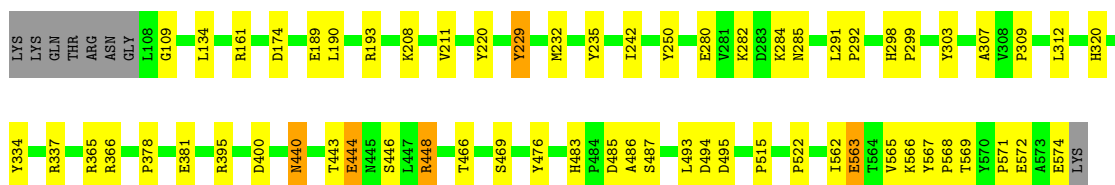
Chain F: 75% 12% 13%



- Molecule 1: Alpha-(1,6)-fucosyltransferase

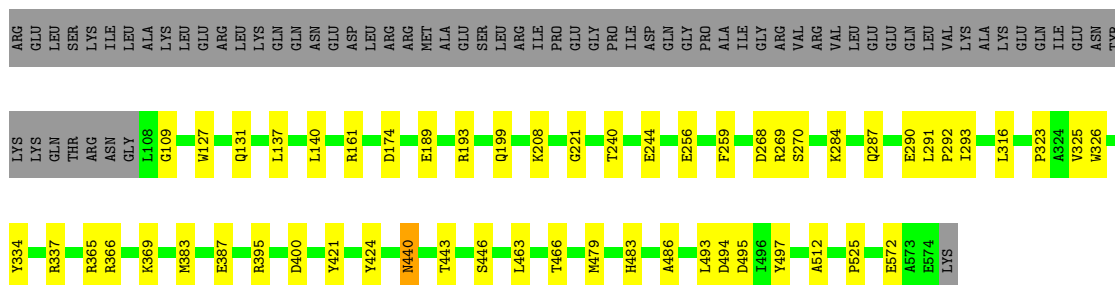
Chain G: 76% 11% 13%





- Molecule 1: Alpha-(1,6)-fucosyltransferase

Chain H:



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

Chain I:



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

Chain K:

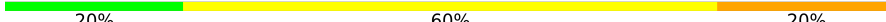


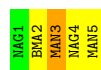
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:

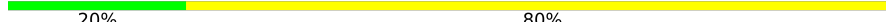


- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  20% 60% 20%



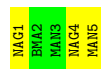
- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  20% 80%

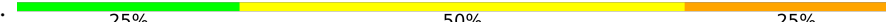


- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  40% 60%



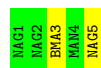
- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  25% 50% 25%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  60% 40%



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	154.02Å 154.02Å 466.92Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	87.84 – 3.20 87.84 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.4 (87.84-3.20) 99.4 (87.84-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.190 , 0.236 0.191 , 0.231	Depositor DCC
R_{free} test set	97028 reflections (94.47%)	wwPDB-VP
Wilson B-factor (Å ²)	64.1	Xtriage
Anisotropy	0.519	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.116 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	31445	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.18 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.4512e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, MAN, NAG, GDP

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	0.20	1/3912 (0.0%)	0.36	1/5304 (0.0%)
1	B	0.16	0/3901	0.33	0/5289
1	C	0.23	0/3901	0.38	3/5289 (0.1%)
1	D	0.12	0/3901	0.32	0/5289
1	E	0.13	0/3901	0.32	0/5289
1	F	0.11	0/3901	0.33	0/5289
1	G	0.11	0/3901	0.34	0/5289
1	H	0.11	0/3901	0.31	0/5289
All	All	0.15	1/31219 (0.0%)	0.34	4/42327 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	567	TYR	CA-C	-5.30	1.46	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	570	TYR	CA-C-N	7.06	126.88	119.05
1	C	570	TYR	C-N-CA	7.06	126.88	119.05
1	A	567	TYR	N-CA-C	6.28	119.24	110.20
1	C	566	LYS	N-CA-C	5.60	117.08	110.97

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	3809	0	3724	34	0
1	B	3801	0	3711	40	0
1	C	3801	0	3711	49	0
1	D	3801	0	3711	29	0
1	E	3801	0	3711	50	0
1	F	3801	0	3711	42	0
1	G	3801	0	3711	56	0
1	H	3801	0	3711	32	0
2	I	75	0	64	1	0
2	K	75	0	64	1	0
2	P	76	0	66	1	0
3	J	62	0	54	2	0
3	L	62	0	54	2	0
3	N	62	0	54	0	0
4	M	51	0	45	2	0
5	O	65	0	57	0	0
6	A	28	0	12	1	0
6	B	28	0	12	0	0
6	C	28	0	12	0	0
6	D	28	0	12	0	0
6	E	28	0	12	1	0
6	F	28	0	12	1	0
6	G	28	0	12	1	0
6	H	28	0	12	1	0
7	A	8	0	4	0	0
7	C	8	0	4	0	0
8	A	34	0	0	2	0
8	B	34	0	0	1	0
8	C	33	0	0	0	0
8	D	31	0	0	1	0
8	E	28	0	0	1	0
8	F	40	0	0	2	0
8	G	25	0	0	3	0
8	H	36	0	0	0	0
All	All	31445	0	30263	322	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 5.

All (322) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:307:ALA:CB	1:G:565:VAL:HG13	1.71	1.21
1:G:307:ALA:HB1	1:G:565:VAL:HG13	1.33	1.05
1:G:307:ALA:CB	1:G:565:VAL:CG1	2.39	0.99
1:E:303:TYR:CD1	1:E:567:TYR:CE1	2.54	0.96
1:E:303:TYR:HD1	1:E:567:TYR:CE1	1.85	0.95
1:F:307:ALA:HB1	1:F:565:VAL:HG23	1.54	0.89
1:H:189:GLU:O	1:H:193:ARG:HG3	1.73	0.89
1:B:303:TYR:CD1	1:B:567:TYR:CE1	2.62	0.87
1:G:569:THR:HG23	1:G:571:PRO:HD3	1.57	0.85
1:G:189:GLU:O	1:G:193:ARG:HG3	1.78	0.84
1:F:307:ALA:HB1	1:F:565:VAL:CG2	2.10	0.81
1:B:303:TYR:CA	1:B:567:TYR:OH	2.29	0.80
1:E:189:GLU:O	1:E:193:ARG:HG3	1.82	0.80
1:B:303:TYR:HA	1:B:567:TYR:OH	1.82	0.79
1:E:303:TYR:HA	1:E:567:TYR:OH	1.83	0.79
1:B:303:TYR:HD1	1:B:567:TYR:CE1	1.99	0.78
1:A:309:PRO:HA	1:A:567:TYR:HD2	1.49	0.77
1:E:235:TYR:O	1:E:568:PRO:HG2	1.85	0.76
1:G:307:ALA:HB3	1:G:565:VAL:CG1	2.17	0.74
1:E:303:TYR:HB3	1:E:567:TYR:HE1	1.53	0.74
1:C:351:LYS:HE3	1:G:285:ASN:HB3	1.69	0.73
1:C:569:THR:HG23	1:C:571:PRO:HD3	1.70	0.72
1:F:307:ALA:CB	1:F:565:VAL:CG2	2.67	0.72
1:A:309:PRO:HA	1:A:567:TYR:CD2	2.26	0.70
1:F:247:ASN:HA	1:F:252:THR:HG23	1.74	0.70
1:B:303:TYR:HB3	1:B:567:TYR:OH	1.93	0.69
1:E:303:TYR:HB3	1:E:567:TYR:CE1	2.28	0.69
1:G:190:LEU:CD1	1:G:574:GLU:OE2	2.41	0.69
1:G:307:ALA:HB3	1:G:565:VAL:HG11	1.75	0.69
1:A:568:PRO:HB2	1:A:570:TYR:CE2	2.29	0.67
1:B:303:TYR:HB3	1:B:567:TYR:CE1	2.30	0.66
1:G:303:TYR:HB3	1:G:567:TYR:HE1	1.60	0.66
1:A:117:ARG:NH2	1:B:317:VAL:O	2.29	0.66
1:C:117:ARG:NH2	1:D:317:VAL:O	2.29	0.66
1:G:235:TYR:O	1:G:568:PRO:HG2	1.96	0.65
1:E:296:SER:HB2	3:J:3:MAN:H5	1.79	0.65
1:C:303:TYR:HB3	1:C:567:TYR:CE1	2.32	0.64
1:F:194:ARG:CZ	1:F:569:THR:O	2.46	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:563:GLU:OE2	1:G:563:GLU:HA	1.96	0.64
1:E:189:GLU:HB3	1:E:193:ARG:NH2	2.14	0.63
1:F:483:HIS:HB2	1:F:486:ALA:HB2	1.81	0.62
1:C:307:ALA:HB1	1:C:565:VAL:HG23	1.82	0.61
1:E:303:TYR:CA	1:E:567:TYR:OH	2.48	0.61
1:C:198:LEU:HD22	1:C:570:TYR:OH	2.01	0.60
1:B:303:TYR:CB	1:B:567:TYR:OH	2.49	0.60
1:H:269:ARG:NH2	1:H:290:GLU:OE2	2.34	0.59
1:H:483:HIS:HB2	1:H:486:ALA:HB2	1.84	0.59
1:C:244:GLU:OE2	1:C:269:ARG:NH1	2.35	0.59
1:B:208:LYS:NZ	1:B:284:LYS:O	2.35	0.59
1:E:466:THR:HG22	1:E:493:LEU:HB2	1.84	0.59
1:C:303:TYR:CD1	1:C:567:TYR:CE1	2.91	0.59
1:G:303:TYR:HB3	1:G:567:TYR:CE1	2.38	0.59
1:E:303:TYR:CD1	1:E:567:TYR:HE1	2.19	0.58
1:C:570:TYR:N	1:C:571:PRO:HD3	2.18	0.58
1:C:466:THR:HG22	1:C:493:LEU:HB2	1.85	0.57
1:E:190:LEU:HA	1:E:193:ARG:NH1	2.19	0.57
1:A:118:ARG:NH1	1:A:166:ASP:OD1	2.38	0.57
1:B:423:ASN:ND2	8:B:802:HOH:O	2.37	0.57
1:E:190:LEU:HA	1:E:193:ARG:CZ	2.35	0.57
1:A:134:LEU:HD23	1:A:137:LEU:HD12	1.87	0.56
1:F:318:ARG:NH1	8:F:804:HOH:O	2.38	0.56
1:F:303:TYR:HE1	1:F:566:LYS:O	1.89	0.56
1:G:466:THR:HG22	1:G:493:LEU:HB2	1.88	0.56
1:F:320:HIS:CE1	1:F:487:SER:HB3	2.41	0.55
1:F:235:TYR:O	1:F:568:PRO:HG2	2.06	0.55
1:C:235:TYR:CE1	1:C:570:TYR:CE2	2.94	0.55
1:D:466:THR:HG22	1:D:493:LEU:HB2	1.87	0.55
1:E:143:ASN:O	1:E:147:ARG:HG3	2.07	0.55
1:B:483:HIS:HB2	1:B:486:ALA:HB2	1.89	0.54
1:C:146:GLN:HG2	1:D:138:LYS:HD2	1.89	0.54
1:A:314:ASP:O	1:A:318:ARG:HG3	2.08	0.54
1:G:303:TYR:HD1	1:G:567:TYR:CE1	2.25	0.54
1:G:303:TYR:CD1	1:G:567:TYR:CE1	2.96	0.54
1:E:118:ARG:NH1	1:E:166:ASP:OD1	2.40	0.54
1:G:307:ALA:HB2	1:G:565:VAL:HG13	1.78	0.54
1:B:303:TYR:HB3	1:B:567:TYR:CZ	2.42	0.54
1:C:194:ARG:HD3	1:C:570:TYR:HB2	1.91	0.53
1:D:199:GLN:HA	1:D:240:THR:HG23	1.89	0.53
1:H:244:GLU:OE2	1:H:269:ARG:NH1	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:232:MET:HG2	1:G:567:TYR:OH	2.09	0.53
1:C:118:ARG:NH1	1:C:166:ASP:OD1	2.41	0.53
1:D:400:ASP:OD1	1:D:400:ASP:N	2.41	0.53
1:C:570:TYR:N	1:C:571:PRO:CD	2.71	0.53
1:E:303:TYR:CD1	1:E:567:TYR:CD1	2.96	0.53
1:F:307:ALA:CB	1:F:565:VAL:HG23	2.31	0.53
1:E:483:HIS:HB2	1:E:486:ALA:HB2	1.91	0.53
1:F:476:TYR:OH	1:F:487:SER:O	2.23	0.53
1:B:466:THR:HG22	1:B:493:LEU:HB2	1.90	0.52
1:B:303:TYR:HB3	1:B:567:TYR:HE1	1.74	0.52
1:C:194:ARG:CZ	1:C:569:THR:O	2.57	0.52
1:H:400:ASP:OD1	1:H:400:ASP:N	2.42	0.52
1:D:260:ARG:NH2	1:D:264:GLU:OE2	2.43	0.52
1:A:466:THR:HG22	1:A:493:LEU:HB2	1.92	0.52
1:G:303:TYR:CE1	1:G:566:LYS:O	2.63	0.52
1:D:244:GLU:OE2	1:D:269:ARG:NH1	2.43	0.52
1:F:466:THR:HG22	1:F:493:LEU:HB2	1.90	0.52
1:A:182:LYS:NZ	8:A:802:HOH:O	2.34	0.52
1:F:469:SER:OG	6:F:600:GDP:O1B	2.25	0.51
1:C:316:LEU:HB3	1:C:323:PRO:HB3	1.92	0.51
1:C:572:GLU:O	1:C:573:ALA:C	2.52	0.51
1:E:572:GLU:OE2	1:E:572:GLU:HA	2.11	0.51
1:A:223:GLN:HE22	1:A:249:ARG:HD2	1.76	0.51
1:D:282:LYS:NZ	8:D:804:HOH:O	2.44	0.51
1:G:307:ALA:HB1	1:G:565:VAL:CG1	2.18	0.51
1:C:400:ASP:N	1:C:400:ASP:OD1	2.42	0.50
1:G:443:THR:HG23	1:G:446:SER:H	1.76	0.50
1:D:338:PRO:HB3	1:D:342:LEU:HD22	1.93	0.50
1:G:469:SER:OG	6:G:600:GDP:O2B	2.29	0.50
1:B:235:TYR:O	1:B:568:PRO:HG2	2.11	0.50
1:B:333:LYS:HB2	1:B:481:THR:HG21	1.93	0.50
1:F:291:LEU:HD12	1:F:292:PRO:HD2	1.93	0.50
1:B:400:ASP:OD1	1:B:400:ASP:N	2.43	0.50
1:B:291:LEU:HD12	1:B:292:PRO:HD2	1.94	0.50
1:B:199:GLN:HA	1:B:240:THR:HG23	1.94	0.50
1:C:512:ALA:HB2	1:C:525:PRO:HD3	1.94	0.50
1:G:208:LYS:NZ	1:G:284:LYS:O	2.35	0.50
1:A:199:GLN:HA	1:A:240:THR:HG23	1.93	0.50
1:H:109:GLY:HA3	1:H:174:ASP:OD2	2.12	0.50
1:G:400:ASP:OD1	1:G:400:ASP:N	2.43	0.49
1:B:223:GLN:HE22	1:B:249:ARG:HD2	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:400:ASP:N	1:E:400:ASP:OD1	2.43	0.49
1:E:303:TYR:CB	1:E:567:TYR:CE1	2.95	0.49
1:A:400:ASP:OD1	1:A:400:ASP:N	2.45	0.49
1:B:208:LYS:HG2	1:B:287:GLN:HA	1.94	0.49
1:E:303:TYR:CB	1:E:567:TYR:HE1	2.20	0.49
1:A:309:PRO:HG2	1:A:312:LEU:HD12	1.94	0.49
1:A:318:ARG:HG2	1:B:110:LYS:HB2	1.95	0.49
1:A:250:TYR:HB3	1:A:447:LEU:HB2	1.94	0.49
1:G:161:ARG:HB3	1:G:395:ARG:HD2	1.93	0.49
1:H:208:LYS:HG2	1:H:287:GLN:HA	1.94	0.49
1:A:303:TYR:HD1	1:A:567:TYR:CE1	2.31	0.49
1:G:365:ARG:NH1	1:G:366:ARG:O	2.46	0.49
1:H:440:ASN:N	1:H:440:ASN:OD1	2.46	0.49
1:A:565:VAL:HB	1:A:567:TYR:CE2	2.47	0.49
1:E:238:GLN:HA	1:E:570:TYR:OH	2.13	0.49
1:G:307:ALA:CB	1:G:565:VAL:HG11	2.31	0.48
1:C:167:LEU:HD13	1:D:120:GLU:HA	1.94	0.48
1:D:320:HIS:CE1	1:D:487:SER:HB3	2.47	0.48
1:H:466:THR:HG22	1:H:493:LEU:HB2	1.95	0.48
1:E:512:ALA:HB2	1:E:525:PRO:HD3	1.96	0.48
1:C:250:TYR:HB3	1:C:447:LEU:HB2	1.95	0.48
1:F:393:LEU:HD22	1:F:397:MET:HE3	1.96	0.48
1:G:440:ASN:OD1	1:G:440:ASN:N	2.46	0.48
1:A:197:TYR:OH	1:A:572:GLU:HG3	2.14	0.48
1:G:303:TYR:HE1	1:G:566:LYS:O	1.96	0.48
1:A:325:VAL:HG22	1:A:497:TYR:CE2	2.48	0.48
1:H:221:GLY:N	6:H:600:GDP:O2A	2.45	0.48
1:F:400:ASP:OD1	1:F:400:ASP:N	2.44	0.48
1:B:296:SER:HB3	4:M:1:NAG:H62	1.95	0.48
1:C:134:LEU:HD23	1:C:137:LEU:HD12	1.94	0.48
1:G:483:HIS:HB2	1:G:486:ALA:HB2	1.95	0.48
1:H:325:VAL:HG22	1:H:497:TYR:CE2	2.49	0.48
1:G:378:PRO:HD2	1:G:381:GLU:HG3	1.96	0.47
1:C:235:TYR:CZ	1:C:570:TYR:CD2	3.03	0.47
1:G:229:TYR:OH	1:G:292:PRO:O	2.30	0.47
1:C:363:HIS:HD2	1:C:472:CYS:HB2	1.79	0.47
1:F:119:ILE:HG23	1:F:163:ILE:HG23	1.97	0.47
1:D:512:ALA:HB2	1:D:525:PRO:HD3	1.97	0.47
1:A:120:GLU:HA	1:B:167:LEU:HD13	1.97	0.46
1:D:515:PRO:HD3	1:D:522:PRO:HB3	1.96	0.46
1:D:226:HIS:CE1	1:D:293:ILE:HA	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:334:TYR:O	1:H:337:ARG:NE	2.44	0.46
1:A:235:TYR:O	1:A:568:PRO:HG2	2.15	0.46
1:D:333:LYS:HB2	1:D:481:THR:HG21	1.98	0.46
1:G:291:LEU:HD12	1:G:292:PRO:HD2	1.97	0.46
1:H:268:ASP:OD2	1:H:270:SER:OG	2.23	0.46
1:H:316:LEU:HB3	1:H:323:PRO:HB3	1.97	0.46
1:E:134:LEU:HD23	1:E:137:LEU:HD12	1.98	0.46
1:G:515:PRO:HD3	1:G:522:PRO:HB3	1.97	0.46
1:C:303:TYR:HD1	1:C:567:TYR:CE1	2.34	0.46
1:A:572:GLU:H	1:A:572:GLU:CD	2.24	0.46
1:C:334:TYR:O	1:C:337:ARG:NE	2.45	0.46
1:F:248:TRP:HB3	1:F:251:ALA:O	2.16	0.46
1:B:414:LEU:HG	1:B:418:LYS:HE2	1.98	0.46
1:C:109:GLY:HA3	1:C:174:ASP:OD2	2.16	0.46
1:C:291:LEU:HD12	1:C:292:PRO:HD2	1.97	0.46
1:D:181:GLU:HA	1:D:484:PRO:HG3	1.98	0.46
1:C:372:THR:HG22	1:H:293:ILE:HD11	1.97	0.46
1:D:443:THR:HG23	1:D:446:SER:H	1.80	0.46
1:F:366:ARG:NH2	1:F:377:HIS:O	2.38	0.45
1:E:320:HIS:CE1	1:E:487:SER:HB3	2.51	0.45
1:A:138:LYS:HB3	1:B:146:GLN:HG2	1.98	0.45
1:E:515:PRO:HB3	1:E:520:GLU:HB2	1.98	0.45
1:G:494:ASP:CG	1:G:495:ASP:H	2.25	0.45
1:F:302:PRO:HB2	1:F:565:VAL:CG1	2.46	0.45
1:F:363:HIS:HD2	1:F:472:CYS:HB2	1.80	0.45
1:C:383:MET:HG3	1:C:421:TYR:CE2	2.52	0.45
1:E:358:PRO:HD2	1:E:403:ARG:HH21	1.82	0.45
1:H:291:LEU:HD12	1:H:292:PRO:HD2	1.98	0.45
1:A:248:TRP:HB3	1:A:251:ALA:O	2.17	0.45
1:E:476:TYR:OH	1:E:487:SER:O	2.34	0.45
1:H:208:LYS:NZ	1:H:284:LYS:O	2.41	0.45
1:B:244:GLU:OE2	1:B:269:ARG:NH1	2.50	0.45
1:C:272:ILE:HG12	1:C:287:GLN:OE1	2.17	0.45
1:F:239:ARG:NH2	1:F:286:VAL:O	2.30	0.45
1:H:189:GLU:O	1:H:193:ARG:CG	2.56	0.45
1:E:303:TYR:CG	1:E:567:TYR:HE1	2.35	0.44
1:H:387:GLU:OE1	1:H:424:TYR:OH	2.25	0.44
1:C:108:LEU:HD12	1:C:109:GLY:H	1.82	0.44
1:C:318:ARG:HG3	1:D:110:LYS:HA	1.99	0.44
1:F:470:GLN:O	1:F:474:VAL:HG23	2.17	0.44
1:E:189:GLU:CB	1:E:193:ARG:NH2	2.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:134:LEU:HD23	1:G:134:LEU:HA	1.85	0.44
1:E:248:TRP:HB3	1:E:251:ALA:O	2.17	0.44
1:F:316:LEU:HD21	1:F:326:TRP:CE3	2.53	0.44
1:G:476:TYR:OH	1:G:487:SER:O	2.34	0.44
1:F:217:GLY:O	8:F:801:HOH:O	2.21	0.44
1:D:498:TYR:CD2	3:L:2:BMA:H62	2.52	0.44
1:G:320:HIS:CE1	1:G:487:SER:HB3	2.53	0.44
2:I:3:BMA:H4	2:I:6:MAN:H5	2.00	0.44
2:K:3:BMA:H62	2:K:6:MAN:H2	1.71	0.44
1:D:242:ILE:HG21	1:D:269:ARG:HB3	2.00	0.44
1:F:137:LEU:HD23	1:F:140:LEU:HD12	1.99	0.44
1:A:291:LEU:HD12	1:A:292:PRO:HD2	2.00	0.44
1:C:303:TYR:CD1	1:C:567:TYR:CD1	3.06	0.44
1:C:483:HIS:HB2	1:C:486:ALA:HB2	1.99	0.44
1:F:250:TYR:HB3	1:F:447:LEU:HB2	2.00	0.44
1:G:572:GLU:CD	1:G:572:GLU:H	2.26	0.44
1:C:400:ASP:OD2	1:C:403:ARG:NH2	2.50	0.43
1:C:463:LEU:HB2	1:C:479:MET:HE1	1.99	0.43
1:H:365:ARG:NH1	1:H:366:ARG:O	2.50	0.43
1:B:179:TRP:CH2	1:B:318:ARG:HD2	2.53	0.43
1:C:310:GLU:HB2	1:C:566:LYS:HA	2.00	0.43
1:E:306:LEU:HB3	1:E:324:ALA:HB1	2.00	0.43
1:F:244:GLU:OE2	1:F:269:ARG:NH1	2.51	0.43
1:B:296:SER:HB2	4:M:2:BMA:H2	1.99	0.43
1:B:568:PRO:HB2	1:B:570:TYR:CE2	2.54	0.43
1:E:297:LEU:O	1:E:298:HIS:ND1	2.51	0.43
1:G:109:GLY:HA3	1:G:174:ASP:OD2	2.18	0.43
1:H:137:LEU:HD23	1:H:140:LEU:HD12	2.00	0.43
1:H:463:LEU:HB2	1:H:479:MET:HE1	2.00	0.43
1:F:554:PRO:HB2	1:F:557:LYS:HG3	2.01	0.43
1:B:303:TYR:C	1:B:567:TYR:OH	2.60	0.43
1:C:235:TYR:CZ	1:C:570:TYR:CE2	3.07	0.43
1:H:494:ASP:CG	1:H:495:ASP:H	2.27	0.43
1:H:256:GLU:HA	1:H:259:PHE:O	2.18	0.43
1:B:498:TYR:CD2	3:J:2:BMA:H62	2.53	0.43
1:B:303:TYR:HA	1:B:567:TYR:CZ	2.54	0.43
1:E:303:TYR:CG	1:E:567:TYR:CE1	3.06	0.43
1:H:512:ALA:HB2	1:H:525:PRO:HD3	2.00	0.43
1:B:494:ASP:CG	1:B:495:ASP:H	2.26	0.43
1:E:190:LEU:CA	1:E:193:ARG:NH1	2.82	0.43
1:G:280:GLU:N	8:G:803:HOH:O	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:309:PRO:HG3	1:G:567:TYR:CD2	2.54	0.43
1:C:351:LYS:HG3	1:G:285:ASN:CG	2.43	0.43
1:E:303:TYR:HD1	1:E:567:TYR:CD1	2.33	0.43
1:F:226:HIS:CE1	1:F:293:ILE:HA	2.54	0.43
1:D:291:LEU:HD12	1:D:292:PRO:HD2	2.01	0.42
1:F:303:TYR:CE1	1:F:566:LYS:O	2.71	0.42
1:G:220:TYR:HB2	1:G:250:TYR:CZ	2.54	0.42
1:G:334:TYR:O	1:G:337:ARG:NE	2.51	0.42
1:A:109:GLY:HA3	1:A:174:ASP:OD2	2.19	0.42
1:A:365:ARG:NE	6:A:600:GDP:O2B	2.39	0.42
1:F:512:ALA:HB2	1:F:525:PRO:HD3	2.01	0.42
1:C:494:ASP:CG	1:C:495:ASP:H	2.26	0.42
1:H:316:LEU:HD21	1:H:326:TRP:CE3	2.54	0.42
1:F:161:ARG:HB3	1:F:395:ARG:HD2	2.02	0.42
1:C:161:ARG:HB3	1:C:395:ARG:HD2	2.01	0.42
1:C:236:GLY:HA2	1:C:568:PRO:HD2	2.01	0.42
1:A:164:MET:HE1	1:B:120:GLU:HG2	2.02	0.42
1:C:325:VAL:HG22	1:C:497:TYR:CE2	2.54	0.42
1:E:316:LEU:HD21	1:E:326:TRP:CE3	2.54	0.42
1:F:147:ARG:HA	1:F:147:ARG:HD2	1.90	0.42
1:G:211:VAL:HG22	1:G:242:ILE:HB	2.01	0.42
1:G:448:ARG:HD3	1:G:448:ARG:C	2.45	0.42
1:A:401:LYS:HE3	1:F:144:GLU:OE1	2.21	0.41
1:D:320:HIS:CD2	1:D:485:ASP:HA	2.56	0.41
1:H:199:GLN:HA	1:H:240:THR:HG23	2.02	0.41
1:B:113:GLU:O	1:B:117:ARG:HG3	2.19	0.41
1:E:109:GLY:HA3	1:E:174:ASP:OD2	2.20	0.41
1:F:307:ALA:HB3	1:F:565:VAL:HG21	2.01	0.41
1:H:383:MET:HG3	1:H:421:TYR:CE2	2.55	0.41
1:A:161:ARG:NH1	8:A:801:HOH:O	2.27	0.41
1:E:226:HIS:CE1	1:E:293:ILE:HA	2.56	0.41
1:E:569:THR:HB	8:E:814:HOH:O	2.19	0.41
1:F:109:GLY:HA3	1:F:174:ASP:OD2	2.20	0.41
2:P:3:BMA:H4	2:P:6:MAN:H5	2.03	0.41
1:B:109:GLY:HA3	1:B:174:ASP:OD2	2.20	0.41
1:D:127:TRP:O	1:D:131:GLN:HG3	2.21	0.41
1:D:383:MET:HG3	1:D:421:TYR:CZ	2.55	0.41
1:E:470:GLN:O	1:E:474:VAL:HG23	2.21	0.41
1:F:309:PRO:HA	1:F:567:TYR:HD2	1.86	0.41
1:B:309:PRO:HG2	1:B:312:LEU:HD12	2.03	0.41
1:D:223:GLN:HE22	1:D:249:ARG:HD2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:494:ASP:CG	1:F:495:ASP:H	2.27	0.41
1:B:512:ALA:HB2	1:B:525:PRO:HD3	2.02	0.41
1:D:250:TYR:HD2	1:D:447:LEU:HB2	1.86	0.41
1:E:258:VAL:HB	1:E:342:LEU:HD22	2.01	0.41
1:E:406:LEU:HB3	1:E:428:SER:HB2	2.01	0.41
1:H:161:ARG:HB3	1:H:395:ARG:HD2	2.03	0.41
1:A:119:ILE:HD13	1:A:167:LEU:HD23	2.02	0.41
1:A:365:ARG:CZ	1:A:369:LYS:HB2	2.51	0.41
1:C:258:VAL:HB	1:C:342:LEU:HD22	2.02	0.41
1:E:250:TYR:HB3	1:E:447:LEU:HB2	2.03	0.41
1:E:309:PRO:HA	1:E:567:TYR:HD2	1.86	0.41
1:G:190:LEU:HD11	1:G:574:GLU:OE2	2.19	0.41
1:G:444:GLU:H	1:G:444:GLU:HG2	1.25	0.41
1:C:383:MET:HG3	1:C:421:TYR:CZ	2.55	0.41
1:E:220:TYR:HB3	6:E:600:GDP:O3'	2.21	0.41
1:E:320:HIS:CD2	1:E:485:ASP:HA	2.56	0.41
1:G:309:PRO:HG2	1:G:312:LEU:HD12	2.01	0.41
1:G:282:LYS:HG3	8:G:823:HOH:O	2.20	0.40
1:H:443:THR:HG23	1:H:446:SER:H	1.85	0.40
1:C:320:HIS:CD2	1:C:485:ASP:HA	2.56	0.40
1:F:444:GLU:OE1	1:F:448:ARG:NH1	2.51	0.40
1:G:298:HIS:HA	1:G:299:PRO:HA	1.95	0.40
1:D:362:VAL:HA	1:D:464:VAL:O	2.21	0.40
1:E:383:MET:HG3	1:E:421:TYR:CZ	2.56	0.40
1:H:127:TRP:O	1:H:131:GLN:HG3	2.21	0.40
1:H:365:ARG:CZ	1:H:369:LYS:HB2	2.51	0.40
1:A:238:GLN:HA	1:A:570:TYR:OH	2.21	0.40
1:C:303:TYR:HB3	1:C:567:TYR:HE1	1.80	0.40
1:D:498:TYR:CG	3:L:2:BMA:H62	2.56	0.40
1:F:256:GLU:HA	1:F:259:PHE:O	2.21	0.40
1:G:220:TYR:HB2	1:G:250:TYR:CE2	2.56	0.40
1:G:232:MET:HE3	1:G:232:MET:HB2	1.97	0.40
1:A:565:VAL:O	1:A:567:TYR:CD2	2.75	0.40
1:B:325:VAL:HG22	1:B:497:TYR:CE2	2.56	0.40
1:D:306:LEU:HB3	1:D:324:ALA:HB1	2.03	0.40
1:E:256:GLU:HA	1:E:259:PHE:O	2.22	0.40
1:G:303:TYR:HD1	1:G:567:TYR:CD1	2.40	0.40
1:G:485:ASP:N	8:G:801:HOH:O	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	466/535 (87%)	452 (97%)	14 (3%)	0	100	100
1	B	465/535 (87%)	452 (97%)	13 (3%)	0	100	100
1	C	465/535 (87%)	449 (97%)	16 (3%)	0	100	100
1	D	465/535 (87%)	450 (97%)	15 (3%)	0	100	100
1	E	465/535 (87%)	451 (97%)	14 (3%)	0	100	100
1	F	465/535 (87%)	452 (97%)	13 (3%)	0	100	100
1	G	465/535 (87%)	452 (97%)	13 (3%)	0	100	100
1	H	465/535 (87%)	452 (97%)	13 (3%)	0	100	100
All	All	3721/4280 (87%)	3610 (97%)	111 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	411/470 (87%)	409 (100%)	2 (0%)	81	85
1	B	410/470 (87%)	407 (99%)	3 (1%)	76	83
1	C	410/470 (87%)	408 (100%)	2 (0%)	81	85
1	D	410/470 (87%)	408 (100%)	2 (0%)	81	85
1	E	410/470 (87%)	409 (100%)	1 (0%)	87	89
1	F	410/470 (87%)	407 (99%)	3 (1%)	76	83

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	410/470 (87%)	404 (98%)	6 (2%)	57	76
1	H	410/470 (87%)	408 (100%)	2 (0%)	81	85
All	All	3281/3760 (87%)	3260 (99%)	21 (1%)	78	84

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	446	SER
1	A	572	GLU
1	B	117	ARG
1	B	440	ASN
1	B	566	LYS
1	C	566	LYS
1	C	569	THR
1	D	440	ASN
1	D	546	LYS
1	E	440	ASN
1	F	440	ASN
1	F	565	VAL
1	F	566	LYS
1	G	229	TYR
1	G	440	ASN
1	G	444	GLU
1	G	448	ARG
1	G	562	ILE
1	G	563	GLU
1	H	440	ASN
1	H	572	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	146	GLN
1	A	223	GLN
1	A	298	HIS
1	A	430	ASN
1	B	159	HIS
1	B	285	ASN
1	B	357	HIS
1	B	391	GLN

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Mol	Chain	Res	Type
1	B	491	HIS
1	B	535	HIS
1	C	215	ASN
1	C	285	ASN
1	C	298	HIS
1	D	285	ASN
1	D	491	HIS
1	E	146	GLN
1	E	285	ASN
1	F	143	ASN
1	F	146	GLN
1	F	159	HIS
1	F	215	ASN
1	F	226	HIS
1	F	285	ASN
1	F	389	HIS
1	F	445	ASN
1	F	459	GLN
1	G	159	HIS
1	G	391	GLN
1	H	215	ASN
1	H	285	ASN

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 ⓘ

42 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	I	1	2,7	14,14,15	0.35	0	17,19,21	1.41	3 (17%)
2	NAG	I	2	2	14,14,15	0.35	0	17,19,21	0.97	1 (5%)
2	BMA	I	3	2	11,11,12	0.33	0	15,15,17	0.79	0
2	MAN	I	4	2	11,11,12	0.75	0	15,15,17	0.86	1 (6%)
2	NAG	I	5	2	14,14,15	0.30	0	17,19,21	0.83	0
2	MAN	I	6	2	11,11,12	0.64	0	15,15,17	1.08	1 (6%)
3	NAG	J	1	3	15,15,15	0.43	0	21,21,21	0.67	0
3	BMA	J	2	3	11,11,12	0.26	0	15,15,17	0.69	0
3	MAN	J	3	3	11,11,12	0.61	0	15,15,17	0.93	1 (6%)
3	NAG	J	4	3	14,14,15	0.34	0	17,19,21	1.12	1 (5%)
3	MAN	J	5	3	11,11,12	0.66	0	15,15,17	1.15	1 (6%)
2	NAG	K	1	2,7	14,14,15	0.31	0	17,19,21	0.97	1 (5%)
2	NAG	K	2	2	14,14,15	0.35	0	17,19,21	1.02	1 (5%)
2	BMA	K	3	2	11,11,12	0.33	0	15,15,17	0.62	0
2	MAN	K	4	2	11,11,12	0.61	0	15,15,17	1.10	1 (6%)
2	NAG	K	5	2	14,14,15	0.33	0	17,19,21	0.87	1 (5%)
2	MAN	K	6	2	11,11,12	0.73	0	15,15,17	0.69	0
3	NAG	L	1	3	15,15,15	0.42	0	21,21,21	0.67	0
3	BMA	L	2	3	11,11,12	0.23	0	15,15,17	0.78	0
3	MAN	L	3	3	11,11,12	0.63	0	15,15,17	0.88	1 (6%)
3	NAG	L	4	3	14,14,15	0.30	0	17,19,21	0.76	1 (5%)
3	MAN	L	5	3	11,11,12	0.62	0	15,15,17	1.19	1 (6%)
4	NAG	M	1	4	15,15,15	0.47	0	21,21,21	1.28	2 (9%)
4	BMA	M	2	4	11,11,12	0.30	0	15,15,17	0.66	0
4	MAN	M	3	4	11,11,12	0.64	0	15,15,17	0.93	1 (6%)
4	NAG	M	4	4	14,14,15	0.31	0	17,19,21	0.72	0
3	NAG	N	1	3	15,15,15	0.46	0	21,21,21	1.33	3 (14%)
3	BMA	N	2	3	11,11,12	0.33	0	15,15,17	0.64	0
3	MAN	N	3	3	11,11,12	0.61	0	15,15,17	0.78	0
3	NAG	N	4	3	14,14,15	0.30	0	17,19,21	0.79	1 (5%)
3	MAN	N	5	3	11,11,12	0.62	0	15,15,17	1.16	1 (6%)
5	NAG	O	1	5	15,15,15	0.44	0	21,21,21	0.73	0
5	NAG	O	2	5	14,14,15	0.42	0	17,19,21	0.70	0
5	BMA	O	3	5	11,11,12	0.41	0	15,15,17	0.80	1 (6%)
5	MAN	O	4	5	11,11,12	0.65	0	15,15,17	0.91	0
5	NAG	O	5	5	14,14,15	0.26	0	17,19,21	0.73	1 (5%)
2	NAG	P	1	2	15,15,15	0.44	0	21,21,21	0.76	0
2	NAG	P	2	2	14,14,15	0.36	0	17,19,21	0.85	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BMA	P	3	2	11,11,12	0.43	0	15,15,17	0.80	1 (6%)
2	MAN	P	4	2	11,11,12	0.61	0	15,15,17	0.95	1 (6%)
2	NAG	P	5	2	14,14,15	0.26	0	17,19,21	0.70	1 (5%)
2	MAN	P	6	2	11,11,12	0.70	0	15,15,17	1.11	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
2	NAG	I	1	2,7	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	BMA	I	3	2	-	0/2/19/22	0/1/1/1
2	MAN	I	4	2	-	0/2/19/22	0/1/1/1
2	NAG	I	5	2	-	0/6/23/26	0/1/1/1
2	MAN	I	6	2	-	2/2/19/22	0/1/1/1
3	NAG	J	1	3	-	2/6/26/26	0/1/1/1
3	BMA	J	2	3	-	0/2/19/22	0/1/1/1
3	MAN	J	3	3	-	0/2/19/22	0/1/1/1
3	NAG	J	4	3	-	0/6/23/26	0/1/1/1
3	MAN	J	5	3	-	1/2/19/22	0/1/1/1
2	NAG	K	1	2,7	-	0/6/23/26	0/1/1/1
2	NAG	K	2	2	-	0/6/23/26	0/1/1/1
2	BMA	K	3	2	-	2/2/19/22	0/1/1/1
2	MAN	K	4	2	-	0/2/19/22	0/1/1/1
2	NAG	K	5	2	-	3/6/23/26	0/1/1/1
2	MAN	K	6	2	-	2/2/19/22	0/1/1/1
3	NAG	L	1	3	-	0/6/26/26	0/1/1/1
3	BMA	L	2	3	-	0/2/19/22	0/1/1/1
3	MAN	L	3	3	-	0/2/19/22	0/1/1/1
3	NAG	L	4	3	-	2/6/23/26	0/1/1/1
3	MAN	L	5	3	-	1/2/19/22	0/1/1/1
4	NAG	M	1	4	-	0/6/26/26	0/1/1/1
4	BMA	M	2	4	-	0/2/19/22	0/1/1/1
4	MAN	M	3	4	-	0/2/19/22	0/1/1/1
4	NAG	M	4	4	-	0/6/23/26	0/1/1/1
3	NAG	N	1	3	-	4/6/26/26	0/1/1/1
3	BMA	N	2	3	-	1/2/19/22	0/1/1/1
3	MAN	N	3	3	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	N	4	3	-	0/6/23/26	0/1/1/1
3	MAN	N	5	3	-	2/2/19/22	1/1/1/1
5	NAG	O	1	5	-	2/6/26/26	0/1/1/1
5	NAG	O	2	5	-	0/6/23/26	0/1/1/1
5	BMA	O	3	5	-	2/2/19/22	0/1/1/1
5	MAN	O	4	5	-	0/2/19/22	0/1/1/1
5	NAG	O	5	5	-	2/6/23/26	0/1/1/1
2	NAG	P	1	2	-	4/6/26/26	0/1/1/1
2	NAG	P	2	2	-	0/6/23/26	0/1/1/1
2	BMA	P	3	2	-	0/2/19/22	0/1/1/1
2	MAN	P	4	2	-	0/2/19/22	0/1/1/1
2	NAG	P	5	2	-	4/6/23/26	0/1/1/1
2	MAN	P	6	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	4	NAG	C1-O5-C5	3.67	117.10	112.19
4	M	1	NAG	O5-C1-C2	3.64	113.17	109.52
3	J	5	MAN	C1-O5-C5	-3.59	107.38	112.19
2	I	1	NAG	C1-O5-C5	3.59	116.99	112.19
3	N	1	NAG	O5-C1-C2	3.58	113.11	109.52
3	L	5	MAN	C1-O5-C5	-3.48	107.53	112.19
2	I	1	NAG	C4-C3-C2	-3.36	106.10	111.02
3	N	1	NAG	C1-C2-C3	3.23	114.95	110.54
4	M	1	NAG	C1-C2-C3	3.23	114.94	110.54
2	P	6	MAN	C1-O5-C5	-3.14	107.97	112.19
2	I	2	NAG	C1-O5-C5	2.90	116.07	112.19
2	P	2	NAG	C1-O5-C5	2.84	116.00	112.19
2	I	6	MAN	C1-O5-C5	-2.82	108.41	112.19
2	K	2	NAG	C1-O5-C5	2.77	115.89	112.19
2	K	5	NAG	C1-O5-C5	2.71	115.81	112.19
2	K	4	MAN	C2-C3-C4	-2.64	106.22	110.86
3	N	1	NAG	C4-C3-C2	2.38	113.87	110.40
2	P	4	MAN	C1-O5-C5	-2.34	109.06	112.19
3	N	4	NAG	C1-O5-C5	2.29	115.25	112.19
5	O	5	NAG	C1-O5-C5	2.29	115.25	112.19
3	J	3	MAN	C2-C3-C4	-2.26	106.89	110.86
4	M	3	MAN	C2-C3-C4	-2.24	106.92	110.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	4	MAN	C1-O5-C5	-2.23	109.19	112.19
3	L	3	MAN	C2-C3-C4	-2.22	106.95	110.86
2	P	5	NAG	C1-O5-C5	2.22	115.16	112.19
5	O	3	BMA	C1-C2-C3	2.20	112.85	109.64
3	N	5	MAN	C6-C5-C4	-2.17	107.69	113.02
2	I	1	NAG	O4-C4-C3	2.14	115.42	110.38
2	K	1	NAG	C4-C3-C2	-2.14	107.89	111.02
2	P	3	BMA	C1-C2-C3	2.12	112.73	109.64
3	L	4	NAG	C1-O5-C5	2.08	114.97	112.19

There are no chirality outliers.

All (38) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	P	1	NAG	C8-C7-N2-C2
2	P	1	NAG	O7-C7-N2-C2
2	K	3	BMA	C4-C5-C6-O6
3	N	1	NAG	O5-C5-C6-O6
2	P	6	MAN	O5-C5-C6-O6
2	I	6	MAN	O5-C5-C6-O6
2	K	3	BMA	O5-C5-C6-O6
2	P	1	NAG	O5-C5-C6-O6
2	P	1	NAG	C4-C5-C6-O6
3	J	1	NAG	C4-C5-C6-O6
3	N	1	NAG	C4-C5-C6-O6
3	J	1	NAG	O5-C5-C6-O6
2	I	6	MAN	C4-C5-C6-O6
3	L	4	NAG	O5-C5-C6-O6
2	P	6	MAN	C4-C5-C6-O6
3	L	4	NAG	C4-C5-C6-O6
5	O	1	NAG	C8-C7-N2-C2
3	J	5	MAN	O5-C5-C6-O6
3	L	5	MAN	O5-C5-C6-O6
2	K	6	MAN	C4-C5-C6-O6
2	K	6	MAN	O5-C5-C6-O6
5	O	3	BMA	C4-C5-C6-O6
5	O	5	NAG	C4-C5-C6-O6
2	P	5	NAG	C4-C5-C6-O6
2	K	5	NAG	C4-C5-C6-O6
2	P	5	NAG	C8-C7-N2-C2
5	O	1	NAG	O7-C7-N2-C2
2	K	5	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	P	5	NAG	O7-C7-N2-C2
5	O	5	NAG	O5-C5-C6-O6
2	P	5	NAG	O5-C5-C6-O6
5	O	3	BMA	O5-C5-C6-O6
2	K	5	NAG	C1-C2-N2-C7
3	N	2	BMA	C4-C5-C6-O6
3	N	5	MAN	O5-C5-C6-O6
3	N	5	MAN	C4-C5-C6-O6
3	N	1	NAG	C8-C7-N2-C2
3	N	1	NAG	O7-C7-N2-C2

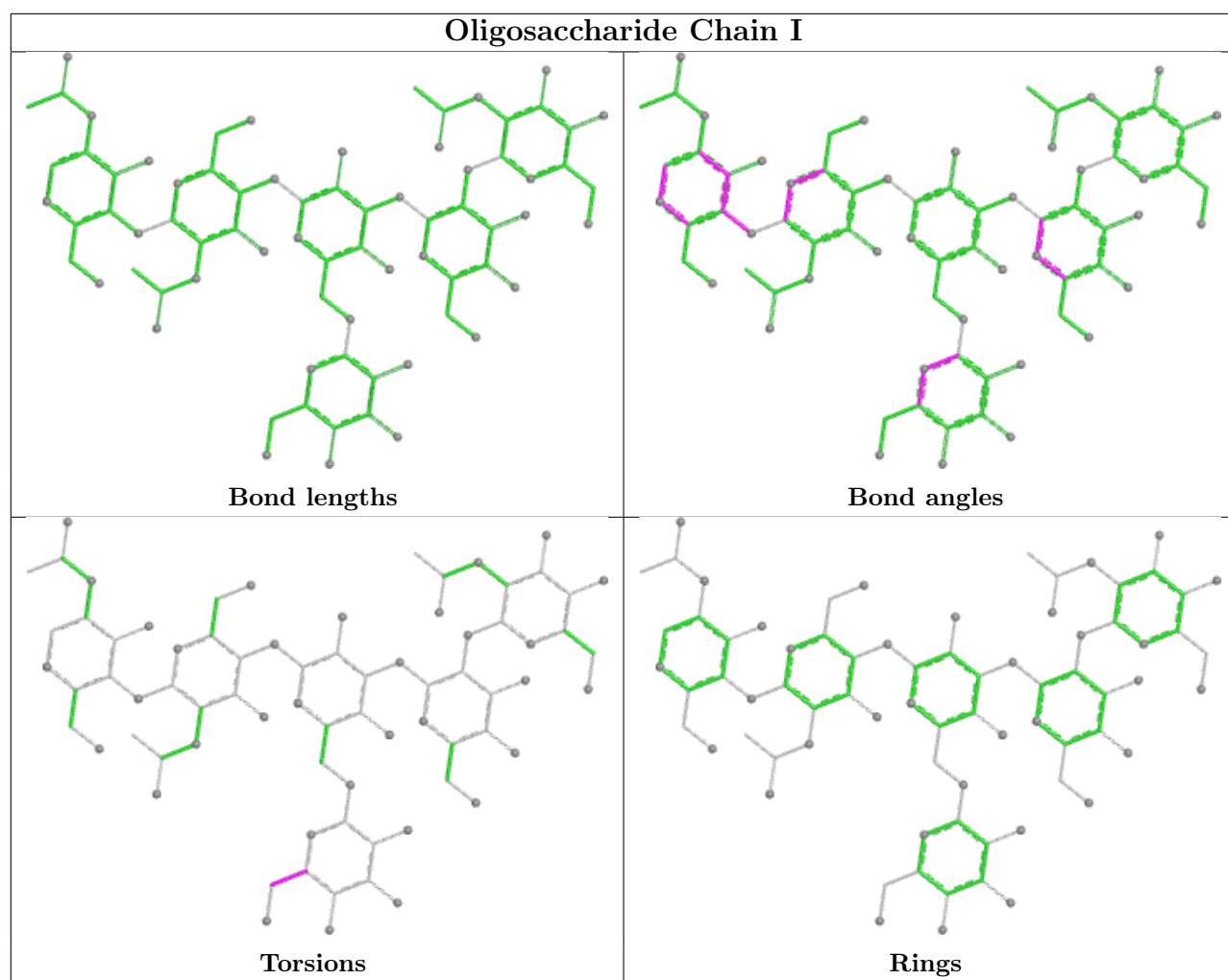
All (1) ring outliers are listed below:

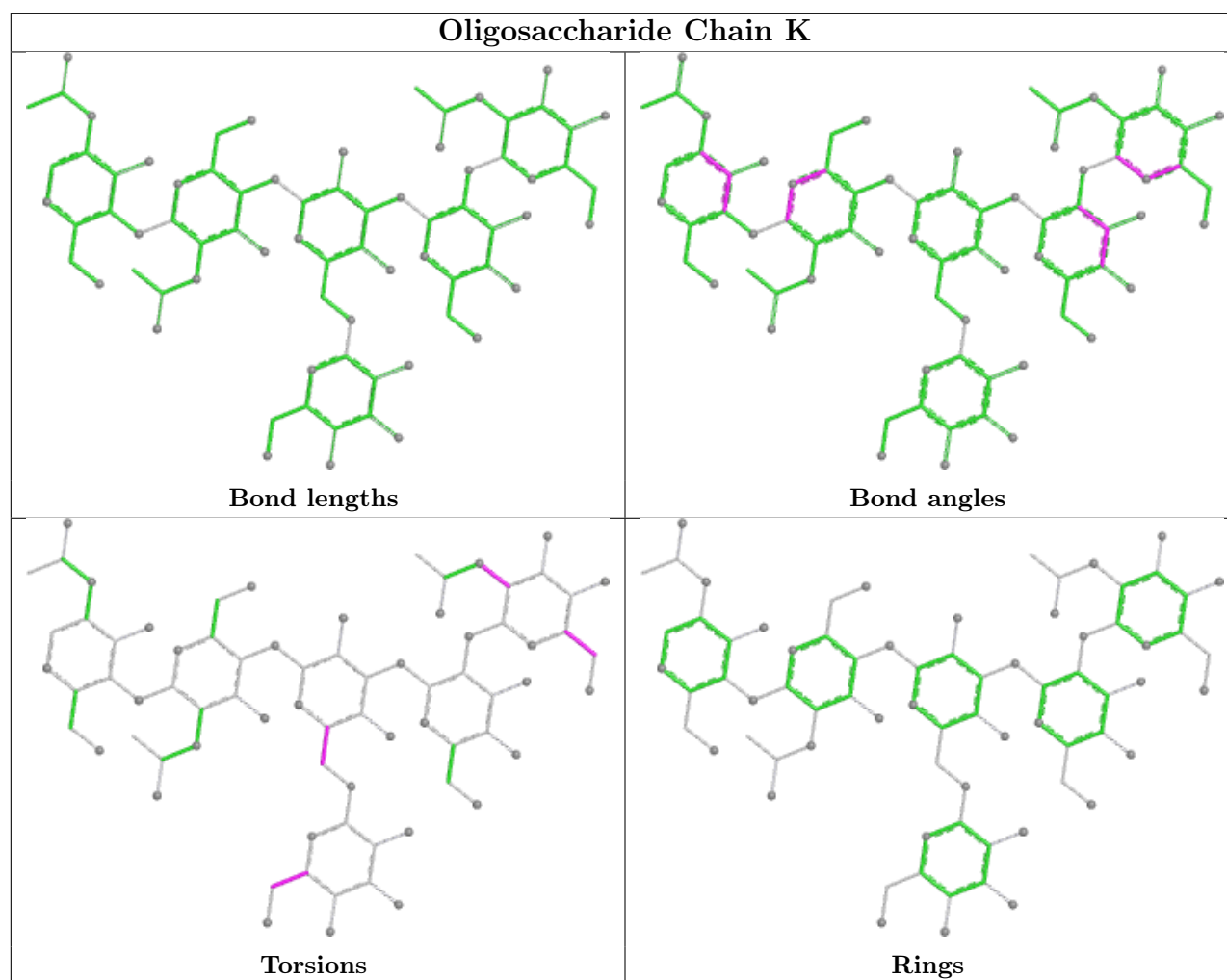
Mol	Chain	Res	Type	Atoms
3	N	5	MAN	C1-C2-C3-C4-C5-O5

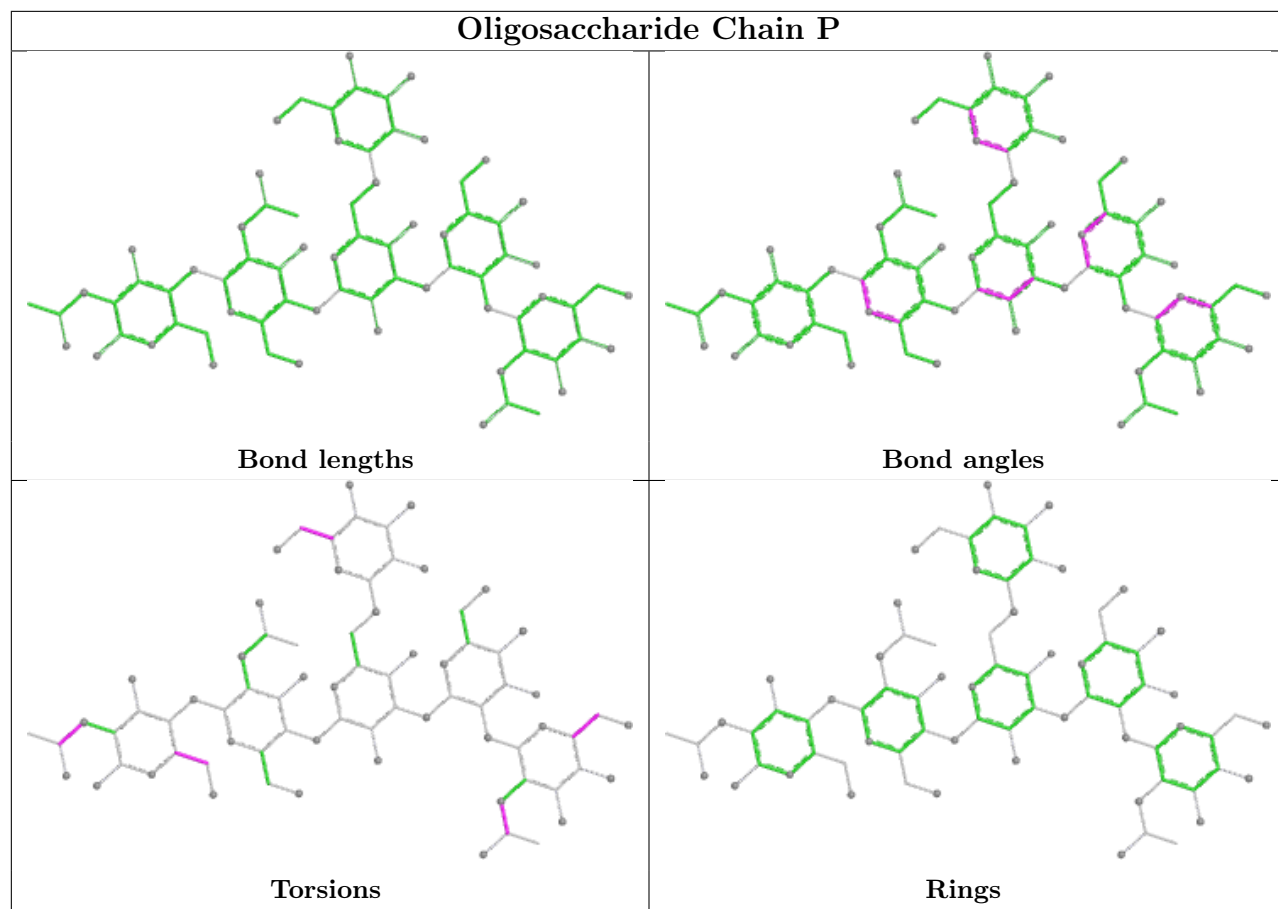
11 monomers are involved in 9 short contacts:

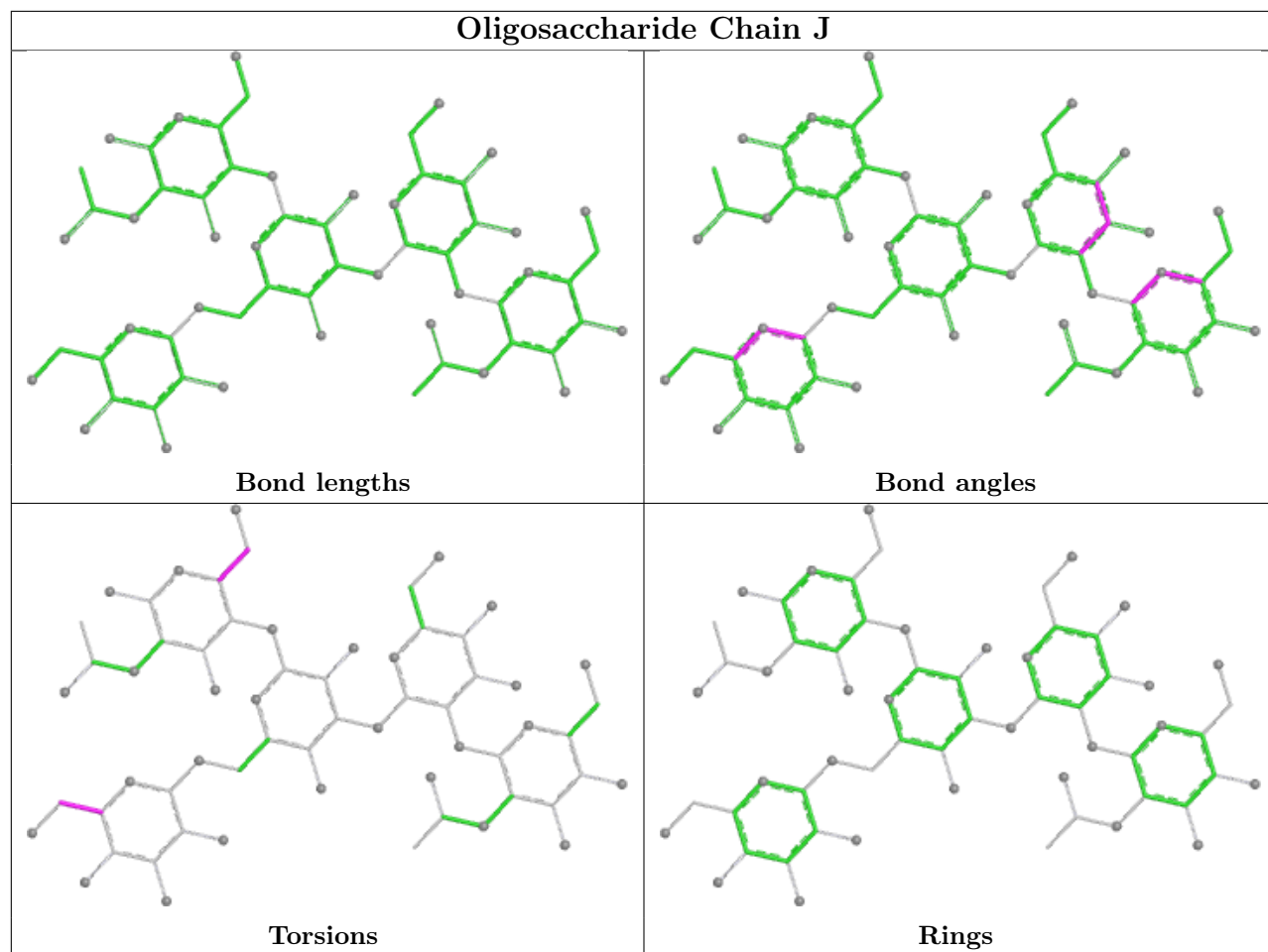
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	M	1	NAG	1	0
2	P	3	BMA	1	0
4	M	2	BMA	1	0
3	J	2	BMA	1	0
3	L	2	BMA	2	0
2	I	6	MAN	1	0
2	I	3	BMA	1	0
3	J	3	MAN	1	0
2	K	3	BMA	1	0
2	P	6	MAN	1	0
2	K	6	MAN	1	0

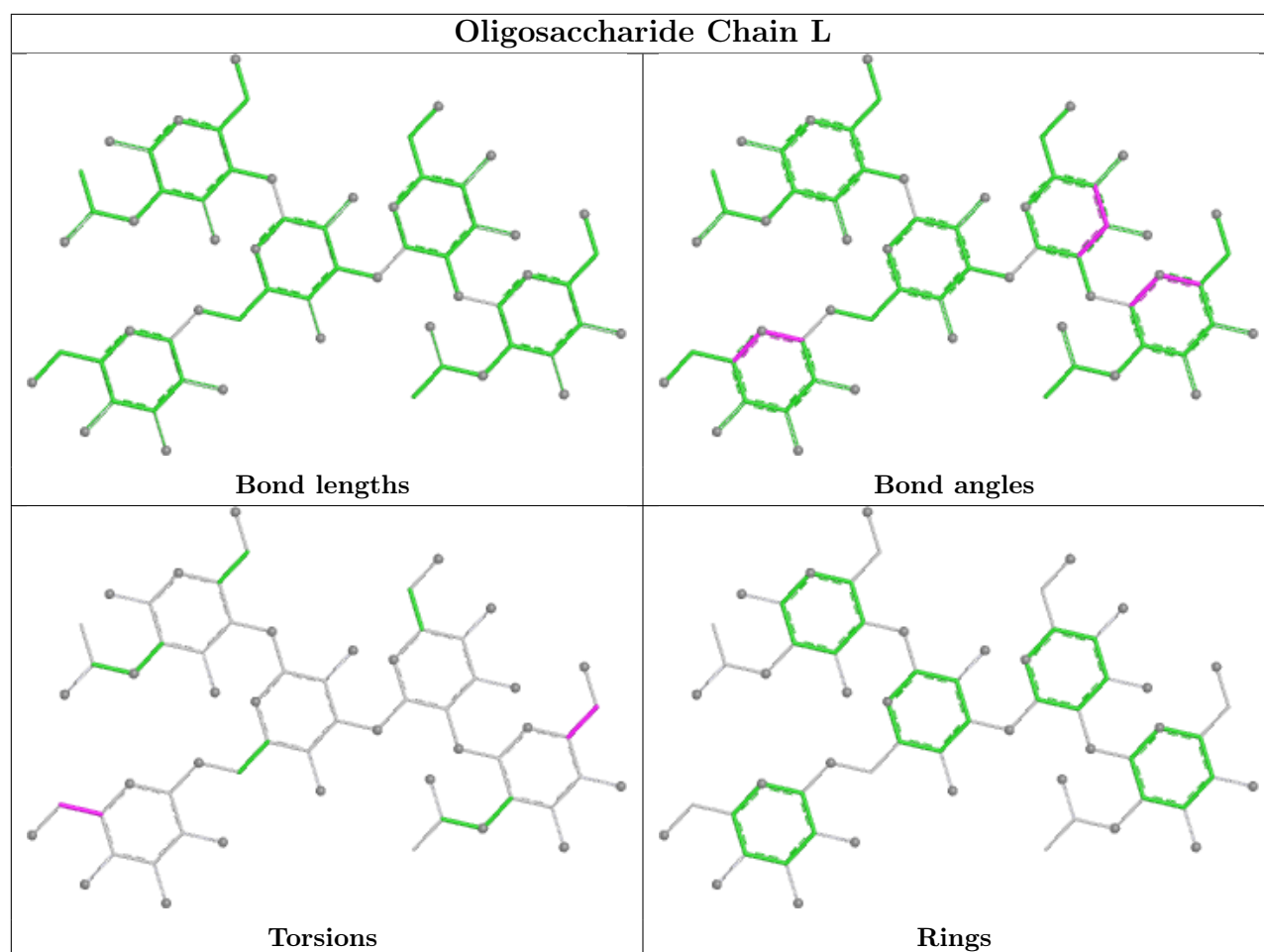
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

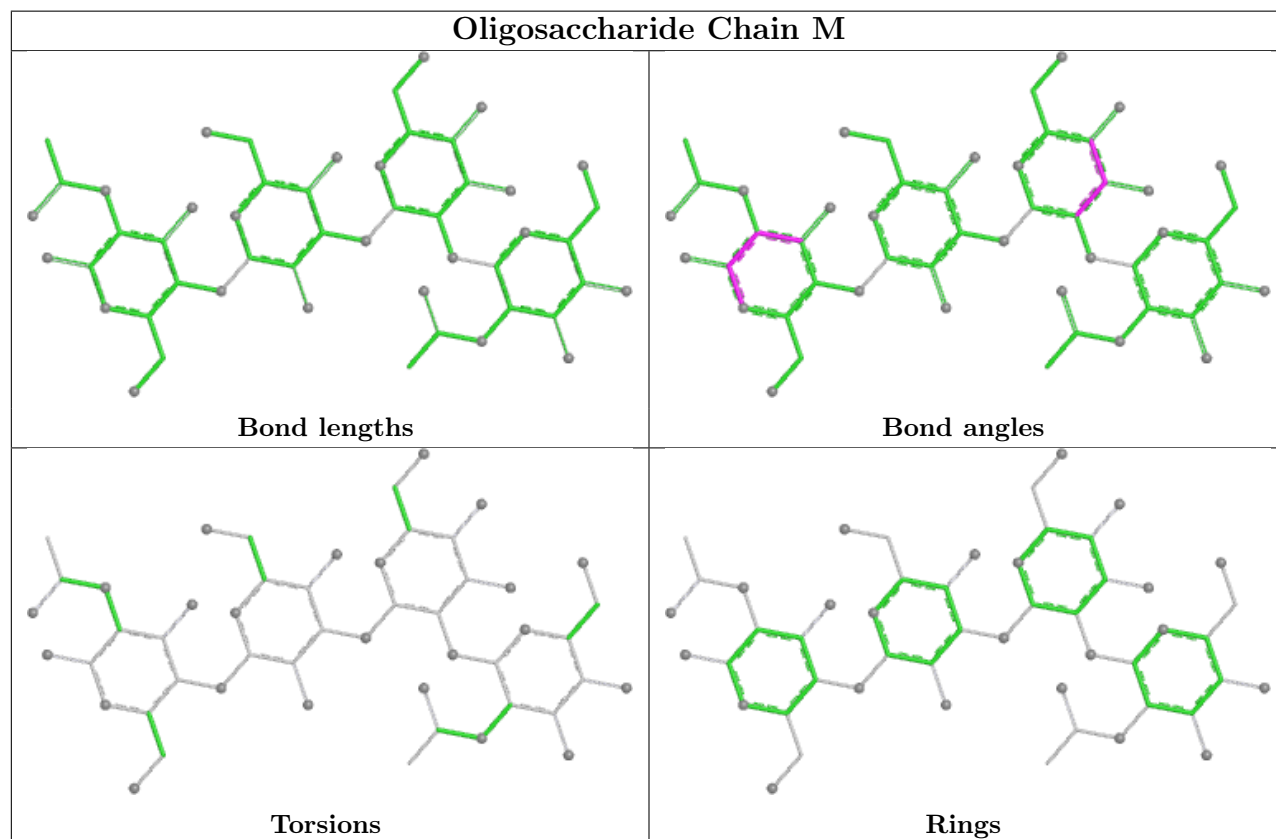
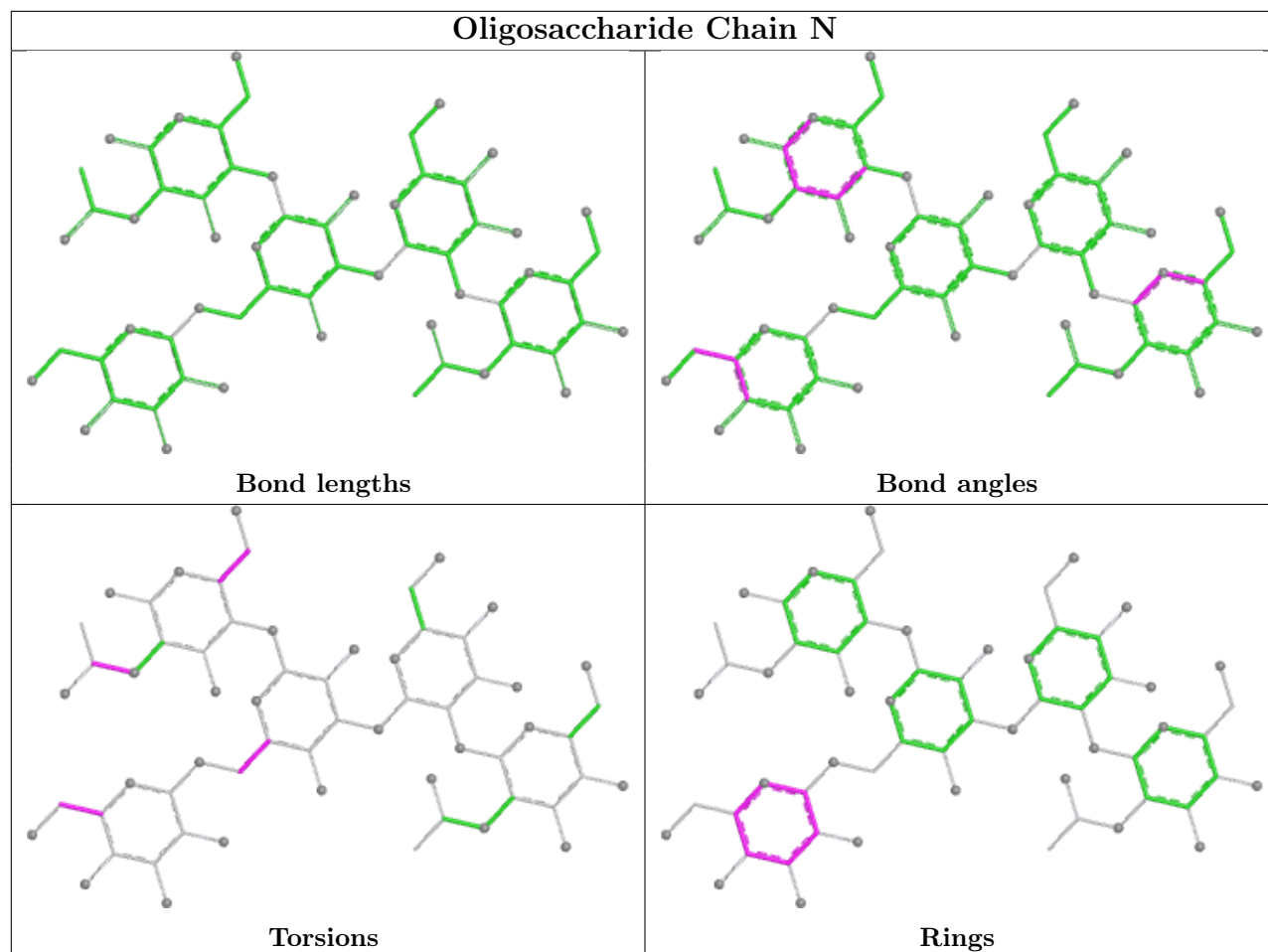


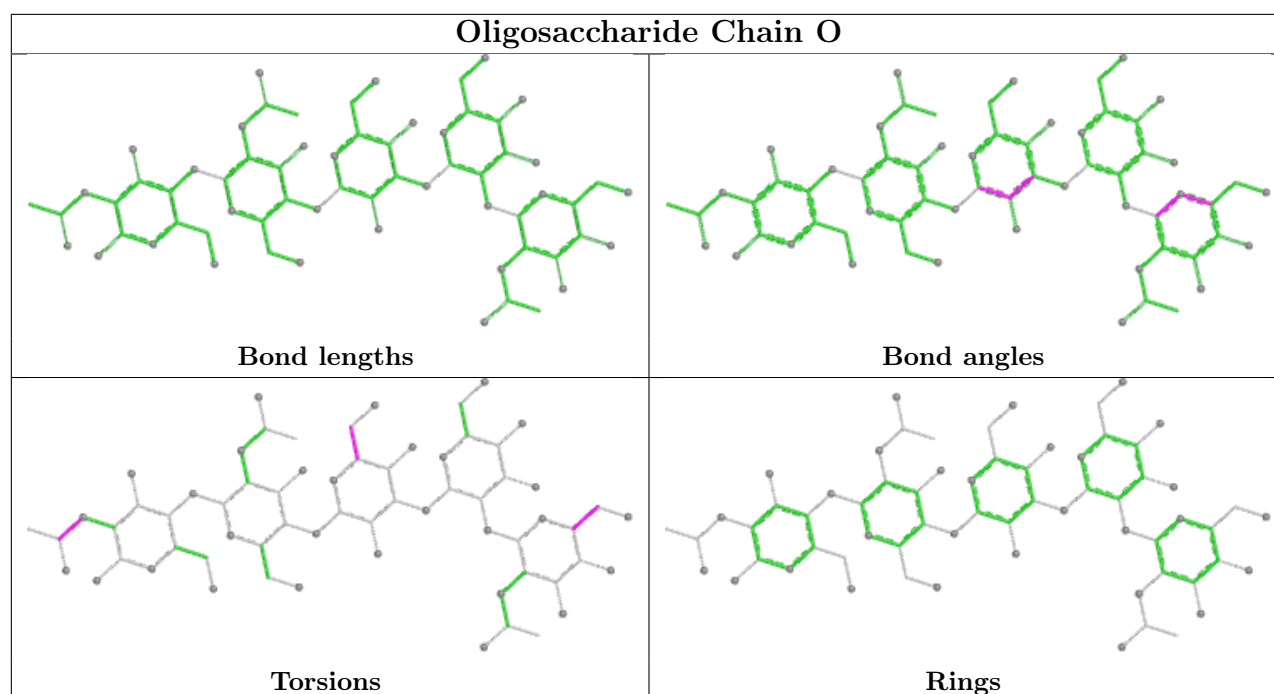












5.6 Ligand geometry [i](#)

10 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
6	GDP	D	600	-	29,30,30	1.13	1 (3%)	45,47,47	1.66	9 (20%)
6	GDP	B	600	-	29,30,30	1.14	1 (3%)	45,47,47	1.64	9 (20%)
6	GDP	H	600	-	29,30,30	1.11	1 (3%)	45,47,47	1.62	9 (20%)
7	ASN	A	700	2	6,7,8	0.48	0	4,8,10	0.15	0
7	ASN	C	700	2	6,7,8	0.48	0	4,8,10	0.17	0
6	GDP	E	600	-	29,30,30	1.15	1 (3%)	45,47,47	1.64	8 (17%)
6	GDP	F	600	-	29,30,30	1.15	2 (6%)	45,47,47	1.63	9 (20%)
6	GDP	C	600	-	29,30,30	1.11	1 (3%)	45,47,47	1.63	9 (20%)
6	GDP	G	600	-	29,30,30	1.11	1 (3%)	45,47,47	1.60	9 (20%)
6	GDP	A	600	-	29,30,30	1.10	1 (3%)	45,47,47	1.65	9 (20%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GDP	D	600	-	-	6/16/32/32	0/3/3/3
6	GDP	B	600	-	-	6/16/32/32	0/3/3/3
6	GDP	H	600	-	-	6/16/32/32	0/3/3/3
7	ASN	A	700	2	-	1/5/6/8	-
7	ASN	C	700	2	-	0/5/6/8	-
6	GDP	E	600	-	-	3/16/32/32	0/3/3/3
6	GDP	F	600	-	-	3/16/32/32	0/3/3/3
6	GDP	C	600	-	-	3/16/32/32	0/3/3/3
6	GDP	G	600	-	-	6/16/32/32	0/3/3/3
6	GDP	A	600	-	-	4/16/32/32	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	E	600	GDP	PB-O1B	3.54	1.61	1.50
6	F	600	GDP	PB-O1B	3.54	1.61	1.50
6	B	600	GDP	PB-O1B	3.54	1.61	1.50
6	D	600	GDP	PB-O1B	3.50	1.61	1.50
6	G	600	GDP	PA-O1A	3.10	1.61	1.50
6	C	600	GDP	PA-O1A	3.10	1.61	1.50
6	A	600	GDP	PA-O1A	3.09	1.61	1.50
6	H	600	GDP	PA-O1A	3.07	1.61	1.50
6	F	600	GDP	PA-O3A	2.17	1.61	1.59

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	600	GDP	C5-C4-N3	-5.24	120.05	128.39
6	B	600	GDP	C5-C4-N3	-5.20	120.11	128.39
6	D	600	GDP	C5-C4-N3	-5.18	120.14	128.39
6	E	600	GDP	C5-C4-N3	-5.16	120.17	128.39
6	A	600	GDP	C5-C4-N3	-5.13	120.22	128.39
6	D	600	GDP	C2-N3-C4	5.13	121.14	112.30
6	A	600	GDP	C2-N3-C4	5.10	121.08	112.30
6	B	600	GDP	C2-N3-C4	5.07	121.03	112.30
6	F	600	GDP	C2-N3-C4	5.07	121.03	112.30
6	E	600	GDP	C2-N3-C4	5.05	121.00	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	600	GDP	C5-C4-N3	-5.05	120.36	128.39
6	C	600	GDP	C2-N3-C4	5.04	120.98	112.30
6	G	600	GDP	C5-C4-N3	-5.02	120.41	128.39
6	H	600	GDP	C5-C4-N3	-5.00	120.43	128.39
6	H	600	GDP	C2-N3-C4	4.99	120.90	112.30
6	G	600	GDP	C2-N3-C4	4.94	120.81	112.30
6	E	600	GDP	N9-C4-N3	3.94	133.83	125.95
6	C	600	GDP	N9-C4-N3	3.91	133.76	125.95
6	B	600	GDP	N9-C4-N3	3.89	133.72	125.95
6	H	600	GDP	N9-C4-N3	3.81	133.57	125.95
6	A	600	GDP	N9-C4-N3	3.72	133.40	125.95
6	G	600	GDP	N9-C4-N3	3.71	133.37	125.95
6	D	600	GDP	N9-C4-N3	3.70	133.35	125.95
6	F	600	GDP	N9-C4-N3	3.64	133.23	125.95
6	F	600	GDP	C8-N7-C5	2.74	109.14	104.26
6	A	600	GDP	C8-N7-C5	2.72	109.11	104.26
6	D	600	GDP	C8-N7-C5	2.72	109.10	104.26
6	C	600	GDP	C8-N7-C5	2.71	109.08	104.26
6	B	600	GDP	C8-N7-C5	2.69	109.04	104.26
6	E	600	GDP	C8-N7-C5	2.67	109.01	104.26
6	H	600	GDP	C8-N7-C5	2.66	109.01	104.26
6	G	600	GDP	C8-N7-C5	2.64	108.96	104.26
6	D	600	GDP	C2-N1-C6	-2.60	120.40	125.11
6	E	600	GDP	C2-N1-C6	-2.56	120.48	125.11
6	B	600	GDP	C2-N1-C6	-2.55	120.48	125.11
6	C	600	GDP	C2-N1-C6	-2.55	120.48	125.11
6	A	600	GDP	C2-N1-C6	-2.53	120.52	125.11
6	F	600	GDP	C2-N1-C6	-2.52	120.53	125.11
6	A	600	GDP	C6-C5-N7	2.51	134.87	130.29
6	H	600	GDP	C2-N1-C6	-2.51	120.56	125.11
6	F	600	GDP	C6-C5-N7	2.51	134.86	130.29
6	D	600	GDP	C6-C5-N7	2.51	134.85	130.29
6	G	600	GDP	C2-N1-C6	-2.48	120.61	125.11
6	H	600	GDP	C6-C5-N7	2.40	134.65	130.29
6	E	600	GDP	O6-C6-C5	-2.39	120.22	126.53
6	G	600	GDP	C6-C5-N7	2.38	134.61	130.29
6	H	600	GDP	O6-C6-C5	-2.36	120.29	126.53
6	B	600	GDP	C6-C5-N7	2.33	134.53	130.29
6	H	600	GDP	C5-C6-N1	2.28	119.06	113.25
6	B	600	GDP	O6-C6-C5	-2.28	120.52	126.53
6	C	600	GDP	O6-C6-C5	-2.28	120.52	126.53
6	E	600	GDP	C5-C6-N1	2.27	119.04	113.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	E	600	GDP	C6-C5-N7	2.27	134.42	130.29
6	D	600	GDP	C5-C6-N1	2.27	119.03	113.25
6	C	600	GDP	C6-C5-N7	2.26	134.40	130.29
6	D	600	GDP	O6-C6-C5	-2.26	120.57	126.53
6	A	600	GDP	C4-C5-N7	-2.24	107.12	110.67
6	B	600	GDP	C5-C6-N1	2.24	118.96	113.25
6	D	600	GDP	C4-C5-N7	-2.23	107.14	110.67
6	A	600	GDP	C5-C6-N1	2.21	118.88	113.25
6	C	600	GDP	C5-C6-N1	2.21	118.87	113.25
6	F	600	GDP	C4-C5-N7	-2.20	107.18	110.67
6	F	600	GDP	O6-C6-C5	-2.20	120.73	126.53
6	G	600	GDP	O6-C6-C5	-2.20	120.73	126.53
6	F	600	GDP	C5-C6-N1	2.19	118.83	113.25
6	G	600	GDP	C5-C6-N1	2.18	118.80	113.25
6	A	600	GDP	O6-C6-C5	-2.17	120.79	126.53
6	C	600	GDP	C4-C5-N7	-2.11	107.33	110.67
6	G	600	GDP	C4-C5-N7	-2.09	107.36	110.67
6	B	600	GDP	C4-C5-N7	-2.07	107.39	110.67
6	H	600	GDP	C4-C5-N7	-2.00	107.49	110.67

There are no chirality outliers.

All (38) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	600	GDP	PA-O3A-PB-O3B
6	A	600	GDP	C5'-O5'-PA-O3A
6	B	600	GDP	C5'-O5'-PA-O1A
6	C	600	GDP	PA-O3A-PB-O3B
6	D	600	GDP	PA-O3A-PB-O3B
6	D	600	GDP	PB-O3A-PA-O5'
6	D	600	GDP	C5'-O5'-PA-O3A
6	D	600	GDP	C5'-O5'-PA-O1A
6	E	600	GDP	C5'-O5'-PA-O3A
6	E	600	GDP	C5'-O5'-PA-O1A
6	G	600	GDP	PA-O3A-PB-O2B
6	G	600	GDP	PA-O3A-PB-O3B
6	G	600	GDP	C5'-O5'-PA-O2A
6	H	600	GDP	PA-O3A-PB-O3B
6	H	600	GDP	PB-O3A-PA-O5'
6	H	600	GDP	C5'-O5'-PA-O2A
7	A	700	ASN	O-C-CA-CB
6	A	600	GDP	PB-O3A-PA-O5'

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Mol	Chain	Res	Type	Atoms
6	B	600	GDP	PB-O3A-PA-O5'
6	C	600	GDP	PB-O3A-PA-O5'
6	E	600	GDP	PB-O3A-PA-O5'
6	F	600	GDP	PB-O3A-PA-O5'
6	G	600	GDP	PB-O3A-PA-O5'
6	B	600	GDP	PA-O3A-PB-O1B
6	B	600	GDP	C5'-O5'-PA-O3A
6	G	600	GDP	C5'-O5'-PA-O3A
6	G	600	GDP	C5'-O5'-PA-O1A
6	H	600	GDP	C5'-O5'-PA-O3A
6	H	600	GDP	PA-O3A-PB-O1B
6	B	600	GDP	PA-O3A-PB-O2B
6	B	600	GDP	PA-O3A-PB-O3B
6	C	600	GDP	PA-O3A-PB-O2B
6	D	600	GDP	PA-O3A-PB-O2B
6	H	600	GDP	PA-O3A-PB-O2B
6	F	600	GDP	PB-O3A-PA-O1A
6	F	600	GDP	PB-O3A-PA-O2A
6	A	600	GDP	PA-O3A-PB-O1B
6	D	600	GDP	PA-O3A-PB-O1B

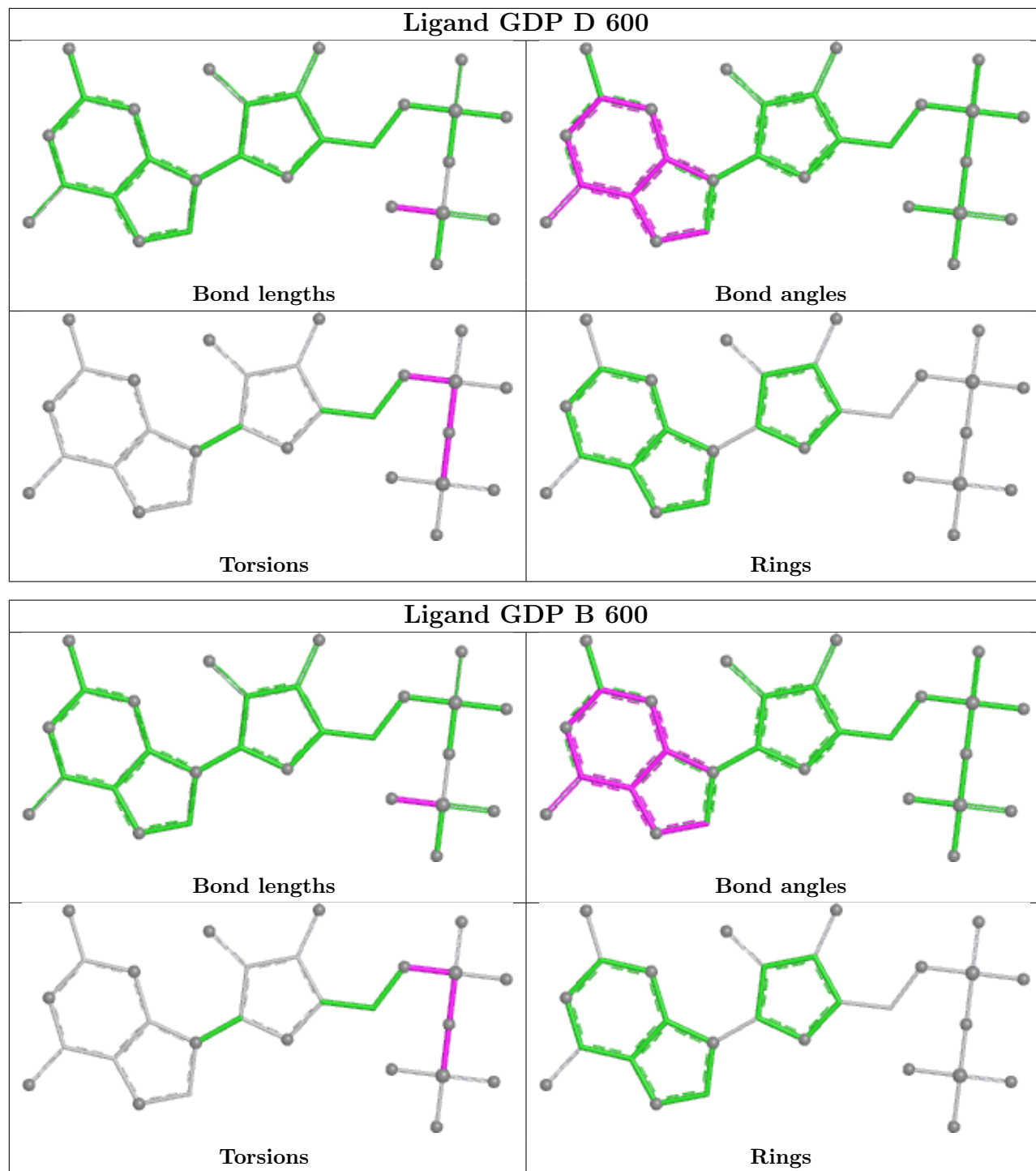
There are no ring outliers.

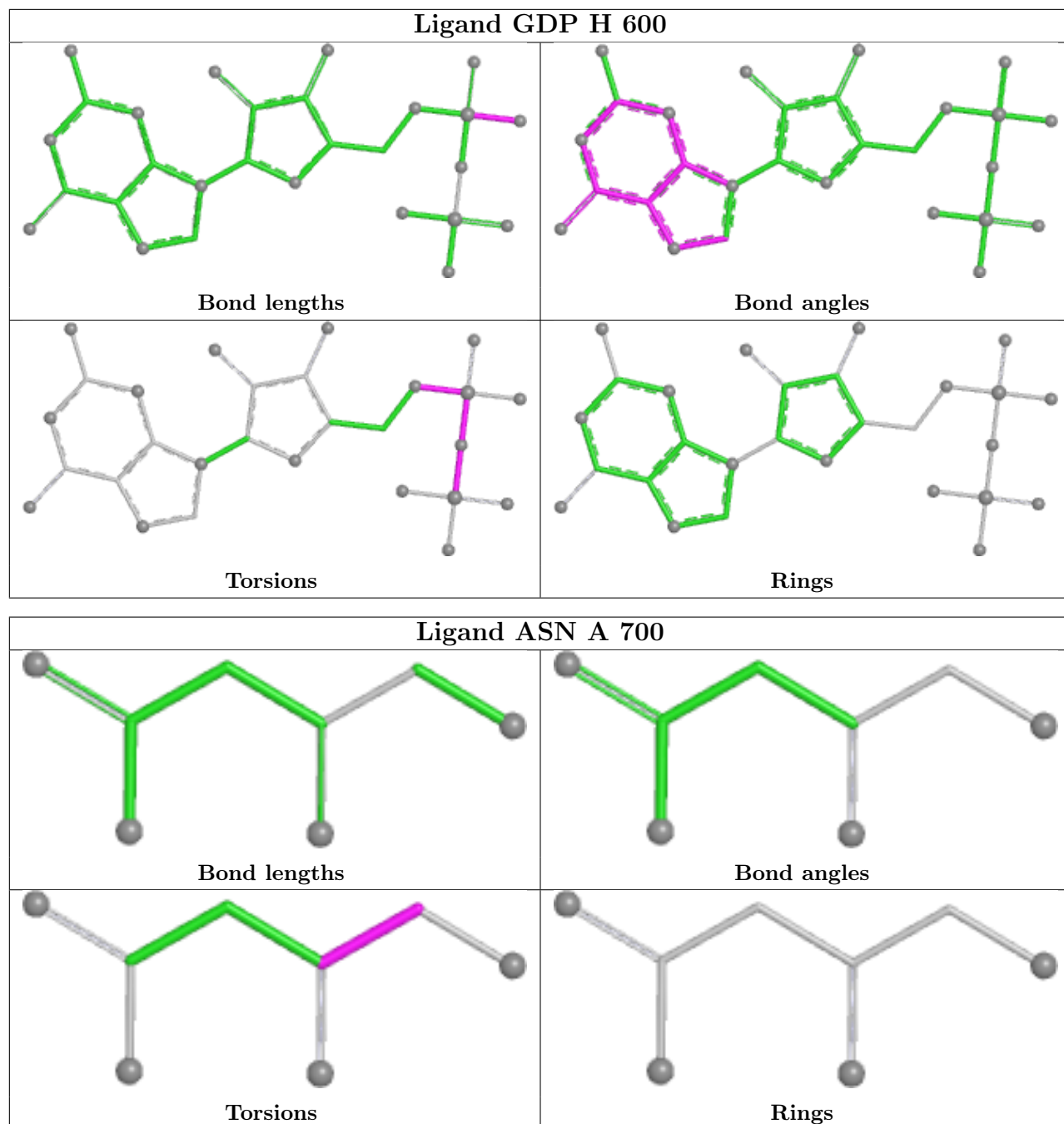
5 monomers are involved in 5 short contacts:

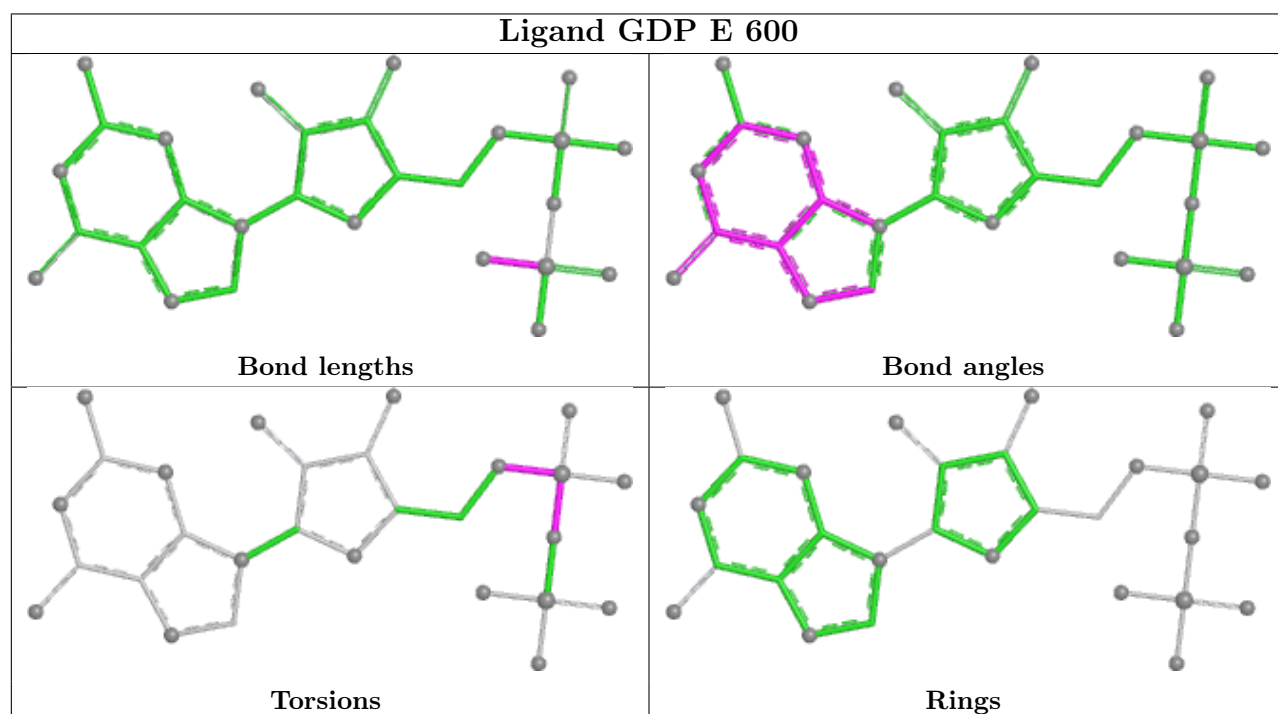
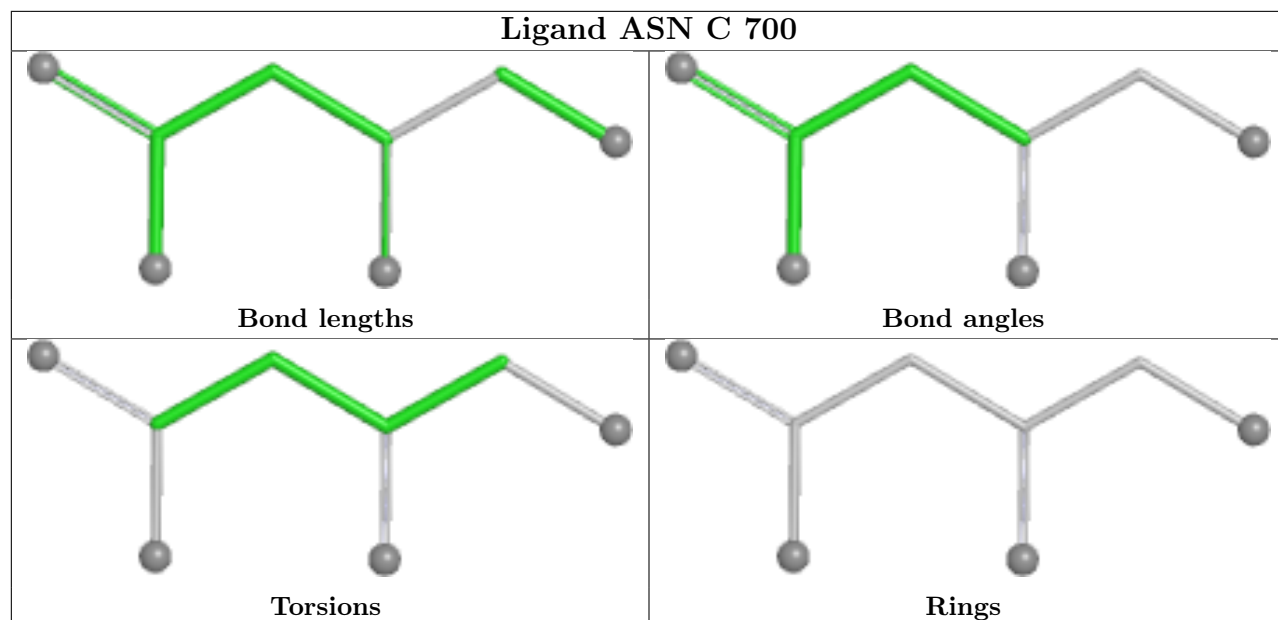
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	H	600	GDP	1	0
6	E	600	GDP	1	0
6	F	600	GDP	1	0
6	G	600	GDP	1	0
6	A	600	GDP	1	0

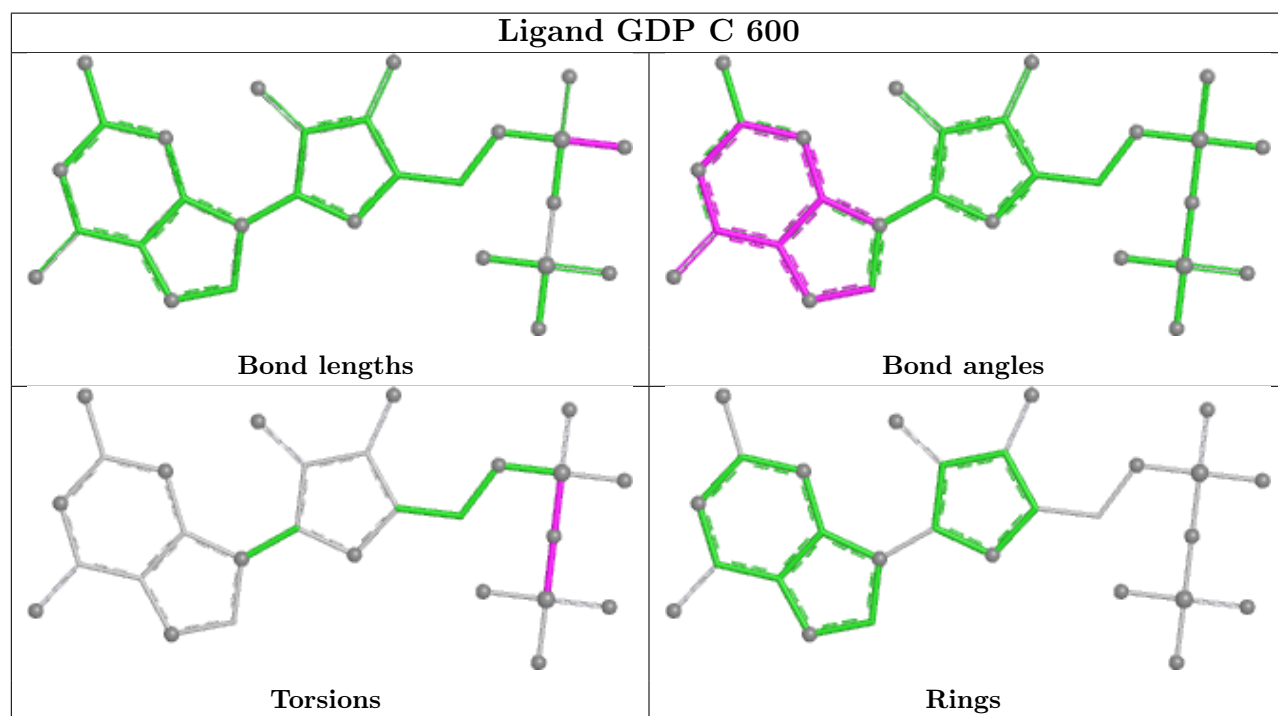
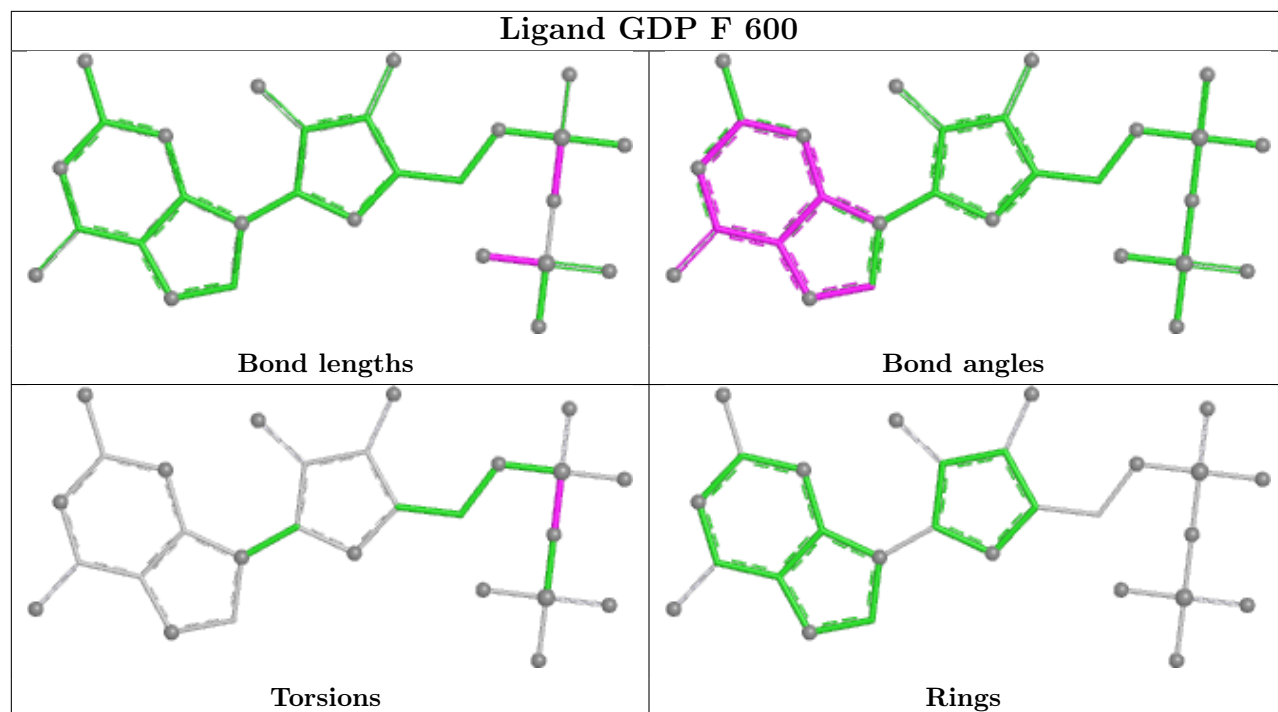
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient

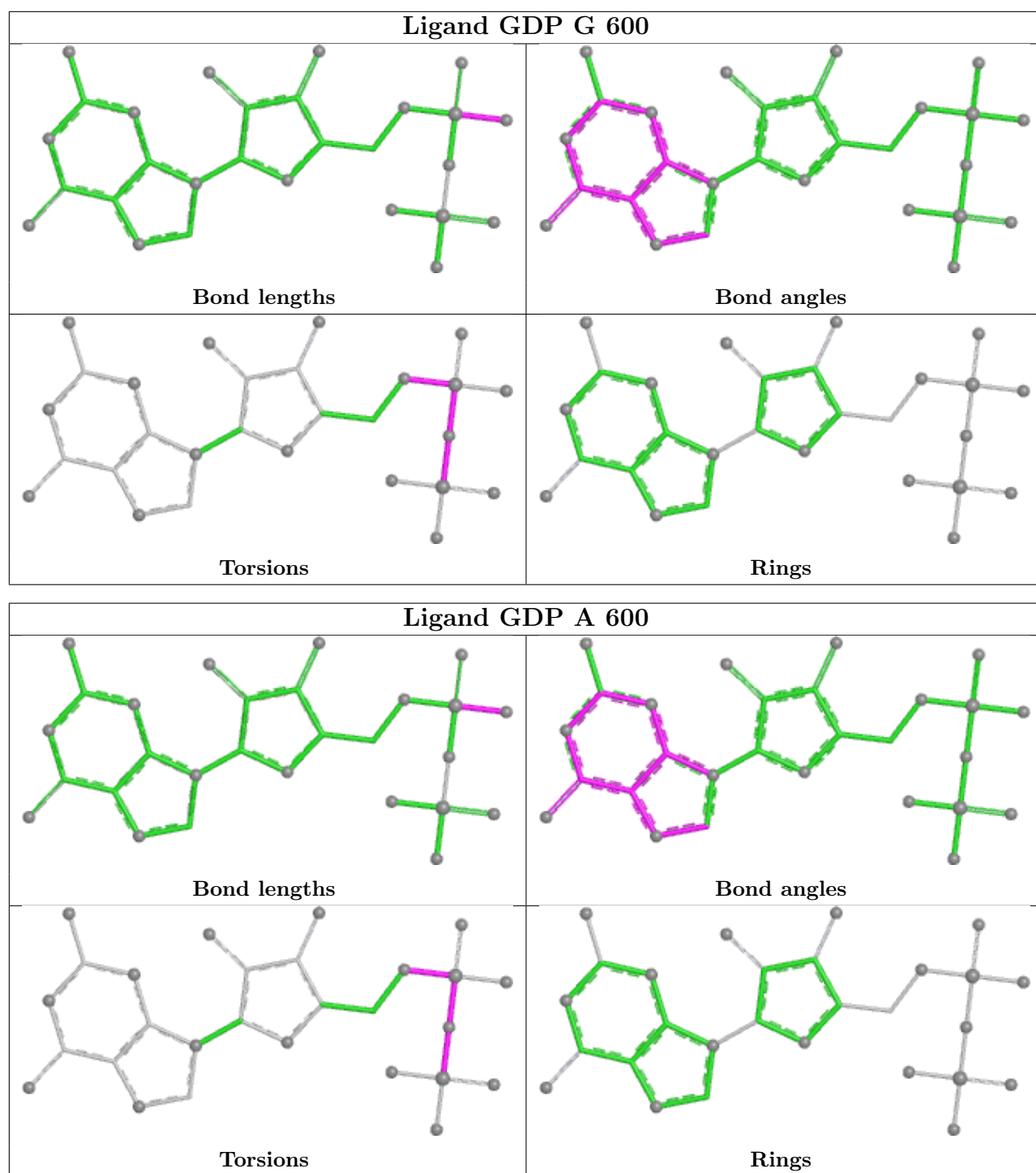
equivalents in the CSD to analyse the geometry.











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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, 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	467/535 (87%)	-1.29	0 100 100	25, 58, 84, 112	1 (0%)
1	B	467/535 (87%)	-1.34	0 100 100	35, 58, 84, 125	0
1	C	467/535 (87%)	-1.26	0 100 100	34, 61, 88, 131	0
1	D	467/535 (87%)	-1.35	0 100 100	35, 58, 80, 116	0
1	E	467/535 (87%)	-1.29	0 100 100	37, 60, 87, 113	0
1	F	467/535 (87%)	-1.32	0 100 100	36, 57, 85, 124	0
1	G	467/535 (87%)	-1.29	0 100 100	34, 60, 88, 127	0
1	H	467/535 (87%)	-1.27	0 100 100	36, 65, 98, 138	0
All	All	3736/4280 (87%)	-1.30	0 100 100	25, 60, 87, 138	1 (0%)

There are no RSRZ outliers to report.

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MAN	I	4	11/12	0.98	0.06	68,73,75,76	0
2	BMA	I	3	11/12	0.99	0.04	60,64,71,77	0
2	NAG	I	2	14/15	0.99	0.03	49,59,70,72	0

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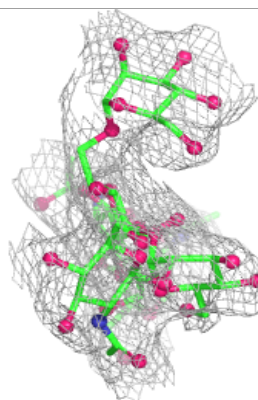
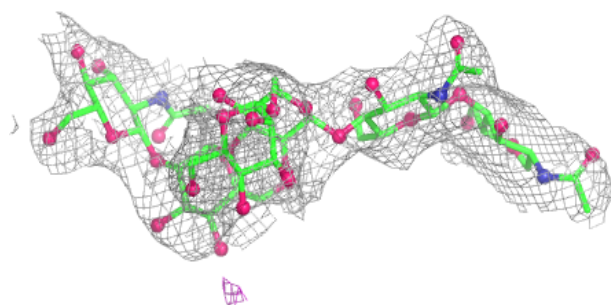
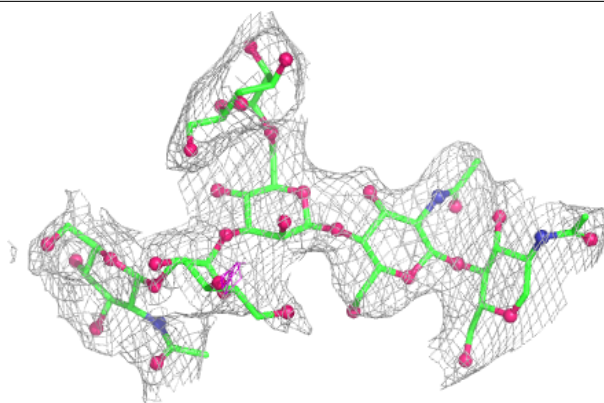
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	I	5	14/15	0.99	0.03	60,64,69,71	0
2	MAN	I	6	11/12	0.99	0.05	78,82,83,84	0
2	NAG	K	1	14/15	0.99	0.04	45,52,56,58	0
2	NAG	K	2	14/15	0.99	0.04	46,49,57,58	0
2	MAN	K	4	11/12	0.99	0.03	59,62,65,66	0
2	NAG	K	5	14/15	0.99	0.04	59,67,69,69	0
2	MAN	K	6	11/12	0.99	0.06	68,73,75,75	0
2	NAG	P	2	14/15	0.99	0.04	71,76,82,84	0
2	MAN	P	4	11/12	0.99	0.04	70,74,75,76	0
2	NAG	P	5	14/15	0.99	0.04	70,73,75,76	0
2	MAN	P	6	11/12	0.99	0.04	74,80,81,81	0
3	NAG	J	1	15/15	0.99	0.04	80,94,97,97	0
3	NAG	J	4	14/15	0.99	0.05	50,59,63,63	0
3	MAN	J	5	11/12	0.99	0.03	66,70,72,73	0
3	MAN	L	3	11/12	0.99	0.04	58,63,75,77	0
3	NAG	L	4	14/15	0.99	0.04	44,54,56,57	0
3	MAN	L	5	11/12	0.99	0.05	81,83,88,92	0
3	NAG	N	1	15/15	0.99	0.07	78,88,91,92	0
3	NAG	N	4	14/15	0.99	0.04	63,66,69,71	0
3	MAN	N	5	11/12	0.99	0.04	62,68,75,76	0
4	NAG	M	1	15/15	0.99	0.05	74,83,88,91	0
4	BMA	M	2	11/12	0.99	0.04	53,59,67,67	0
4	NAG	M	4	14/15	0.99	0.05	63,69,72,74	0
5	NAG	O	1	15/15	0.99	0.05	88,101,107,108	0
5	BMA	O	3	11/12	0.99	0.04	70,79,80,82	0
5	MAN	O	4	11/12	0.99	0.04	66,69,72,72	0
5	NAG	O	5	14/15	0.99	0.04	48,58,64,67	0
3	MAN	N	3	11/12	1.00	0.03	67,71,74,75	0
2	NAG	I	1	14/15	1.00	0.03	54,60,62,63	0
2	NAG	P	1	15/15	1.00	0.04	61,67,74,77	0
3	NAG	L	1	15/15	1.00	0.04	80,88,89,89	0
3	BMA	L	2	11/12	1.00	0.04	66,69,74,78	0
4	MAN	M	3	11/12	1.00	0.03	66,70,84,87	0
2	BMA	K	3	11/12	1.00	0.03	61,62,69,75	0
2	BMA	P	3	11/12	1.00	0.03	67,70,72,75	0
5	NAG	O	2	14/15	1.00	0.03	73,85,89,92	0
3	BMA	J	2	11/12	1.00	0.03	61,66,70,72	0
3	MAN	J	3	11/12	1.00	0.03	61,67,73,75	0
3	BMA	N	2	11/12	1.00	0.03	56,65,69,70	0

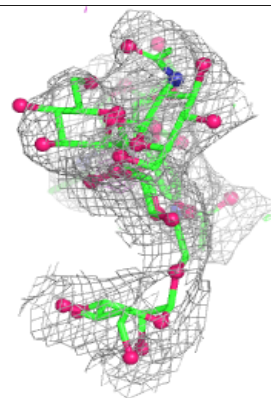
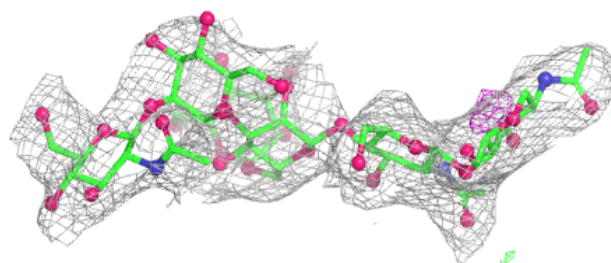
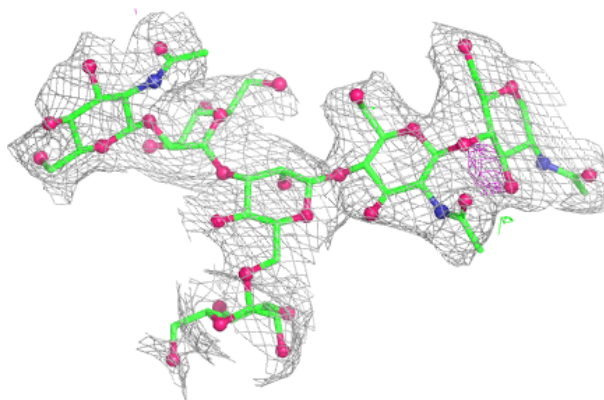
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain I:

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

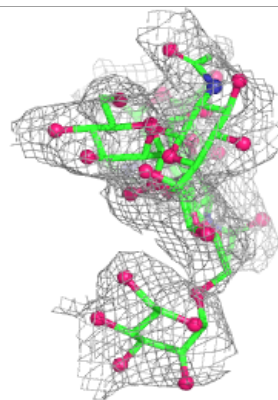
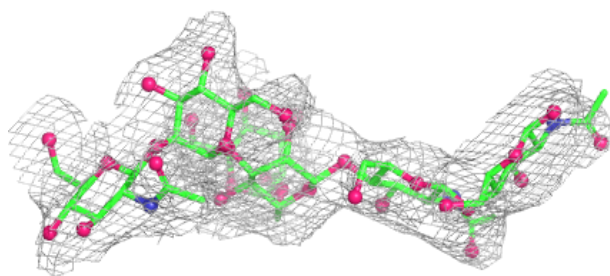
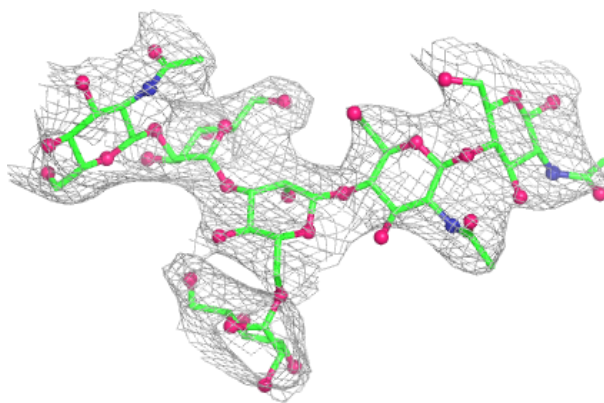
**Electron density around Chain K:**

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

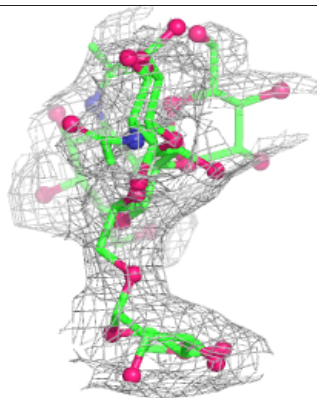
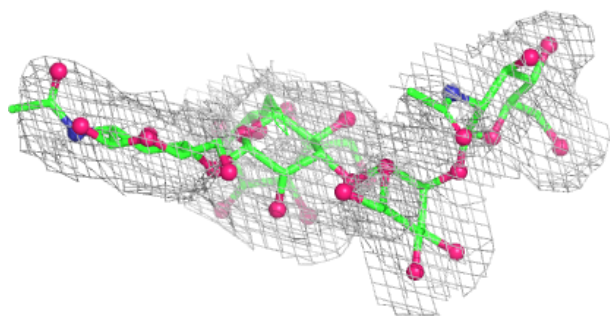
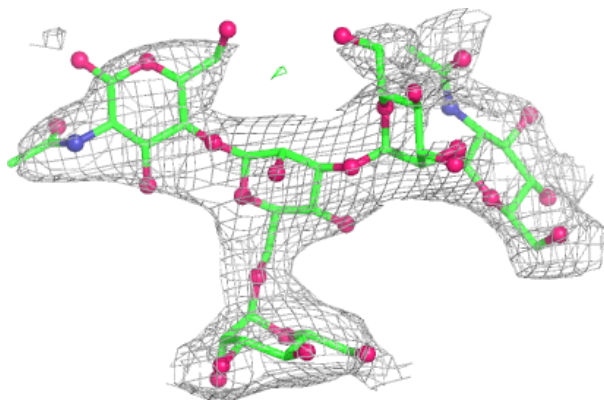


Electron density around Chain P:

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

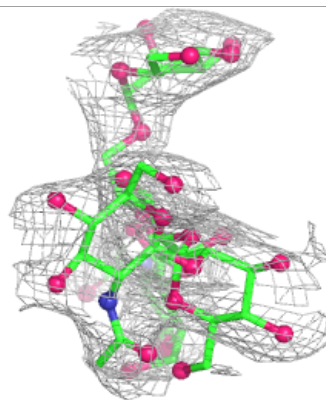
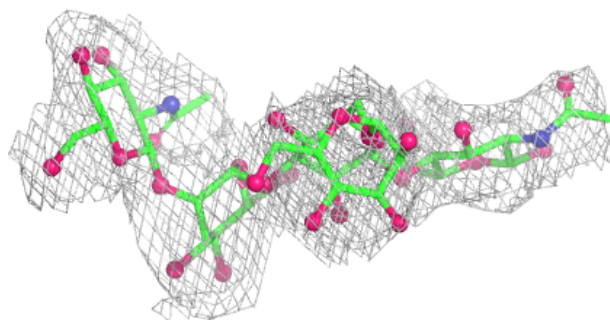
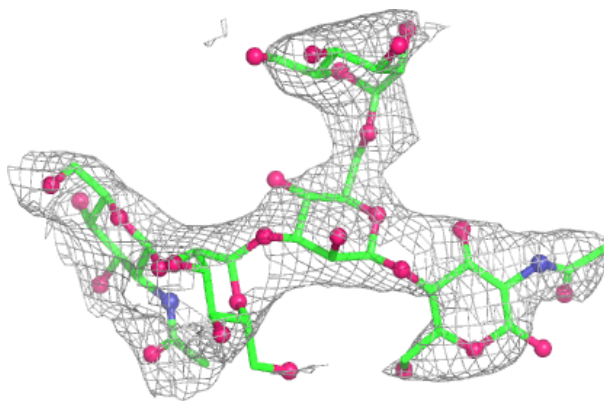
**Electron density around Chain J:**

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

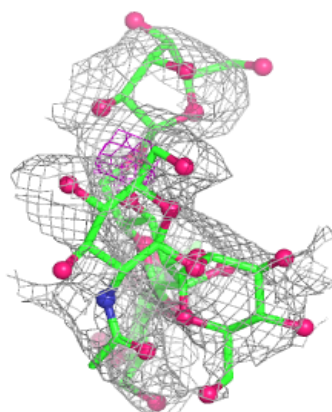
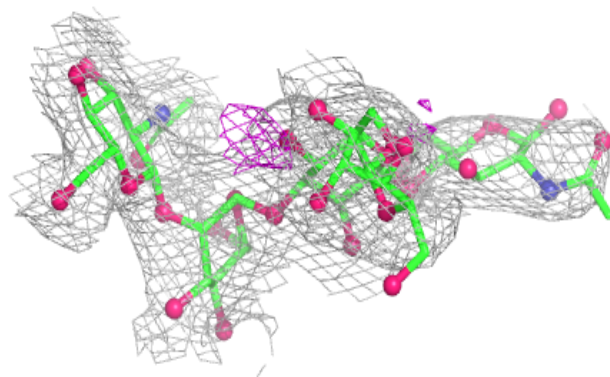
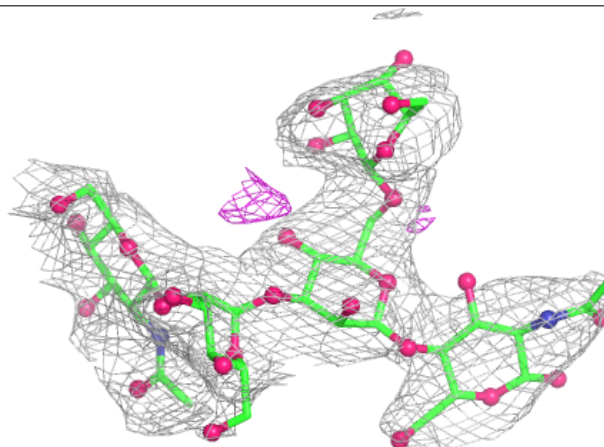


Electron density around Chain L:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

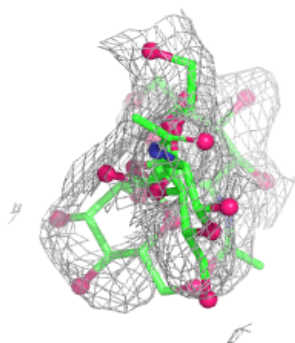
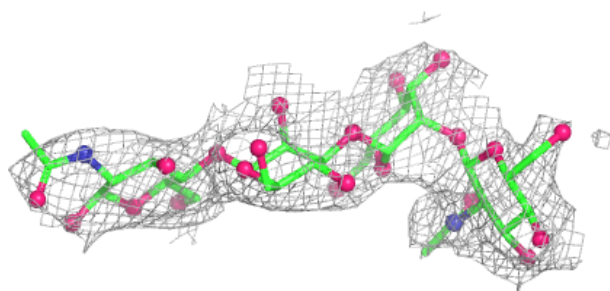
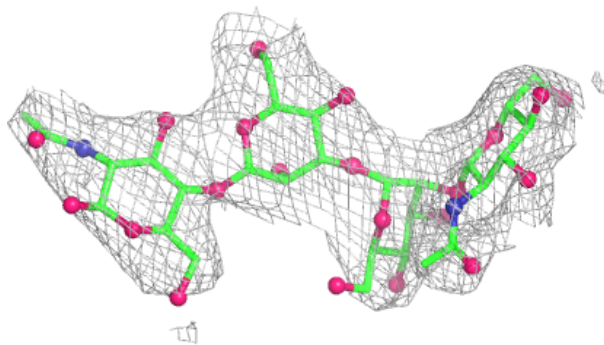
**Electron density around Chain N:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

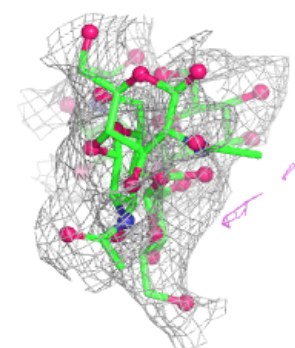
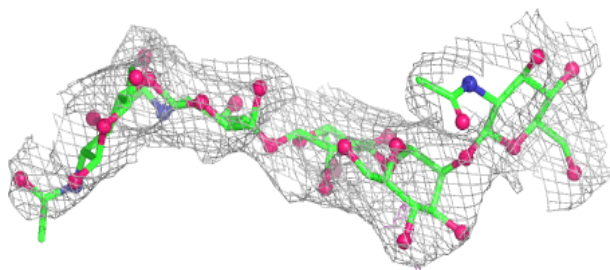
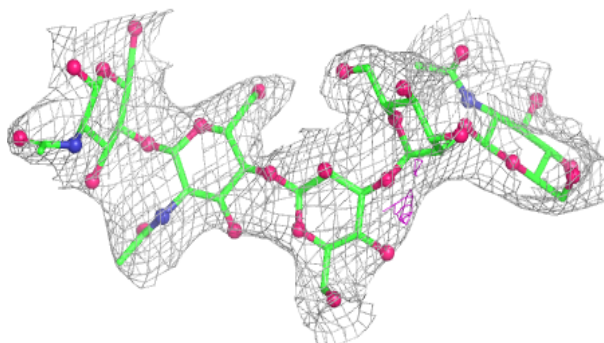


Electron density around Chain M:

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

**Electron density around Chain O:**

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

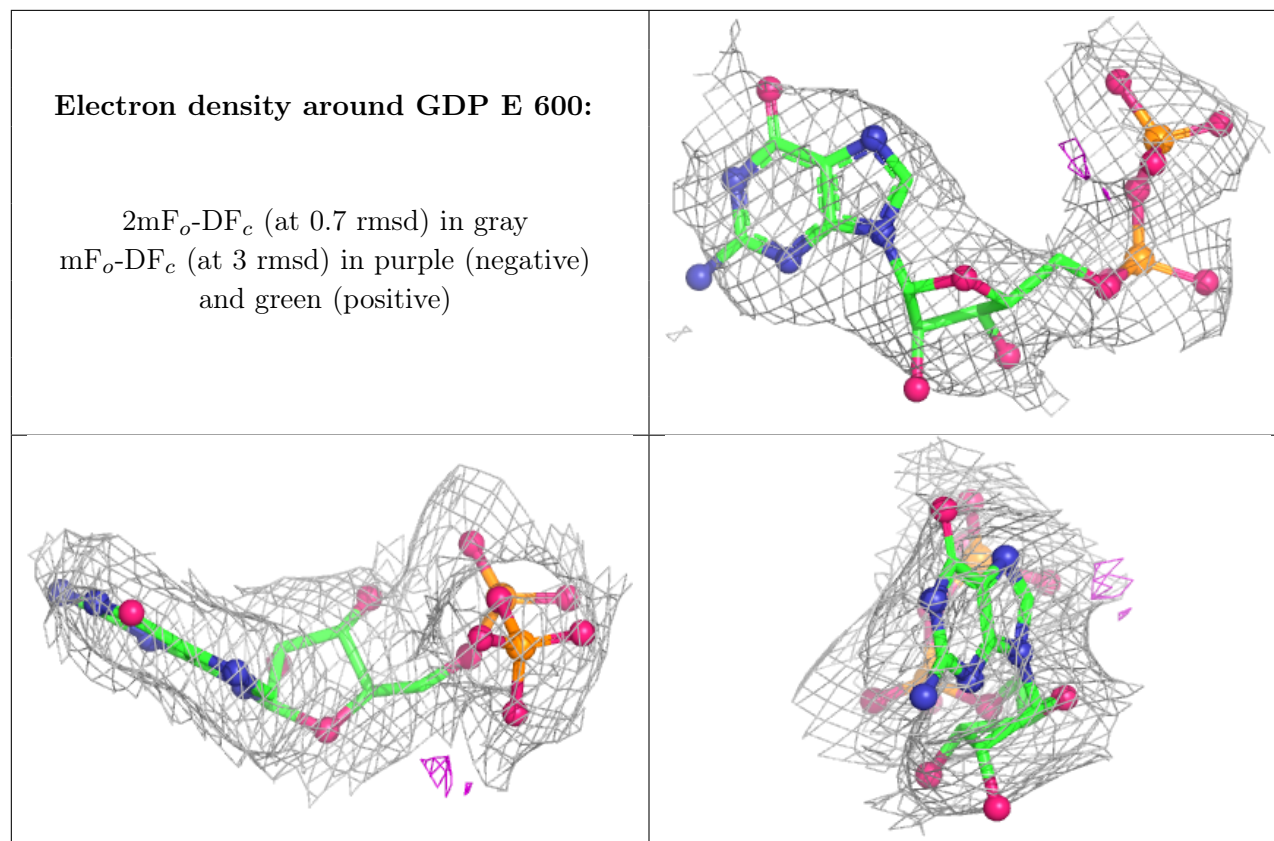


6.4 Ligands ⓘ

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, 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.

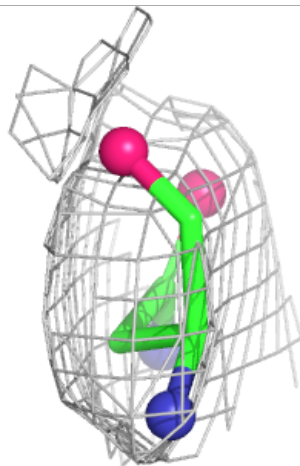
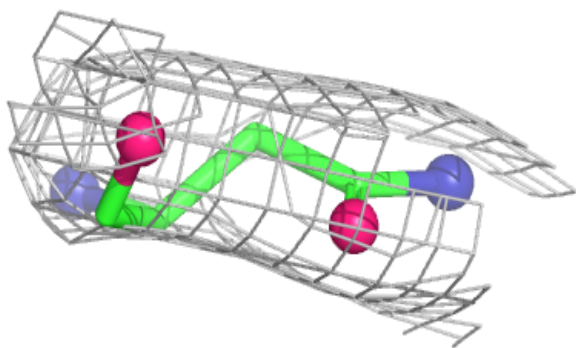
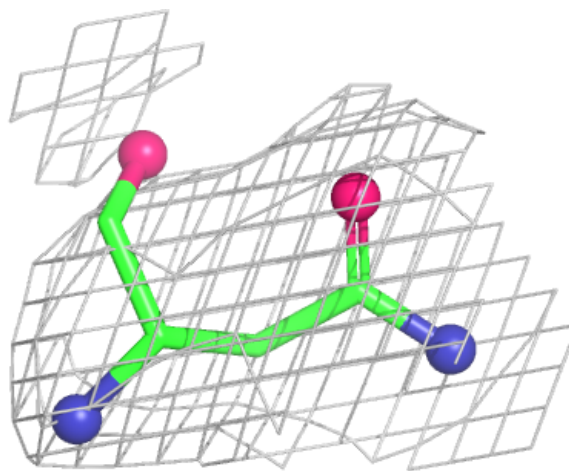
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	GDP	E	600	28/28	0.99	0.04	53,65,68,69	0
7	ASN	C	700	8/9	0.99	0.04	59,66,75,77	0
6	GDP	C	600	28/28	1.00	0.03	50,61,67,69	0
6	GDP	D	600	28/28	1.00	0.03	51,57,61,64	0
6	GDP	A	600	28/28	1.00	0.03	54,61,64,66	0
6	GDP	F	600	28/28	1.00	0.03	40,62,71,72	0
6	GDP	G	600	28/28	1.00	0.03	57,64,67,69	0
6	GDP	H	600	28/28	1.00	0.03	68,74,79,80	0
7	ASN	A	700	8/9	1.00	0.04	64,68,78,81	0
6	GDP	B	600	28/28	1.00	0.03	48,58,63,64	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.



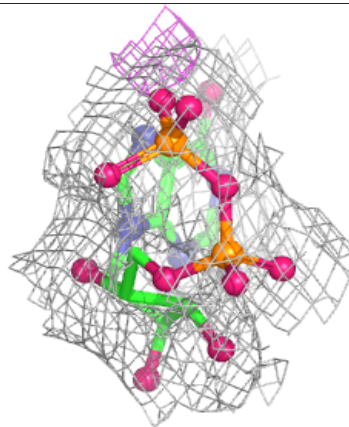
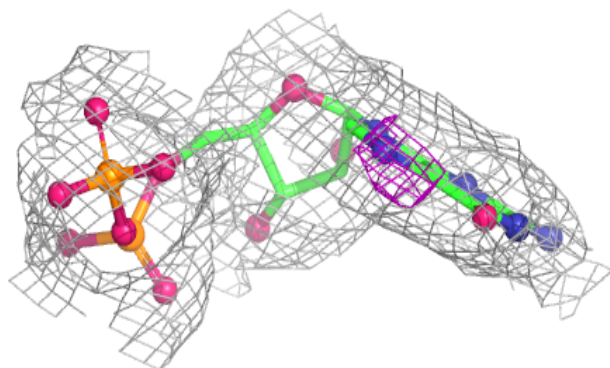
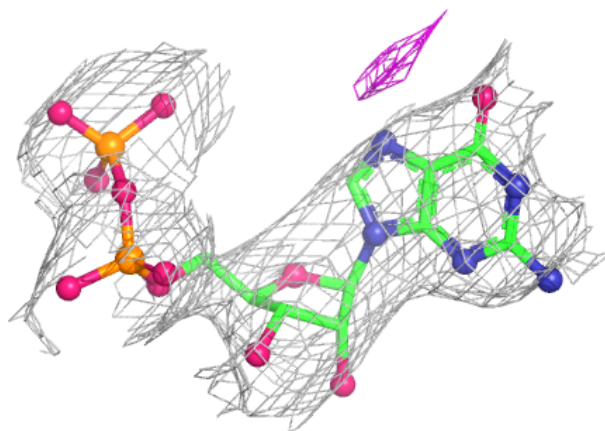
Electron density around ASN C 700:

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



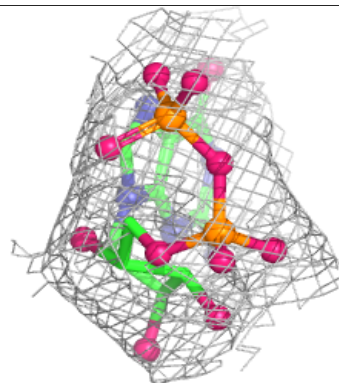
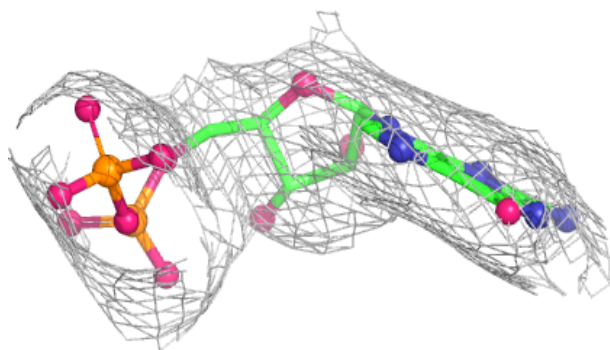
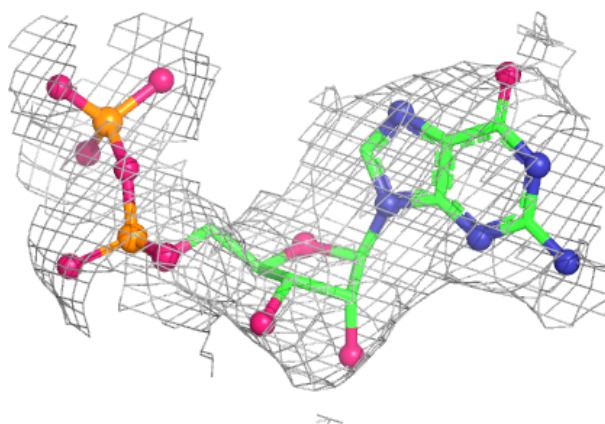
Electron density around GDP C 600:

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

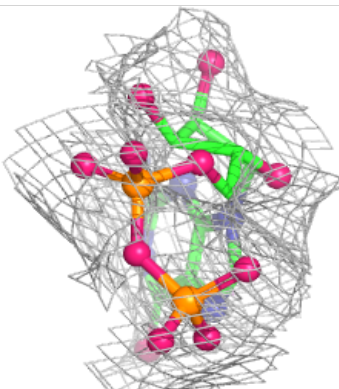
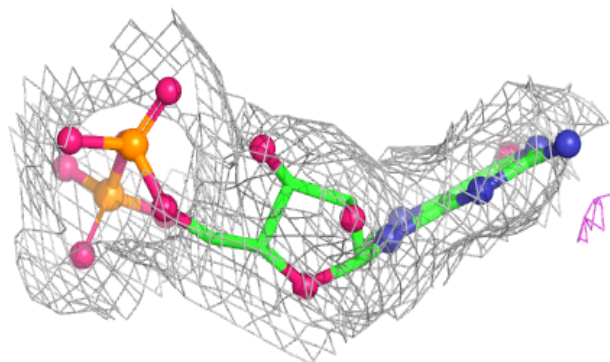
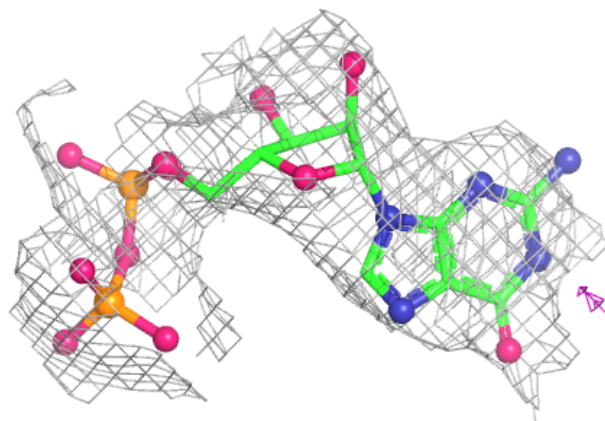


Electron density around GDP D 600:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

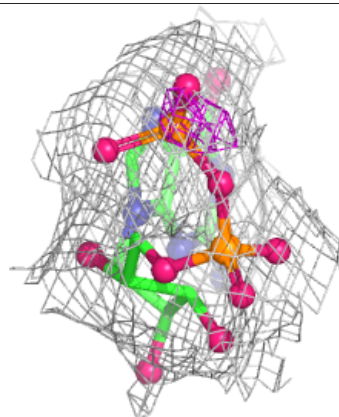
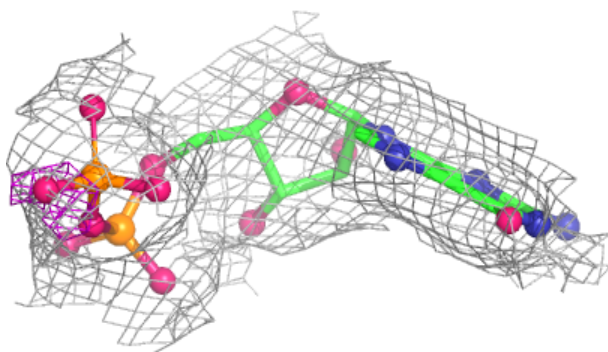
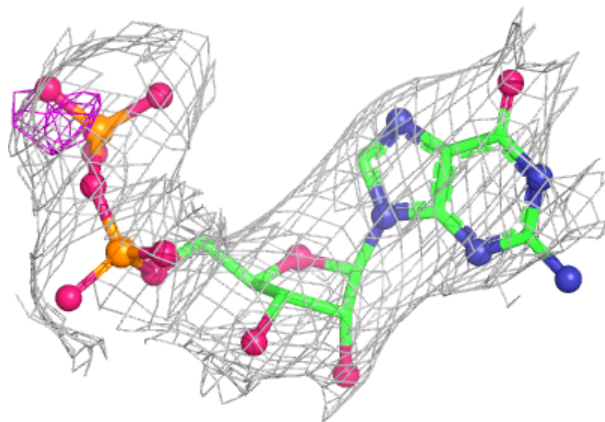
**Electron density around GDP A 600:**

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



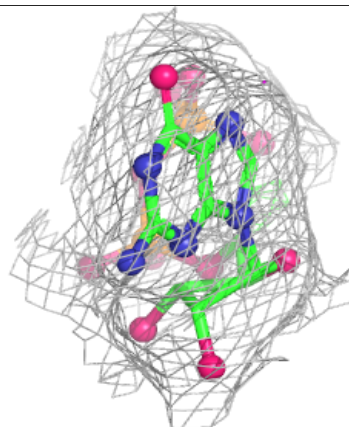
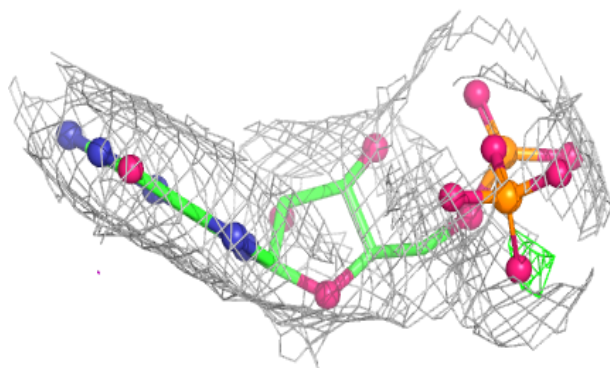
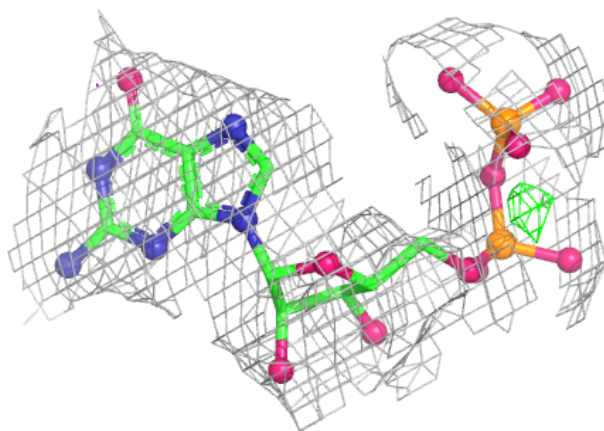
Electron density around GDP F 600:

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



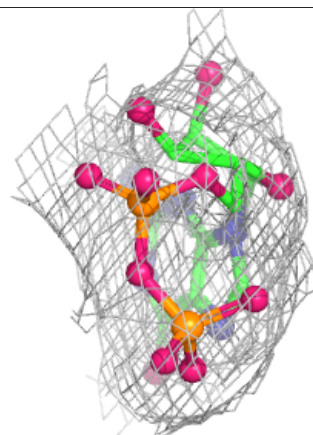
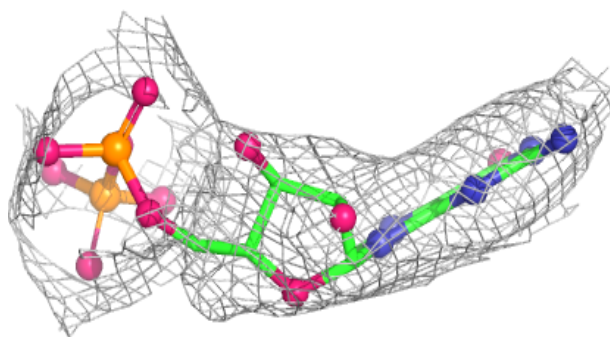
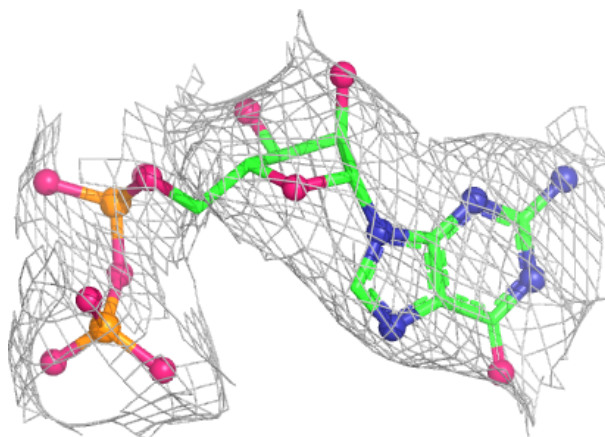
Electron density around GDP G 600:

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



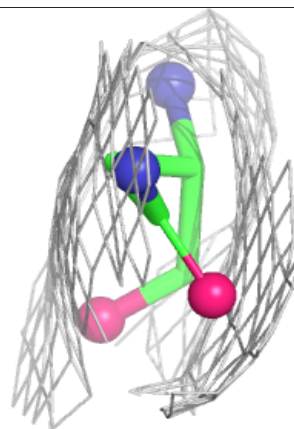
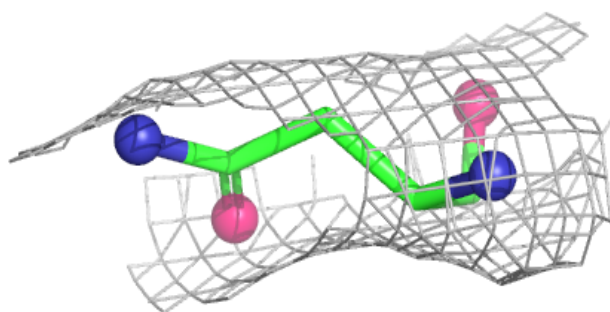
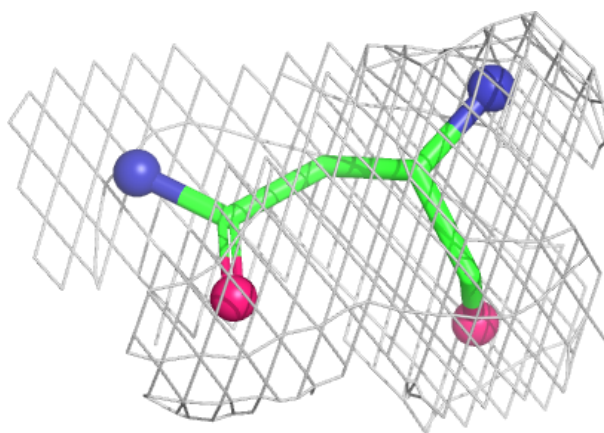
Electron density around GDP H 600:

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

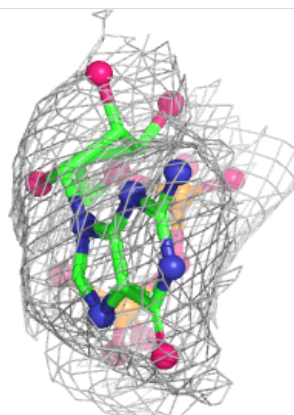
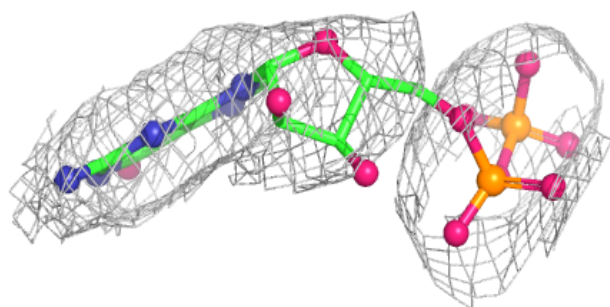
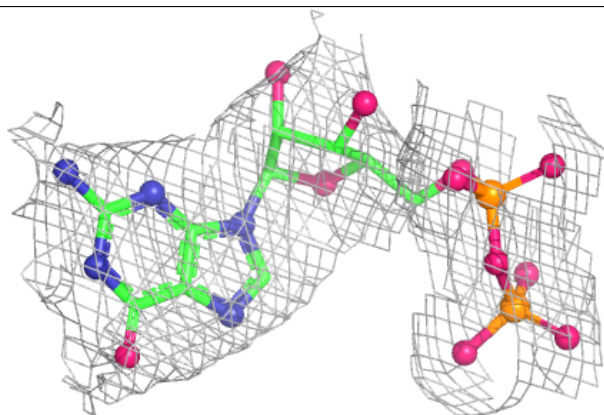


Electron density around ASN A 700:

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

**Electron density around GDP B 600:**

$2mF_o-DF_c$ (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.