



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 12:55 PM UTC

PDB ID : 7ZYF / pdb_00007zyf
Title : Insulin regulated aminopeptidase (IRAP) in complex with a nanomolar alpha hydroxy beta amino acid based inhibitor.
Authors : Mpakali, A.; Stratikos, E.; Giastas, P.; Papakyriakou, A.
Deposited on : 2022-05-24
Resolution : 2.81 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

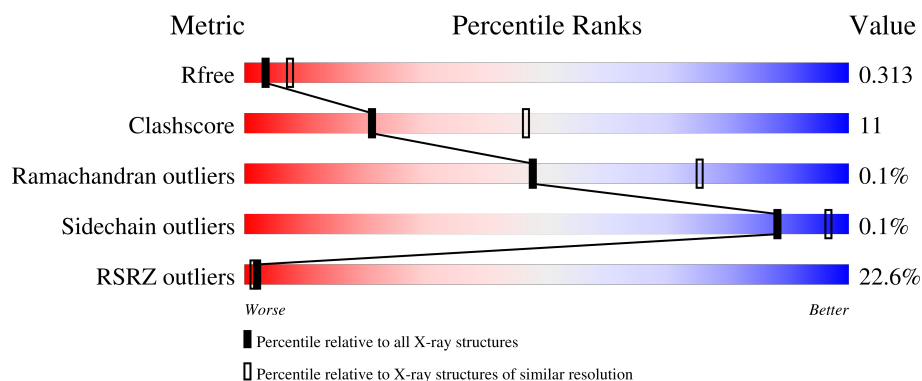
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	871	18% 77% 21%
1	B	871	26% 69% 29%
2	C	4	50% 50%
3	D	2	100%
3	F	2	100%

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Mol	Chain	Length	Quality of chain
3	G	2	 100%
3	K	2	 50% 50%
4	E	3	 33% 67%
4	J	3	 100%
5	H	5	 20% 80%
5	I	5	 40% 60%

2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 14588 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 Leucyl-cystinyl aminopeptidase, pregnancy serum form.

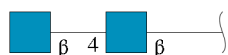
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	852	Total	C	N	O	S	0	4	0
			6910	4481	1123	1279	27			
1	B	854	Total	C	N	O	S	0	2	0
			6882	4457	1122	1277	26			

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

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



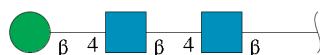
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

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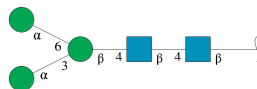
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	K	2	Total	C	N	O	0	0	0
			28	16	2	10			

- 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			
4	J	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is an oligosaccharide called 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
5	H	5	Total	C	N	O	0	0	0
			61	34	2	25			
5	I	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn).

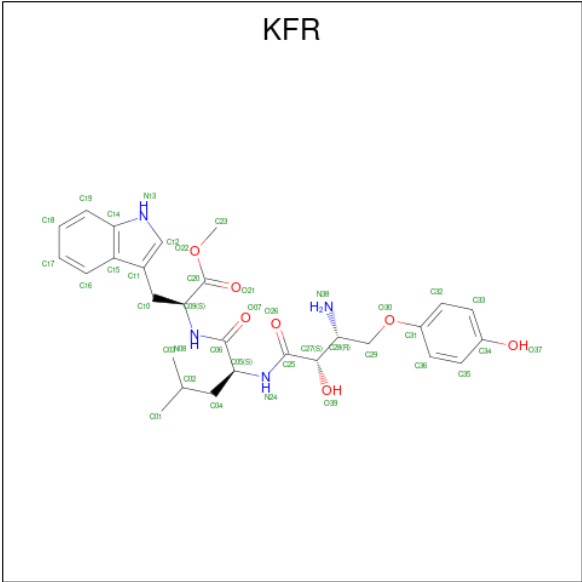
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Zn	0	0
			1	1		
6	B	1	Total	Zn	0	0
			1	1		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



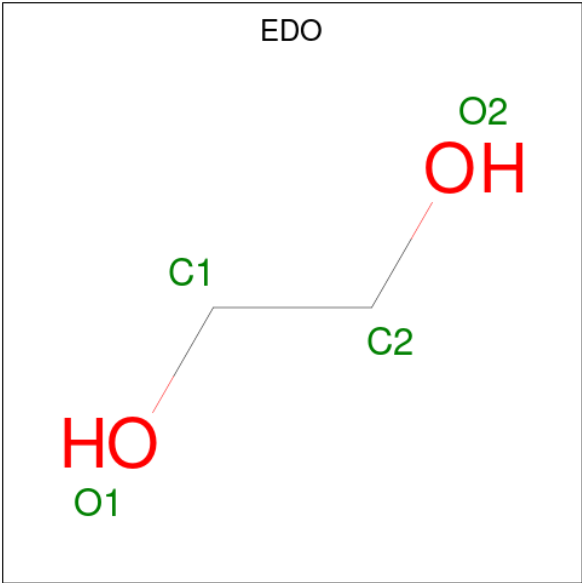
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1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

- Molecule 8 is methyl (2S)-2-[[[(2S)-2-[[[(2S,3R)-3-azanyl-2-oxidanyl-4-(4-oxidanylphenoxy)butanoyl]amino]-4-methyl-pentanoyl]amino]-3-(1H-indol-3-yl)propanoate (CCD ID: KFR) (formula: C₂₈H₃₆N₄O₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			39	28	4	7		
8	B	1	Total	C	N	O	0	1
			78	56	8	14		

- Molecule 9 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



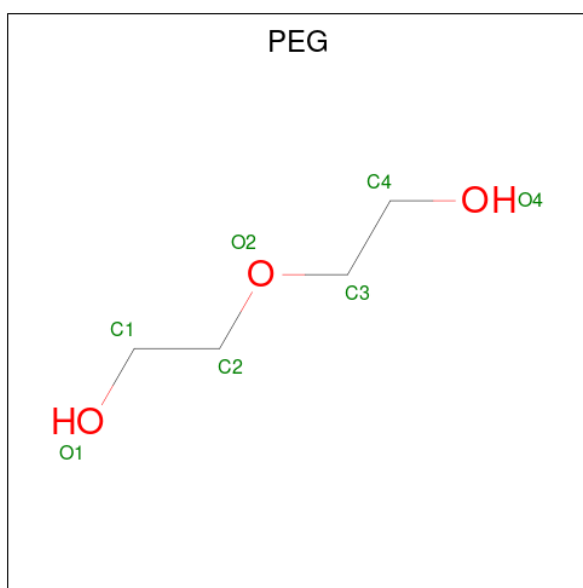
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		

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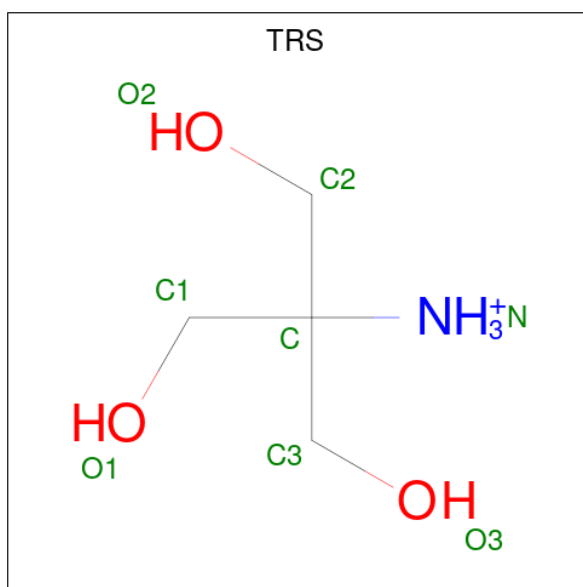
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			4	2	2		
9	B	1	Total	C	O	0	0
			4	2	2		
9	B	1	Total	C	O	0	0
			4	2	2		
9	B	1	Total	C	O	0	0
			4	2	2		
9	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 10 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

- Molecule 11 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	B	1	Total	C	N	O	0	0
			8	4	1	3		

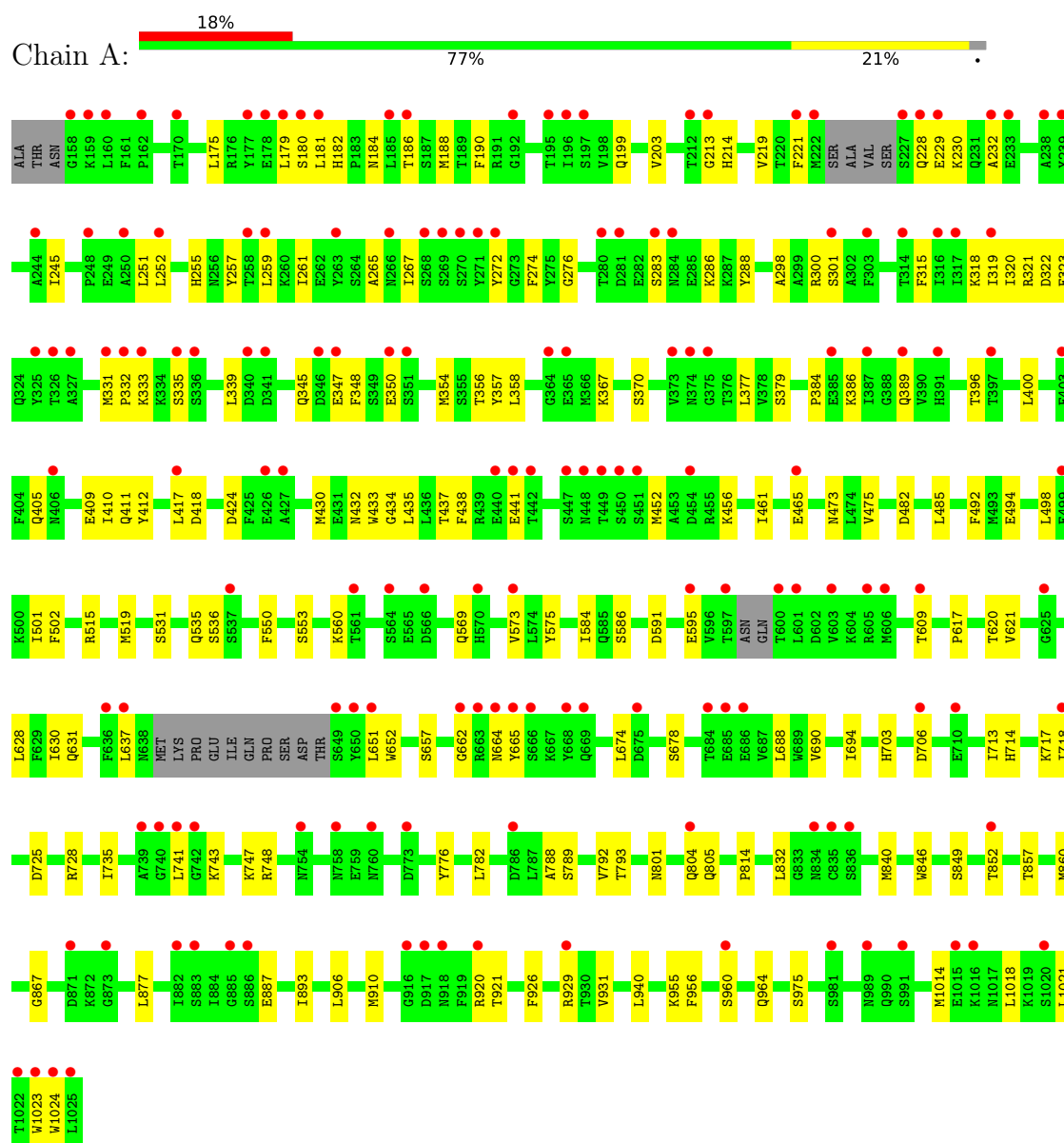
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	51	Total	O	0	0
			51	51		
12	B	53	Total	O	0	0
			53	53		

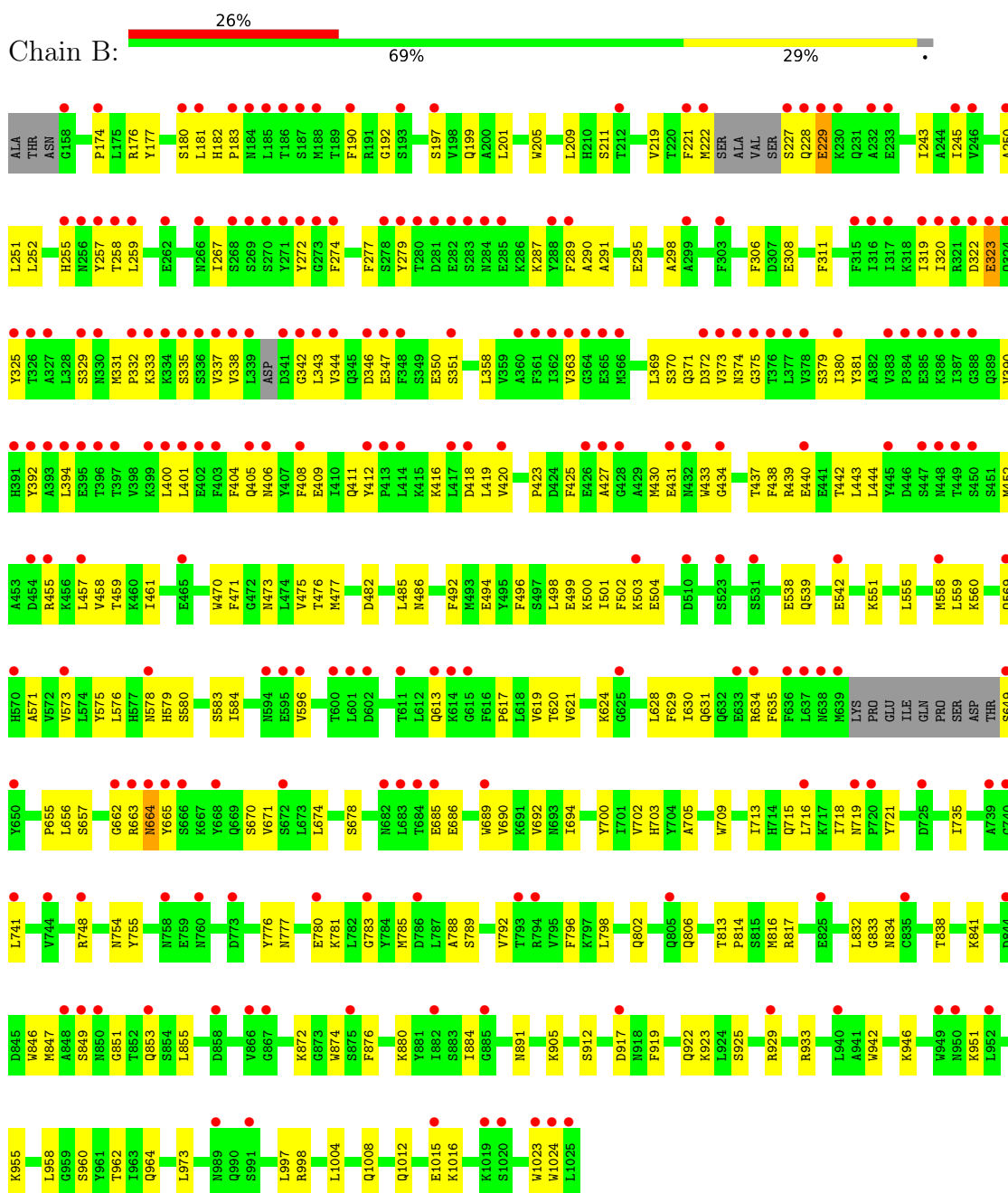
3 Residue-property plots

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.

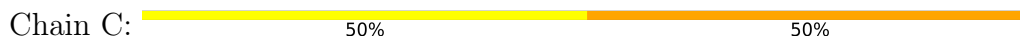
- Molecule 1: Leucyl-cystinyl aminopeptidase, pregnancy serum form



- Molecule 1: Leucyl-cystinyl aminopeptidase, pregnancy serum form



- Molecule 2: 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



MAG1
MAG2
EMA3
MAV4

- Molecule 3: 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)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%



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

Chain G:  100%



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

Chain K:  50% 50%



- 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:  33% 67%



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

Chain J:  100%



- Molecule 5: 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 H:  20% 80%



- Molecule 5: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain I:



MAG1	MAG2	EMJ3	MAT4	MAT5
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.48Å 120.05Å 141.25Å 90.00° 103.28° 90.00°	Depositor
Resolution (Å)	137.47 – 2.81 137.47 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.5 (137.47-2.81) 99.6 (137.47-2.81)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.82Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660, PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.271 , 0.314 0.271 , 0.313	Depositor DCC
R_{free} test set	2883 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	35.3	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 59.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.004 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	14588	wwPDB-VP
Average B, all atoms (Å ²)	39.0	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 28.40 % 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 1.8538e-03. 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, EDO, ZN, PEG, KFR, TRS, BMA, NAG

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	0.21	0/7091	0.50	0/9615
1	B	0.24	0/7057	0.56	4/9576 (0.0%)
All	All	0.23	0/14148	0.53	4/19191 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	227	SER	CA-C-N	7.36	135.59	121.54
1	B	227	SER	C-N-CA	7.36	135.59	121.54
1	B	664	ASN	N-CA-C	-5.38	104.83	113.28
1	B	374	ASN	CB-CA-C	-5.25	102.51	111.02

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	6910	0	6805	130	0
1	B	6882	0	6717	180	0
2	C	50	0	43	4	0
3	D	28	0	25	0	0
3	F	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	28	0	25	0	0
3	K	28	0	25	2	0
4	E	39	0	34	0	0
4	J	39	0	34	0	0
5	H	61	0	52	1	0
5	I	61	0	52	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	56	0	52	4	0
7	B	98	0	89	8	0
8	A	39	0	0	0	0
8	B	78	0	0	4	0
9	A	12	0	18	0	0
9	B	16	0	24	0	0
10	A	14	0	20	0	0
10	B	7	0	10	0	0
11	B	8	0	12	0	0
12	A	51	0	0	0	0
12	B	53	0	0	1	0
All	All	14588	0	14062	315	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 11.

The worst 5 of 315 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:441:GLU:HG2	1:A:887:GLU:HB3	1.46	0.97
1:B:735:ILE:HD11	1:B:748:ARG:HG3	1.51	0.92
1:B:833:GLY:HA3	3:K:1:NAG:H82	1.55	0.88
1:B:847:MET:HE1	1:B:872:LYS:HG2	1.62	0.82
1:A:245:ILE:HD13	1:A:259:LEU:HD21	1.62	0.81

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	848/871 (97%)	821 (97%)	27 (3%)	0	100	100
1	B	848/871 (97%)	812 (96%)	34 (4%)	2 (0%)	43	71
All	All	1696/1742 (97%)	1633 (96%)	61 (4%)	2 (0%)	48	75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	229	GLU
1	B	323	GLU

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	756/780 (97%)	755 (100%)	1 (0%)	88	96
1	B	745/780 (96%)	745 (100%)	0	100	100
All	All	1501/1560 (96%)	1500 (100%)	1 (0%)	88	96

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	283	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1001	GLN
1	B	733	ASN
1	B	324	GLN
1	B	965	ASN
1	B	715	GLN

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 ⓘ

28 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	1,2	14,14,15	0.59	0	17,19,21	1.45	2 (11%)
2	NAG	C	2	2	14,14,15	1.02	1 (7%)	17,19,21	0.92	1 (5%)
2	BMA	C	3	2	11,11,12	0.81	0	15,15,17	1.73	5 (33%)
2	MAN	C	4	2	11,11,12	0.54	0	15,15,17	0.96	2 (13%)
3	NAG	D	1	3,1	14,14,15	0.48	0	17,19,21	0.54	0
3	NAG	D	2	3	14,14,15	0.33	0	17,19,21	0.43	0
4	NAG	E	1	4,1	14,14,15	0.48	0	17,19,21	0.77	1 (5%)
4	NAG	E	2	4	14,14,15	0.31	0	17,19,21	0.58	0
4	BMA	E	3	4	11,11,12	1.12	1 (9%)	15,15,17	1.15	2 (13%)
3	NAG	F	1	3,1	14,14,15	0.57	0	17,19,21	0.71	0
3	NAG	F	2	3	14,14,15	0.54	0	17,19,21	0.53	0
3	NAG	G	1	3,1	14,14,15	0.55	0	17,19,21	0.83	0
3	NAG	G	2	3	14,14,15	0.44	0	17,19,21	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	H	1	5,1	14,14,15	0.29	0	17,19,21	0.81	0
5	NAG	H	2	5	14,14,15	0.20	0	17,19,21	0.53	0
5	BMA	H	3	5	11,11,12	1.46	2 (18%)	15,15,17	1.18	1 (6%)
5	MAN	H	4	5	11,11,12	0.79	1 (9%)	15,15,17	1.50	3 (20%)
5	MAN	H	5	5	11,11,12	0.90	1 (9%)	15,15,17	1.17	2 (13%)
5	NAG	I	1	5,1	14,14,15	0.37	0	17,19,21	0.55	0
5	NAG	I	2	5	14,14,15	0.39	0	17,19,21	0.54	0
5	BMA	I	3	5	11,11,12	0.62	0	15,15,17	0.77	0
5	MAN	I	4	5	11,11,12	0.72	0	15,15,17	1.01	2 (13%)
5	MAN	I	5	5	11,11,12	0.84	0	15,15,17	1.01	2 (13%)
4	NAG	J	1	4,1	14,14,15	0.33	0	17,19,21	0.77	0
4	NAG	J	2	4	14,14,15	0.46	0	17,19,21	0.69	0
4	BMA	J	3	4	11,11,12	0.60	0	15,15,17	0.77	0
3	NAG	K	1	3,1	14,14,15	0.36	0	17,19,21	0.70	0
3	NAG	K	2	3	14,14,15	0.63	0	17,19,21	0.45	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	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	3/6/23/26	0/1/1/1
2	BMA	C	3	2	-	2/2/19/22	0/1/1/1
2	MAN	C	4	2	-	2/2/19/22	0/1/1/1
3	NAG	D	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	3/6/23/26	0/1/1/1
4	NAG	E	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	1/6/23/26	0/1/1/1
4	BMA	E	3	4	-	0/2/19/22	0/1/1/1
3	NAG	F	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	NAG	G	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	G	2	3	-	1/6/23/26	0/1/1/1
5	NAG	H	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	H	2	5	-	3/6/23/26	0/1/1/1
5	BMA	H	3	5	-	1/2/19/22	0/1/1/1
5	MAN	H	4	5	-	1/2/19/22	0/1/1/1

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

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	3	BMA	C2-C3	3.73	1.58	1.52
2	C	2	NAG	O5-C1	-3.51	1.37	1.43
5	H	4	MAN	C1-C2	2.21	1.57	1.52
5	H	3	BMA	O3-C3	2.16	1.48	1.43
4	E	3	BMA	C4-C3	2.05	1.57	1.52

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	O4-C4-C5	-4.29	98.75	109.32
5	H	4	MAN	C1-O5-C5	4.15	117.75	112.19
2	C	3	BMA	C1-O5-C5	3.79	117.27	112.19
5	H	3	BMA	O3-C3-C2	3.46	117.12	110.05
5	H	5	MAN	C1-O5-C5	3.30	116.61	112.19

There are no chirality outliers.

5 of 37 torsion outliers are listed below:

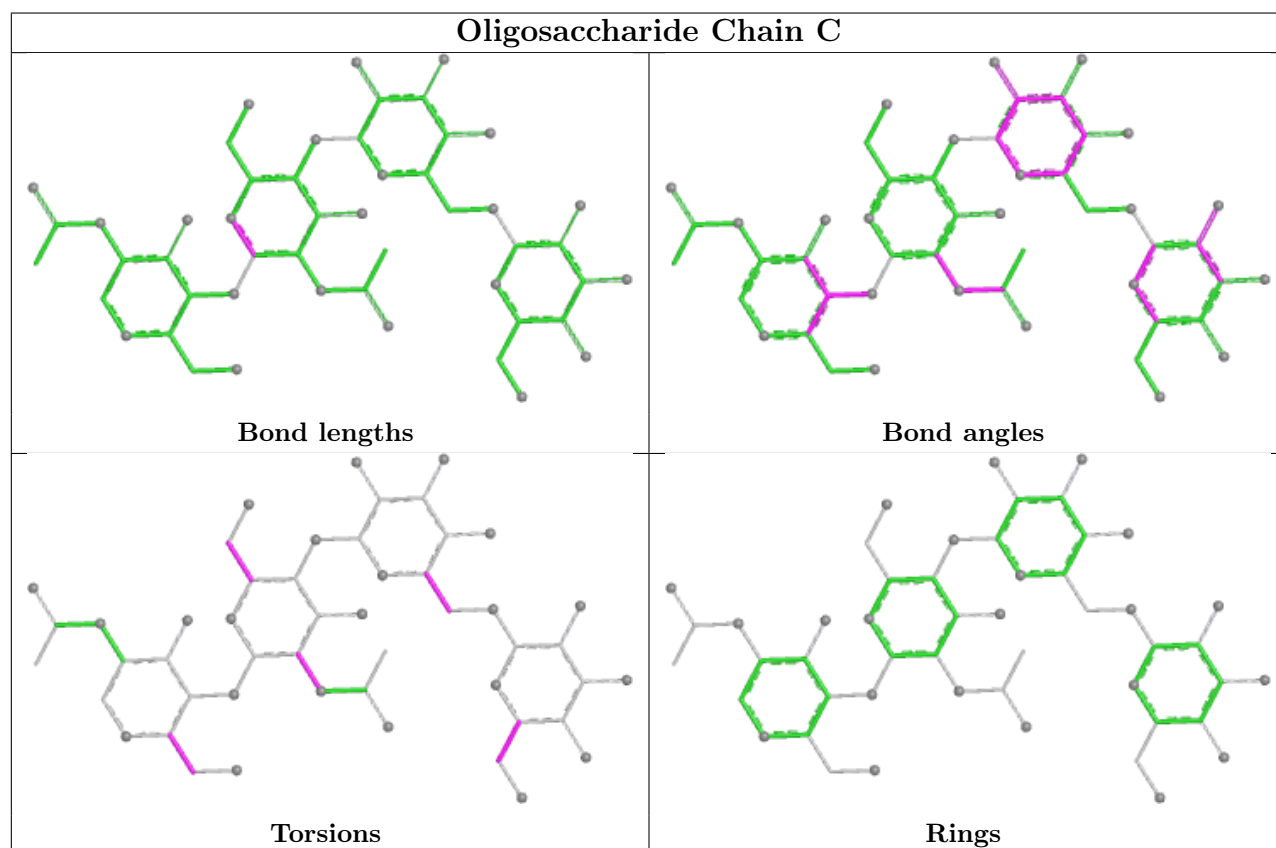
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	C3-C2-N2-C7
5	H	2	NAG	C1-C2-N2-C7
5	I	2	NAG	O5-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6

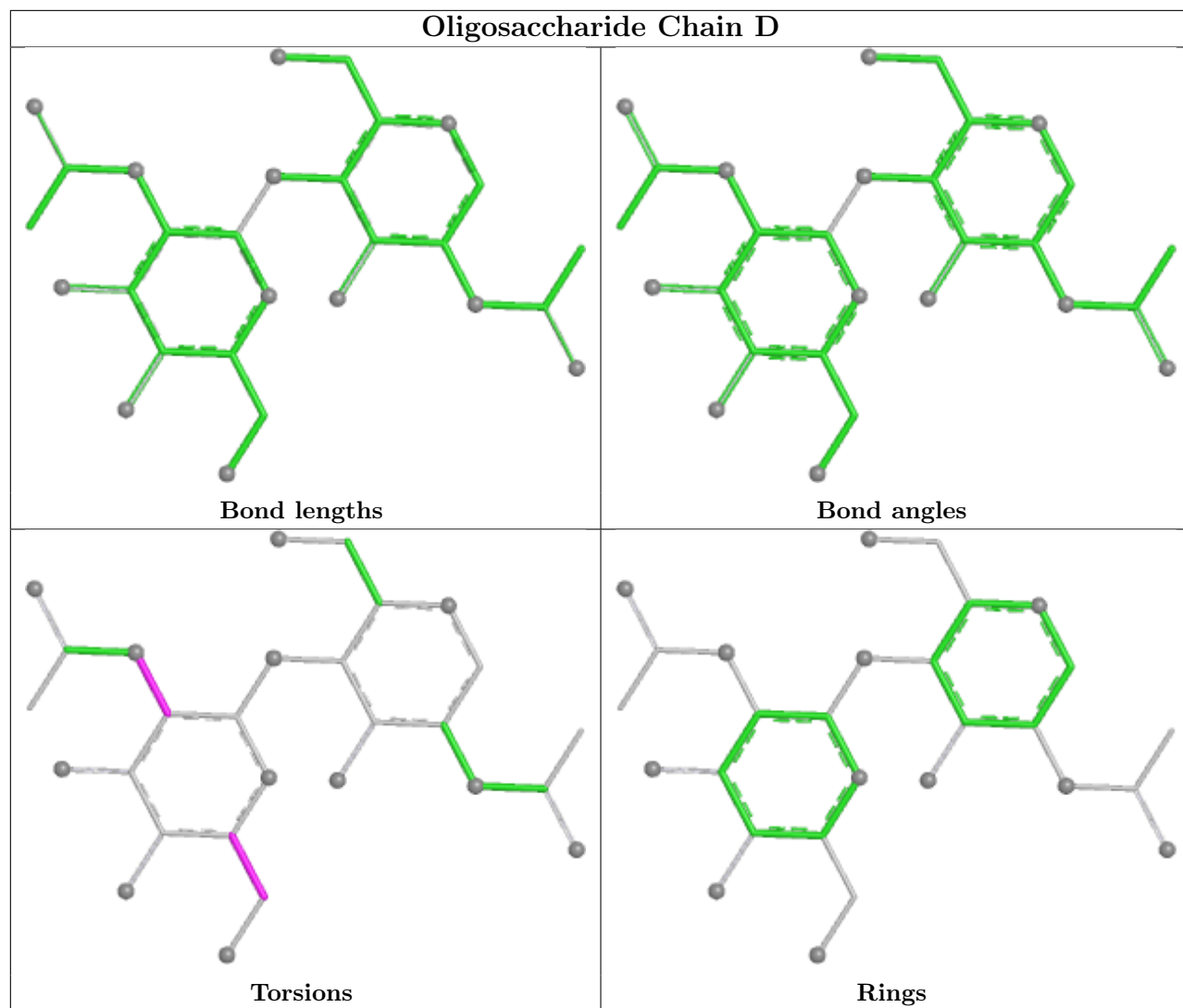
There are no ring outliers.

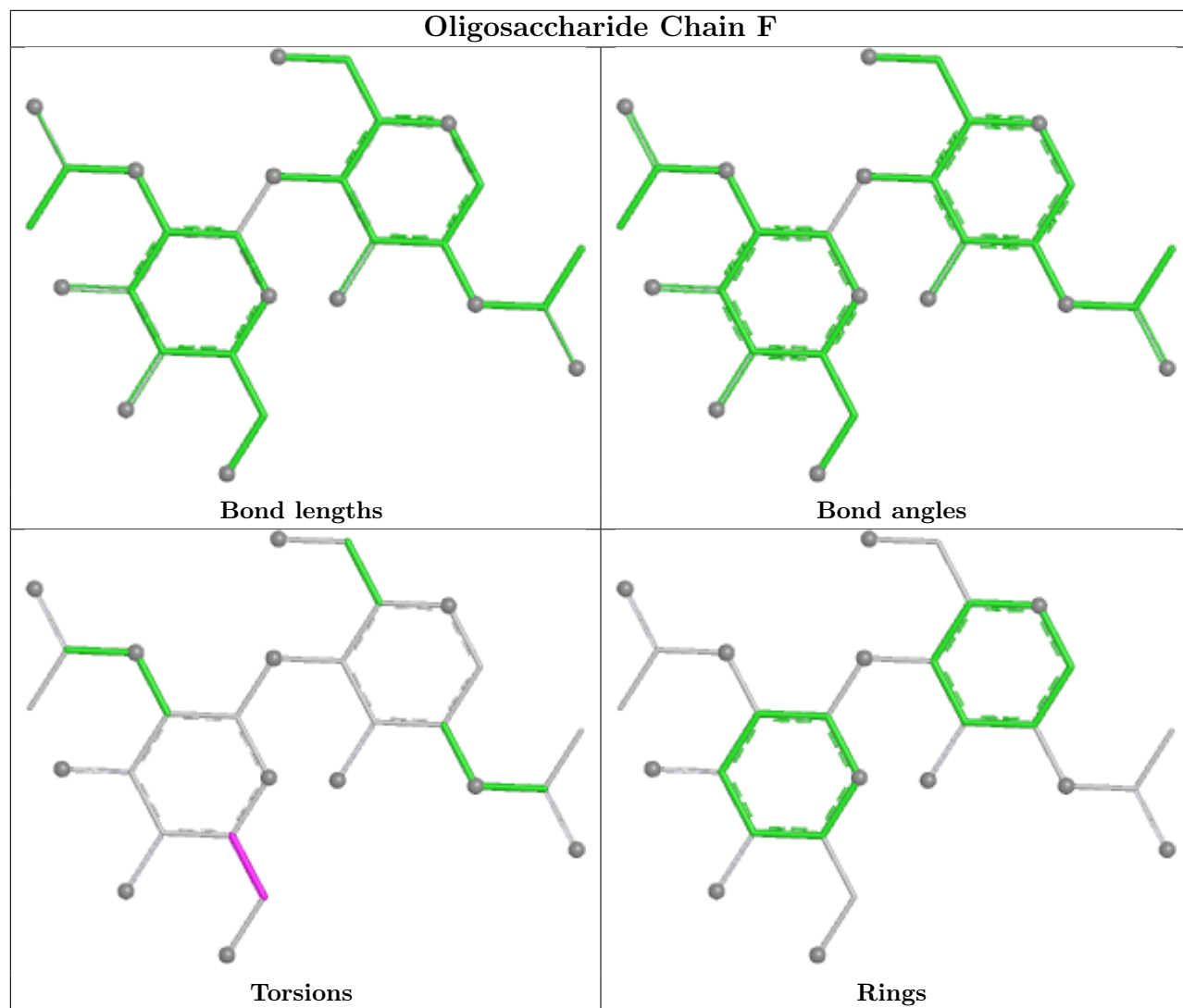
5 monomers are involved in 8 short contacts:

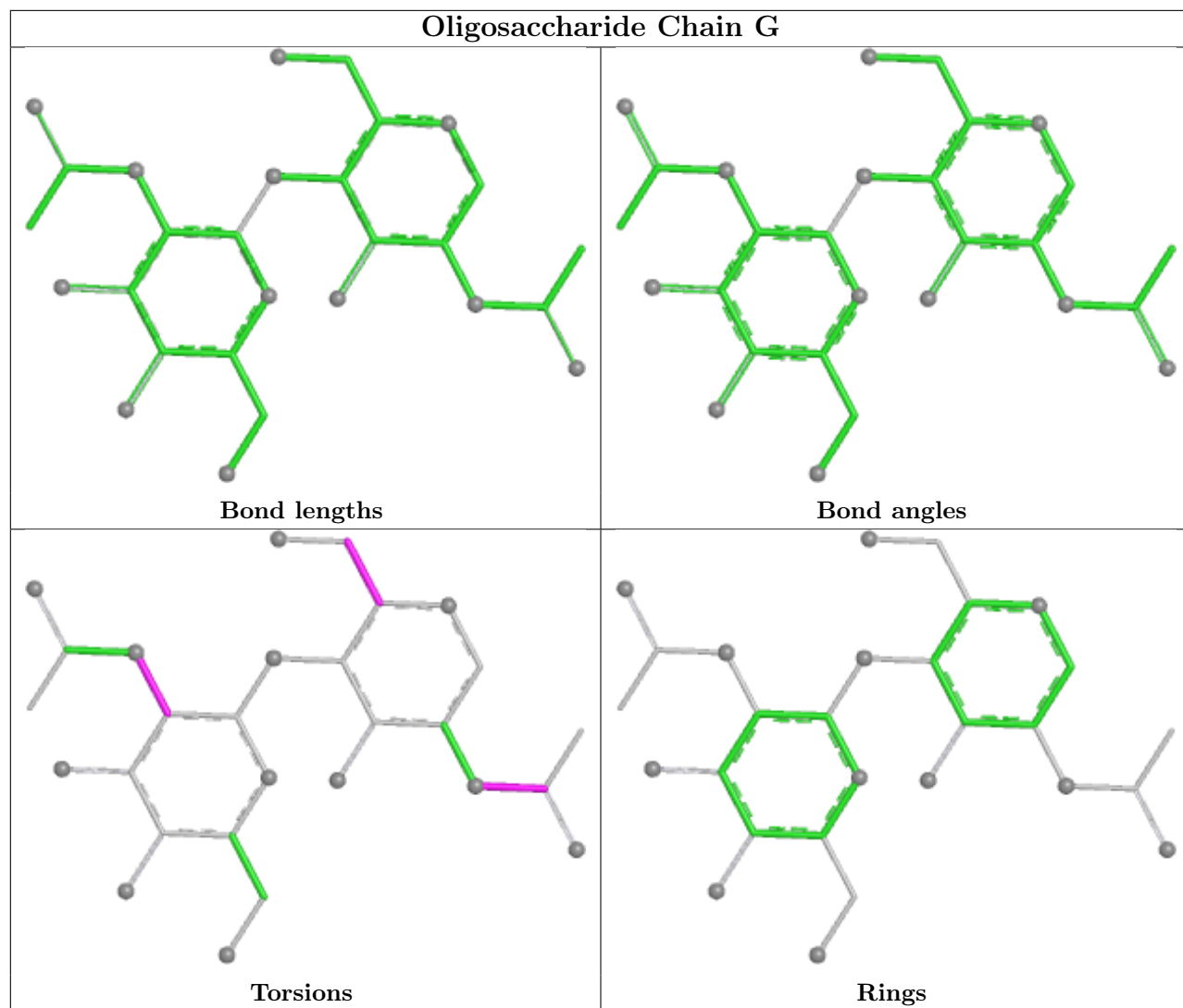
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	K	1	NAG	2	0
5	H	1	NAG	1	0
5	I	1	NAG	1	0
2	C	2	NAG	4	0
2	C	3	BMA	1	0

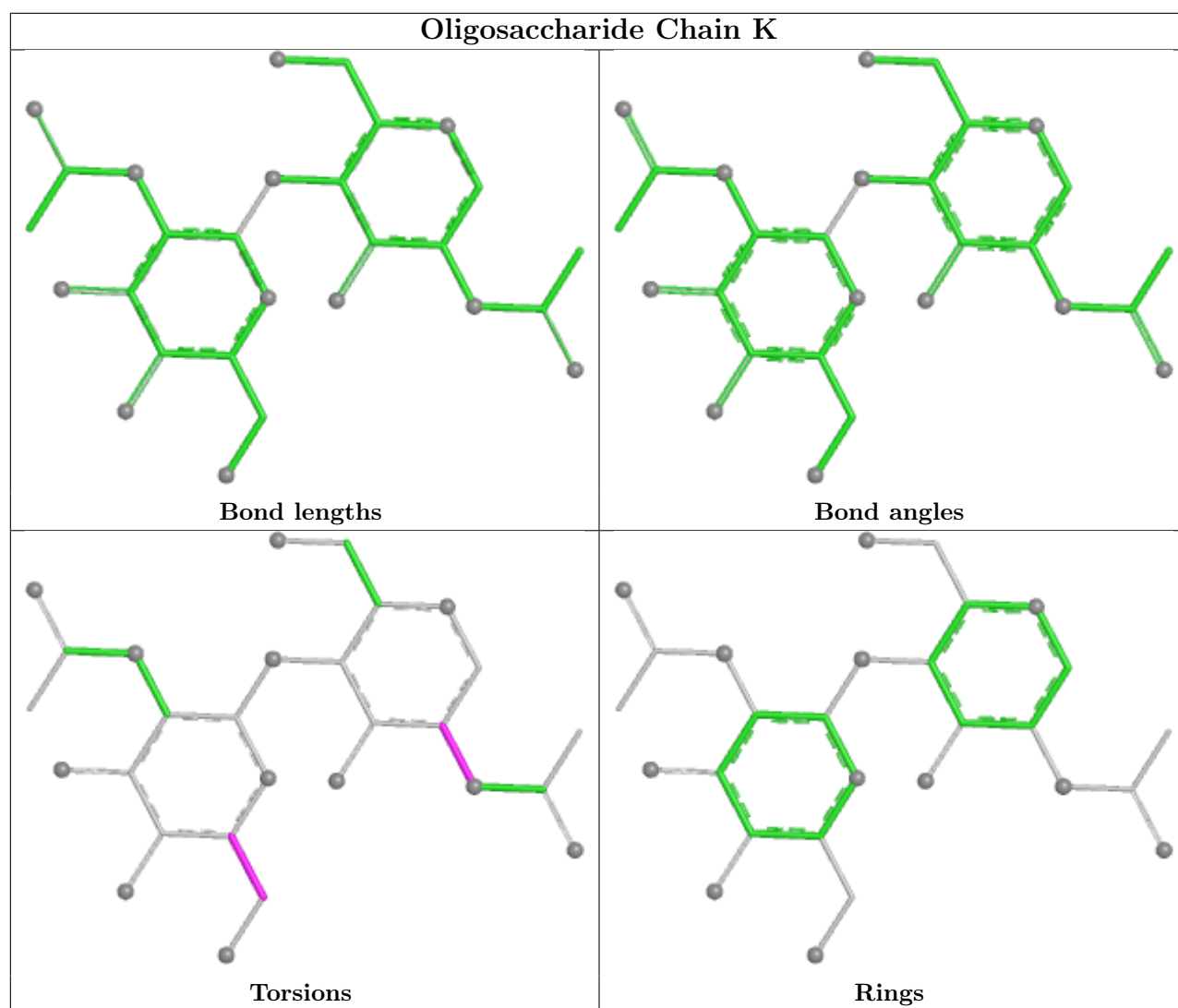
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

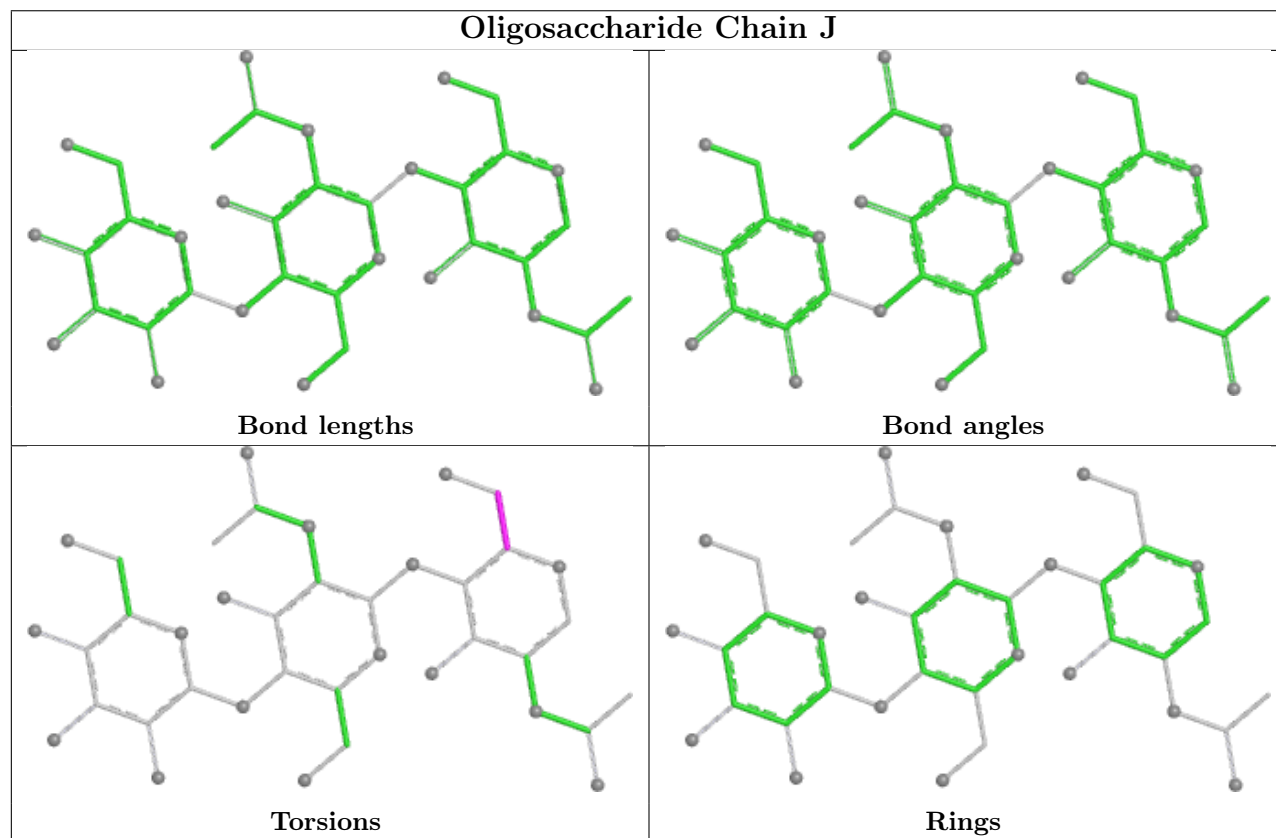
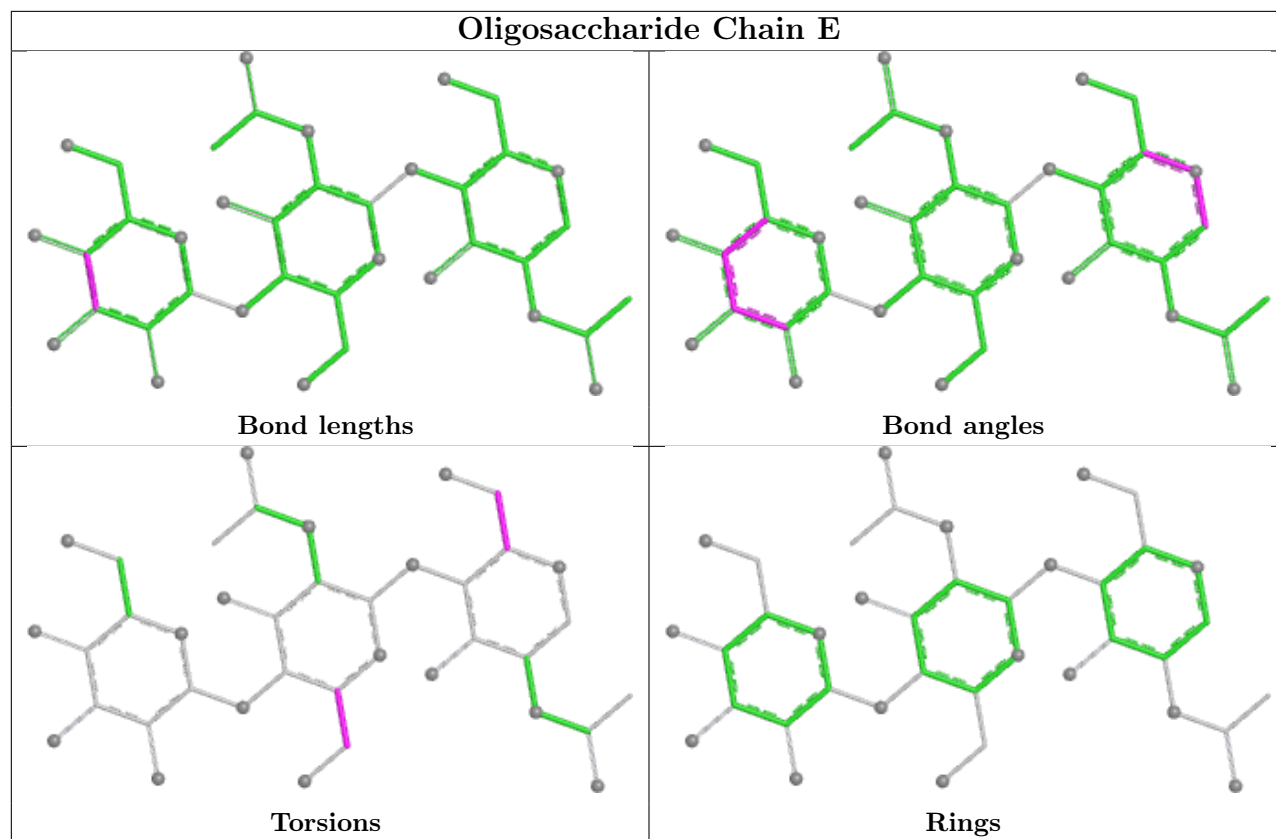


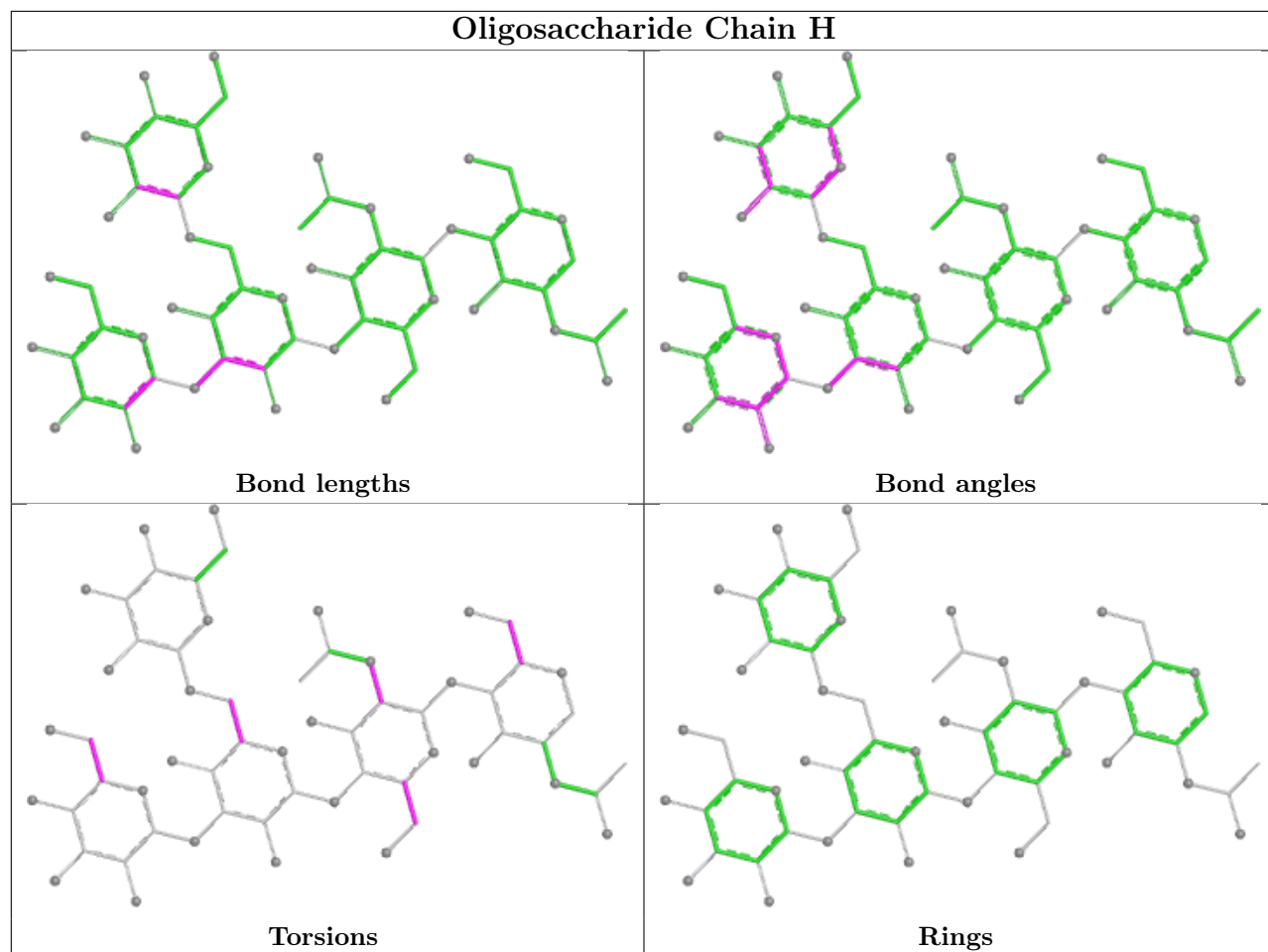


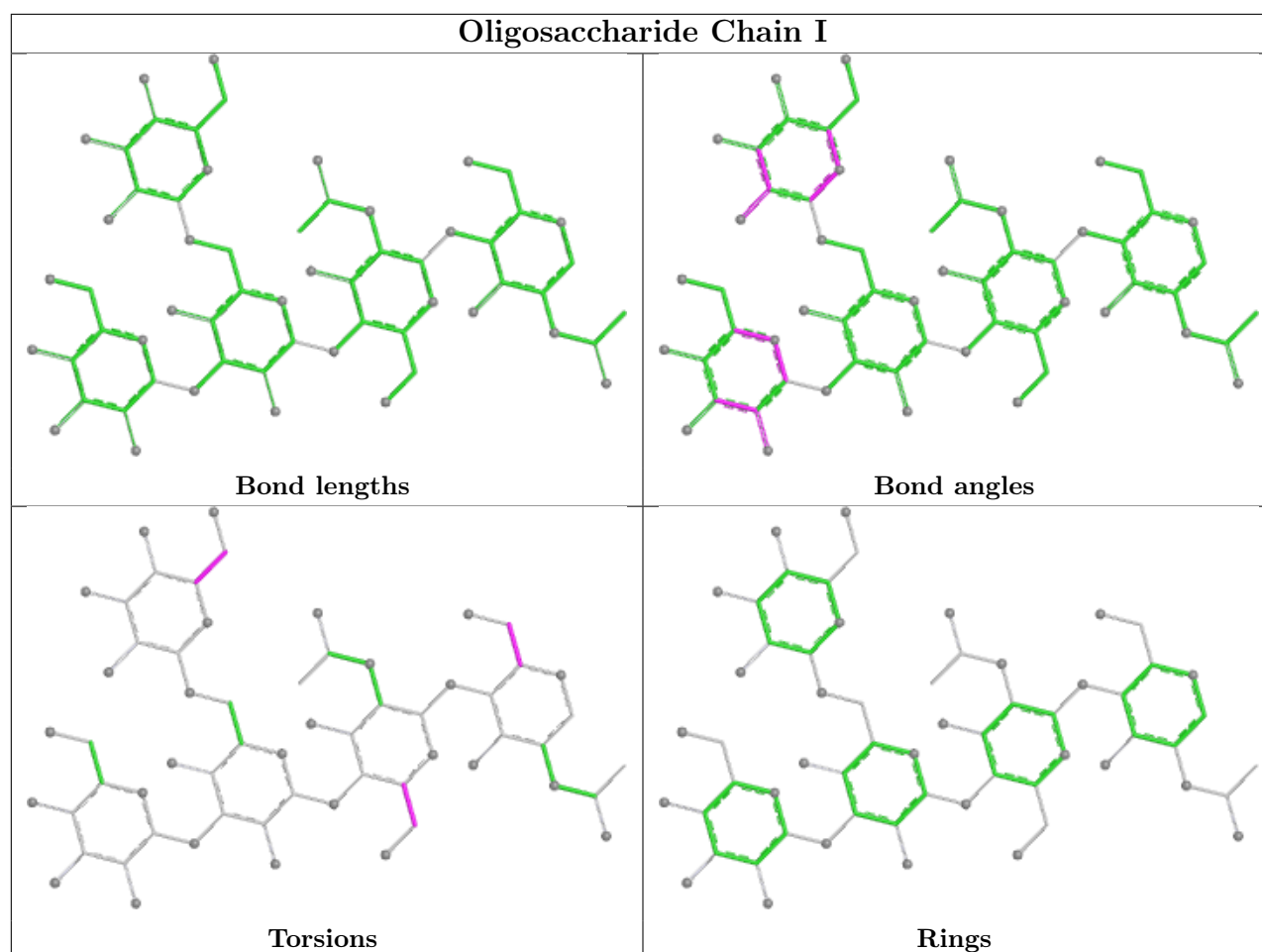












5.6 Ligand geometry [i](#)

Of 27 ligands modelled in this entry, 2 are monoatomic - leaving 25 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$
9	EDO	A	1108	-	3,3,3	0.42	0	2,2,2	0.39	0
7	NAG	B	1803	1	14,14,15	1.10	2 (14%)	17,19,21	0.84	1 (5%)
9	EDO	B	1808	-	3,3,3	0.42	0	2,2,2	0.43	0
9	EDO	A	1109	-	3,3,3	0.43	0	2,2,2	0.39	0
7	NAG	B	1813	1,7	14,14,15	0.48	0	17,19,21	0.86	1 (5%)
9	EDO	B	1805	-	3,3,3	0.43	0	2,2,2	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NAG	B	1811	7	14,14,15	0.63	1 (7%)	17,19,21	0.94	0
9	EDO	A	1107	-	3,3,3	0.45	0	2,2,2	0.34	0
9	EDO	B	1806	-	3,3,3	0.44	0	2,2,2	0.40	0
10	PEG	A	1111	-	6,6,6	0.50	0	5,5,5	0.28	0
9	EDO	B	1804	-	3,3,3	0.43	0	2,2,2	0.41	0
8	KFR	A	1106	6	41,41,41	1.73	8 (19%)	53,56,56	1.07	3 (5%)
7	NAG	B	1814	1	14,14,15	0.72	1 (7%)	17,19,21	1.74	1 (5%)
11	TRS	B	1807	-	7,7,7	0.33	0	9,9,9	0.35	0
7	NAG	B	1810	1,7	14,14,15	0.79	0	17,19,21	1.41	1 (5%)
7	NAG	A	1105	1	14,14,15	0.31	0	17,19,21	0.62	0
7	NAG	A	1103	1	14,14,15	0.36	0	17,19,21	0.48	0
7	NAG	A	1102	1	14,14,15	0.19	0	17,19,21	0.57	0
8	KFR	B	1815[A]	6	41,41,41	1.76	8 (19%)	53,56,56	1.03	4 (7%)
7	NAG	B	1802	1	14,14,15	1.01	1 (7%)	17,19,21	0.80	1 (5%)
10	PEG	B	1809	-	6,6,6	0.51	0	5,5,5	0.28	0
7	NAG	B	1812	7	14,14,15	0.66	1 (7%)	17,19,21	0.70	0
7	NAG	A	1104	1	14,14,15	0.87	1 (7%)	17,19,21	0.89	0
8	KFR	B	1815[B]	6	41,41,41	1.77	8 (19%)	53,56,56	1.05	4 (7%)
10	PEG	A	1110	-	6,6,6	0.51	0	5,5,5	0.30	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
9	EDO	A	1108	-	-	0/1/1/1	-
7	NAG	B	1803	1	-	2/6/23/26	0/1/1/1
9	EDO	B	1808	-	-	0/1/1/1	-
9	EDO	A	1109	-	-	0/1/1/1	-
7	NAG	B	1813	1,7	-	1/6/23/26	0/1/1/1
9	EDO	B	1805	-	-	0/1/1/1	-
7	NAG	B	1811	7	-	1/6/23/26	0/1/1/1
9	EDO	A	1107	-	-	1/1/1/1	-
9	EDO	B	1806	-	-	1/1/1/1	-
10	PEG	A	1111	-	-	3/4/4/4	-
9	EDO	B	1804	-	-	1/1/1/1	-
8	KFR	A	1106	6	-	15/39/48/48	0/3/3/3
7	NAG	B	1814	1	-	3/6/23/26	0/1/1/1
11	TRS	B	1807	-	-	0/9/9/9	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	B	1810	1,7	-	4/6/23/26	0/1/1/1
7	NAG	A	1105	1	-	3/6/23/26	0/1/1/1
7	NAG	A	1103	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1102	1	-	0/6/23/26	0/1/1/1
8	KFR	B	1815[A]	6	-	13/39/48/48	0/3/3/3
7	NAG	B	1802	1	-	2/6/23/26	0/1/1/1
10	PEG	B	1809	-	-	0/4/4/4	-
7	NAG	B	1812	7	-	1/6/23/26	0/1/1/1
7	NAG	A	1104	1	-	3/6/23/26	0/1/1/1
8	KFR	B	1815[B]	6	-	9/39/48/48	0/3/3/3
10	PEG	A	1110	-	-	0/4/4/4	-

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	1815[B]	KFR	C06-N08	5.66	1.46	1.34
8	B	1815[A]	KFR	C06-N08	5.60	1.46	1.34
8	B	1815[B]	KFR	C25-N24	5.50	1.45	1.34
8	A	1106	KFR	C06-N08	5.49	1.45	1.34
8	B	1815[A]	KFR	C25-N24	5.40	1.45	1.34

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	1814	NAG	C2-N2-C7	6.52	131.64	122.90
7	B	1810	NAG	C1-O5-C5	4.56	118.30	112.19
8	B	1815[A]	KFR	O22-C20-C09	3.49	120.36	111.49
8	B	1815[B]	KFR	O22-C20-C09	3.48	120.34	111.49
8	A	1106	KFR	O22-C20-C09	3.30	119.89	111.49

There are no chirality outliers.

5 of 65 torsion outliers are listed below:

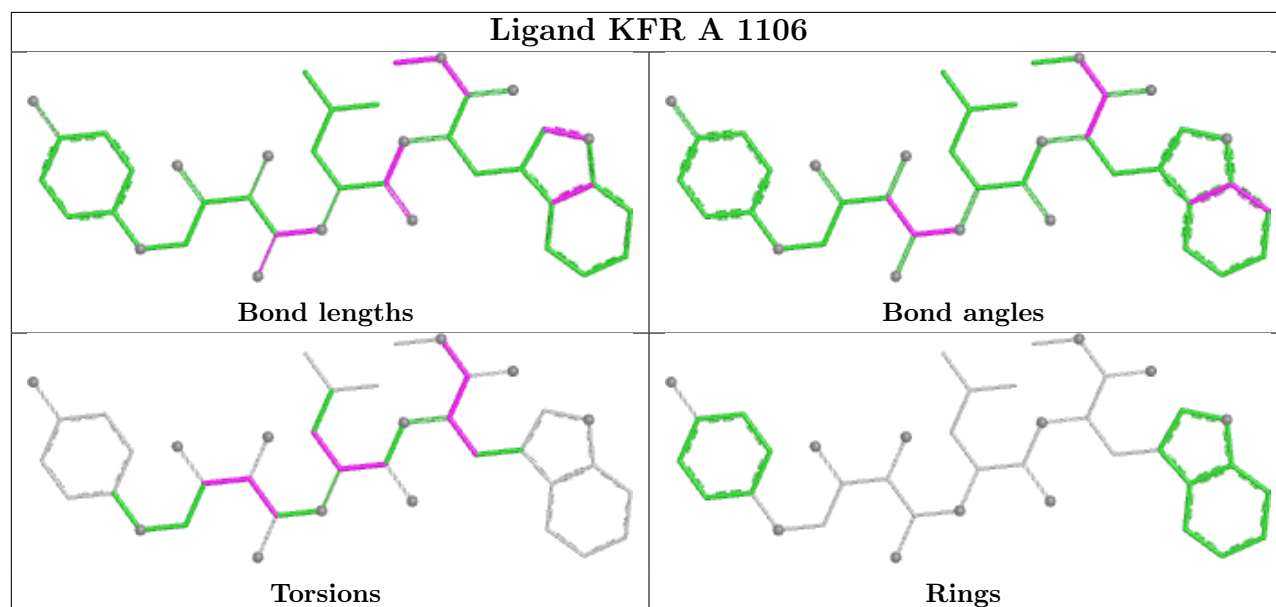
Mol	Chain	Res	Type	Atoms
7	A	1104	NAG	C1-C2-N2-C7
7	A	1105	NAG	C1-C2-N2-C7
7	B	1813	NAG	C1-C2-N2-C7
7	B	1814	NAG	C3-C2-N2-C7
8	A	1106	KFR	C25-C27-C28-C29

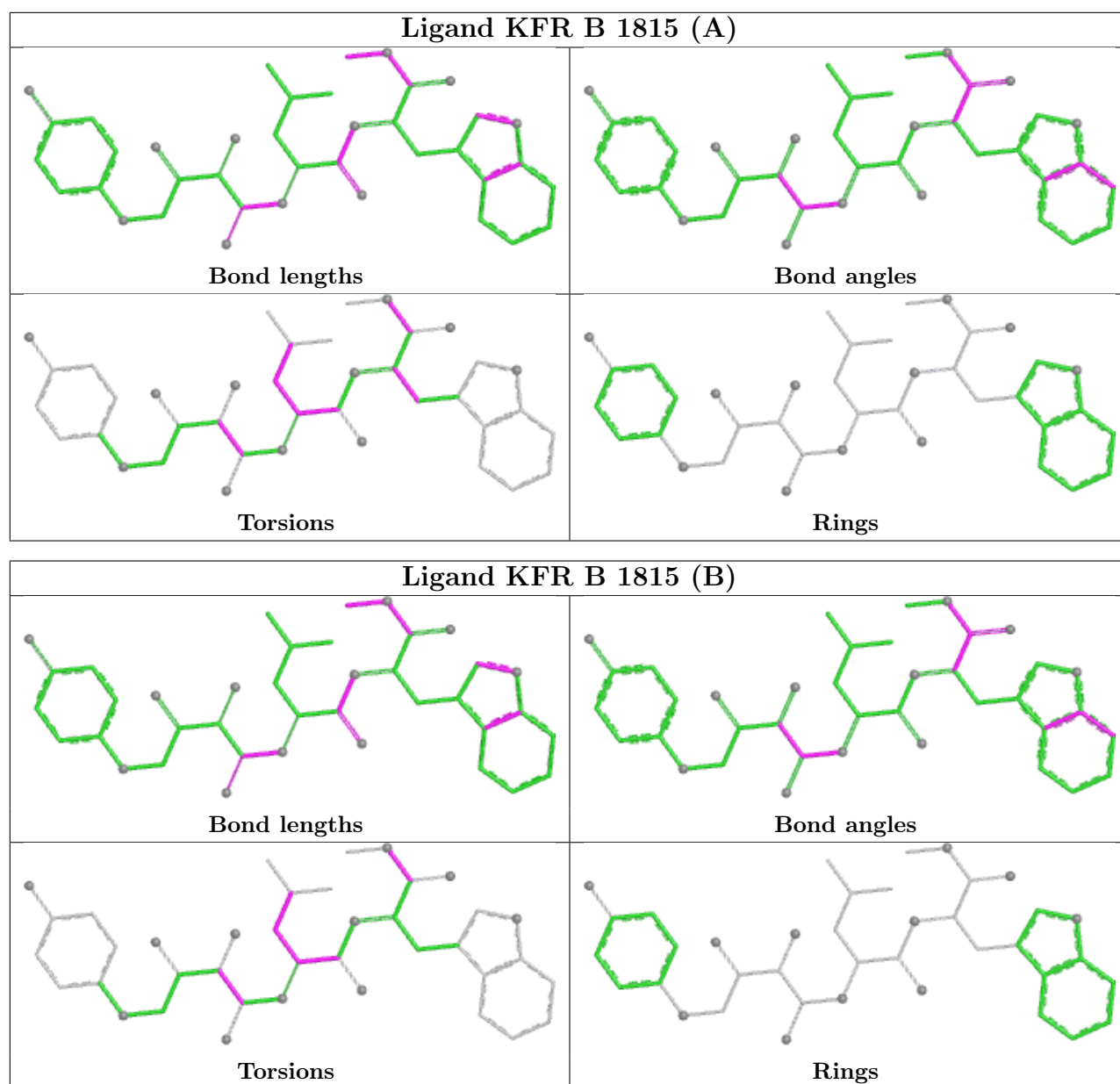
There are no ring outliers.

9 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	1813	NAG	2	0
7	B	1814	NAG	3	0
7	B	1810	NAG	2	0
7	A	1103	NAG	1	0
7	A	1102	NAG	1	0
8	B	1815[A]	KFR	2	0
7	B	1802	NAG	1	0
7	A	1104	NAG	2	0
8	B	1815[B]	KFR	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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	852/871 (97%)	1.17	160 (18%) 3 2	8, 33, 62, 96	4 (0%)
1	B	854/871 (98%)	1.44	225 (26%) 1 1	7, 39, 68, 115	2 (0%)
All	All	1706/1742 (97%)	1.31	385 (22%) 2 1	7, 36, 66, 115	6 (0%)

The worst 5 of 385 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	639	MET	8.0
1	B	222	MET	7.3
1	A	665	TYR	6.8
1	B	638	ASN	6.7
1	A	917	ASP	6.2

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
4	BMA	J	3	11/12	0.19	0.26	92,94,97,98	0
5	BMA	H	3	11/12	0.36	0.24	90,97,100,100	0
5	MAN	H	5	11/12	0.39	0.24	83,87,89,89	0
5	NAG	H	2	14/15	0.46	0.26	105,113,117,118	0
4	NAG	J	2	14/15	0.49	0.26	85,90,91,91	0

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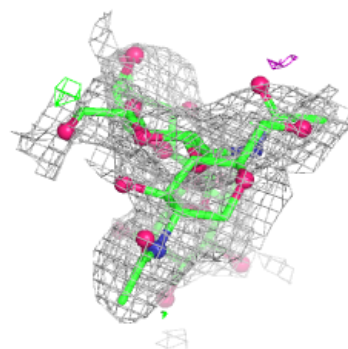
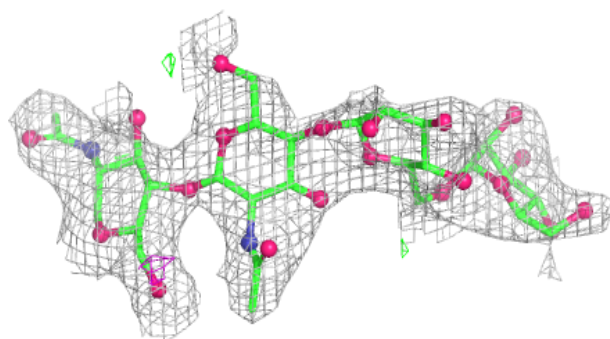
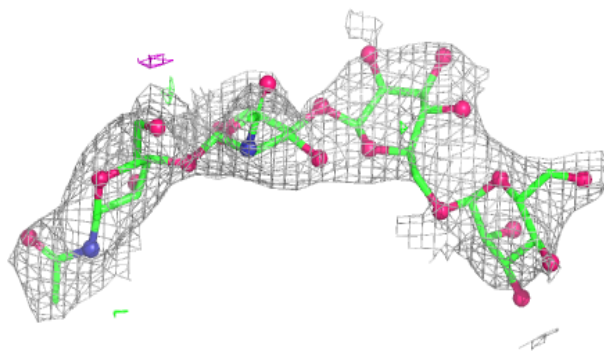
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MAN	H	4	11/12	0.50	0.26	80,82,84,86	0
3	NAG	F	2	14/15	0.51	0.25	60,68,69,72	0
3	NAG	G	2	14/15	0.51	0.26	82,87,89,93	0
3	NAG	G	1	14/15	0.52	0.35	104,109,116,117	0
3	NAG	F	1	14/15	0.52	0.23	79,87,89,94	0
4	BMA	E	3	11/12	0.52	0.20	73,75,76,77	0
3	NAG	K	1	14/15	0.53	0.28	121,128,131,131	0
5	NAG	I	2	14/15	0.55	0.23	47,60,62,65	0
5	NAG	H	1	14/15	0.57	0.27	91,103,108,114	0
2	BMA	C	3	11/12	0.57	0.22	76,79,81,82	0
5	MAN	I	4	11/12	0.57	0.18	83,86,88,88	0
5	MAN	I	5	11/12	0.57	0.25	68,71,73,75	0
3	NAG	K	2	14/15	0.65	0.24	76,89,94,101	0
5	NAG	I	1	14/15	0.67	0.26	63,67,75,76	0
5	BMA	I	3	11/12	0.68	0.19	68,69,73,78	0
4	NAG	J	1	14/15	0.68	0.21	76,83,89,92	0
2	NAG	C	2	14/15	0.68	0.22	67,71,74,79	0
3	NAG	D	2	14/15	0.70	0.16	59,65,70,71	0
4	NAG	E	2	14/15	0.73	0.20	68,73,76,78	0
2	MAN	C	4	11/12	0.77	0.18	64,68,71,71	0
4	NAG	E	1	14/15	0.78	0.19	39,48,52,60	0
2	NAG	C	1	14/15	0.82	0.16	42,47,54,61	0
3	NAG	D	1	14/15	0.84	0.12	37,43,49,53	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.

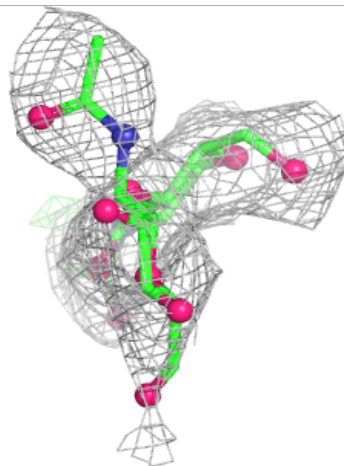
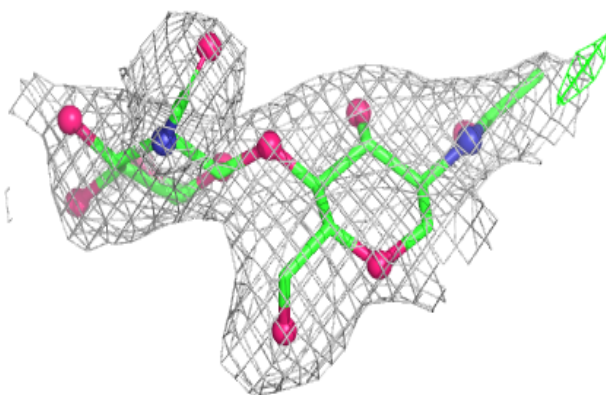
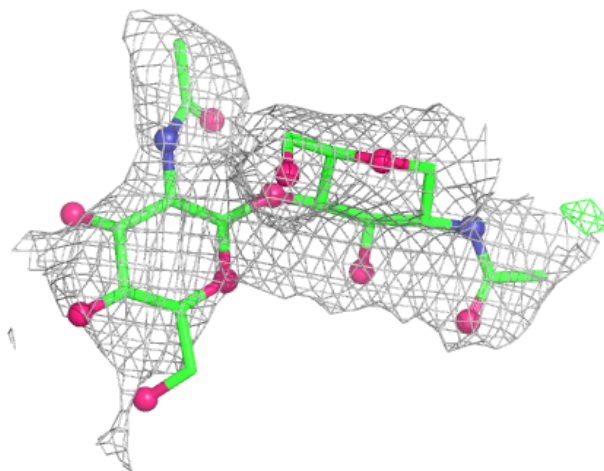
Electron density around Chain C:

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



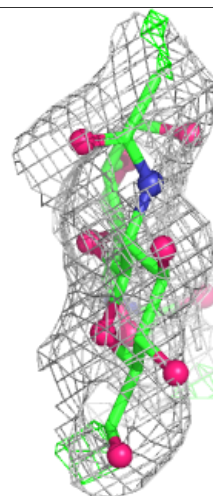
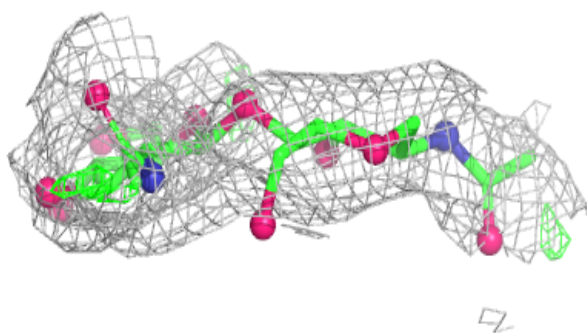
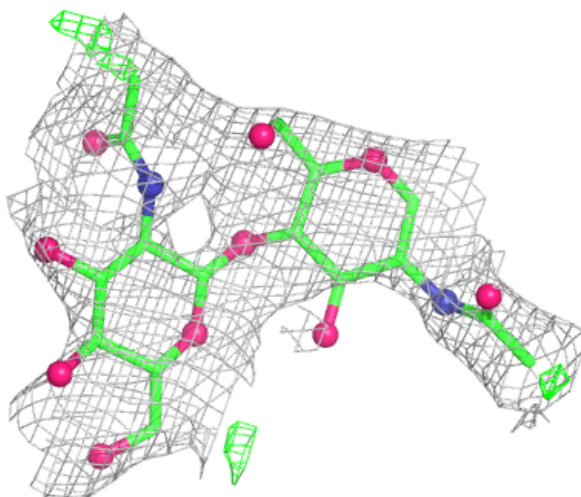
Electron density around Chain D:

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



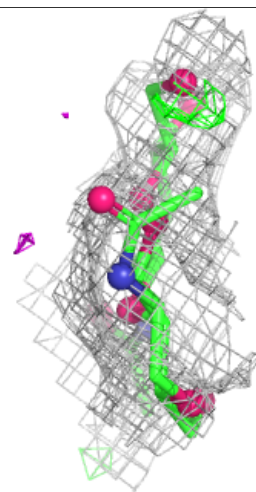
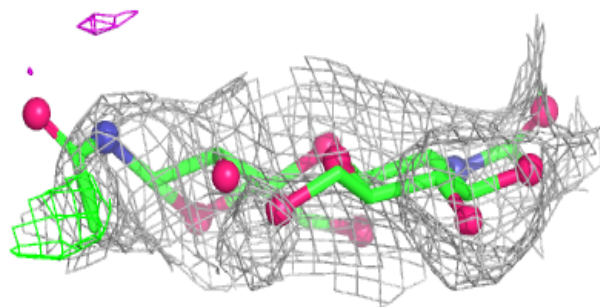
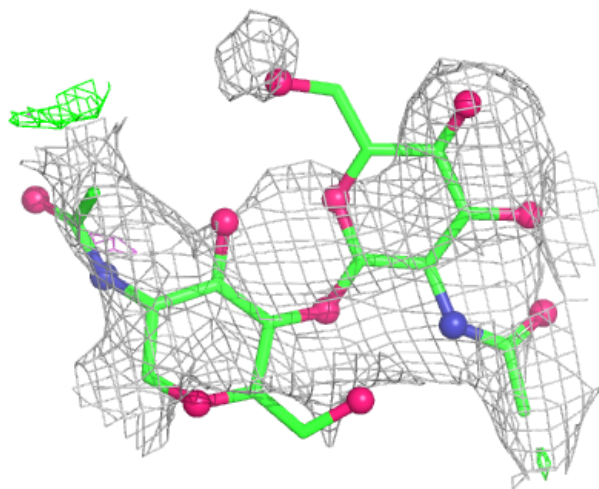
Electron density around Chain F:

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



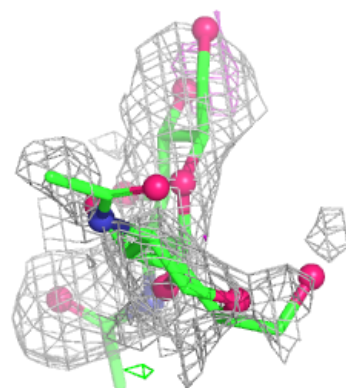
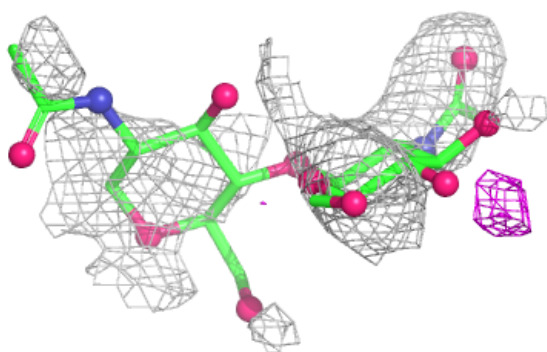
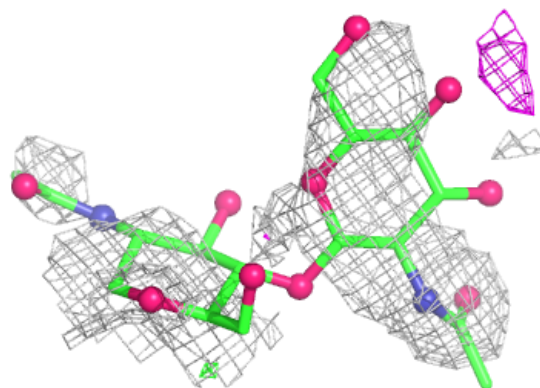
Electron density around Chain G:

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

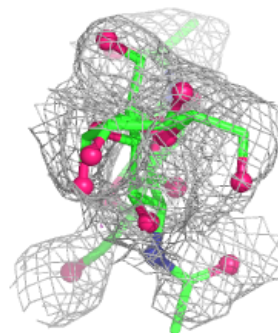
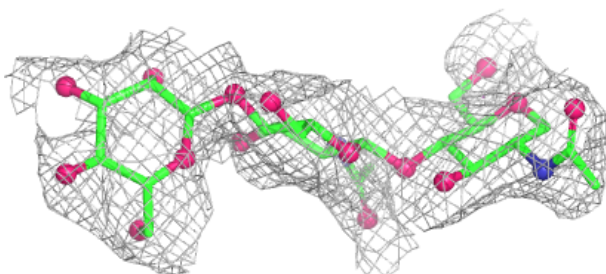
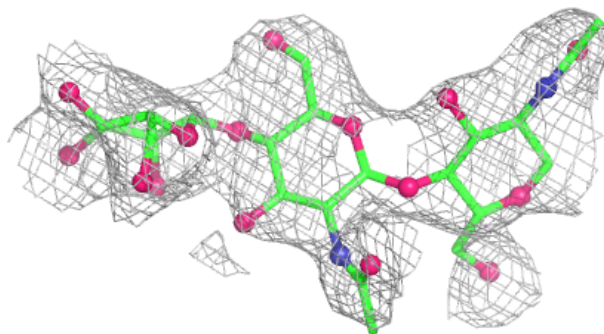


Electron density around Chain K:

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

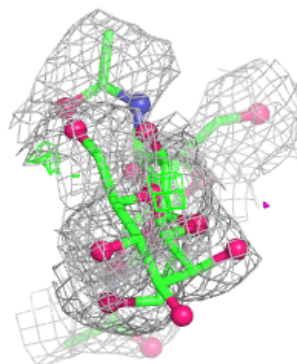
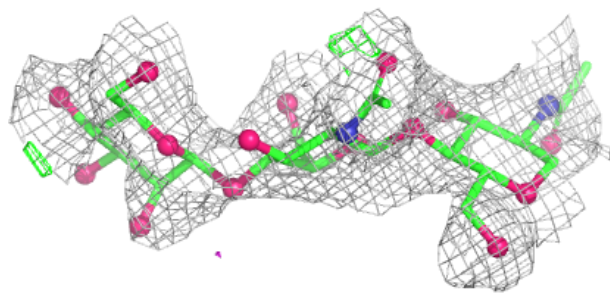
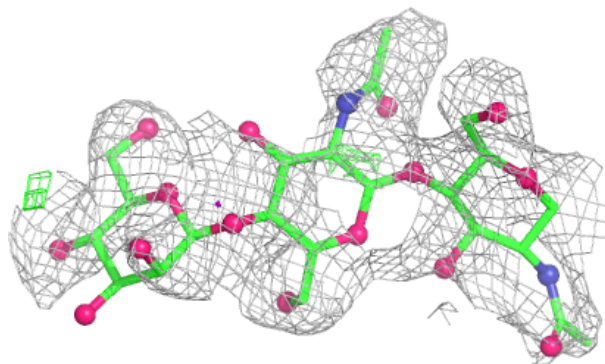
**Electron density around Chain E:**

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



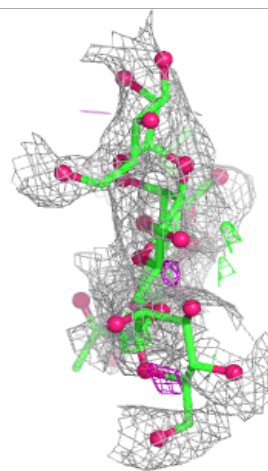
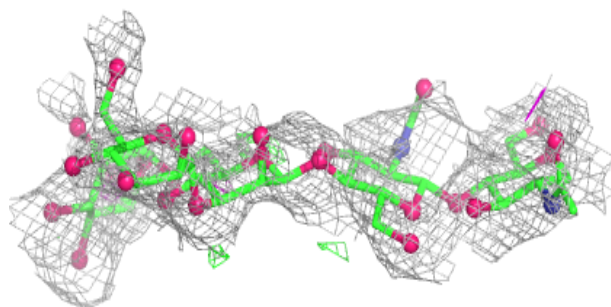
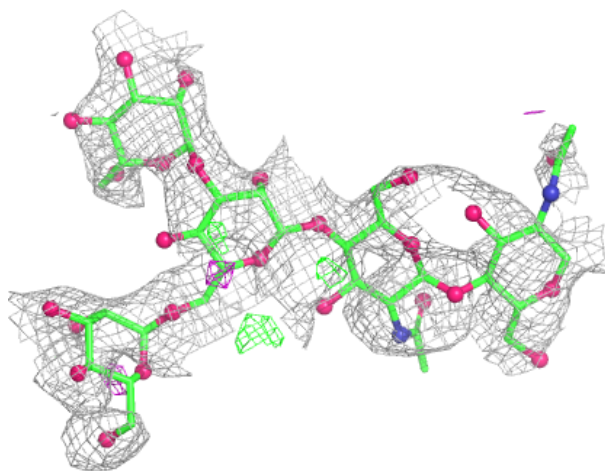
Electron density around Chain J:

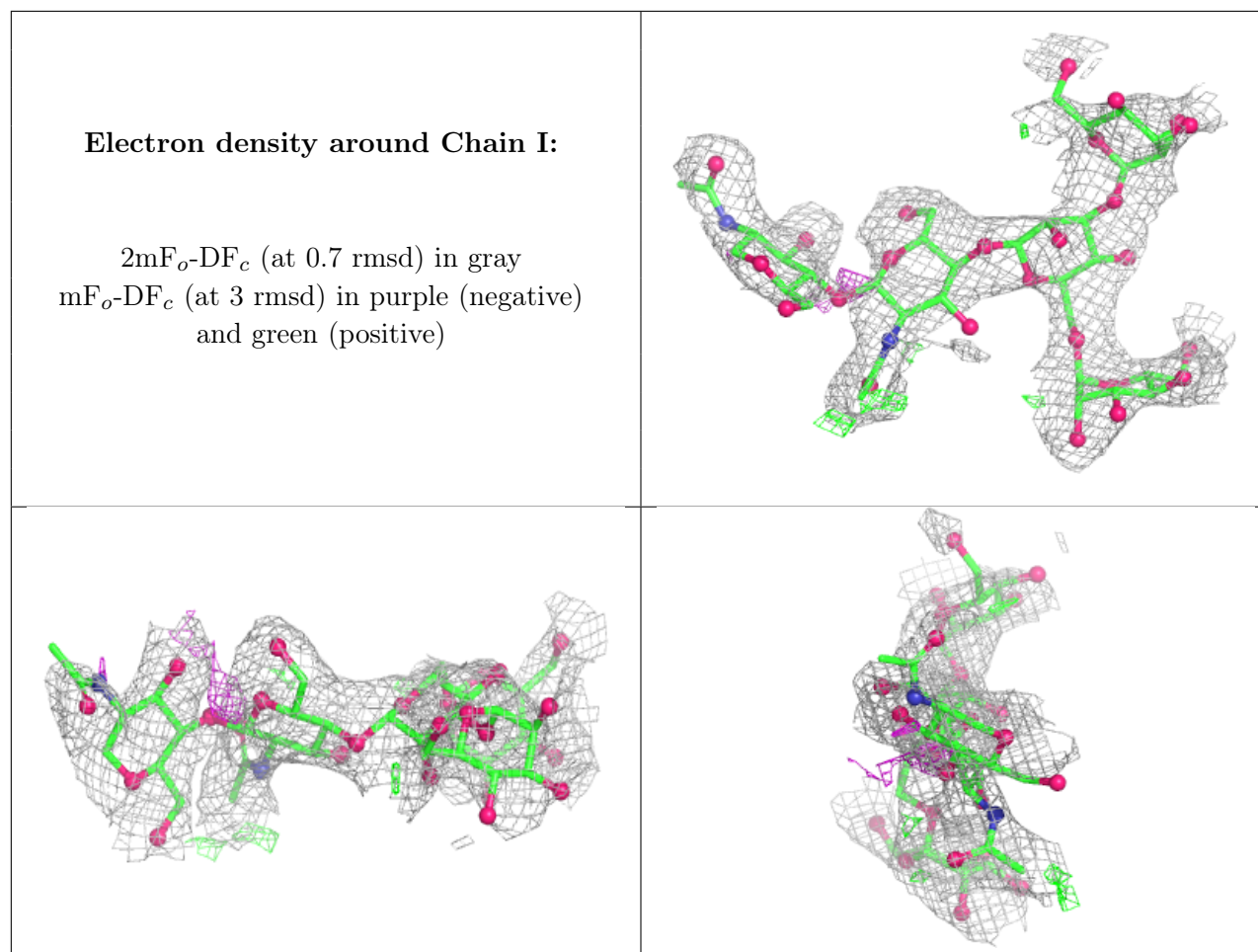
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





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
10	PEG	B	1809	7/7	0.47	0.27	55,58,62,63	0
10	PEG	A	1111	7/7	0.54	0.25	50,53,56,58	0
7	NAG	B	1812	14/15	0.57	0.22	82,85,88,88	0
7	NAG	B	1811	14/15	0.59	0.25	62,67,74,77	0
7	NAG	B	1803	14/15	0.60	0.26	70,74,75,76	0
7	NAG	B	1810	14/15	0.60	0.32	77,84,88,88	0
7	NAG	B	1802	14/15	0.60	0.27	78,84,86,86	0
7	NAG	A	1105	14/15	0.62	0.19	78,89,91,91	0
7	NAG	B	1814	14/15	0.62	0.23	74,77,82,89	0
7	NAG	A	1103	14/15	0.63	0.20	64,65,66,68	0
7	NAG	B	1813	14/15	0.67	0.19	63,66,72,78	0

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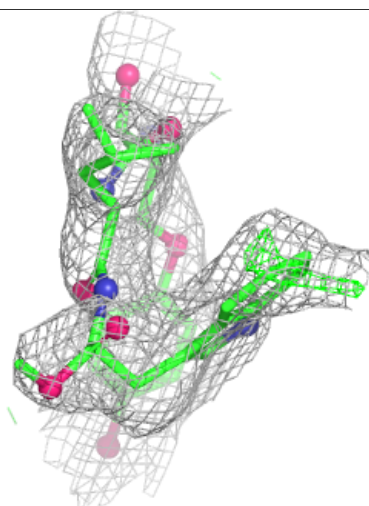
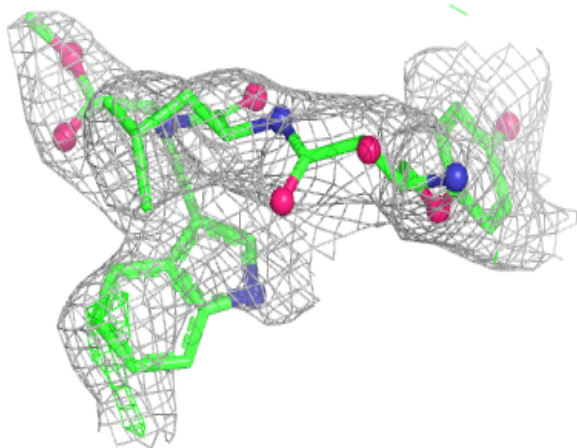
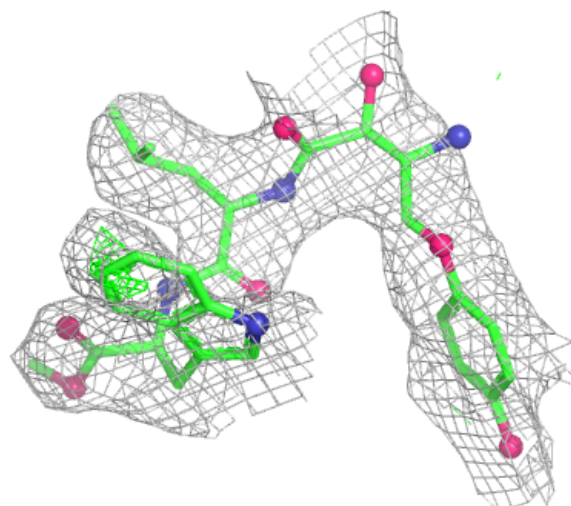
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	NAG	A	1104	14/15	0.68	0.20	65,68,73,73	0
7	NAG	A	1102	14/15	0.70	0.17	68,73,76,76	0
11	TRS	B	1807	8/8	0.71	0.17	34,36,37,37	0
9	EDO	B	1806	4/4	0.72	0.17	42,43,43,43	0
8	KFR	B	1815[A]	39/39	0.75	0.23	36,42,49,52	39
8	KFR	B	1815[B]	39/39	0.75	0.23	37,42,48,50	39
10	PEG	A	1110	7/7	0.79	0.20	34,37,40,41	0
9	EDO	A	1107	4/4	0.79	0.21	44,45,45,45	0
9	EDO	B	1805	4/4	0.81	0.14	36,37,37,38	0
9	EDO	B	1804	4/4	0.81	0.14	31,32,33,33	0
9	EDO	B	1808	4/4	0.84	0.16	40,41,41,41	0
9	EDO	A	1109	4/4	0.85	0.12	31,32,32,33	0
8	KFR	A	1106	39/39	0.86	0.16	34,43,55,58	0
9	EDO	A	1108	4/4	0.87	0.17	40,40,41,41	0
6	ZN	B	1801	1/1	1.00	0.02	29,29,29,29	0
6	ZN	A	1101	1/1	1.00	0.02	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.

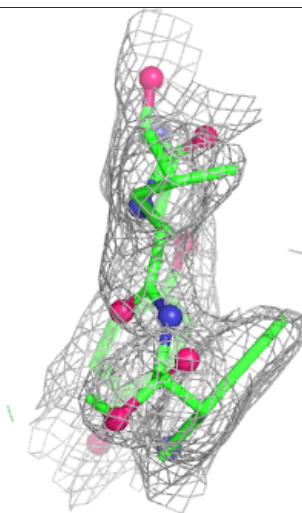
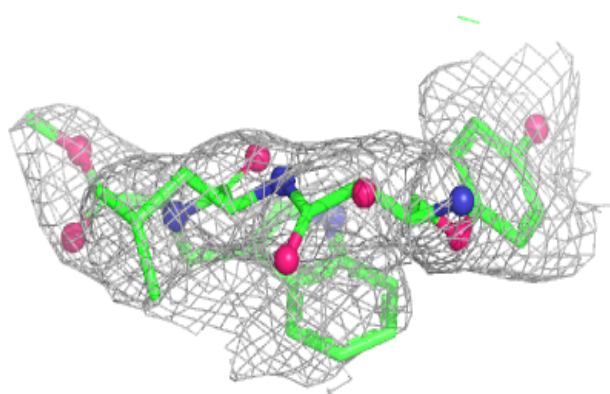
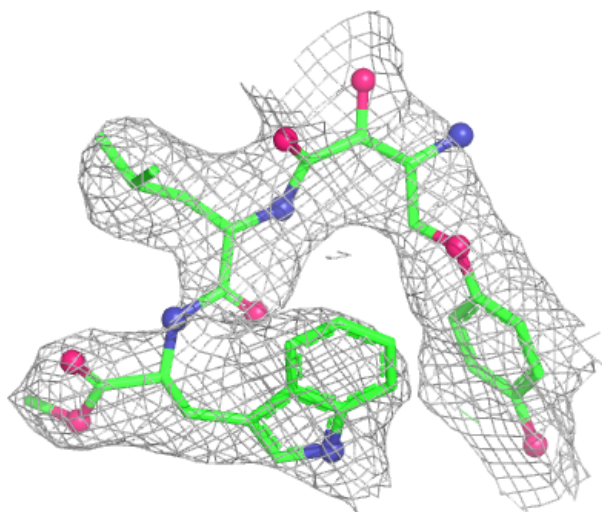
Electron density around KFR B 1815 (A):

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



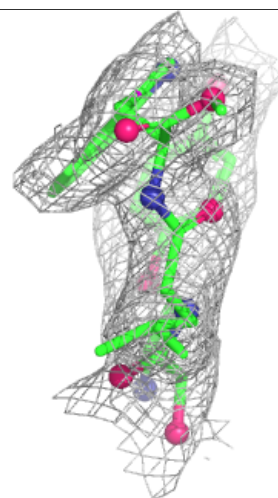
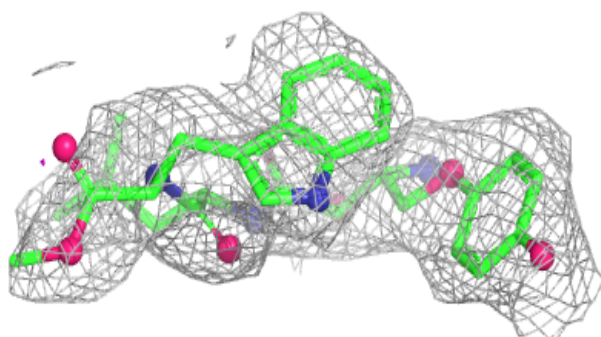
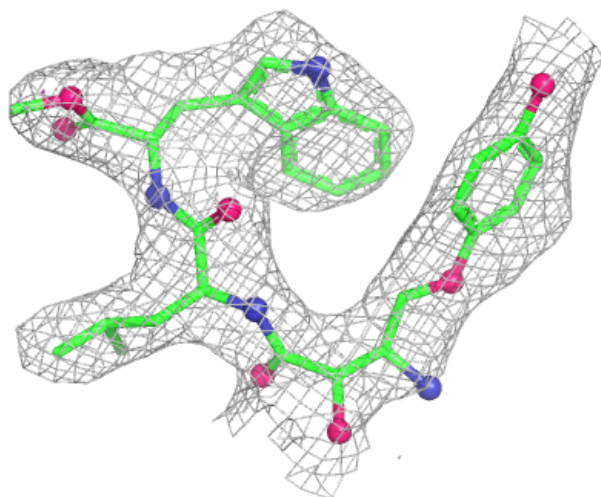
Electron density around KFR B 1815 (B):

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



Electron density around KFR A 1106:

$2mF_o-DF_c$ (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.