



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:58 PM UTC

PDB ID : 5AM9 / pdb_00005am9
Title : Crystal structure of the Angiotensin-1 converting enzyme N-domain in complex with amyloid-beta 10-16
Authors : Masuyer, G.; Larmuth, K.M.; Douglas, R.G.; Sturrock, E.D.; Acharya, K.R.
Deposited on : 2015-03-10
Resolution : 1.80 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

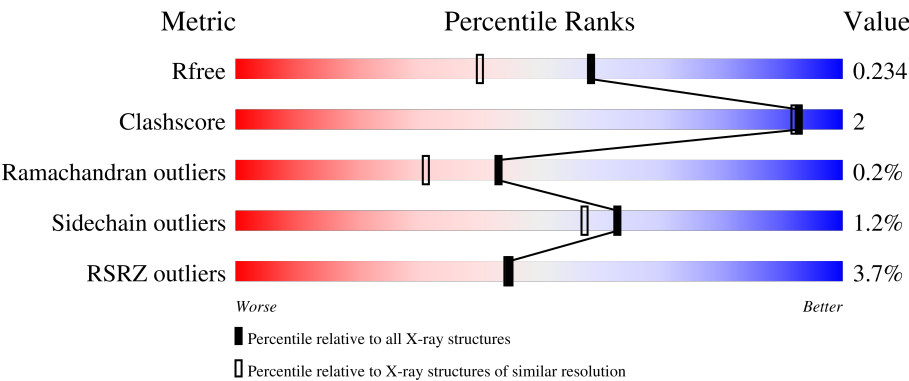
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	629	2% 92%
1	B	629	% 91% 5%
1	C	629	2% 92%
1	D	629	8% 92%
2	E	2	100%

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Mol	Chain	Length	Quality of chain
3	F	2	 100%
3	I	2	 100%
3	L	2	 100%
3	O	2	 100%
4	G	3	 67% 33%
4	J	3	 100%
5	H	2	 100%
5	K	2	 100%
5	N	2	 50% 50%
6	M	4	 25% 75%
6	P	4	 75% 25%

2 Entry composition

There are 17 unique types of molecules in this entry. The entry contains 22178 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 ANGIOTENSIN-CONVERTING ENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	603	Total	C	N	O	S	0	1	0
			4924	3164	844	897	19			
1	B	605	Total	C	N	O	S	0	3	0
			4962	3188	855	900	19			
1	C	611	Total	C	N	O	S	0	3	0
			5006	3214	858	914	20			
1	D	609	Total	C	N	O	S	0	3	0
			4995	3207	859	910	19			

There are 35 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLN	ASN	engineered mutation	UNP P12821
A	25	GLN	ASN	engineered mutation	UNP P12821
A	82	GLN	ASN	engineered mutation	UNP P12821
A	117	GLN	ASN	engineered mutation	UNP P12821
A	131	GLN	ASN	engineered mutation	UNP P12821
A	289	GLN	ASN	engineered mutation	UNP P12821
A	545	ARG	GLN	engineered mutation	UNP P12821
A	576	LEU	PRO	engineered mutation	UNP P12821
A	629	LEU	ARG	engineered mutation	UNP P12821
B	9	GLN	ASN	engineered mutation	UNP P12821
B	25	GLN	ASN	engineered mutation	UNP P12821
B	82	GLN	ASN	engineered mutation	UNP P12821
B	117	GLN	ASN	engineered mutation	UNP P12821
B	289	GLN	ASN	engineered mutation	UNP P12821
B	545	ARG	GLN	engineered mutation	UNP P12821
B	576	LEU	PRO	engineered mutation	UNP P12821
B	629	LEU	ARG	engineered mutation	UNP P12821
C	9	GLN	ASN	engineered mutation	UNP P12821
C	25	GLN	ASN	engineered mutation	UNP P12821
C	82	GLN	ASN	engineered mutation	UNP P12821
C	117	GLN	ASN	engineered mutation	UNP P12821

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Chain	Residue	Modelled	Actual	Comment	Reference
C	131	GLN	ASN	engineered mutation	UNP P12821
C	289	GLN	ASN	engineered mutation	UNP P12821
C	545	ARG	GLN	engineered mutation	UNP P12821
C	576	LEU	PRO	engineered mutation	UNP P12821
C	629	LEU	ARG	engineered mutation	UNP P12821
D	9	GLN	ASN	engineered mutation	UNP P12821
D	25	GLN	ASN	engineered mutation	UNP P12821
D	82	GLN	ASN	engineered mutation	UNP P12821
D	117	GLN	ASN	engineered mutation	UNP P12821
D	131	GLN	ASN	engineered mutation	UNP P12821
D	289	GLN	ASN	engineered mutation	UNP P12821
D	545	ARG	GLN	engineered mutation	UNP P12821
D	576	LEU	PRO	engineered mutation	UNP P12821
D	629	LEU	ARG	engineered mutation	UNP P12821

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



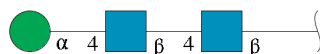
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	0	0	0
			24	14	1	9			

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



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

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



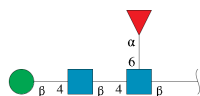
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	J	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.



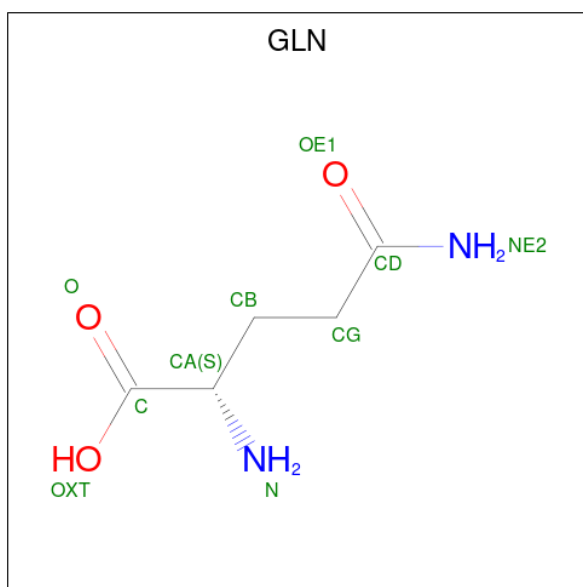
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	H	2	Total	C	N	O	0	0	0
			24	14	1	9			
5	K	2	Total	C	N	O	0	0	0
			24	14	1	9			
5	N	2	Total	C	N	O	0	0	0
			24	14	1	9			

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



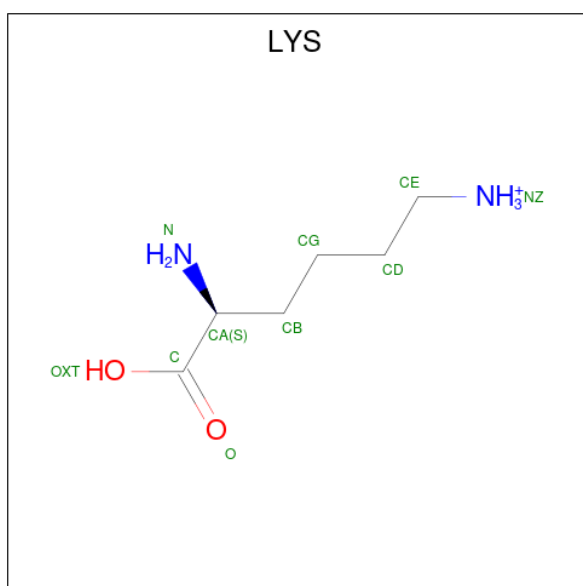
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	M	4	Total	C	N	O	0	0	0
			49	28	2	19			
6	P	4	Total	C	N	O	0	0	0
			49	28	2	19			

- Molecule 7 is GLUTAMINE (CCD ID: GLN) (formula: C₅H₁₀N₂O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			9	5	2	2		
7	D	1	Total	C	N	O	0	0
			9	5	2	2		

- Molecule 8 is LYSINE (CCD ID: LYS) (formula: $C_6H_{15}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			10	6	2	2		
8	D	1	Total	C	N	O	0	0
			10	6	2	2		

- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1	Total 1	Zn 1	0	0
9	B	1	Total 1	Zn 1	0	0
9	C	1	Total 1	Zn 1	0	0
9	D	1	Total 1	Zn 1	0	0

- Molecule 10 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	1	Total 1	Cl 1	0	0
10	B	1	Total 1	Cl 1	0	0
10	C	1	Total 1	Cl 1	0	0
10	D	1	Total 1	Cl 1	0	0

- Molecule 11 is SODIUM ION (CCD ID: NA) (formula: Na).

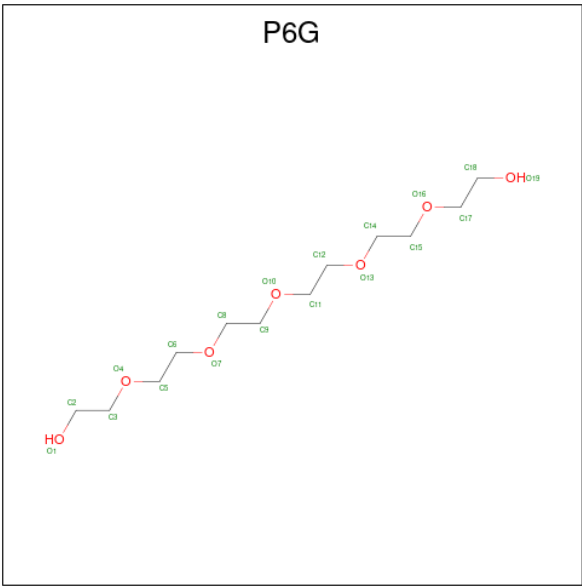
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	1	Total 1	Na 1	0	0
11	D	1	Total 1	Na 1	0	0

- Molecule 12 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



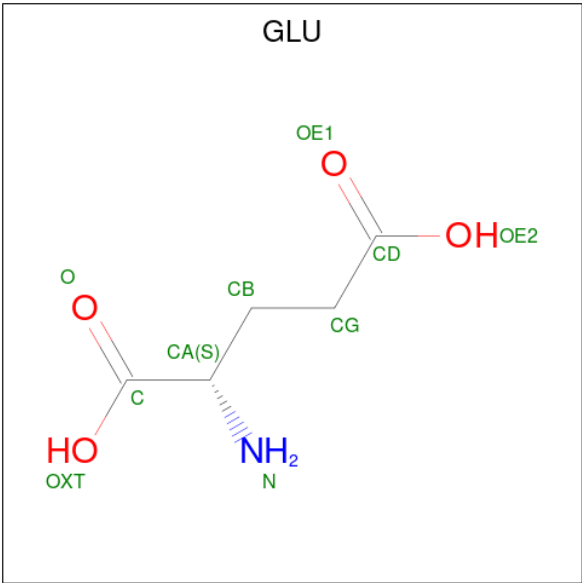
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	A	1	Total	C	O	0	0
			7	4	3		
12	A	1	Total	C	O	0	0
			7	4	3		
12	A	1	Total	C	O	0	0
			7	4	3		
12	B	1	Total	C	O	0	0
			7	4	3		
12	B	1	Total	C	O	0	0
			7	4	3		
12	B	1	Total	C	O	0	0
			7	4	3		
12	C	1	Total	C	O	0	0
			7	4	3		
12	C	1	Total	C	O	0	0
			7	4	3		
12	C	1	Total	C	O	0	0
			7	4	3		
12	C	1	Total	C	O	0	0
			7	4	3		
12	D	1	Total	C	O	0	0
			7	4	3		
12	D	1	Total	C	O	0	0
			7	4	3		
12	D	1	Total	C	O	0	0
			7	4	3		
12	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 13 is HEXAETHYLENE GLYCOL (CCD ID: P6G) (formula: C₁₂H₂₆O₇).



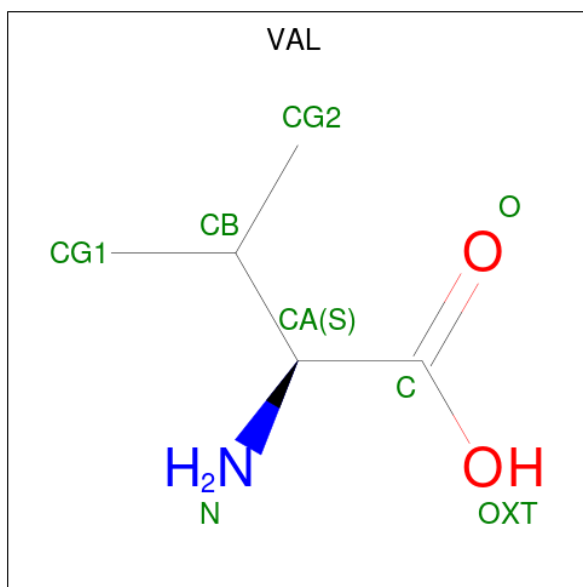
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	A	1	Total	C	O	0	0
			19	12	7		
13	B	1	Total	C	O	0	0
			19	12	7		
13	B	1	Total	C	O	0	0
			19	12	7		
13	D	1	Total	C	O	0	0
			19	12	7		

- Molecule 14 is GLUTAMIC ACID (CCD ID: GLU) (formula: C₅H₉NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
14	B	1	Total	C	N	O	0	0
			9	5	1	3		
14	C	1	Total	C	N	O	0	0
			9	5	1	3		

- Molecule 15 is VALINE (CCD ID: VAL) (formula: C₅H₁₁NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
15	B	1	Total	C	N	O	0	0
			8	5	1	2		
15	C	1	Total	C	N	O	0	0
			8	5	1	2		

- Molecule 16 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	B	1	Total	Ca	0	0
			1	1		
16	C	1	Total	Ca	0	0
			1	1		

- Molecule 17 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	A	367	Total	O	0	0
			367	367		

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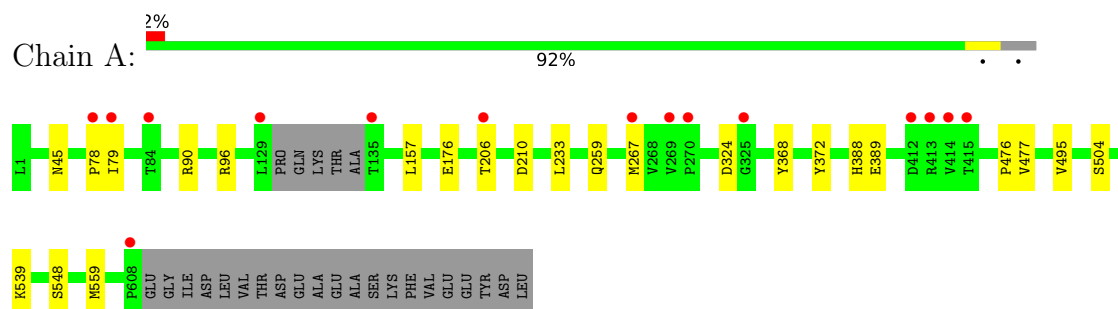
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	A	3	Total 3	O 3	0	0
17	B	407	Total 407	O 407	0	0
17	B	1	Total 1	O 1	0	0
17	C	465	Total 465	O 465	0	0
17	C	2	Total 2	O 2	0	0
17	D	403	Total 403	O 403	0	0
17	D	1	Total 1	O 1	0	0

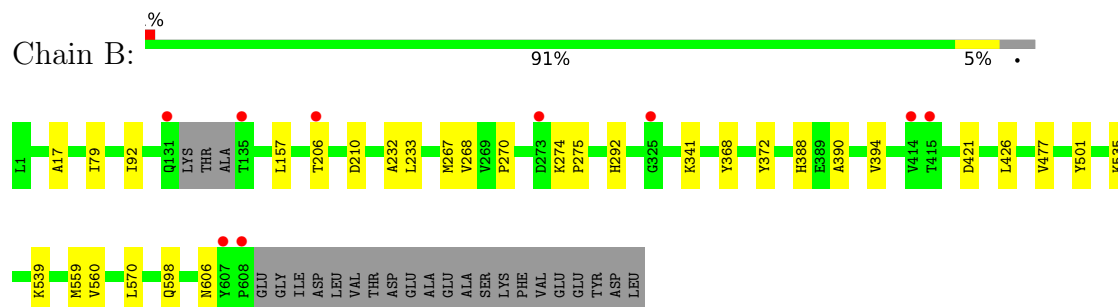
3 Residue-property plots [i](#)

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

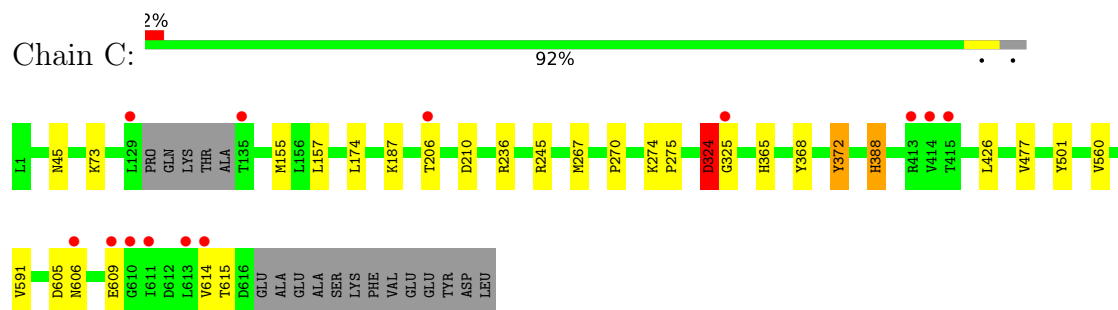
• Molecule 1: ANGIOTENSIN-CONVERTING ENZYME



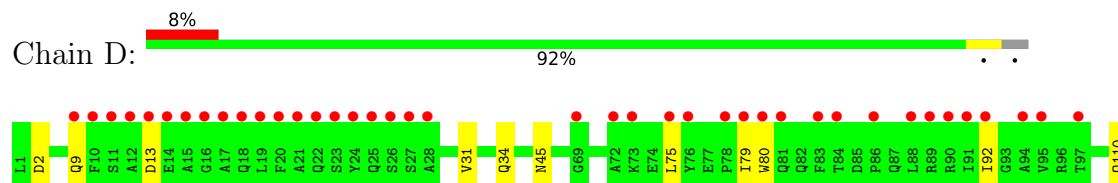
• Molecule 1: ANGIOTENSIN-CONVERTING ENZYME

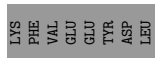


• Molecule 1: ANGIOTENSIN-CONVERTING ENZYME



• Molecule 1: ANGIOTENSIN-CONVERTING ENZYME





- NAG1
FUL2

- NAG1
NAG2

- NAG1
NAG2

- NAG1
NAG2

- NAG1
NAG2

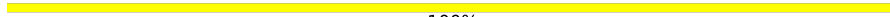
- NAG1
NAG2
MAN3

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

Chain J:  100%

MAG1
MAG2
MAN3

- Molecule 5: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  100%

MAG1
FUC2

- Molecule 5: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  100%

MAG1
FUC2

- Molecule 5: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  50% 50%

MAG1
FUC2

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

Chain M:  25% 75%

MAG1
MAG2
BMA3
FUC4

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

Chain P:  75% 25%

MAG1
MAG2
BMA3
FUC4

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	73.35Å 101.80Å 113.95Å 85.04° 85.55° 81.88°	Depositor
Resolution (Å)	113.27 – 1.80 113.27 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.0 (113.27-1.80) 97.0 (113.27-1.80)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.197 , 0.229 0.204 , 0.234	Depositor DCC
R_{free} test set	14480 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.439	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	22178	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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 36.56 % 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 4.9036e-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: P6G, NA, PEG, CA, CL, FUL, MAN, NAG, BMA, FUC, ZN

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	0.73	0/5083	0.87	0/6926
1	B	0.74	0/5124	0.85	1/6978 (0.0%)
1	C	0.78	0/5167	0.87	2/7036 (0.0%)
1	D	0.78	0/5157	0.86	1/7023 (0.0%)
All	All	0.76	0/20531	0.86	4/27963 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	324	ASP	N-CA-C	7.29	121.09	112.93
1	C	605	ASP	N-CA-C	6.28	118.68	111.02
1	B	79	ILE	N-CA-C	5.61	119.03	113.53
1	D	239	ASP	N-CA-C	5.38	117.90	111.71

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	324	ASP	Peptide

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	4924	0	4688	11	0
1	B	4962	0	4737	13	0
1	C	5006	0	4780	21	0
1	D	4995	0	4775	13	0
2	E	24	0	22	0	0
3	F	28	0	25	0	0
3	I	28	0	25	0	0
3	L	28	0	25	0	0
3	O	28	0	25	0	0
4	G	39	0	34	0	0
4	J	39	0	34	0	0
5	H	24	0	22	0	0
5	K	24	0	22	0	0
5	N	24	0	22	0	0
6	M	49	0	43	0	0
6	P	49	0	43	0	0
7	A	9	0	7	0	0
7	D	9	0	7	1	0
8	A	10	0	13	2	0
8	D	10	0	13	0	0
9	A	1	0	0	0	0
9	B	1	0	0	0	0
9	C	1	0	0	0	0
9	D	1	0	0	0	0
10	A	1	0	0	0	0
10	B	1	0	0	0	0
10	C	1	0	0	0	0
10	D	1	0	0	0	0
11	A	1	0	0	0	0
11	D	1	0	0	0	0
12	A	21	0	30	0	0
12	B	21	0	30	0	0
12	C	28	0	40	2	0
12	D	28	0	40	0	0
13	A	19	0	26	0	0
13	B	38	0	52	1	0
13	D	19	0	26	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	B	9	0	5	0	0
14	C	9	0	5	0	0
15	B	8	0	9	1	0
15	C	8	0	9	1	0
16	B	1	0	0	0	0
16	C	1	0	0	0	0
17	A	370	0	0	3	0
17	B	408	0	0	0	0
17	C	467	0	0	3	0
17	D	404	0	0	2	0
All	All	22178	0	19634	61	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 2.

All (61) 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:C:155[A]:MET:HA	1:C:155[A]:MET:HE2	1.42	1.01
1:C:155[A]:MET:HE2	1:C:155[A]:MET:CA	2.06	0.74
1:C:365:HIS:HD1	1:C:388:HIS:CD2	2.08	0.71
1:C:206:THR:HG23	1:C:210:ASP:OD2	1.94	0.68
1:C:365:HIS:HD1	1:C:388:HIS:HD2	1.39	0.66
1:B:501:TYR:CD1	15:B:912:VAL:HG12	2.31	0.66
1:C:155[A]:MET:HA	1:C:155[A]:MET:CE	2.19	0.65
1:A:157:LEU:HD11	1:A:477:VAL:HG13	1.79	0.65
8:A:916:LYS:HE2	17:A:2293:HOH:O	1.97	0.64
1:A:206:THR:HG23	1:A:210:ASP:OD2	1.98	0.63
1:C:501:TYR:CD1	15:C:912:VAL:HG12	2.38	0.59
1:C:157:LEU:HD11	1:C:477:VAL:HG13	1.83	0.59
1:D:157:LEU:HD11	1:D:477:VAL:HG13	1.85	0.58
1:B:157:LEU:HD11	1:B:477:VAL:HG13	1.86	0.57
12:C:1201:PEG:H41	17:C:2379:HOH:O	2.04	0.56
1:B:206:THR:HG23	1:B:210:ASP:OD2	2.06	0.56
1:A:539:LYS:HE3	1:A:559:MET:O	2.07	0.54
1:B:270:PRO:HD3	1:B:426:LEU:HD22	1.89	0.53
1:D:176:GLU:OE1	17:D:2144:HOH:O	2.18	0.53
1:B:233:LEU:HD23	1:B:267:MET:HE1	1.91	0.52
1:B:17:ALA:HB1	1:B:92:ILE:HD11	1.92	0.52
1:C:155[A]:MET:HE1	1:C:615:THR:HG23	1.92	0.52
1:C:270:PRO:HD3	1:C:426:LEU:HD22	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:236:ARG:HD2	1:C:267:MET:HE3	1.93	0.51
1:D:331:HIS:CE1	7:D:915:GLN:HG3	2.46	0.50
1:A:233:LEU:HD23	1:A:267:MET:HE1	1.95	0.49
1:D:13:ASP:HA	1:D:79:ILE:HD13	1.95	0.48
1:D:268:VAL:HG12	1:D:426:LEU:HD11	1.95	0.48
1:C:426:LEU:O	1:C:426:LEU:HG	2.09	0.48
1:A:389:GLU:HB2	1:A:504:SER:HB2	1.96	0.47
1:B:390:ALA:O	1:B:394:VAL:HG23	2.14	0.47
1:C:155[A]:MET:CE	1:C:614:VAL:HG23	2.45	0.47
1:B:570:LEU:C	1:B:570:LEU:HD23	2.40	0.47
1:C:606:ASN:O	1:C:609:GLU:O	2.33	0.46
1:C:73:LYS:NZ	17:C:2086:HOH:O	2.44	0.46
1:A:90:ARG:NH2	17:A:2086:HOH:O	2.48	0.46
1:C:324:ASP:N	1:C:325:GLY:HA2	2.30	0.46
1:D:274:LYS:HB3	1:D:275:PRO:CD	2.46	0.45
1:C:274:LYS:HB3	1:C:275:PRO:CD	2.47	0.45
1:D:110:GLN:OE1	17:D:2077:HOH:O	2.21	0.45
1:D:228:PHE:CZ	13:D:1201:P6G:H142	2.51	0.45
1:D:324:ASP:OD1	1:D:326:ARG:HB2	2.16	0.45
1:B:292:HIS:CE1	13:B:1203:P6G:H51	2.52	0.45
1:A:495:VAL:HG12	1:A:495:VAL:O	2.18	0.44
1:B:232:ALA:CB	1:B:268:VAL:HG12	2.47	0.44
1:C:245:ARG:HG2	1:C:591:VAL:HG11	2.00	0.43
12:C:1201:PEG:C4	17:C:2379:HOH:O	2.65	0.43
1:D:31:VAL:O	1:D:34:GLN:HG3	2.19	0.43
1:B:274:LYS:HB3	1:B:275:PRO:HD2	2.00	0.43
1:A:259:GLN:HE21	8:A:916:LYS:NZ	2.17	0.43
1:D:268:VAL:CG1	1:D:426:LEU:HD11	2.49	0.42
1:B:539:LYS:HE3	1:B:559:MET:O	2.19	0.42
1:C:245:ARG:HG2	1:C:591:VAL:CG1	2.50	0.42
1:A:176:GLU:OE1	17:A:2156:HOH:O	2.21	0.42
1:A:157:LEU:HD13	1:A:476:PRO:HB2	2.02	0.42
1:C:236:ARG:CD	1:C:267:MET:HE3	2.50	0.41
1:A:90:ARG:NE	1:A:548:SER:O	2.54	0.41
1:B:270:PRO:HD3	1:B:426:LEU:CD2	2.50	0.41
1:D:9[A]:GLN:NE2	1:D:75:LEU:HD21	2.35	0.41
1:C:372:TYR:OH	1:C:388:HIS:HE1	2.04	0.41
1:D:80:TRP:HB2	1:D:92:ILE:HG21	2.03	0.41

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	600/629 (95%)	589 (98%)	8 (1%)	3 (0%)	24	14
1	B	604/629 (96%)	594 (98%)	10 (2%)	0	100	100
1	C	610/629 (97%)	599 (98%)	10 (2%)	1 (0%)	43	31
1	D	608/629 (97%)	594 (98%)	13 (2%)	1 (0%)	43	31
All	All	2422/2516 (96%)	2376 (98%)	41 (2%)	5 (0%)	43	31

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	45	ASN
1	D	45	ASN
1	A	45	ASN
1	A	78	PRO
1	A	324	ASP

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	517/541 (96%)	512 (99%)	5 (1%)	68	64
1	B	521/541 (96%)	512 (98%)	9 (2%)	53	45
1	C	527/541 (97%)	521 (99%)	6 (1%)	65	60
1	D	526/541 (97%)	520 (99%)	6 (1%)	65	60
All	All	2091/2164 (97%)	2065 (99%)	26 (1%)	63	57

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	ILE
1	A	96	ARG
1	A	368	TYR
1	A	372	TYR
1	A	388	HIS
1	B	341	LYS
1	B	368	TYR
1	B	372	TYR
1	B	388	HIS
1	B	421	ASP
1	B	535	LYS
1	B	560	VAL
1	B	598	GLN
1	B	606	ASN
1	C	174	LEU
1	C	187	LYS
1	C	368	TYR
1	C	372	TYR
1	C	388	HIS
1	C	560	VAL
1	D	2	ASP
1	D	187	LYS
1	D	372	TYR
1	D	377	VAL
1	D	388	HIS
1	D	414	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	289	GLN
1	A	371	GLN
1	B	87	GLN
1	B	263	ASN
1	B	371	GLN
1	C	22	GLN
1	C	117	GLN
1	C	213	HIS
1	C	224	ASN
1	C	388	HIS

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Mol	Chain	Res	Type
1	C	568	GLN
1	C	582	GLN
1	C	585	ASN
1	D	18	GLN
1	D	22	GLN
1	D	82	GLN
1	D	109	GLN
1	D	110	GLN
1	D	224	ASN
1	D	263	ASN
1	D	585	ASN
1	D	586	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

30 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	E	1	2,1	14,14,15	0.46	0	17,19,21	1.06	1 (5%)
2	FUL	E	2	2	10,10,11	0.53	0	14,14,16	1.18	2 (14%)
3	NAG	F	1	1,3	14,14,15	0.65	0	17,19,21	0.99	1 (5%)
3	NAG	F	2	3	14,14,15	0.50	0	17,19,21	1.32	1 (5%)
4	NAG	G	1	1,4	14,14,15	0.42	0	17,19,21	0.78	0
4	NAG	G	2	4	14,14,15	0.40	0	17,19,21	0.87	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	G	3	4	11,11,12	0.57	0	15,15,17	0.98	2 (13%)
5	NAG	H	1	5,1	14,14,15	0.76	0	17,19,21	1.52	4 (23%)
5	FUC	H	2	5	10,10,11	0.49	0	14,14,16	1.40	1 (7%)
3	NAG	I	1	1,3	14,14,15	0.50	0	17,19,21	1.20	1 (5%)
3	NAG	I	2	3	14,14,15	0.62	0	17,19,21	1.20	2 (11%)
4	NAG	J	1	1,4	14,14,15	0.43	0	17,19,21	1.00	1 (5%)
4	NAG	J	2	4	14,14,15	0.50	0	17,19,21	1.14	2 (11%)
4	MAN	J	3	4	11,11,12	0.79	0	15,15,17	2.10	5 (33%)
5	NAG	K	1	5,1	14,14,15	0.69	0	17,19,21	1.63	3 (17%)
5	FUC	K	2	5	10,10,11	0.72	0	14,14,16	1.72	2 (14%)
3	NAG	L	1	1,3	14,14,15	0.45	0	17,19,21	1.34	2 (11%)
3	NAG	L	2	3	14,14,15	0.50	0	17,19,21	1.22	2 (11%)
6	NAG	M	1	6,1	14,14,15	0.55	0	17,19,21	1.02	0
6	NAG	M	2	6	14,14,15	0.48	0	17,19,21	1.19	1 (5%)
6	BMA	M	3	6	11,11,12	0.52	0	15,15,17	0.99	1 (6%)
6	FUC	M	4	6	10,10,11	0.57	0	14,14,16	1.18	1 (7%)
5	NAG	N	1	5,1	14,14,15	0.65	0	17,19,21	1.01	0
5	FUC	N	2	5	10,10,11	0.66	0	14,14,16	1.67	3 (21%)
3	NAG	O	1	1,3	14,14,15	0.50	0	17,19,21	1.04	1 (5%)
3	NAG	O	2	3	14,14,15	0.45	0	17,19,21	1.42	4 (23%)
6	NAG	P	1	6,1	14,14,15	0.69	0	17,19,21	0.95	0
6	NAG	P	2	6	14,14,15	0.37	0	17,19,21	0.94	1 (5%)
6	BMA	P	3	6	11,11,12	0.52	0	15,15,17	0.77	0
6	FUC	P	4	6	10,10,11	0.43	0	14,14,16	0.75	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
2	NAG	E	1	2,1	-	0/6/23/26	0/1/1/1
2	FUL	E	2	2	-	-	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
4	NAG	G	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
4	MAN	G	3	4	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	H	1	5,1	-	2/6/23/26	0/1/1/1
5	FUC	H	2	5	-	-	0/1/1/1
3	NAG	I	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	I	2	3	-	2/6/23/26	0/1/1/1
4	NAG	J	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1
4	MAN	J	3	4	-	0/2/19/22	0/1/1/1
5	NAG	K	1	5,1	-	2/6/23/26	0/1/1/1
5	FUC	K	2	5	-	-	0/1/1/1
3	NAG	L	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	L	2	3	-	0/6/23/26	0/1/1/1
6	NAG	M	1	6,1	-	0/6/23/26	0/1/1/1
6	NAG	M	2	6	-	0/6/23/26	0/1/1/1
6	BMA	M	3	6	-	0/2/19/22	0/1/1/1
6	FUC	M	4	6	-	-	0/1/1/1
5	NAG	N	1	5,1	-	2/6/23/26	0/1/1/1
5	FUC	N	2	5	-	-	0/1/1/1
3	NAG	O	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	O	2	3	-	1/6/23/26	0/1/1/1
6	NAG	P	1	6,1	-	0/6/23/26	0/1/1/1
6	NAG	P	2	6	-	0/6/23/26	0/1/1/1
6	BMA	P	3	6	-	0/2/19/22	0/1/1/1
6	FUC	P	4	6	-	-	0/1/1/1

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	K	2	FUC	C1-C2-C3	4.80	116.63	109.64
3	F	2	NAG	C1-O5-C5	4.34	118.00	112.19
5	N	2	FUC	C1-C2-C3	4.05	115.54	109.64
4	J	3	MAN	C1-O5-C5	3.97	117.51	112.19
3	L	1	NAG	O5-C1-C2	-3.75	105.49	111.29
4	J	3	MAN	C1-C2-C3	3.66	114.98	109.64
4	J	3	MAN	C3-C4-C5	3.61	116.78	110.23
5	K	1	NAG	C4-C3-C2	3.48	116.11	111.02
5	K	1	NAG	C1-C2-N2	-3.33	105.18	110.43
5	H	1	NAG	O7-C7-N2	-3.24	116.26	121.98
5	K	2	FUC	O5-C1-C2	3.19	118.41	110.79
4	J	3	MAN	O5-C5-C6	3.17	113.84	107.66
4	J	3	MAN	C2-C3-C4	3.15	116.40	110.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	P	2	NAG	C1-C2-N2	-3.14	105.49	110.43
6	M	2	NAG	C2-N2-C7	3.10	127.06	122.90
2	E	2	FUL	C1-O5-C5	3.07	120.20	112.97
3	I	1	NAG	C1-O5-C5	2.96	116.16	112.19
3	I	2	NAG	C4-C3-C2	2.95	115.35	111.02
3	O	2	NAG	C1-C2-N2	-2.94	105.81	110.43
5	H	1	NAG	C2-N2-C7	-2.86	119.06	122.90
6	M	4	FUC	O2-C2-C3	2.81	115.97	110.15
3	O	2	NAG	O5-C5-C4	-2.78	104.06	110.83
5	H	1	NAG	C8-C7-N2	2.78	120.73	116.12
5	H	2	FUC	C1-O5-C5	2.75	119.45	112.97
3	O	1	NAG	C1-O5-C5	2.63	115.72	112.19
3	L	1	NAG	C1-O5-C5	2.38	115.38	112.19
5	N	2	FUC	O5-C5-C6	2.32	112.43	107.40
4	G	3	MAN	C1-O5-C5	2.30	115.26	112.19
4	J	1	NAG	C2-N2-C7	2.25	125.92	122.90
3	O	2	NAG	O5-C5-C6	2.23	112.00	107.66
3	F	1	NAG	O5-C1-C2	-2.21	107.88	111.29
4	G	3	MAN	C1-C2-C3	2.18	112.81	109.64
6	M	3	BMA	C2-C3-C4	-2.13	107.12	110.86
3	O	2	NAG	C4-C3-C2	2.13	114.14	111.02
5	K	1	NAG	O3-C3-C4	-2.10	105.42	110.38
3	L	2	NAG	C2-N2-C7	2.10	125.71	122.90
4	J	2	NAG	C2-N2-C7	2.09	125.70	122.90
5	N	2	FUC	O3-C3-C4	-2.08	105.47	110.38
4	J	2	NAG	C1-C2-N2	-2.08	107.16	110.43
2	E	2	FUL	O5-C5-C4	2.08	113.29	109.55
3	I	2	NAG	C2-N2-C7	2.06	125.66	122.90
3	L	2	NAG	O7-C7-N2	2.04	125.58	121.98
5	H	1	NAG	O5-C5-C6	2.02	111.59	107.66
2	E	1	NAG	C8-C7-N2	2.00	119.44	116.12

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	K	1	NAG	O5-C5-C6-O6
5	N	1	NAG	O5-C5-C6-O6
5	H	1	NAG	O5-C5-C6-O6
4	G	3	MAN	C4-C5-C6-O6
5	H	1	NAG	C4-C5-C6-O6
5	K	1	NAG	C4-C5-C6-O6

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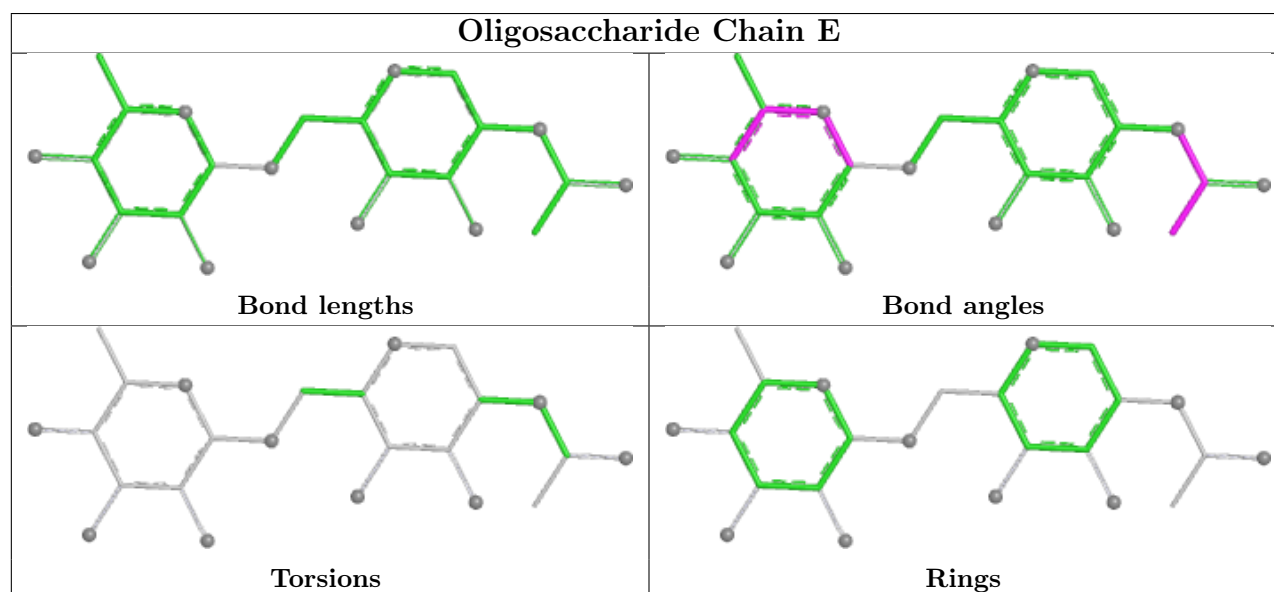
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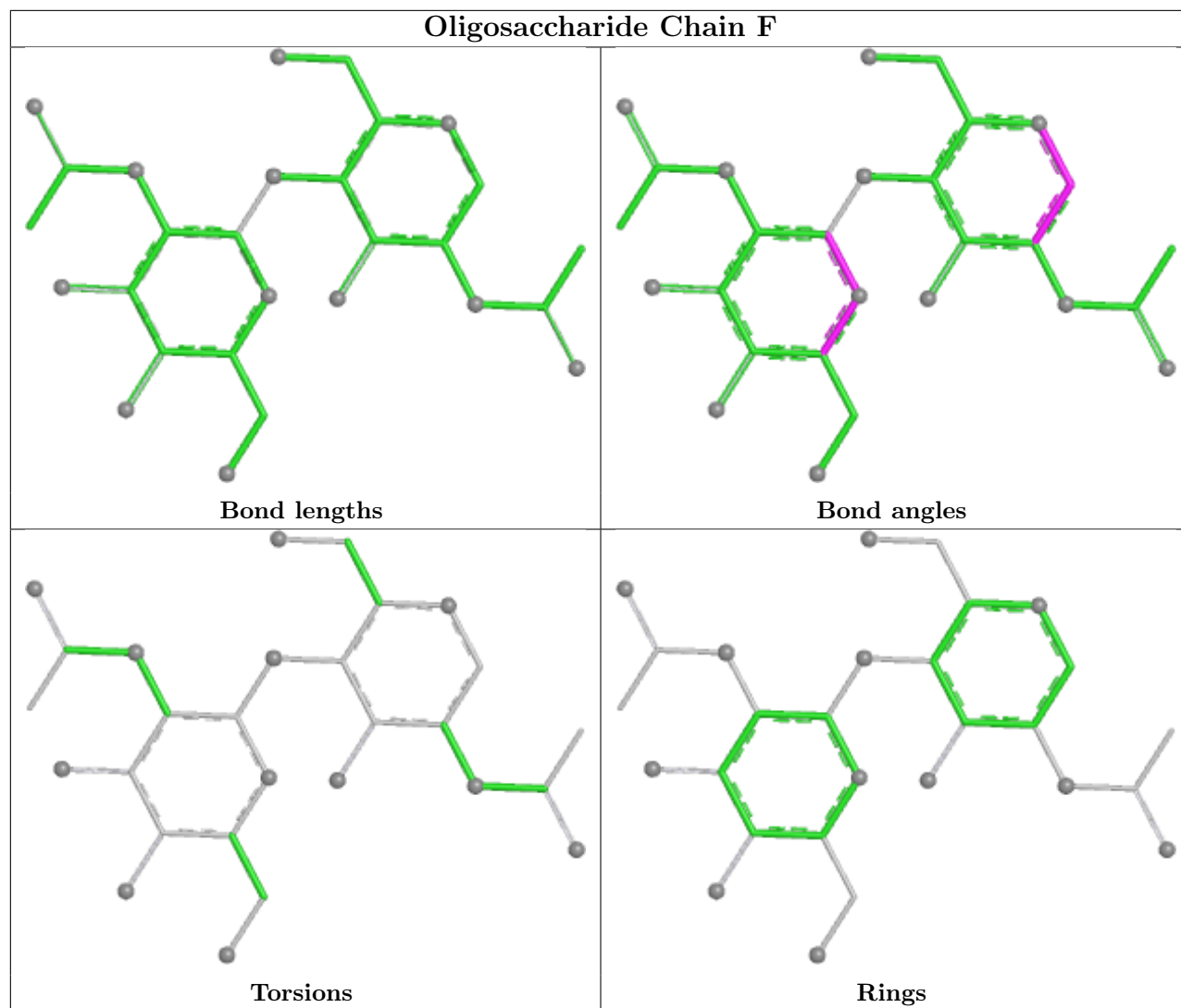
Mol	Chain	Res	Type	Atoms
4	G	3	MAN	O5-C5-C6-O6
5	N	1	NAG	C4-C5-C6-O6
3	I	2	NAG	C4-C5-C6-O6
3	I	2	NAG	O5-C5-C6-O6
3	I	1	NAG	C4-C5-C6-O6
3	O	2	NAG	C1-C2-N2-C7
3	I	1	NAG	O5-C5-C6-O6

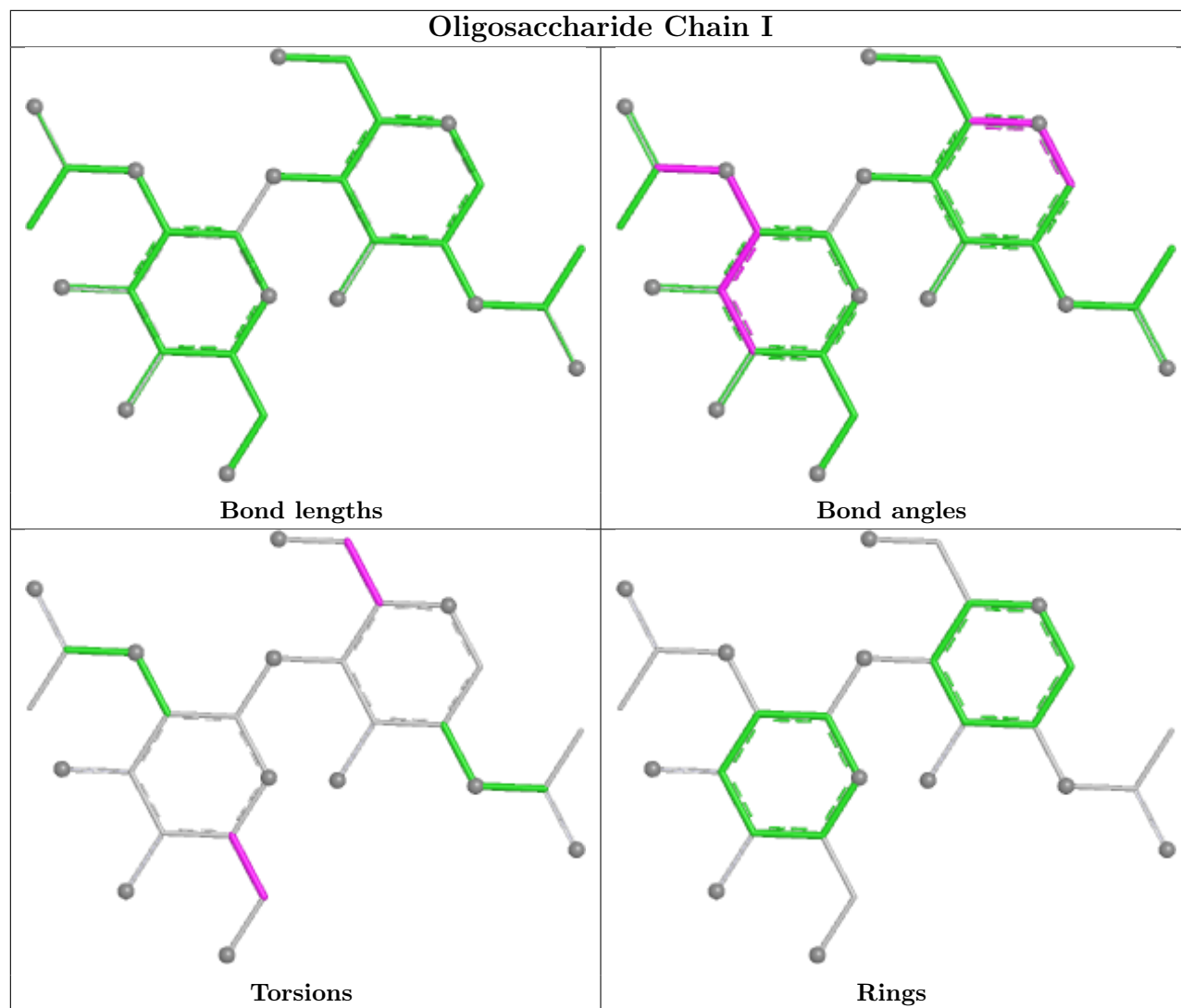
There are no ring outliers.

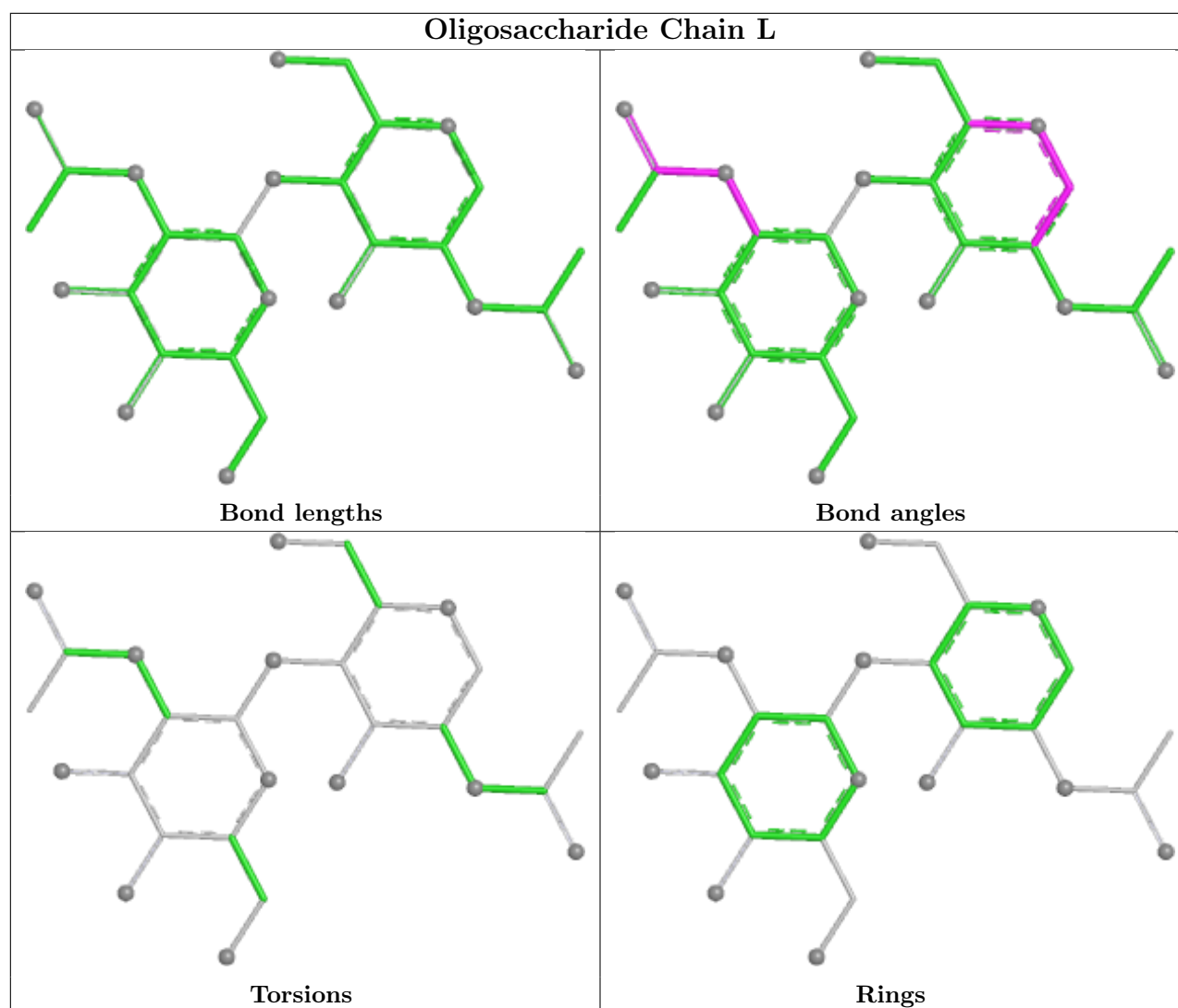
No monomer is involved in short contacts.

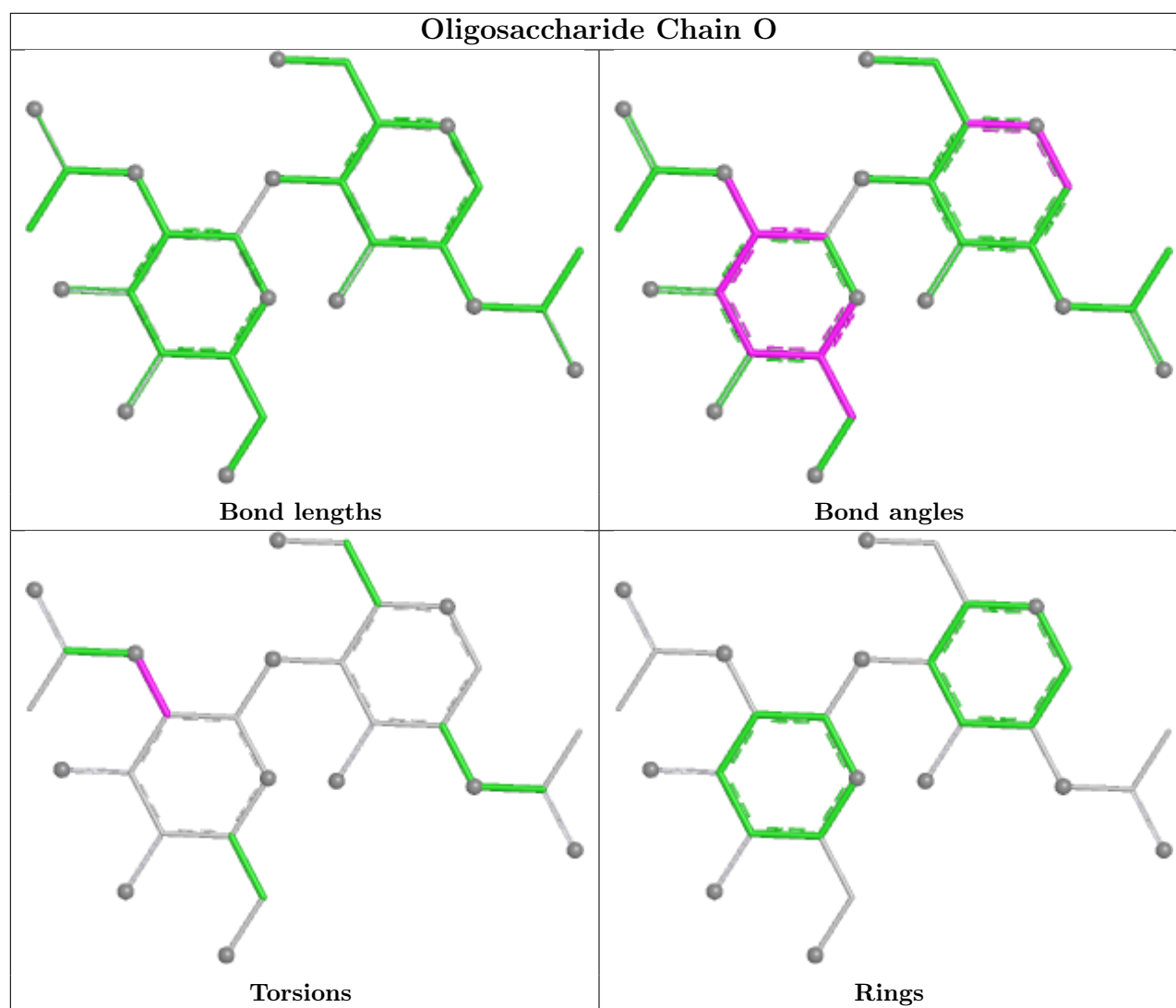
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

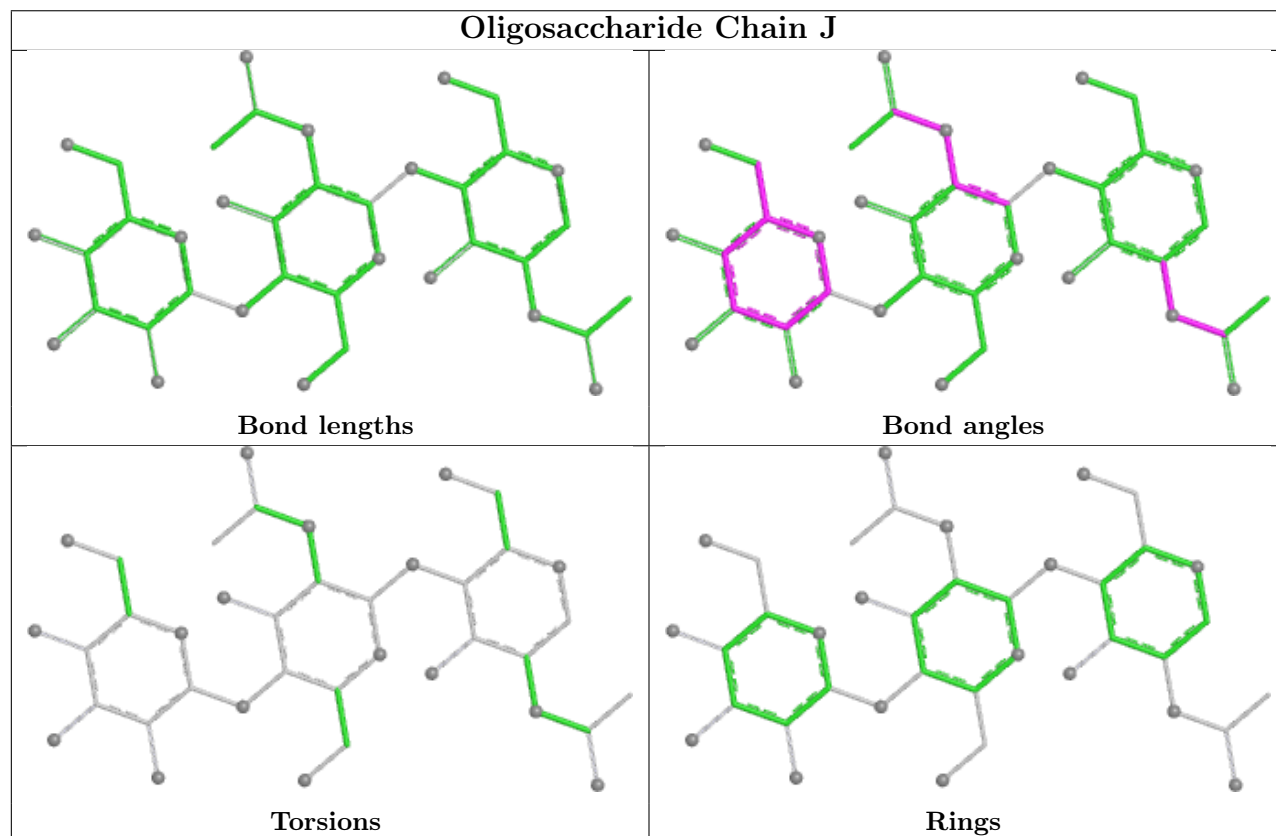
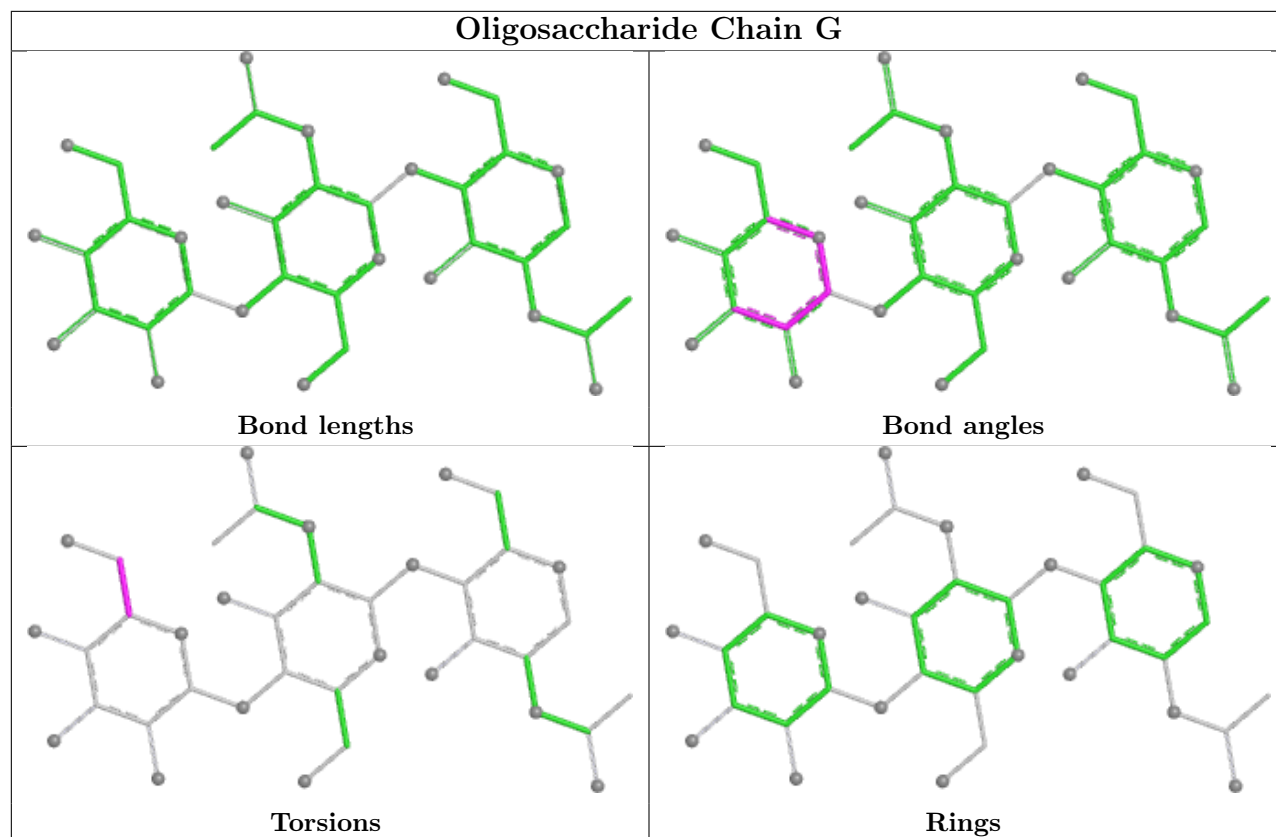


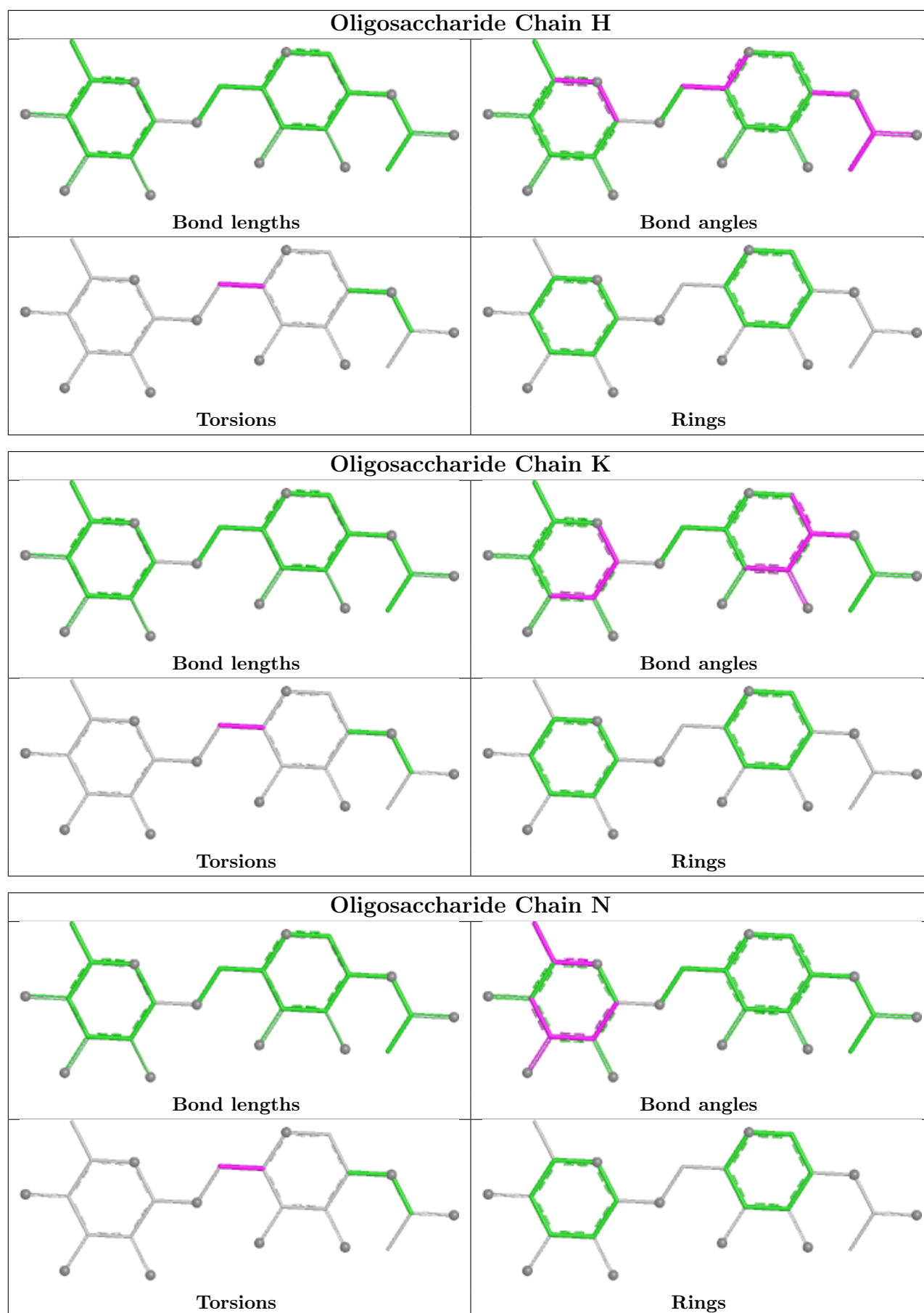


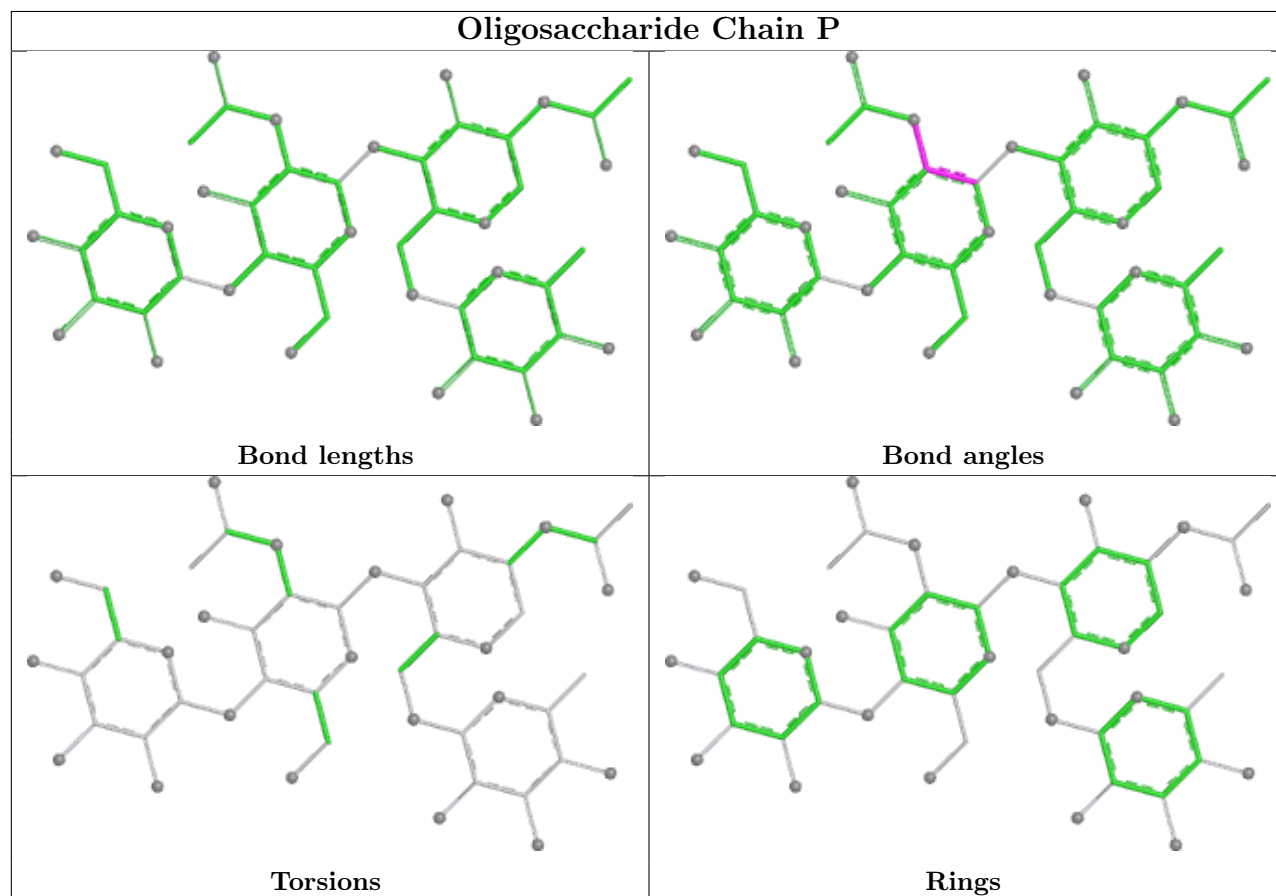
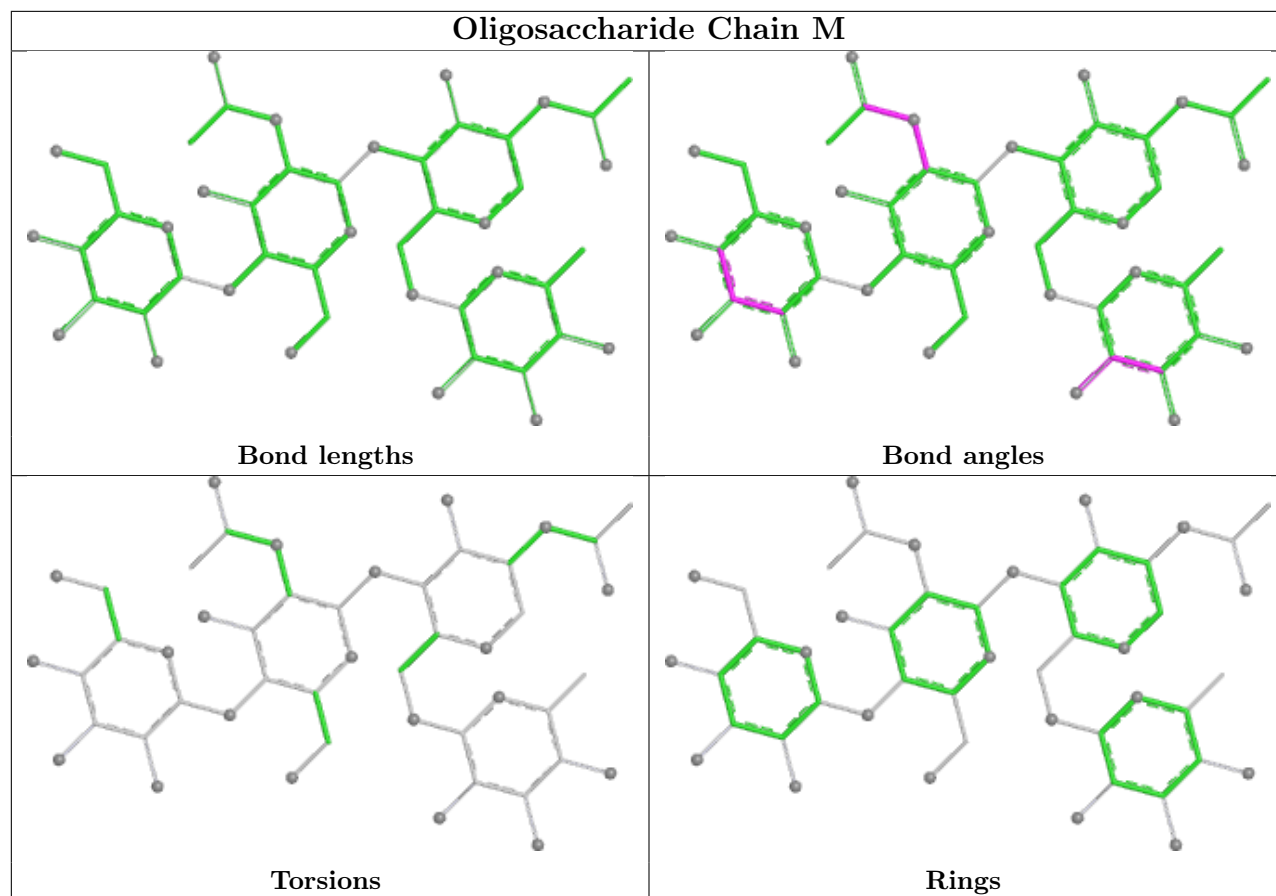












5.6 Ligand geometry

Of 38 ligands modelled in this entry, 12 are monoatomic - leaving 26 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$
13	P6G	D	1201	-	18,18,18	0.72	0	17,17,17	1.11	2 (11%)
12	PEG	D	1202	-	6,6,6	0.49	0	5,5,5	0.60	0
14	GLU	B	911	-	7,8,9	1.11	0	4,9,11	1.19	0
12	PEG	B	1201	-	6,6,6	0.52	0	5,5,5	0.30	0
13	P6G	A	1202	-	18,18,18	0.50	0	17,17,17	0.41	0
14	GLU	C	911	-	7,8,9	1.05	0	4,9,11	1.06	0
12	PEG	B	1205	-	6,6,6	0.43	0	5,5,5	0.44	0
12	PEG	D	1200	-	6,6,6	0.39	0	5,5,5	0.39	0
15	VAL	B	912	-	6,7,7	1.06	1 (16%)	8,9,9	1.12	1 (12%)
12	PEG	C	1201	-	6,6,6	0.55	0	5,5,5	0.66	0
15	VAL	C	912	-	6,7,7	0.85	0	8,9,9	1.11	1 (12%)
8	LYS	A	916	-	8,9,9	0.84	0	7,10,10	1.19	1 (14%)
12	PEG	C	1200	-	6,6,6	0.44	0	5,5,5	0.24	0
12	PEG	A	1201	-	6,6,6	0.47	0	5,5,5	0.51	0
12	PEG	A	1203	-	6,6,6	0.42	0	5,5,5	0.49	0
7	GLN	D	915	-	7,8,9	1.08	0	4,9,11	0.82	0
12	PEG	A	1200	-	6,6,6	0.44	0	5,5,5	0.37	0
8	LYS	D	916	-	8,9,9	1.08	0	7,10,10	0.87	0
12	PEG	C	1202	-	6,6,6	0.53	0	5,5,5	0.36	0
12	PEG	C	1203	-	6,6,6	0.52	0	5,5,5	0.77	0
13	P6G	B	1204	-	18,18,18	0.45	0	17,17,17	0.38	0
12	PEG	D	1203	-	6,6,6	0.44	0	5,5,5	0.43	0
13	P6G	B	1203	-	18,18,18	0.61	0	17,17,17	0.65	0
12	PEG	B	1202	-	6,6,6	0.66	0	5,5,5	0.62	0
7	GLN	A	915	-	7,8,9	0.68	0	4,9,11	0.43	0
12	PEG	D	1204	-	6,6,6	0.48	0	5,5,5	0.41	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
13	P6G	D	1201	-	-	4/16/16/16	-
12	PEG	D	1202	-	-	1/4/4/4	-
14	GLU	B	911	-	-	2/6/7/9	-
12	PEG	B	1201	-	-	1/4/4/4	-
13	P6G	A	1202	-	-	8/16/16/16	-
14	GLU	C	911	-	-	2/6/7/9	-
12	PEG	B	1205	-	-	0/4/4/4	-
12	PEG	D	1200	-	-	1/4/4/4	-
15	VAL	B	912	-	-	0/8/8/8	-
12	PEG	C	1201	-	-	2/4/4/4	-
15	VAL	C	912	-	-	0/8/8/8	-
8	LYS	A	916	-	-	2/9/9/9	-
12	PEG	C	1200	-	-	3/4/4/4	-
12	PEG	A	1201	-	-	2/4/4/4	-
12	PEG	A	1203	-	-	1/4/4/4	-
7	GLN	D	915	-	-	2/6/7/9	-
12	PEG	A	1200	-	-	2/4/4/4	-
8	LYS	D	916	-	-	1/9/9/9	-
12	PEG	C	1202	-	-	3/4/4/4	-
12	PEG	C	1203	-	-	3/4/4/4	-
13	P6G	B	1204	-	-	9/16/16/16	-
12	PEG	D	1203	-	-	3/4/4/4	-
13	P6G	B	1203	-	-	9/16/16/16	-
12	PEG	B	1202	-	-	3/4/4/4	-
7	GLN	A	915	-	-	2/6/7/9	-
12	PEG	D	1204	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	B	912	VAL	OXT-C	-2.52	1.22	1.30

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	916	LYS	OXT-C-O	-2.82	117.69	124.08
15	B	912	VAL	OXT-C-O	-2.43	118.57	124.08
13	D	1201	P6G	O16-C17-C18	2.35	120.47	110.11
15	C	912	VAL	OXT-C-O	-2.25	118.98	124.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	D	1201	P6G	O10-C9-C8	2.06	119.72	110.35

There are no chirality outliers.

All (70) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
13	D	1201	P6G	C8-C9-O10-C11
13	B	1204	P6G	O13-C14-C15-O16
13	B	1204	P6G	O4-C5-C6-O7
13	B	1204	P6G	O7-C8-C9-O10
13	A	1202	P6G	O4-C5-C6-O7
13	A	1202	P6G	O7-C8-C9-O10
13	B	1203	P6G	O4-C5-C6-O7
12	A	1201	PEG	O2-C3-C4-O4
12	D	1203	PEG	O2-C3-C4-O4
13	B	1204	P6G	O10-C11-C12-O13
13	B	1203	P6G	O10-C11-C12-O13
12	C	1200	PEG	O2-C3-C4-O4
12	C	1202	PEG	O2-C3-C4-O4
13	B	1203	P6G	O1-C2-C3-O4
13	B	1203	P6G	O16-C17-C18-O19
13	D	1201	P6G	O1-C2-C3-O4
12	B	1201	PEG	O1-C1-C2-O2
13	B	1204	P6G	O1-C2-C3-O4
13	B	1203	P6G	O7-C8-C9-O10
13	B	1203	P6G	C8-C9-O10-C11
12	B	1202	PEG	O1-C1-C2-O2
13	A	1202	P6G	O16-C17-C18-O19
13	A	1202	P6G	O10-C11-C12-O13
12	C	1200	PEG	O1-C1-C2-O2
12	C	1202	PEG	O1-C1-C2-O2
12	C	1203	PEG	O1-C1-C2-O2
12	D	1200	PEG	O1-C1-C2-O2
12	D	1203	PEG	O1-C1-C2-O2
12	D	1204	PEG	O1-C1-C2-O2
12	D	1202	PEG	O1-C1-C2-O2
7	A	915	GLN	OE1-CD-CG-CB
7	A	915	GLN	NE2-CD-CG-CB
13	B	1204	P6G	C15-C14-O13-C12
13	A	1202	P6G	C8-C9-O10-C11
12	C	1201	PEG	O1-C1-C2-O2
8	A	916	LYS	CG-CD-CE-NZ

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Mol	Chain	Res	Type	Atoms
12	D	1204	PEG	O2-C3-C4-O4
13	A	1202	P6G	O1-C2-C3-O4
13	A	1202	P6G	C18-C17-O16-C15
7	D	915	GLN	OE1-CD-CG-CB
12	C	1201	PEG	C1-C2-O2-C3
12	C	1203	PEG	C1-C2-O2-C3
13	B	1203	P6G	C2-C3-O4-C5
13	B	1203	P6G	C11-C12-O13-C14
12	A	1200	PEG	O1-C1-C2-O2
12	B	1202	PEG	O2-C3-C4-O4
12	B	1202	PEG	C1-C2-O2-C3
12	C	1203	PEG	C4-C3-O2-C2
12	D	1204	PEG	C1-C2-O2-C3
13	D	1201	P6G	O13-C14-C15-O16
12	D	1204	PEG	C4-C3-O2-C2
12	A	1203	PEG	C1-C2-O2-C3
7	D	915	GLN	NE2-CD-CG-CB
12	A	1200	PEG	C1-C2-O2-C3
13	B	1204	P6G	C8-C9-O10-C11
13	B	1204	P6G	C6-C5-O4-C3
12	C	1200	PEG	C1-C2-O2-C3
13	B	1203	P6G	O13-C14-C15-O16
13	D	1201	P6G	O10-C11-C12-O13
12	D	1203	PEG	C4-C3-O2-C2
13	B	1204	P6G	C11-C12-O13-C14
12	A	1201	PEG	C4-C3-O2-C2
14	B	911	GLU	OE2-CD-CG-CB
14	C	911	GLU	OE2-CD-CG-CB
12	C	1202	PEG	C4-C3-O2-C2
14	B	911	GLU	OE1-CD-CG-CB
8	D	916	LYS	OXT-C-CA-N
13	A	1202	P6G	C5-C6-O7-C8
8	A	916	LYS	OXT-C-CA-N
14	C	911	GLU	OE1-CD-CG-CB

There are no ring outliers.

7 monomers are involved in 9 short contacts:

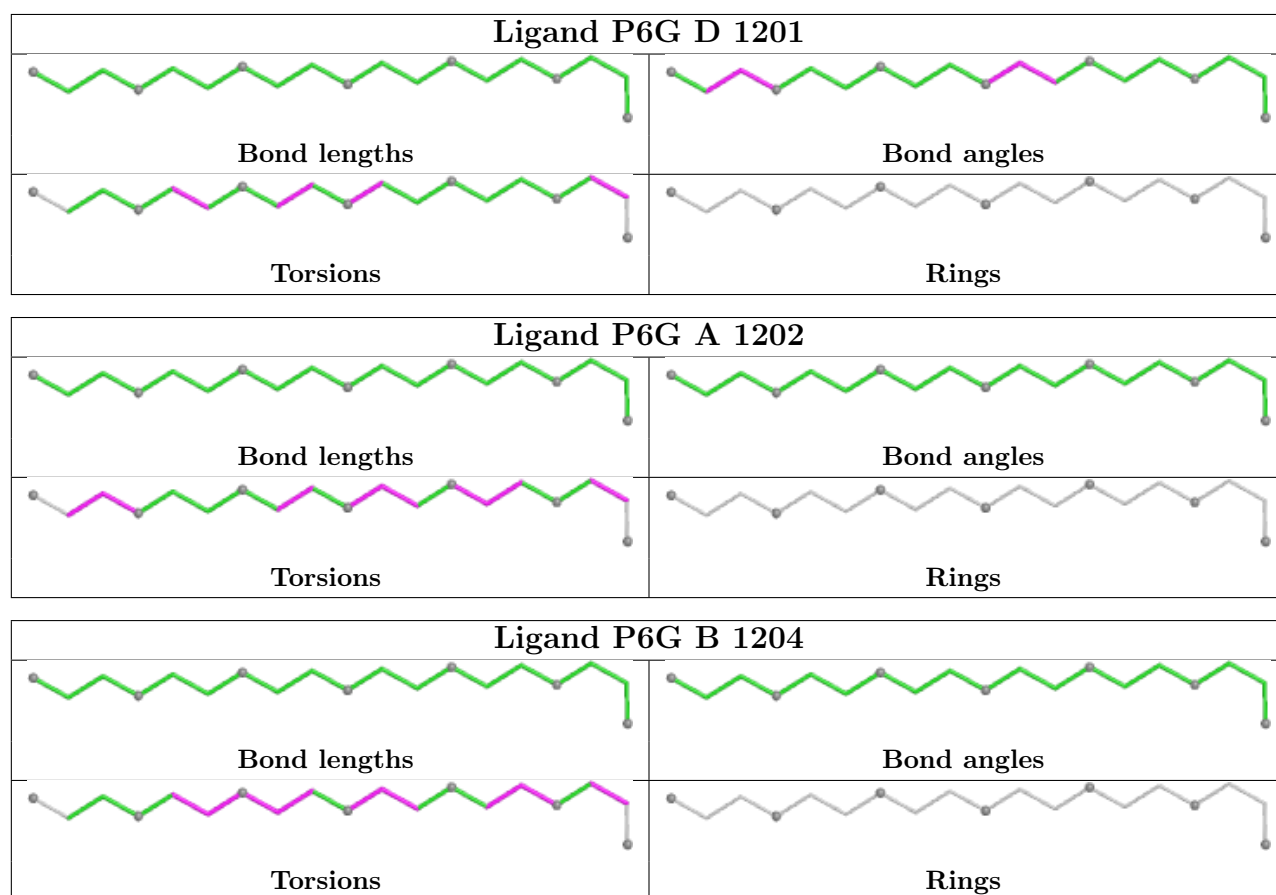
Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	D	1201	P6G	1	0
15	B	912	VAL	1	0
12	C	1201	PEG	2	0

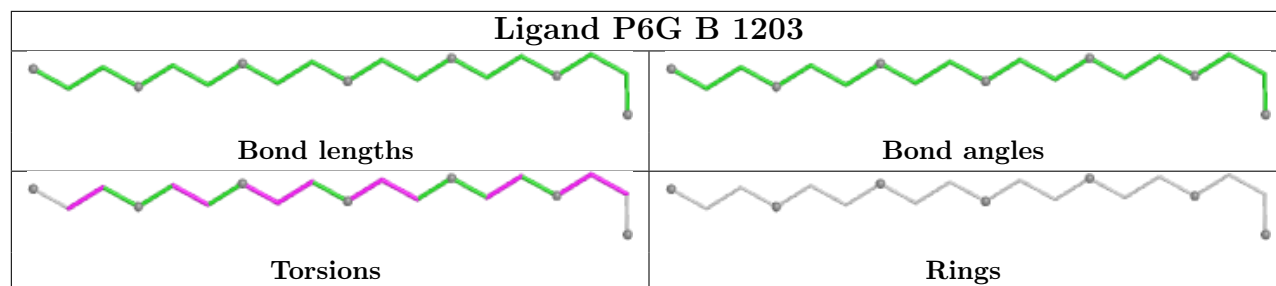
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
15	C	912	VAL	1	0
8	A	916	LYS	2	0
7	D	915	GLN	1	0
13	B	1203	P6G	1	0

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	603/629 (95%)	0.07	15 (2%) 58 58	16, 27, 46, 70	1 (0%)
1	B	605/629 (96%)	-0.11	9 (1%) 72 72	11, 24, 41, 67	3 (0%)
1	C	611/629 (97%)	-0.19	13 (2%) 63 64	12, 22, 39, 60	3 (0%)
1	D	609/629 (96%)	0.11	52 (8%) 16 14	15, 24, 50, 79	3 (0%)
All	All	2428/2516 (96%)	-0.03	89 (3%) 45 45	11, 24, 45, 79	10 (0%)

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	79	ILE	7.2
1	D	83	PHE	6.1
1	C	414	VAL	5.9
1	A	415	THR	5.2
1	D	325	GLY	5.1
1	D	20	PHE	5.0
1	D	88	LEU	4.9
1	D	78	PRO	4.8
1	A	414	VAL	4.7
1	D	76	TYR	4.6
1	D	95	VAL	4.6
1	D	80	TRP	4.6
1	C	135	THR	4.5
1	D	24	TYR	4.5
1	D	18	GLN	4.3
1	D	134	ALA	4.2
1	D	414	VAL	4.1
1	A	135	THR	4.1
1	D	17	ALA	4.0
1	D	15	ALA	3.9
1	D	11	SER	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	12	ALA	3.9
1	A	325	GLY	3.9
1	B	135	THR	3.8
1	A	129	LEU	3.8
1	D	21	ALA	3.8
1	D	19	LEU	3.8
1	D	133	THR	3.7
1	D	27	SER	3.7
1	D	84	THR	3.6
1	D	72	ALA	3.6
1	D	16	GLY	3.6
1	C	415	THR	3.5
1	C	325	GLY	3.5
1	D	25	GLN	3.5
1	A	270	PRO	3.4
1	D	10	PHE	3.4
1	D	14	GLU	3.3
1	B	415	THR	3.3
1	B	607	TYR	3.2
1	B	131	GLN	3.1
1	A	78	PRO	3.1
1	D	612	ASP	3.0
1	D	75	LEU	3.0
1	B	325	GLY	3.0
1	D	135	THR	2.9
1	C	206	THR	2.8
1	D	91	ILE	2.7
1	D	92	ILE	2.7
1	D	26	SER	2.7
1	D	69	GLY	2.6
1	C	606	ASN	2.6
1	D	607	TYR	2.6
1	D	9[A]	GLN	2.6
1	D	23	SER	2.6
1	C	413	ARG	2.5
1	D	86	PRO	2.5
1	B	206	THR	2.5
1	A	608	PRO	2.5
1	A	79	ILE	2.5
1	A	412	ASP	2.5
1	D	94	ALA	2.4
1	D	415	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	13	ASP	2.4
1	A	269	VAL	2.4
1	C	129	LEU	2.4
1	D	90	ARG	2.3
1	D	609	GLU	2.3
1	D	413	ARG	2.3
1	D	22	GLN	2.3
1	B	414	VAL	2.3
1	D	97	THR	2.3
1	D	453	ARG	2.2
1	C	613	LEU	2.2
1	A	413	ARG	2.2
1	A	206	THR	2.2
1	B	273	ASP	2.2
1	C	610	GLY	2.2
1	D	28	ALA	2.2
1	A	267	MET	2.1
1	C	614	VAL	2.1
1	C	609	GLU	2.1
1	D	81	GLN	2.1
1	A	84	THR	2.1
1	D	73	LYS	2.1
1	D	377	VAL	2.0
1	C	611	ILE	2.0
1	B	608	PRO	2.0
1	D	89	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	FUC	N	2	10/11	0.53	0.27	57,62,65,66	0
3	NAG	I	2	14/15	0.55	0.16	65,71,75,79	0

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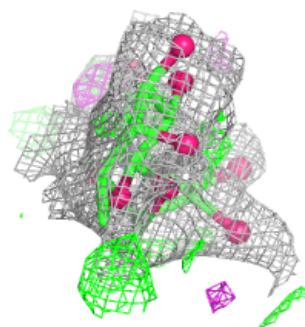
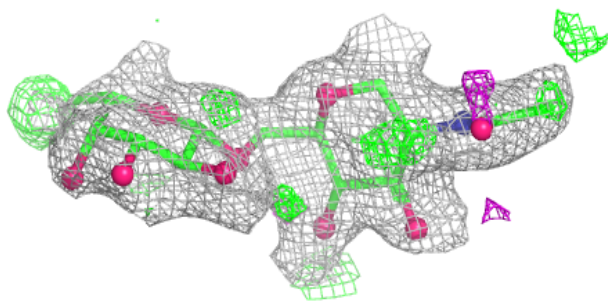
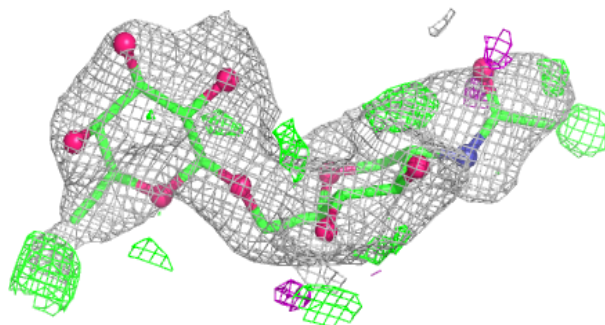
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FUL	E	2	10/11	0.55	0.21	63,66,66,68	0
4	MAN	J	3	11/12	0.62	0.17	64,66,69,69	0
5	FUC	K	2	10/11	0.65	0.23	58,63,67,68	0
5	FUC	H	2	10/11	0.65	0.24	58,64,67,68	0
3	NAG	F	2	14/15	0.66	0.14	62,68,70,70	0
3	NAG	O	2	14/15	0.66	0.14	63,69,74,75	0
6	BMA	P	3	11/12	0.67	0.14	65,67,70,72	0
4	MAN	G	3	11/12	0.71	0.13	70,74,76,76	0
6	BMA	M	3	11/12	0.74	0.14	46,48,53,54	0
2	NAG	E	1	14/15	0.76	0.15	44,50,58,58	0
3	NAG	L	2	14/15	0.77	0.14	49,54,63,67	0
4	NAG	G	2	14/15	0.79	0.14	67,70,72,75	0
3	NAG	O	1	14/15	0.79	0.13	41,45,51,58	0
5	NAG	N	1	14/15	0.81	0.14	38,44,49,52	0
5	NAG	H	1	14/15	0.81	0.13	36,42,50,50	0
5	NAG	K	1	14/15	0.82	0.13	39,45,50,52	0
6	FUC	P	4	10/11	0.82	0.14	45,48,51,51	0
4	NAG	G	1	14/15	0.84	0.12	62,63,67,67	0
3	NAG	I	1	14/15	0.85	0.11	40,44,50,58	0
3	NAG	F	1	14/15	0.85	0.11	39,41,47,54	0
6	FUC	M	4	10/11	0.86	0.11	42,43,46,47	0
6	NAG	P	2	14/15	0.86	0.12	46,52,56,61	0
4	NAG	J	2	14/15	0.88	0.11	52,55,57,60	0
6	NAG	M	2	14/15	0.89	0.12	38,41,47,50	0
4	NAG	J	1	14/15	0.89	0.10	46,48,52,54	0
3	NAG	L	1	14/15	0.91	0.09	34,37,40,46	0
6	NAG	P	1	14/15	0.92	0.09	34,38,42,42	0
6	NAG	M	1	14/15	0.92	0.09	32,33,36,39	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.

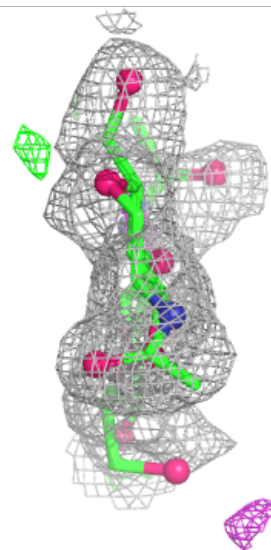
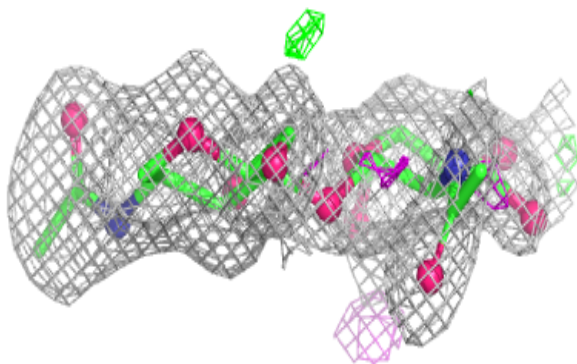
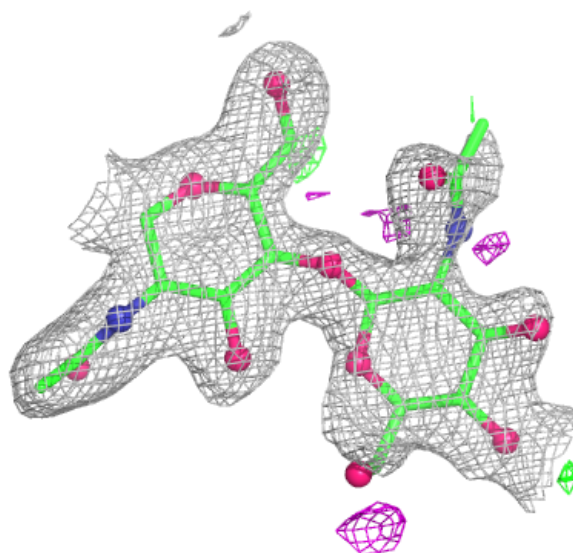
Electron density around Chain E:

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



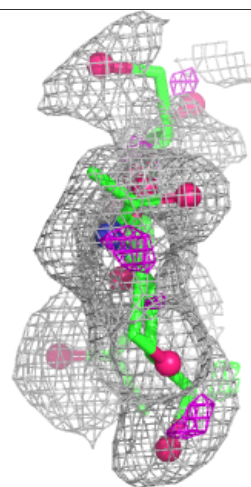
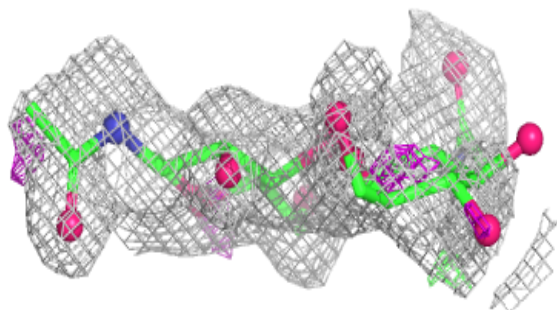
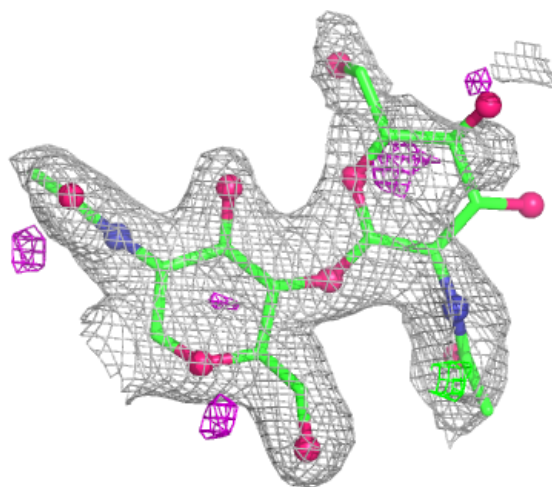
Electron density around Chain F:

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



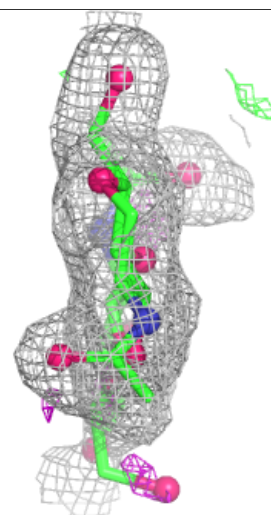
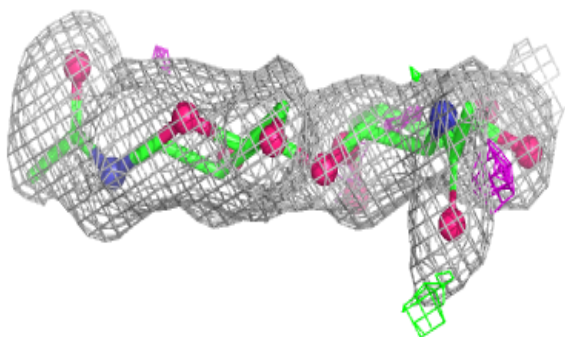
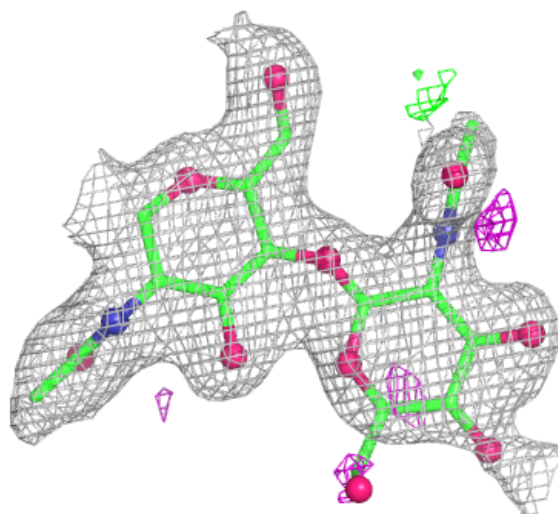
Electron density around Chain I:

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



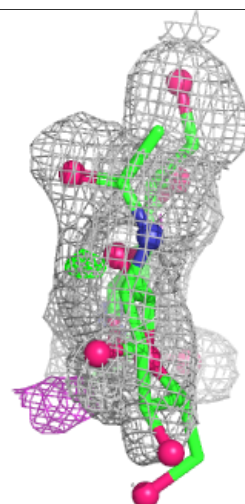
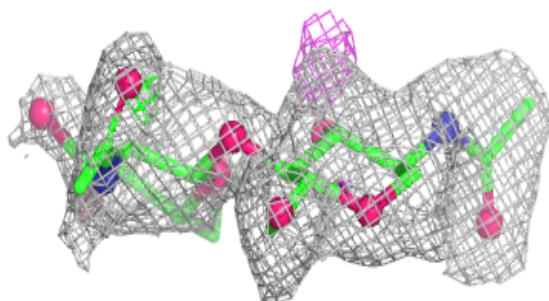
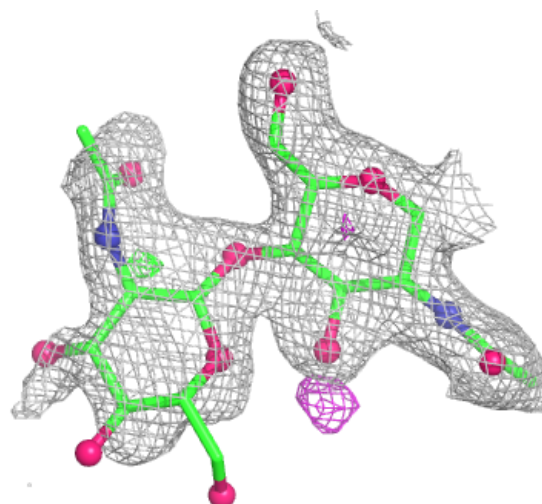
Electron density around Chain L:

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



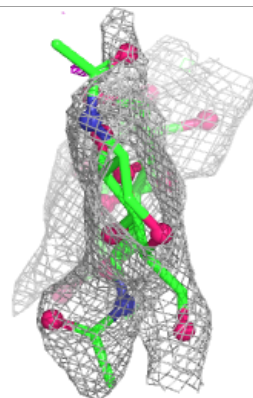
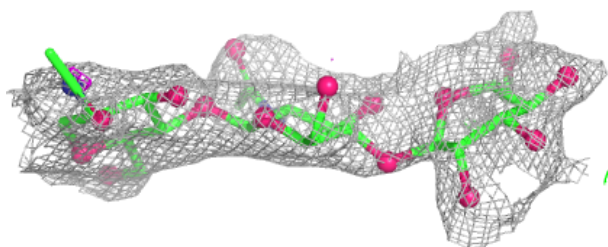
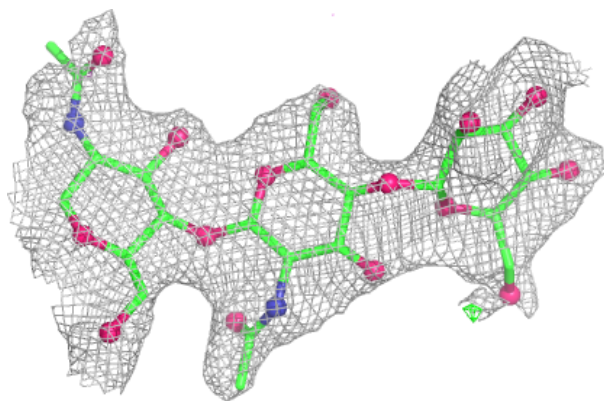
Electron density around Chain O:

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

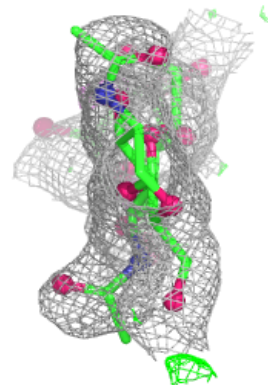
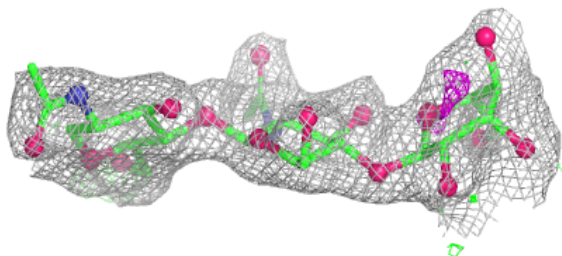
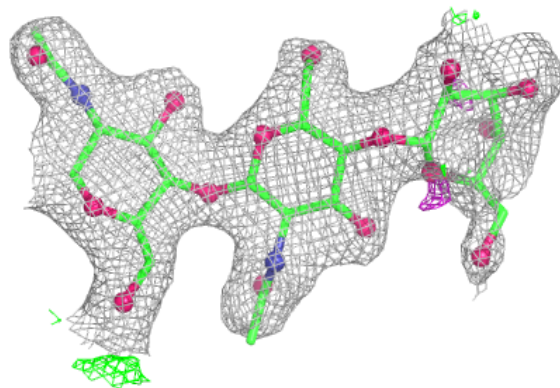


Electron density around Chain G:

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

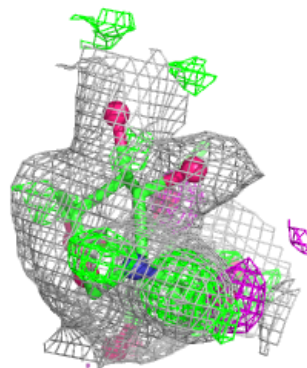
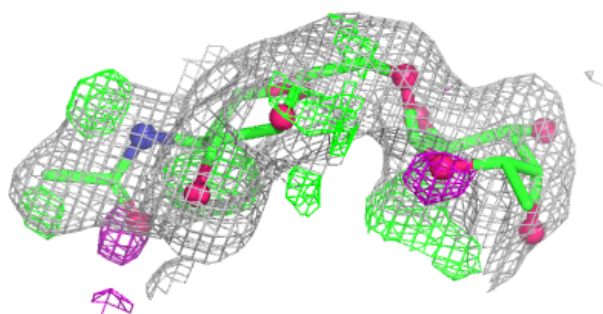
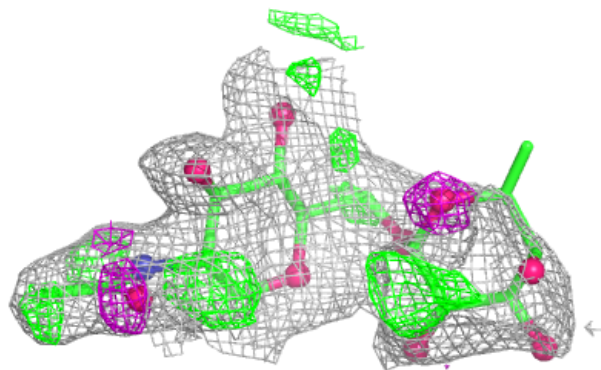
**Electron density around Chain J:**

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

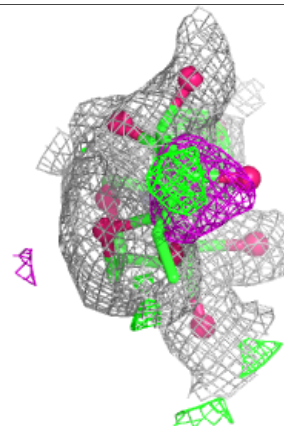
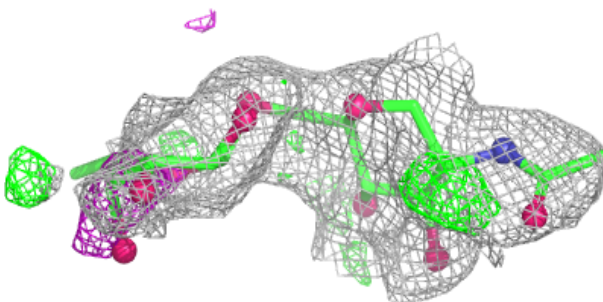
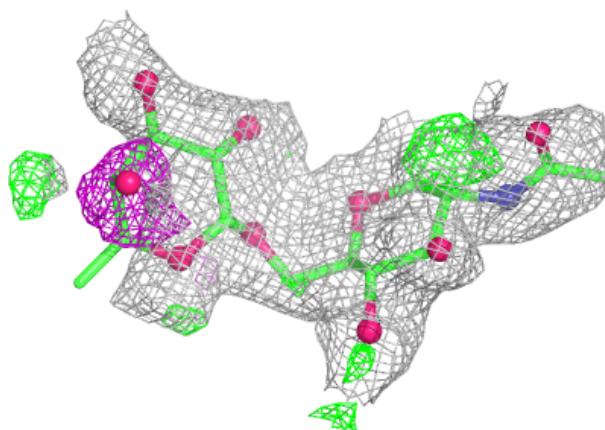


Electron density around Chain H:

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

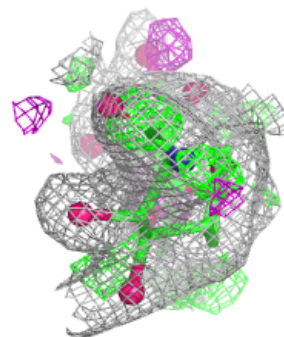
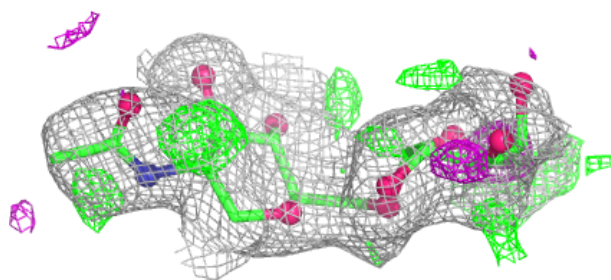
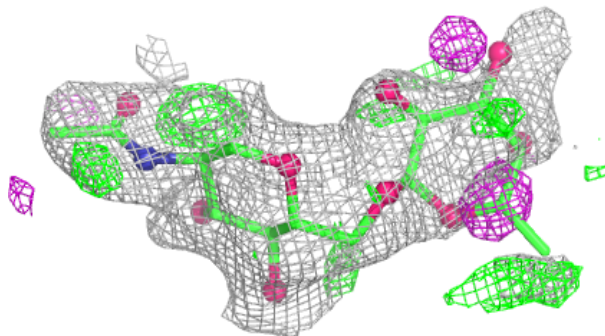
**Electron density around Chain K:**

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



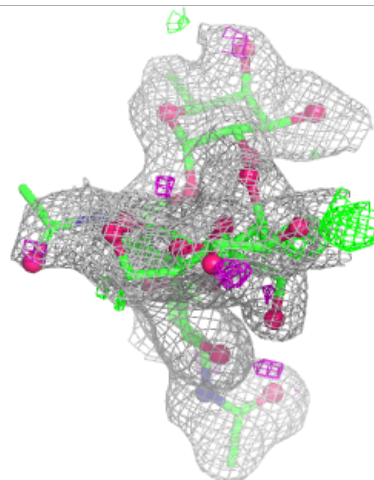
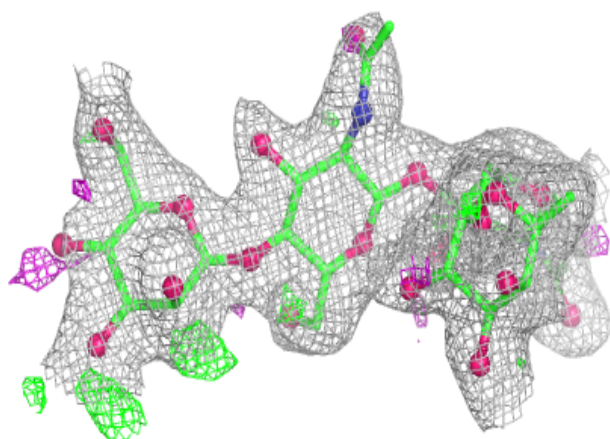
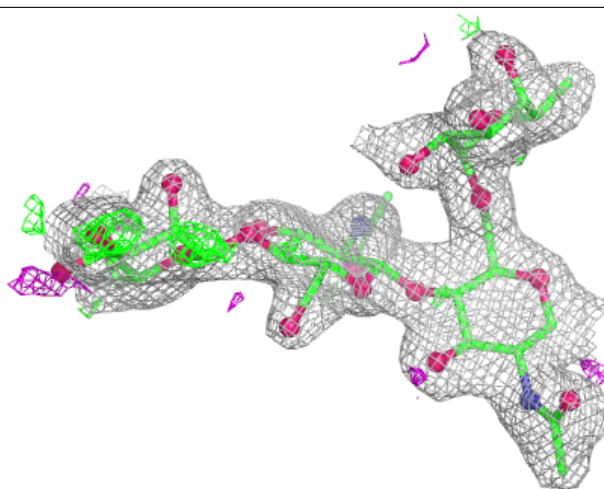
Electron density around Chain N:

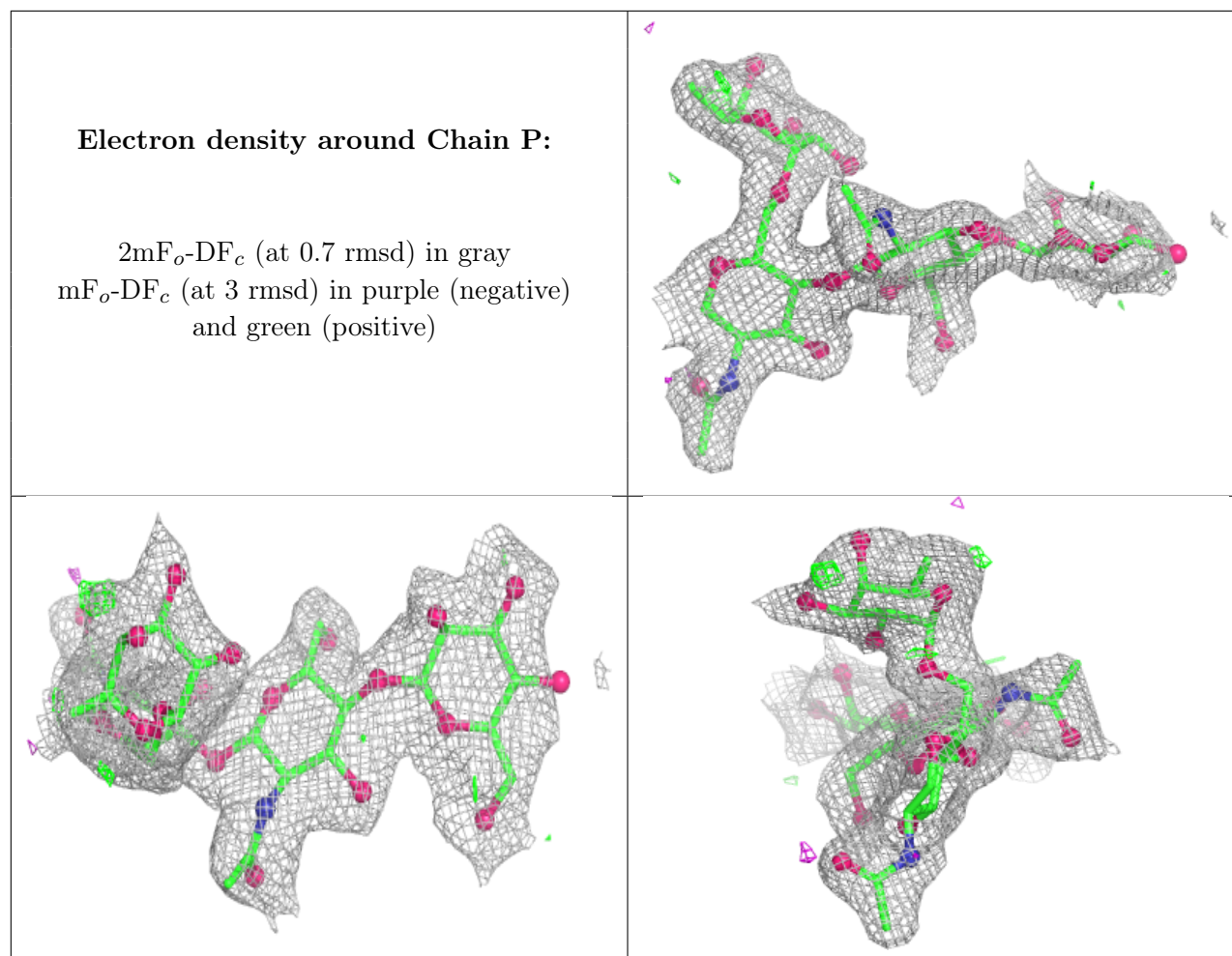
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain M:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
12	PEG	D	1203	7/7	0.75	0.24	57,59,60,62	0
12	PEG	A	1201	7/7	0.80	0.17	40,43,46,47	0
12	PEG	C	1202	7/7	0.82	0.19	51,52,56,56	0
12	PEG	D	1200	7/7	0.84	0.14	56,57,58,59	0
12	PEG	D	1202	7/7	0.84	0.16	51,53,53,54	0
12	PEG	B	1201	7/7	0.84	0.19	53,54,56,57	0
13	P6G	B	1204	19/19	0.84	0.17	50,57,62,64	0
12	PEG	C	1203	7/7	0.86	0.16	40,41,43,45	0
8	LYS	D	916	10/10	0.87	0.12	21,25,33,33	0
12	PEG	B	1202	7/7	0.87	0.16	43,44,46,47	0

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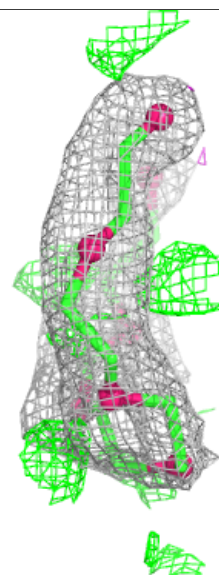
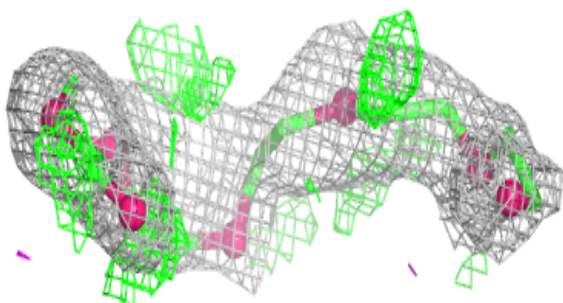
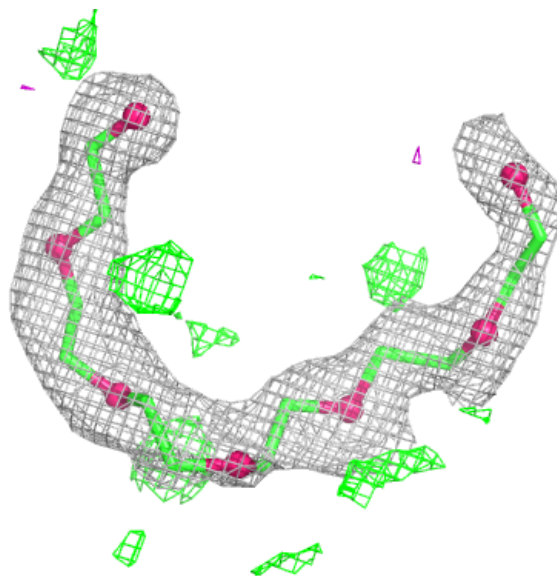
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	GLN	D	915	9/10	0.88	0.12	23,27,32,33	0
12	PEG	D	1204	7/7	0.88	0.16	52,53,53,54	0
8	LYS	A	916	10/10	0.88	0.11	23,25,34,34	0
12	PEG	B	1205	7/7	0.89	0.12	56,57,58,60	0
12	PEG	C	1201	7/7	0.89	0.13	41,42,44,45	0
13	P6G	B	1203	19/19	0.89	0.14	38,41,52,55	0
12	PEG	A	1200	7/7	0.89	0.13	53,53,58,58	0
13	P6G	D	1201	19/19	0.89	0.13	29,36,49,51	0
13	P6G	A	1202	19/19	0.90	0.13	37,43,49,50	0
12	PEG	A	1203	7/7	0.90	0.11	40,41,43,45	0
12	PEG	C	1200	7/7	0.92	0.10	46,46,50,51	0
7	GLN	A	915	9/10	0.92	0.08	23,26,27,28	0
14	GLU	B	911	9/10	0.93	0.08	22,23,26,27	0
15	VAL	B	912	8/8	0.95	0.07	19,22,23,25	0
14	GLU	C	911	9/10	0.96	0.07	18,21,23,24	0
11	NA	D	1003	1/1	0.96	0.07	40,40,40,40	0
15	VAL	C	912	8/8	0.97	0.07	19,21,23,23	0
11	NA	A	1003	1/1	0.98	0.04	30,30,30,30	0
10	CL	B	1002	1/1	0.99	0.03	19,19,19,19	0
10	CL	D	1002	1/1	0.99	0.03	20,20,20,20	0
16	CA	B	1003	1/1	0.99	0.05	41,41,41,41	0
9	ZN	D	1001	1/1	1.00	0.01	19,19,19,19	0
10	CL	A	1002	1/1	1.00	0.01	21,21,21,21	0
9	ZN	A	1001	1/1	1.00	0.01	20,20,20,20	0
10	CL	C	1002	1/1	1.00	0.01	17,17,17,17	0
9	ZN	B	1001	1/1	1.00	0.01	18,18,18,18	0
9	ZN	C	1001	1/1	1.00	0.02	17,17,17,17	0
16	CA	C	1003	1/1	1.00	0.05	25,25,25,25	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.

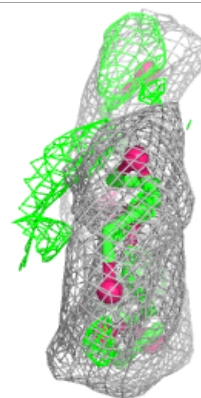
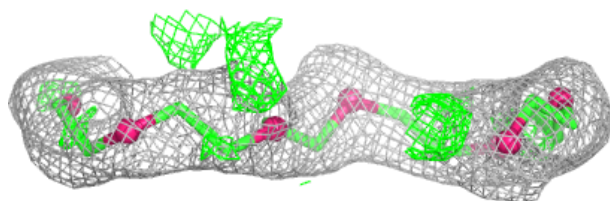
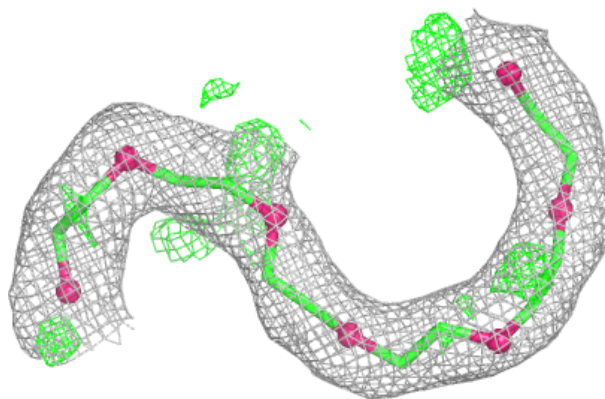
Electron density around P6G B 1204:

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



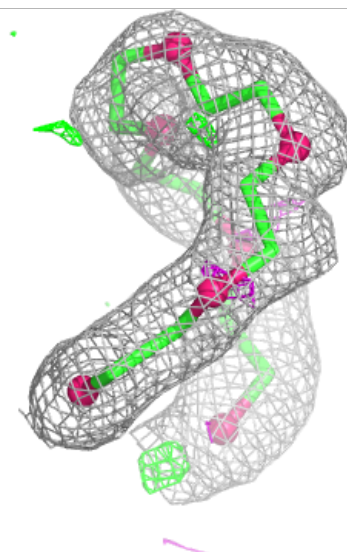
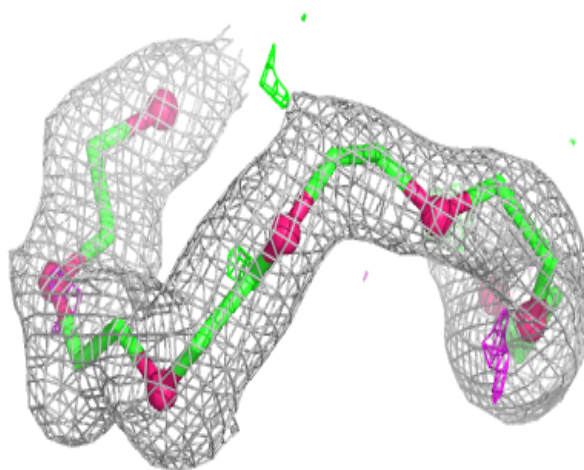
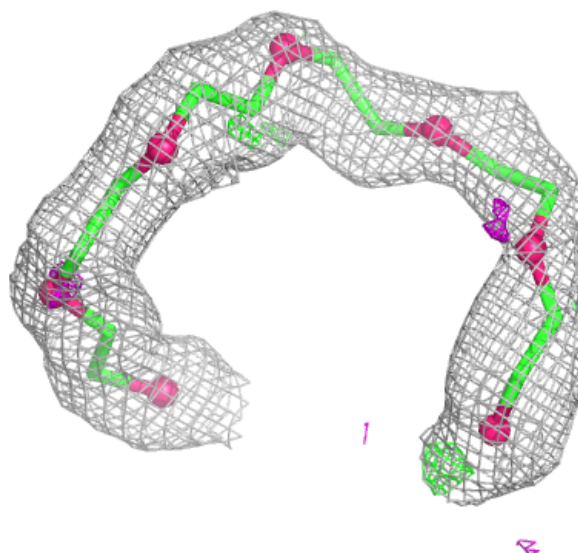
Electron density around P6G B 1203:

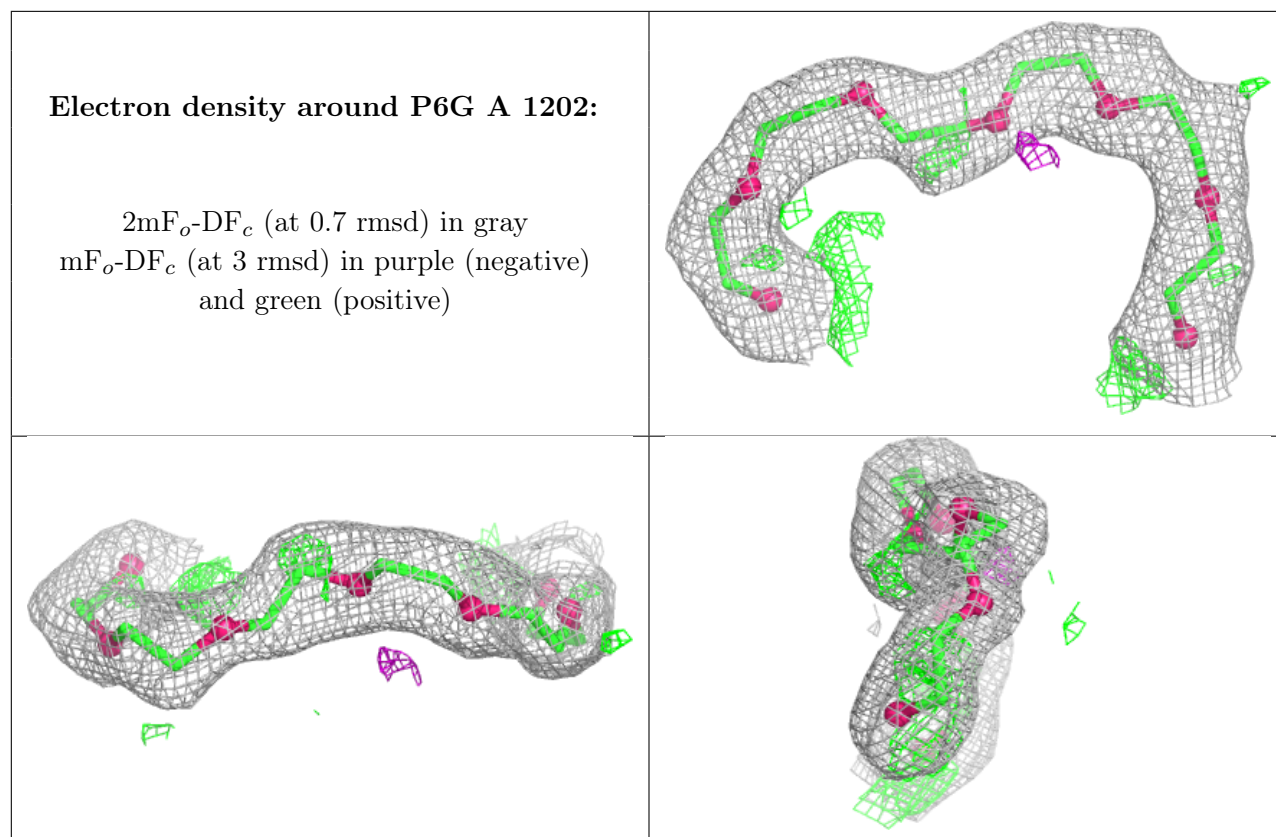
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around P6G D 1201:

$2mF_o-DF_c$ (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.