



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 9, 2024 – 01:52 pm GMT

PDB ID : 4P4E
Title : X-ray structure of human glutamate carboxypeptidase II (GCPII) in complex with a phosphoramidate inhibitor MP1D
Authors : Barinka, C.
Deposited on : 2014-03-12
Resolution : 1.67 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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

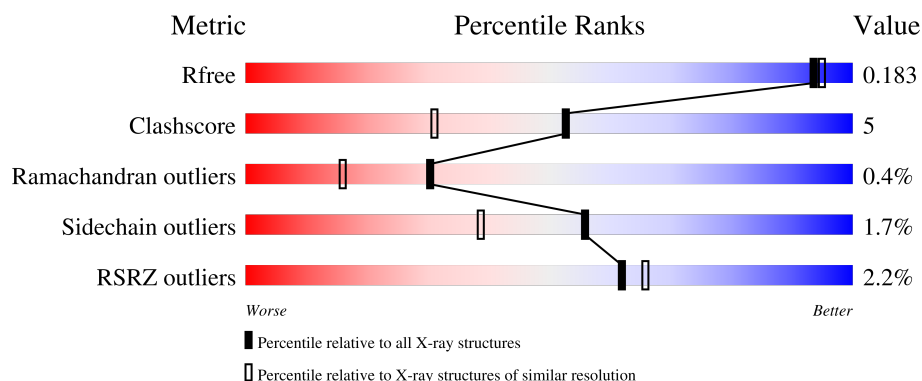
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.67 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}	164625	8422 (1.70-1.66)
Clashscore	180529	1005 (1.68-1.68)
Ramachandran outliers	177936	9065 (1.70-1.66)
Sidechain outliers	177891	9064 (1.70-1.66)
RSRZ outliers	164620	8421 (1.70-1.66)

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	707	<div> <div>2%</div> <div>87%</div> <div>10%</div> <div>..</div> </div>
2	B	2	<div> <div>50%</div> <div>50%</div> </div>
3	C	3	<div> <div>33%</div> <div>67%</div> </div>
3	D	3	<div> <div>33%</div> <div>67%</div> </div>
4	E	3	<div> <div>100%</div> </div>

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Mol	Chain	Length	Quality of chain
5	F	4	 25% 75%

2 Entry composition [i](#)

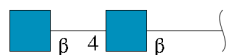
There are 11 unique types of molecules in this entry. The entry contains 6739 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glutamate carboxypeptidase 2.

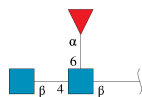
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	694	Total	C	N	O	S	0	75	0
			5928	3801	992	1109	26			

- Molecule 2 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
2	B	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



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

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



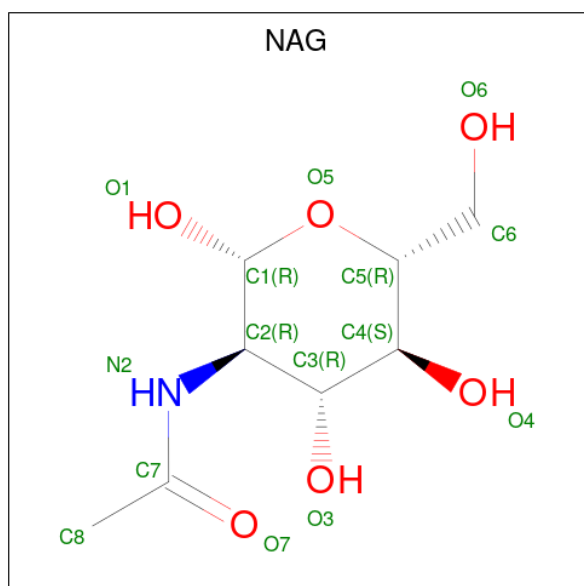
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is an oligosaccharide called 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	F	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 7 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	Zn	0	0
			2	2		

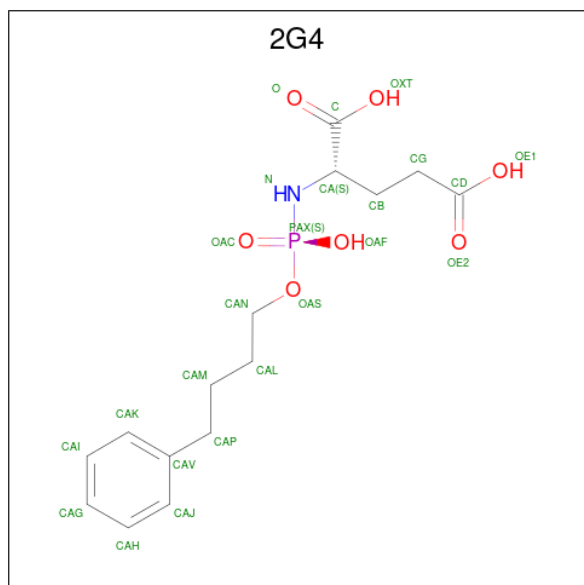
- Molecule 8 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Ca	0	0
			1	1		

- Molecule 9 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1	Total	Cl	0	0
			1	1		

- Molecule 10 is N-[(S)-hydroxy(4-phenylbutoxy)phosphoryl]-L-glutamic acid (three-letter code: 2G4) (formula: C₁₅H₂₂NO₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	A	1	Total	C	N	O	P	0	0
			24	15	1	7	1		

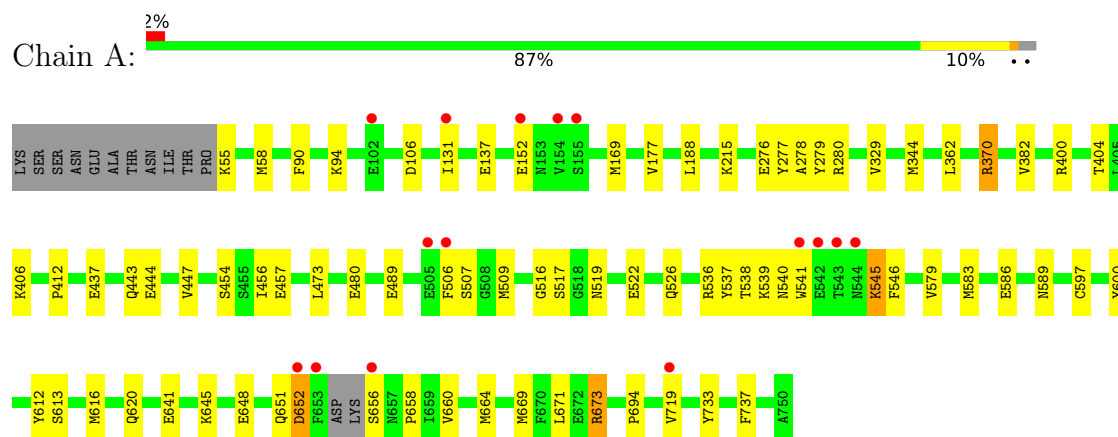
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	562	Total 562	O 562	0	0

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: Glutamate carboxypeptidase 2



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



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



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



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

Chain E:  100%

MAG1
MAG2
BMA3

- Molecule 5: 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 F:  25% 75%

MAG1
MAG2
BMA3
MAN4

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	101.75Å 130.68Å 159.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.29 – 1.67 28.29 – 1.67	Depositor EDS
% Data completeness (in resolution range)	96.6 (28.29-1.67) 96.6 (28.29-1.67)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 1.67Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.154 , 0.183 0.154 , 0.183	Depositor DCC
R_{free} test set	1194 reflections (1.01%)	wwPDB-VP
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6739	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% of the height of the origin peak. No significant pseudotranslation is detected.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, 2G4, FUC, CL, CA, BMA, ZN, MAN

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	0.93	2/6245 (0.0%)	0.83	2/8450 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	737	PHE	CD2-CE2	5.19	1.49	1.39
1	A	733	TYR	CD1-CE1	5.13	1.47	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	370	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	A	673	ARG	NE-CZ-NH1	6.16	123.38	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5928	0	5782	59	1
2	B	28	0	25	1	0
3	C	38	0	34	0	0
3	D	38	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	39	0	34	0	0
5	F	50	0	43	0	0
6	A	28	0	26	0	0
7	A	2	0	0	0	0
8	A	1	0	0	0	0
9	A	1	0	0	0	0
10	A	24	0	19	0	0
11	A	562	0	0	13	2
All	All	6739	0	5997	60	2

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 (60) 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:A:651[B]:GLN:O	1:A:651[B]:GLN:NE2	1.65	1.27
1:A:660[A]:VAL:O	1:A:664[A]:MET:HG2	1.43	1.17
1:A:651[B]:GLN:O	1:A:651[B]:GLN:CD	1.88	1.11
1:A:658[A]:PRO:HB3	11:A:1344:HOH:O	1.62	1.00
1:A:412:PRO:HA	1:A:589[B]:ASN:HD21	1.34	0.87
1:A:215[B]:LYS:NZ	11:A:902:HOH:O	2.13	0.80
1:A:597[B]:CYS:SG	1:A:671:LEU:HD22	2.21	0.80
1:A:597[B]:CYS:SG	1:A:671:LEU:CD2	2.76	0.73
1:A:437[A]:GLU:OE1	11:A:901:HOH:O	2.07	0.73
1:A:641:GLU:HG3	11:A:1265:HOH:O	1.90	0.71
1:A:400:ARG:O	1:A:404[B]:THR:HG23	1.91	0.71
1:A:362:LEU:CD1	1:A:406:LYS:HD2	2.23	0.69
1:A:131[A]:ILE:HG22	1:A:137:GLU:HG2	1.75	0.69
1:A:276[B]:GLU:HG2	1:A:277:TYR:N	2.10	0.67
1:A:370:ARG:HD2	1:A:669[B]:MET:HE1	1.77	0.66
1:A:447:VAL:HG22	1:A:669[B]:MET:CE	2.26	0.65
1:A:719[A]:VAL:HG12	1:A:719[A]:VAL:O	1.98	0.64
1:A:90:PHE:CE2	1:A:94:LYS:HE2	2.37	0.59
1:A:480[B]:GLU:O	1:A:480[B]:GLU:HG3	2.03	0.58
1:A:719[A]:VAL:O	1:A:719[A]:VAL:CG1	2.52	0.58
1:A:516:GLY:O	1:A:526[B]:GLN:NE2	2.34	0.58
1:A:719[B]:VAL:HG22	11:A:1347:HOH:O	2.05	0.57
1:A:412:PRO:HA	1:A:589[B]:ASN:ND2	2.15	0.56
1:A:447:VAL:HG22	1:A:669[B]:MET:HE2	1.88	0.55
1:A:177:VAL:HG12	1:A:188:LEU:HD11	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:LEU:HD11	1:A:406:LYS:HD2	1.87	0.55
1:A:480[B]:GLU:O	1:A:480[B]:GLU:CG	2.54	0.55
1:A:152:GLU:OE2	11:A:903:HOH:O	2.19	0.52
1:A:447:VAL:HG22	1:A:669[B]:MET:HE1	1.90	0.52
1:A:579:VAL:O	1:A:583[B]:MET:HG2	2.09	0.52
1:A:545:LYS:HE3	1:A:546:PHE:CZ	2.48	0.49
1:A:169:MET:HA	1:A:344:MET:O	2.14	0.48
1:A:58[B]:MET:SD	1:A:586:GLU:HG2	2.54	0.48
1:A:215[B]:LYS:HE2	11:A:958:HOH:O	2.14	0.48
1:A:489:GLU:H	1:A:489:GLU:CD	2.18	0.47
1:A:106:ASP:OD1	1:A:406:LYS:HE3	2.14	0.47
1:A:620:GLN:NE2	11:A:920:HOH:O	2.47	0.47
1:A:506:PHE:CB	1:A:509:MET:HG3	2.45	0.47
1:A:188:LEU:HD21	1:A:329[A]:VAL:HG11	1.97	0.46
1:A:612:TYR:CZ	1:A:616:MET:HG3	2.50	0.46
1:A:651[B]:GLN:O	1:A:651[B]:GLN:CG	2.58	0.45
1:A:278:ALA:HB3	1:A:280[A]:ARG:CZ	2.47	0.44
1:A:517:SER:OG	1:A:522:GLU:OE2	2.30	0.44
1:A:517:SER:HB2	1:A:694:PRO:HG3	1.98	0.44
1:A:645[A]:LYS:HD2	11:A:953:HOH:O	2.17	0.44
1:A:454[B]:SER:OG	1:A:457[B]:GLU:OE2	2.35	0.44
1:A:719[B]:VAL:HG13	11:A:1343:HOH:O	2.16	0.44
1:A:613[B]:SER:OG	11:A:904:HOH:O	2.20	0.44
1:A:443[A]:GLN:HG3	1:A:444:GLU:CD	2.38	0.44
1:A:648:GLU:O	1:A:652:ASP:HB2	2.17	0.44
11:A:1404:HOH:O	2:B:2:NAG:H81	2.17	0.43
1:A:276[B]:GLU:HB3	11:A:908:HOH:O	2.18	0.43
1:A:456[B]:ILE:O	1:A:456[B]:ILE:HG13	2.19	0.43
1:A:473:LEU:HG	1:A:583[B]:MET:SD	2.59	0.42
1:A:188:LEU:CD2	1:A:329[A]:VAL:HG11	2.49	0.42
1:A:656[B]:SER:O	1:A:658[B]:PRO:HD3	2.19	0.42
1:A:457[A]:GLU:HG2	1:A:538:THR:HA	2.02	0.41
1:A:506:PHE:HB3	1:A:509:MET:HG3	2.02	0.41
1:A:669[B]:MET:HE2	1:A:669[B]:MET:HB2	1.55	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:A:910:HOH:O	11:A:1150:HOH:O[2_565]	1.83	0.37
1:A:656[B]:SER:O	11:A:1344:HOH:O[4_566]	2.10	0.10

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	760/707 (108%)	738 (97%)	19 (2%)	3 (0%)	30 16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	541	TRP
1	A	382	VAL
1	A	652	ASP

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	662/603 (110%)	652 (98%)	10 (2%)	60 44

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	LYS
1	A	507	SER
1	A	519	ASN
1	A	536	ARG
1	A	537	TYR
1	A	539[A]	LYS
1	A	540	ASN
1	A	545	LYS

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Mol	Chain	Res	Type
1	A	600	TYR
1	A	673	ARG

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

Mol	Chain	Res	Type
1	A	540	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 ⓘ

15 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	B	1	1,2	14,14,15	0.54	0	17,19,21	1.40	2 (11%)
2	NAG	B	2	2	14,14,15	0.78	0	17,19,21	1.35	2 (11%)
3	NAG	C	1	3,1	14,14,15	0.75	1 (7%)	17,19,21	1.36	4 (23%)
3	NAG	C	2	3	14,14,15	0.56	0	17,19,21	1.42	2 (11%)
3	FUC	C	3	3	10,10,11	0.69	0	14,14,16	1.08	0
3	NAG	D	1	3,1	14,14,15	0.61	0	17,19,21	0.83	0
3	NAG	D	2	3	14,14,15	0.51	0	17,19,21	1.05	1 (5%)
3	FUC	D	3	3	10,10,11	0.79	0	14,14,16	1.22	2 (14%)
4	NAG	E	1	1,4	14,14,15	0.96	1 (7%)	17,19,21	1.02	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	E	2	4	14,14,15	0.53	0	17,19,21	1.28	2 (11%)
4	BMA	E	3	4	11,11,12	0.86	0	15,15,17	1.09	1 (6%)
5	NAG	F	1	5,1	14,14,15	0.65	0	17,19,21	1.54	4 (23%)
5	NAG	F	2	5	14,14,15	0.49	0	17,19,21	1.42	3 (17%)
5	BMA	F	3	5	11,11,12	0.59	0	15,15,17	0.95	0
5	MAN	F	4	5	11,11,12	0.68	0	15,15,17	1.46	2 (13%)

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	B	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	2/6/23/26	0/1/1/1
3	NAG	C	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	C	2	3	-	2/6/23/26	0/1/1/1
3	FUC	C	3	3	-	-	0/1/1/1
3	NAG	D	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	1/6/23/26	0/1/1/1
3	FUC	D	3	3	-	-	0/1/1/1
4	NAG	E	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	2/6/23/26	0/1/1/1
4	BMA	E	3	4	-	2/2/19/22	0/1/1/1
5	NAG	F	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	F	2	5	-	2/6/23/26	0/1/1/1
5	BMA	F	3	5	-	0/2/19/22	0/1/1/1
5	MAN	F	4	5	-	0/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1	NAG	O7-C7	2.96	1.29	1.23
3	C	1	NAG	O5-C1	-2.06	1.40	1.43

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	4	MAN	O5-C5-C6	3.61	112.87	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	NAG	O5-C5-C6	3.56	112.79	107.20
4	E	2	NAG	C8-C7-N2	3.39	121.84	116.10
5	F	2	NAG	C1-O5-C5	3.10	116.40	112.19
5	F	1	NAG	O5-C5-C6	2.99	111.88	107.20
2	B	2	NAG	C2-N2-C7	2.95	127.10	122.90
5	F	1	NAG	O5-C1-C2	-2.88	106.75	111.29
4	E	1	NAG	C2-N2-C7	2.72	126.78	122.90
5	F	1	NAG	O4-C4-C5	-2.66	102.68	109.30
4	E	3	BMA	O5-C1-C2	-2.57	106.80	110.77
2	B	2	NAG	C4-C3-C2	2.44	114.60	111.02
3	D	2	NAG	O4-C4-C5	2.44	115.35	109.30
2	B	1	NAG	C2-N2-C7	-2.44	119.43	122.90
3	C	1	NAG	C1-C2-N2	-2.41	106.37	110.49
5	F	2	NAG	C8-C7-N2	2.41	120.17	116.10
3	C	1	NAG	C1-O5-C5	-2.40	108.94	112.19
3	D	3	FUC	O2-C2-C3	2.37	114.89	110.14
5	F	1	NAG	C2-N2-C7	-2.28	119.66	122.90
3	C	1	NAG	O3-C3-C2	-2.24	104.83	109.47
4	E	2	NAG	O7-C7-C8	-2.22	117.93	122.06
5	F	4	MAN	O3-C3-C2	2.21	114.22	109.99
3	C	2	NAG	C1-O5-C5	2.18	115.15	112.19
3	D	3	FUC	O5-C1-C2	-2.16	107.44	110.77
2	B	1	NAG	C4-C3-C2	2.14	114.16	111.02
3	C	1	NAG	C8-C7-N2	2.09	119.63	116.10
5	F	2	NAG	O4-C4-C5	-2.06	104.19	109.30

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2	NAG	O5-C5-C6-O6
4	E	3	BMA	O5-C5-C6-O6
4	E	3	BMA	C4-C5-C6-O6
3	C	2	NAG	C4-C5-C6-O6
2	B	2	NAG	C8-C7-N2-C2
2	B	2	NAG	O7-C7-N2-C2
3	C	1	NAG	C8-C7-N2-C2
3	C	1	NAG	O7-C7-N2-C2
4	E	1	NAG	C8-C7-N2-C2
4	E	1	NAG	O7-C7-N2-C2
4	E	2	NAG	C8-C7-N2-C2
4	E	2	NAG	O7-C7-N2-C2

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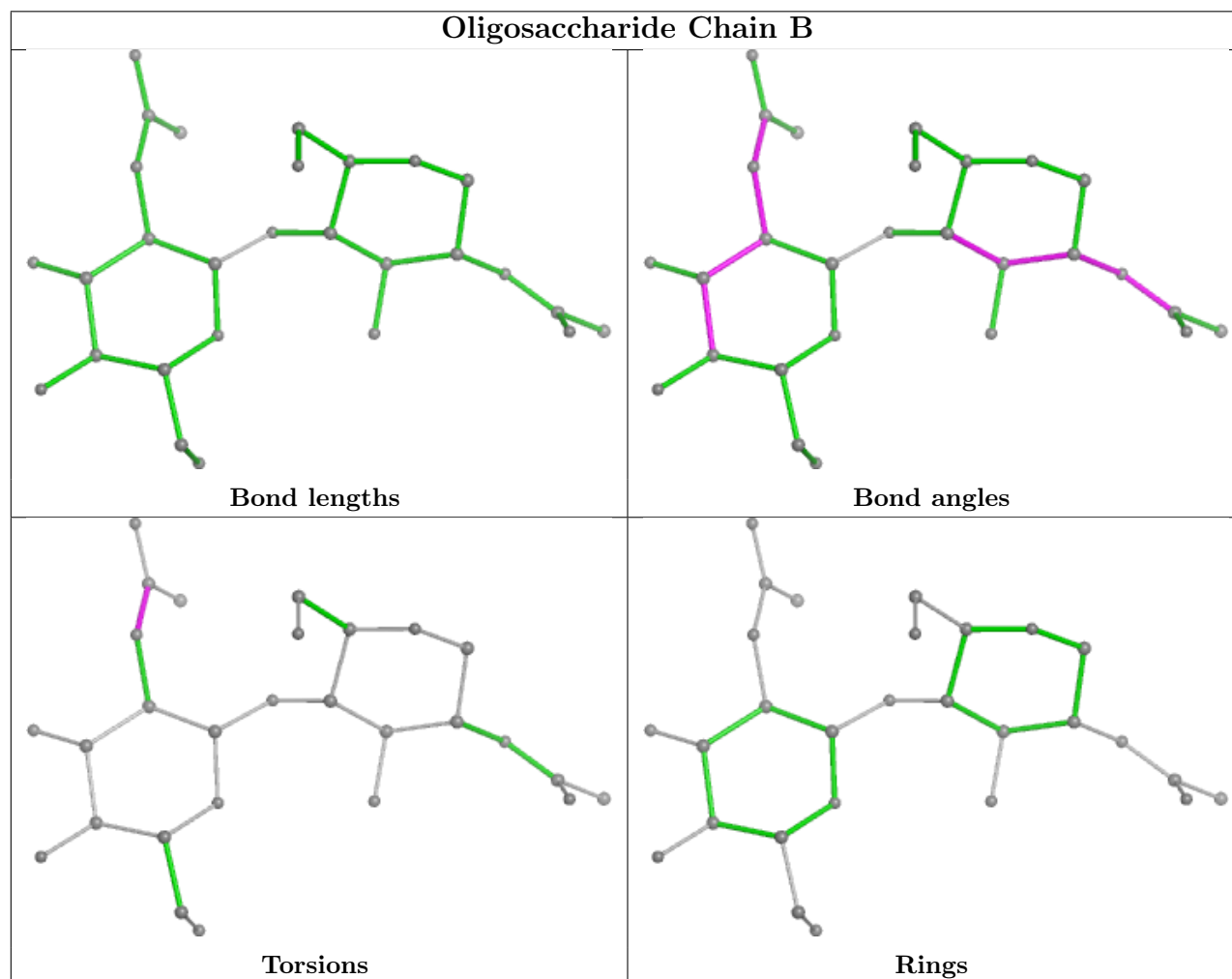
Mol	Chain	Res	Type	Atoms
5	F	2	NAG	C8-C7-N2-C2
5	F	2	NAG	O7-C7-N2-C2
5	F	1	NAG	C4-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6

There are no ring outliers.

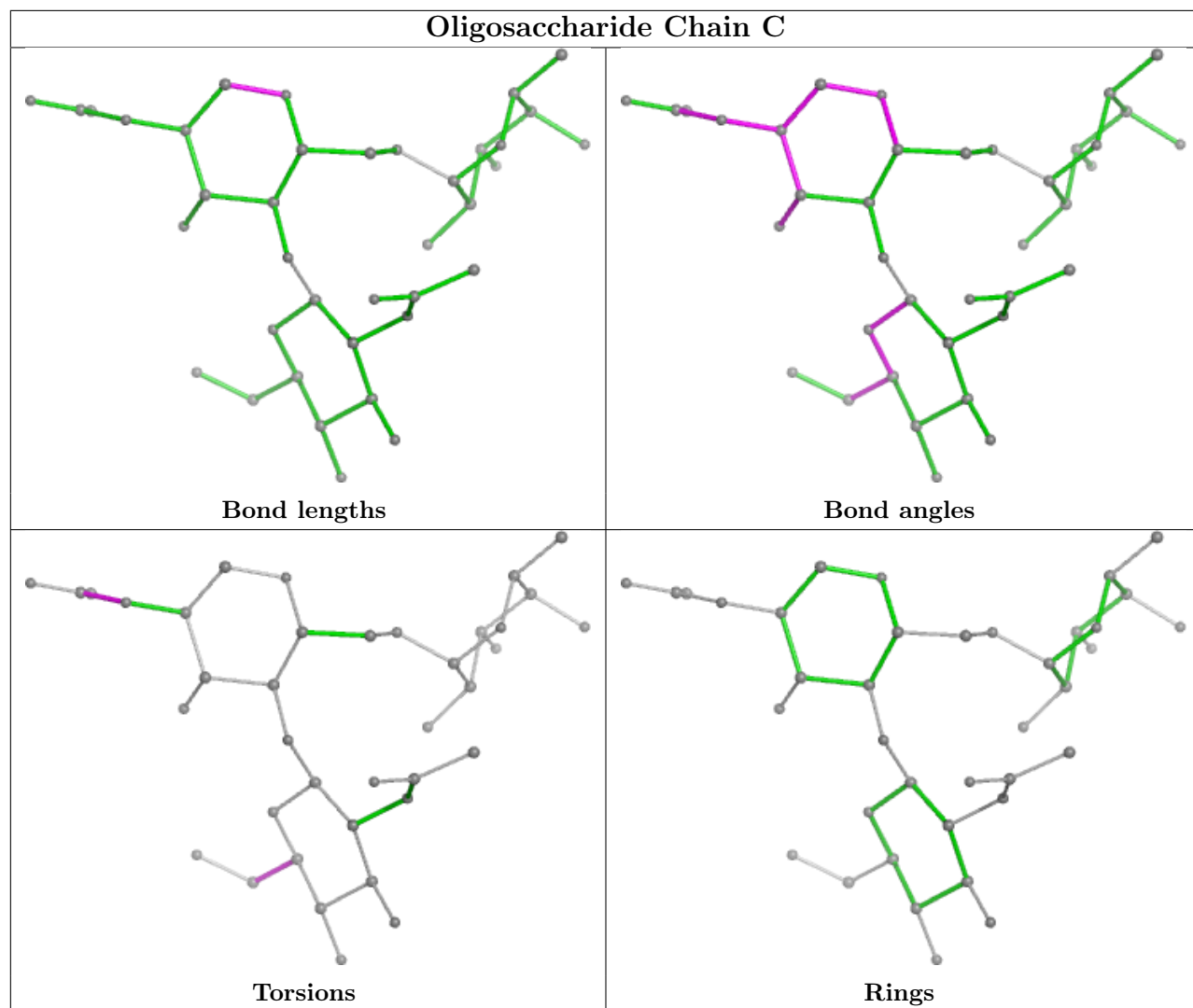
1 monomer is involved in 1 short contact:

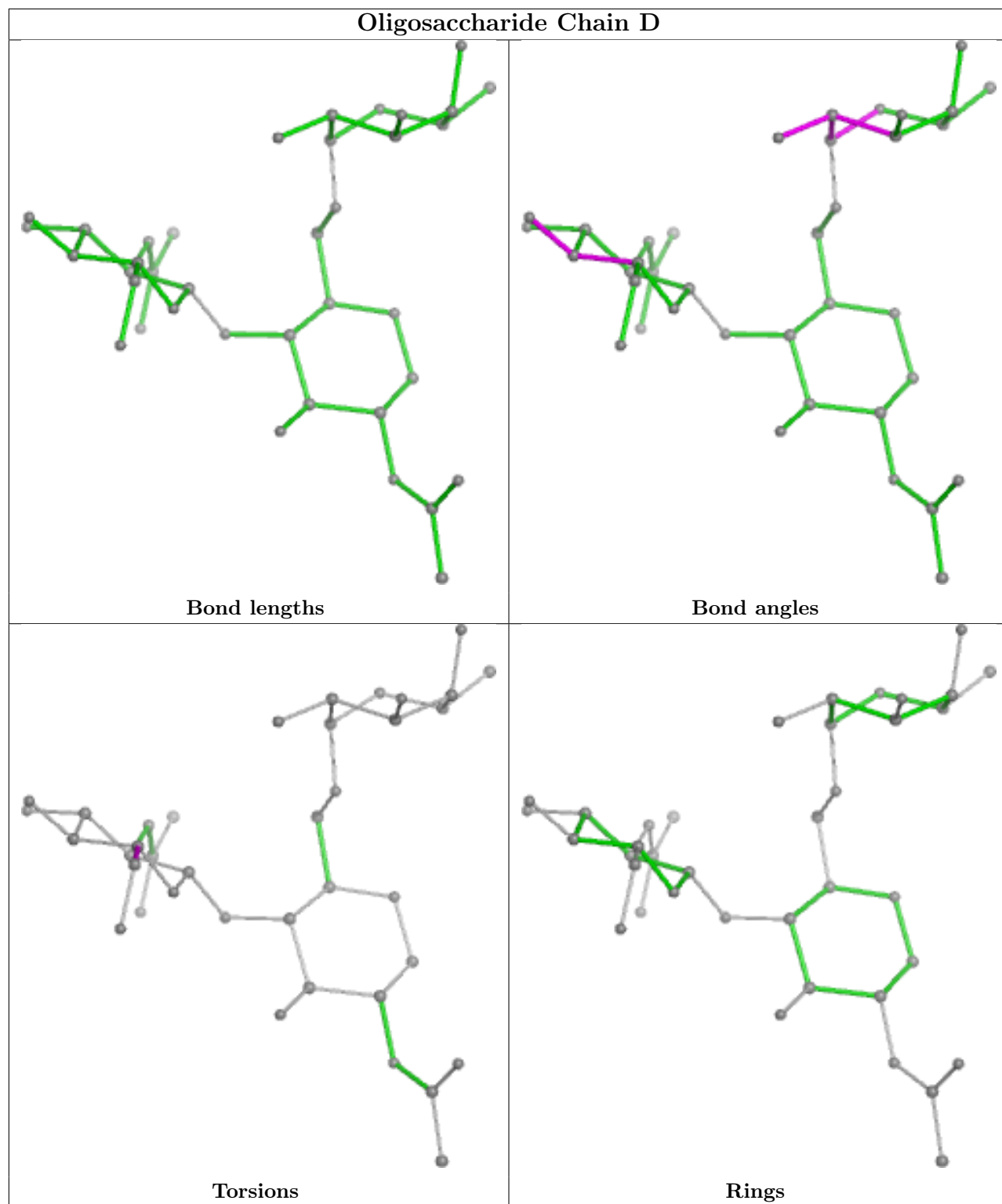
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	NAG	1	0

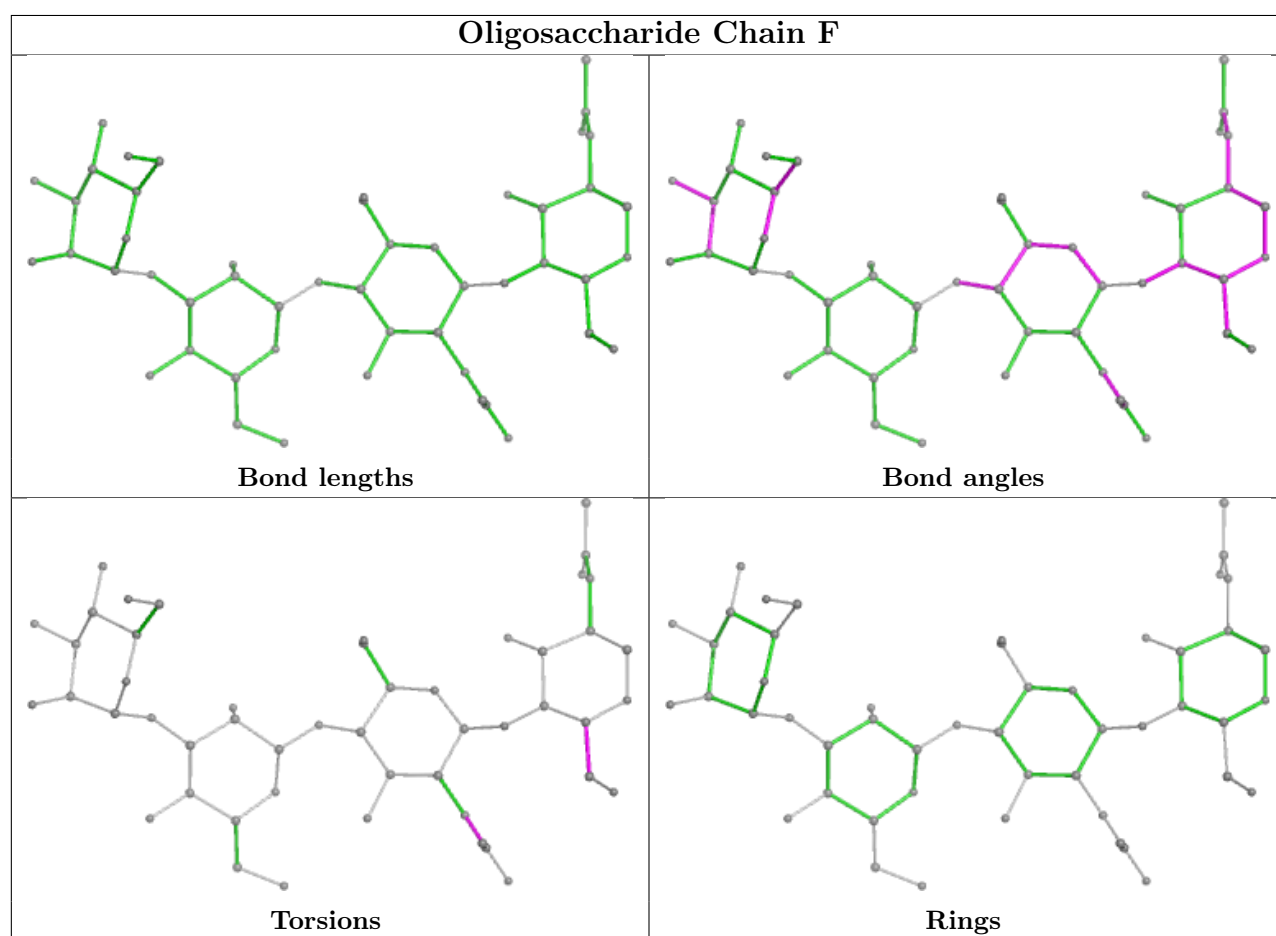
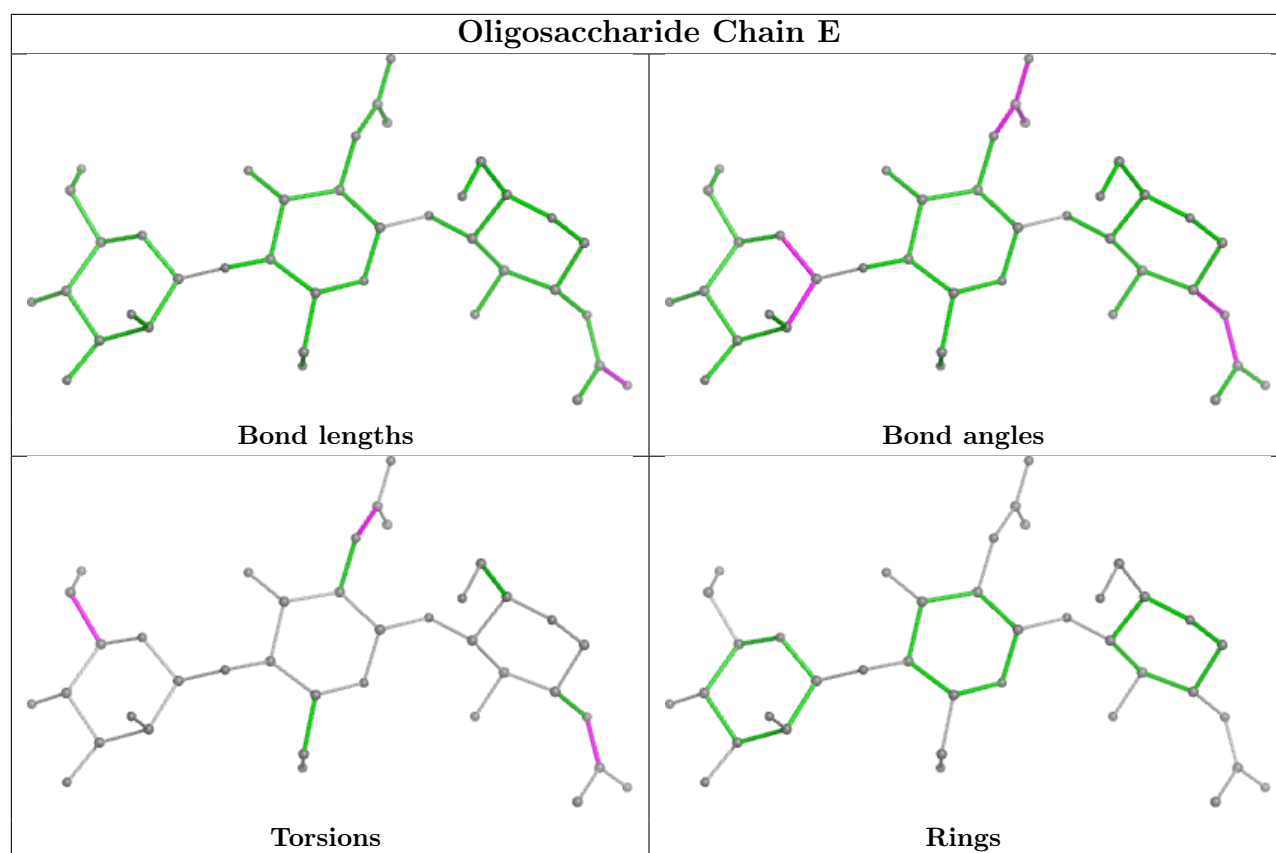
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



Oligosaccharide Chain C







5.6 Ligand geometry

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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
10	2G4	A	822	7	24,24,24	1.96	5 (20%)	30,31,31	1.25	3 (10%)
6	NAG	A	809	1	14,14,15	0.52	0	17,19,21	2.37	3 (17%)
6	NAG	A	810	1	14,14,15	0.84	0	17,19,21	1.82	4 (23%)

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
10	2G4	A	822	7	-	1/22/23/23	0/1/1/1
6	NAG	A	809	1	-	0/6/23/26	0/1/1/1
6	NAG	A	810	1	-	2/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	A	822	2G4	PAX-OAC	6.25	1.56	1.46
10	A	822	2G4	PAX-N	4.11	1.66	1.61
10	A	822	2G4	PAX-OAF	-2.51	1.50	1.56
10	A	822	2G4	CG-CD	2.27	1.55	1.50
10	A	822	2G4	CA-N	2.20	1.50	1.47

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	809	NAG	C1-O5-C5	8.11	123.18	112.19
6	A	810	NAG	O5-C5-C6	4.22	113.82	107.20
10	A	822	2G4	PAX-N-CA	-3.26	116.60	123.83
6	A	810	NAG	C3-C4-C5	-3.12	104.67	110.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	810	NAG	C1-O5-C5	-2.92	108.23	112.19
6	A	809	NAG	C6-C5-C4	-2.75	106.56	113.00
6	A	810	NAG	C1-C2-N2	-2.33	106.51	110.49
10	A	822	2G4	CG-CB-CA	-2.29	108.88	113.16
10	A	822	2G4	CAK-CAV-CAJ	2.19	121.61	118.17
6	A	809	NAG	O5-C5-C6	2.07	110.45	107.20

There are no chirality outliers.

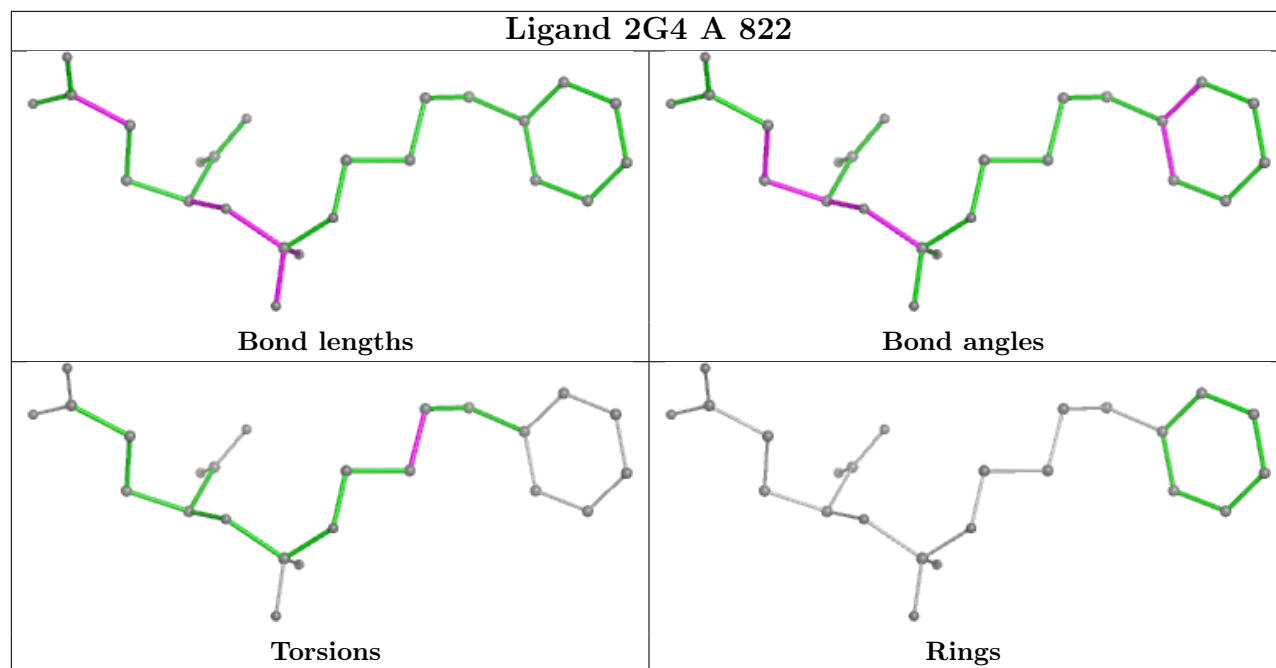
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	810	NAG	O5-C5-C6-O6
6	A	810	NAG	C4-C5-C6-O6
10	A	822	2G4	CAN-CAL-CAM-CAP

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 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	694/707 (98%)	-0.25	15 (2%) 62 66	10, 26, 48, 70	72 (10%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	541	TRP	5.0
1	A	656[A]	SER	4.2
1	A	719[A]	VAL	4.0
1	A	152	GLU	3.7
1	A	544	ASN	3.3
1	A	542	GLU	3.1
1	A	543	THR	2.8
1	A	506	PHE	2.8
1	A	154	VAL	2.7
1	A	102[A]	GLU	2.6
1	A	505	GLU	2.6
1	A	652	ASP	2.6
1	A	131[A]	ILE	2.5
1	A	653	PHE	2.3
1	A	155[A]	SER	2.1

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

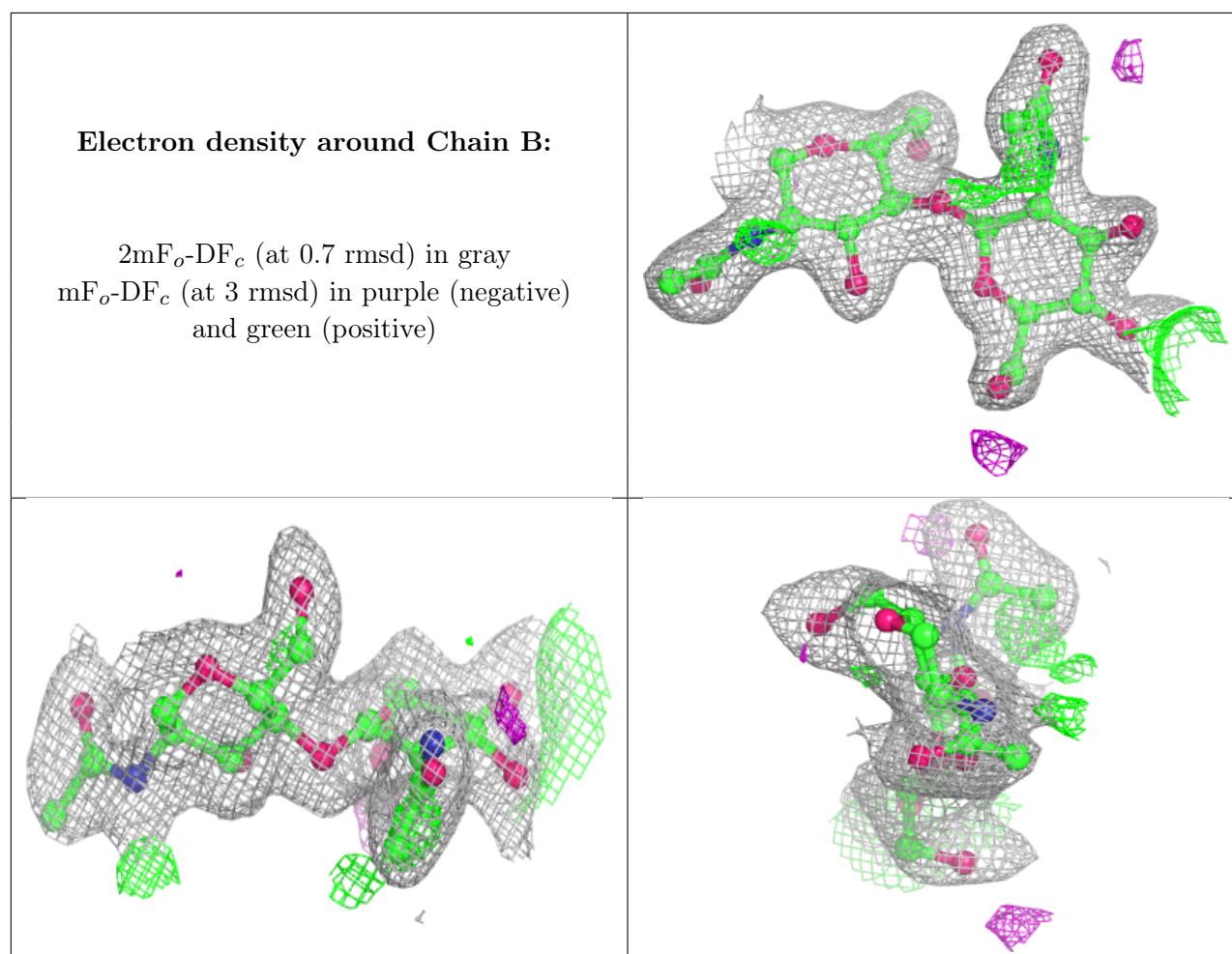
There are no non-standard protein/DNA/RNA residues in this entry.

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.

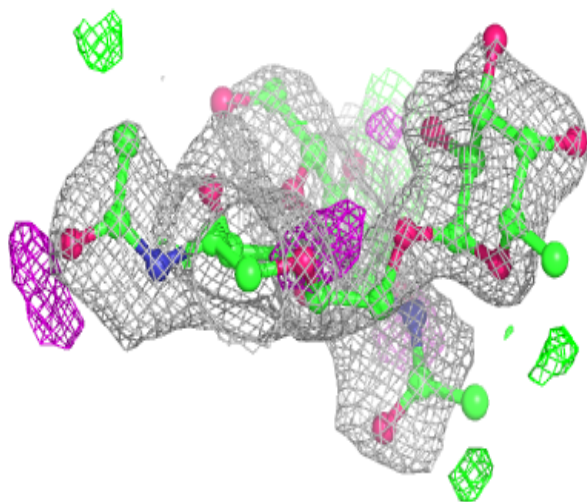
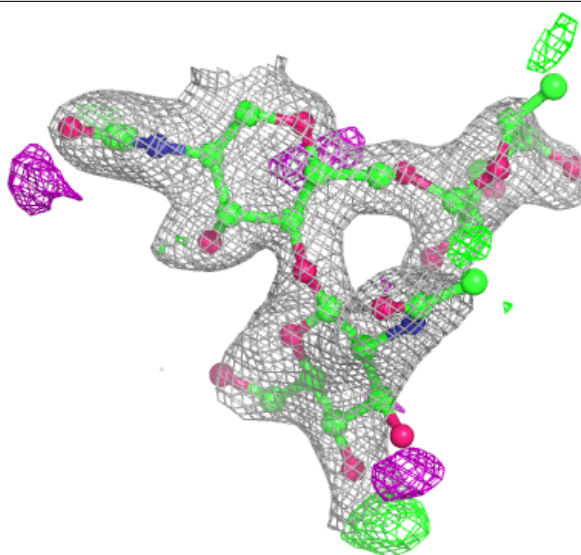
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	C	2	14/15	0.61	0.16	72,76,78,79	0
4	BMA	E	3	11/12	0.66	0.16	61,66,70,71	0
2	NAG	B	2	14/15	0.79	0.13	44,51,59,60	0
3	NAG	D	2	14/15	0.80	0.13	59,65,70,71	0
3	FUC	D	3	10/11	0.81	0.17	56,62,64,66	0
3	NAG	C	1	14/15	0.83	0.12	48,55,67,73	0
5	NAG	F	2	14/15	0.84	0.15	34,40,50,53	0
3	FUC	C	3	10/11	0.85	0.13	64,68,69,69	0
4	NAG	E	2	14/15	0.85	0.13	37,45,53,53	0
5	MAN	F	4	11/12	0.85	0.12	47,50,54,55	0
5	BMA	F	3	11/12	0.90	0.10	39,43,46,47	0
3	NAG	D	1	14/15	0.93	0.09	45,49,56,59	0
4	NAG	E	1	14/15	0.94	0.08	29,34,41,41	0
2	NAG	B	1	14/15	0.94	0.08	35,43,49,52	0
5	NAG	F	1	14/15	0.96	0.07	23,29,38,44	0

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



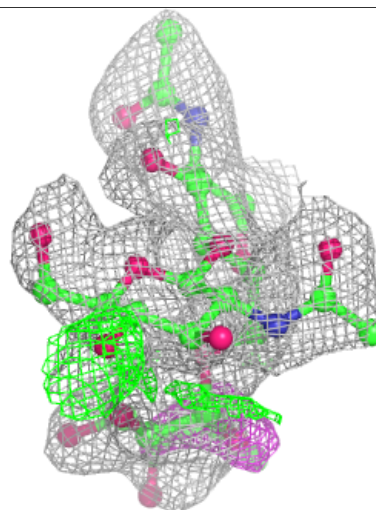
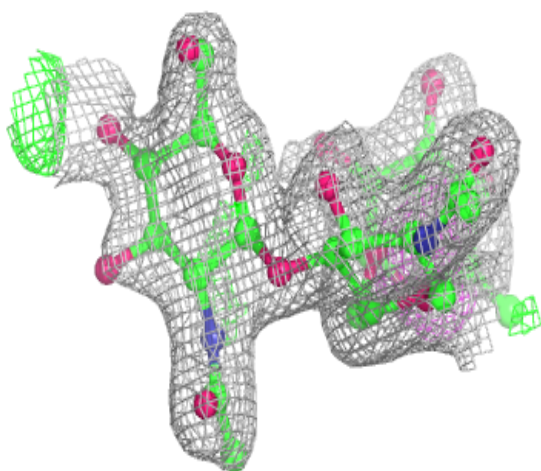
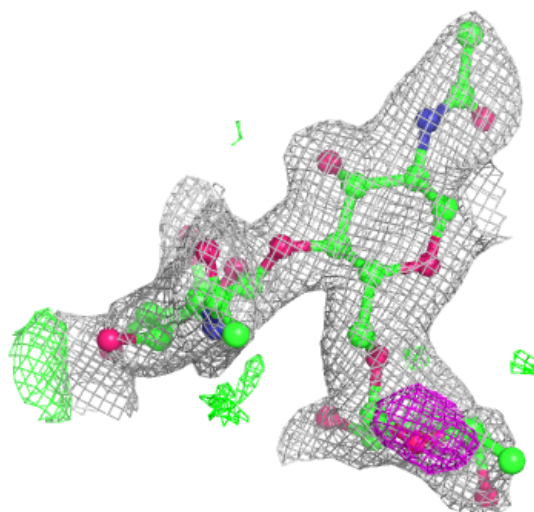
Electron density around Chain C:

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



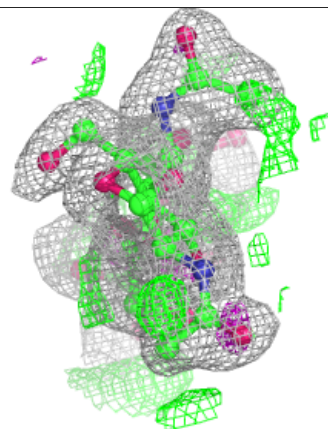
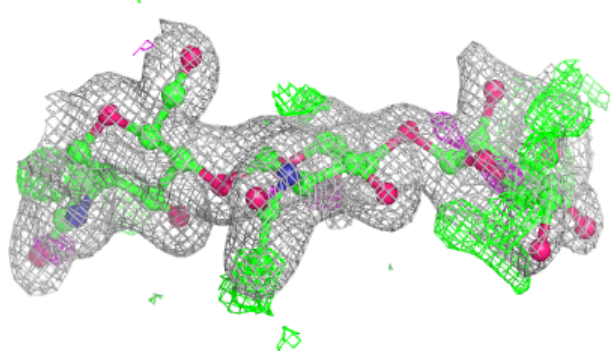
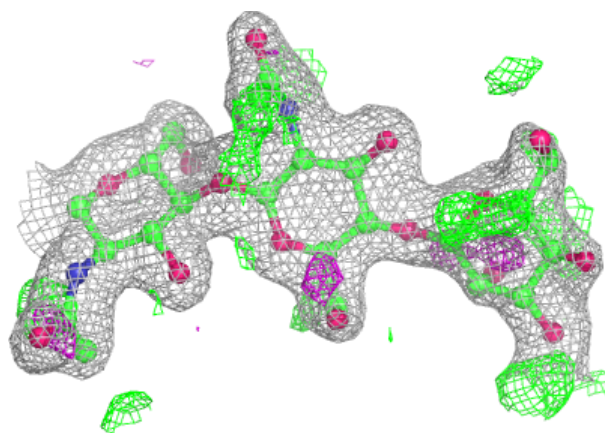
Electron density around Chain D:

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

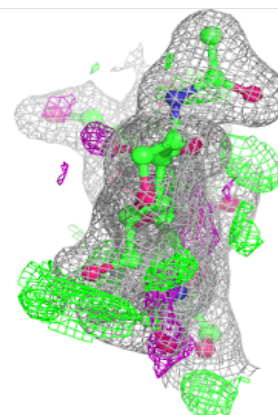
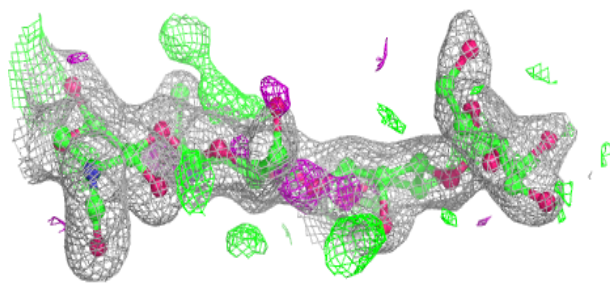
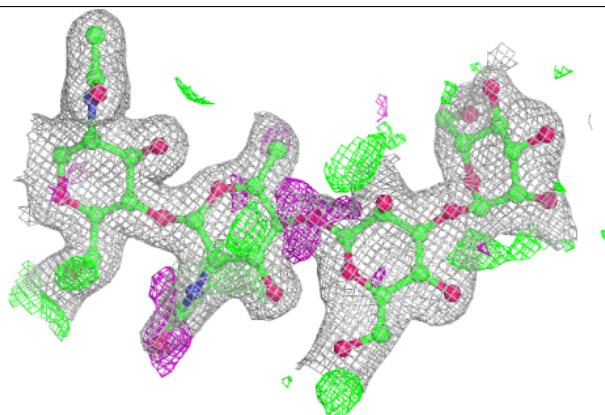


Electron density around Chain E:

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

**Electron density around Chain F:**

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



6.4 Ligands [i](#)

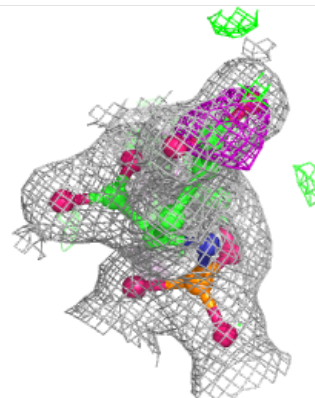
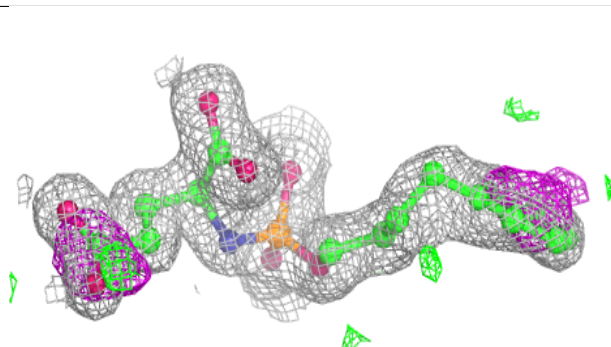
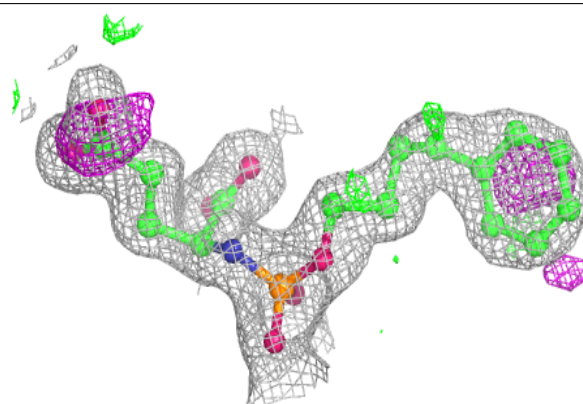
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	NAG	A	809	14/15	0.77	0.14	69,75,77,77	0
6	NAG	A	810	14/15	0.88	0.10	33,44,51,54	0
10	2G4	A	822	24/24	0.98	0.08	20,24,43,45	0
9	CL	A	821	1/1	0.99	0.07	28,28,28,28	0
8	CA	A	820	1/1	1.00	0.01	18,18,18,18	0
7	ZN	A	818	1/1	1.00	0.01	20,20,20,20	0
7	ZN	A	819	1/1	1.00	0.01	21,21,21,21	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 2G4 A 822:

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.