



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 22, 2025 – 03:12 AM EDT

PDB ID : 7LFS / pdb\_00007lfs  
Title : Crystal structure of the epidermal growth factor receptor extracellular region with A265V mutation in complex with epiregulin  
Authors : Hu, C.; Leche II, C.A.; Stayrook, S.E.; Ferguson, K.M.; Lemmon, M.A.  
Deposited on : 2021-01-18  
Resolution : 3.50 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](mailto: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>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0rc1
EDS	:	<b>FAILED</b>
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.42

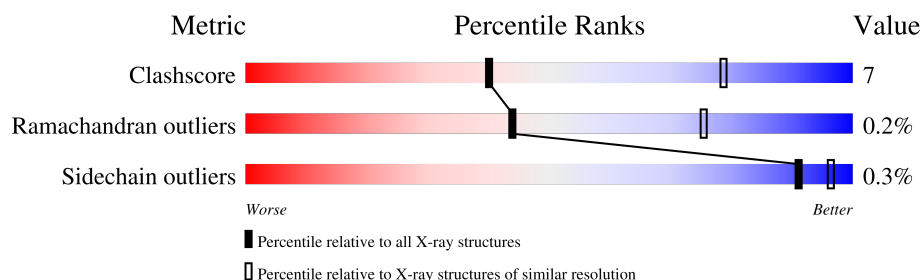
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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))
Clashscore	180529	1045 (3.54-3.46)
Ramachandran outliers	177936	1032 (3.54-3.46)
Sidechain outliers	177891	1033 (3.54-3.46)






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  $\geq 3$ , 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  $\leq 5\%$ .

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	507	86% 13% .
1	B	507	84% 15% .
1	C	507	81% 18% .
1	D	507	81% 17% .
2	E	48	92% 8%
2	F	48	90% . 6%
2	G	48	85% 10% .
2	H	48	83% 8% 8%

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Mol	Chain	Length	Quality of chain
3	I	4	 75% 25%
3	K	4	 25% 25% 50%
3	M	4	 25% 75%
4	J	5	 60% 40%
5	L	2	 50% 50%

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16443 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 Isoform 4 of Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	500	Total	C	N	O	S	0	0	0
			3643	2251	640	712	40			
1	B	503	Total	C	N	O	S	0	0	0
			3793	2349	663	739	42			
1	C	500	Total	C	N	O	S	0	0	0
			3663	2269	635	718	41			
1	D	498	Total	C	N	O	S	0	0	0
			3620	2254	615	709	42			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	265	VAL	ALA	engineered mutation	UNP P00533
A	502	HIS	-	expression tag	UNP P00533
A	503	HIS	-	expression tag	UNP P00533
A	504	HIS	-	expression tag	UNP P00533
A	505	HIS	-	expression tag	UNP P00533
A	506	HIS	-	expression tag	UNP P00533
A	507	HIS	-	expression tag	UNP P00533
B	265	VAL	ALA	engineered mutation	UNP P00533
B	502	HIS	-	expression tag	UNP P00533
B	503	HIS	-	expression tag	UNP P00533
B	504	HIS	-	expression tag	UNP P00533
B	505	HIS	-	expression tag	UNP P00533
B	506	HIS	-	expression tag	UNP P00533
B	507	HIS	-	expression tag	UNP P00533
C	265	VAL	ALA	engineered mutation	UNP P00533
C	502	HIS	-	expression tag	UNP P00533
C	503	HIS	-	expression tag	UNP P00533
C	504	HIS	-	expression tag	UNP P00533
C	505	HIS	-	expression tag	UNP P00533
C	506	HIS	-	expression tag	UNP P00533
C	507	HIS	-	expression tag	UNP P00533

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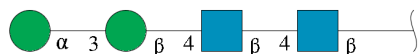
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Chain	Residue	Modelled	Actual	Comment	Reference
D	265	VAL	ALA	engineered mutation	UNP P00533
D	502	HIS	-	expression tag	UNP P00533
D	503	HIS	-	expression tag	UNP P00533
D	504	HIS	-	expression tag	UNP P00533
D	505	HIS	-	expression tag	UNP P00533
D	506	HIS	-	expression tag	UNP P00533
D	507	HIS	-	expression tag	UNP P00533

- Molecule 2 is a protein called Proepiregulin.

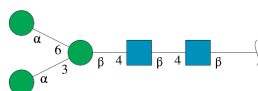
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	48	Total	C	N	O	S	0	0	0
			371	229	63	71	8			
2	F	45	Total	C	N	O	S	0	0	0
			347	215	58	66	8			
2	G	46	Total	C	N	O	S	0	0	0
			351	220	56	67	8			
2	H	44	Total	C	N	O	S	0	0	0
			332	206	52	66	8			

- Molecule 3 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
3	I	4	Total	C	N	O	0	0	0
			50	28	2	20			
3	K	4	Total	C	N	O	0	0	0
			50	28	2	20			
3	M	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 4 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.



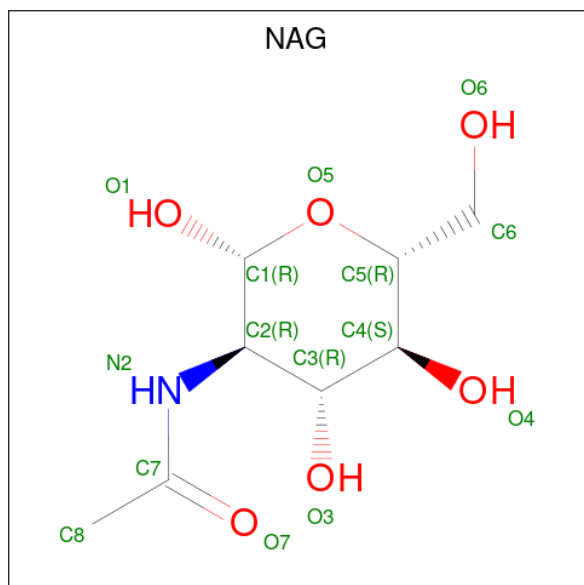
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	J	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 5 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
5	L	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: 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	B	1	Total	C	N	O	0	0
			14	8	1	5		

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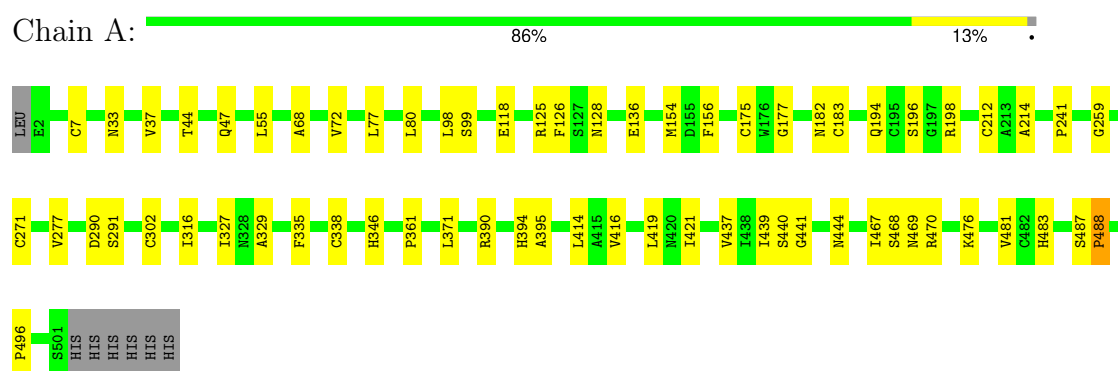
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			14	8	1	5		
6	C	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		

### 3 Residue-property plots

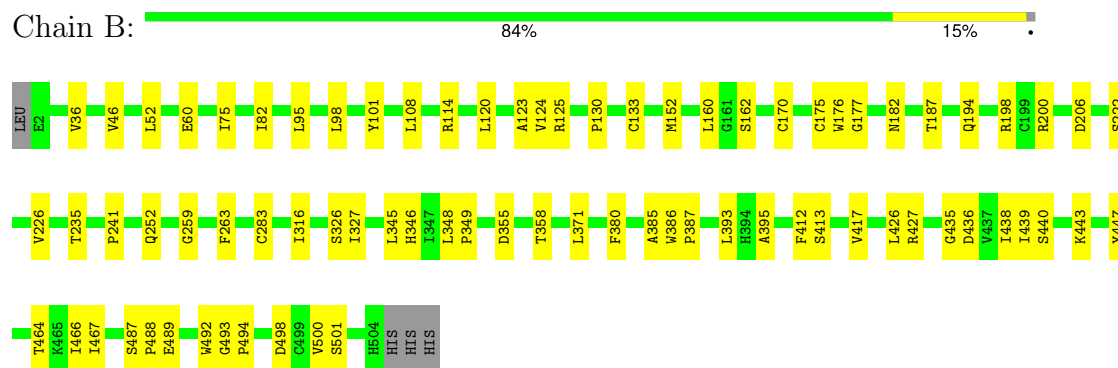
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS failed to run properly.

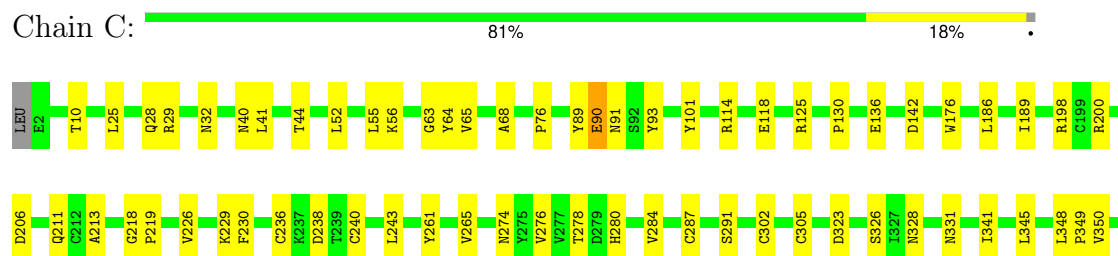
- Molecule 1: Isoform 4 of Epidermal growth factor receptor



- Molecule 1: Isoform 4 of Epidermal growth factor receptor



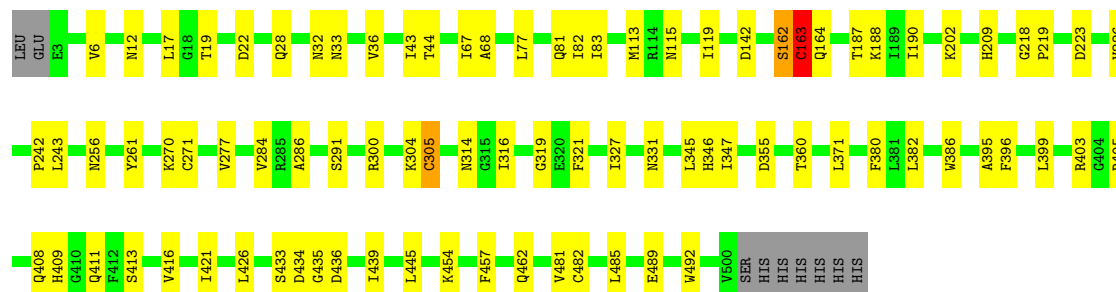
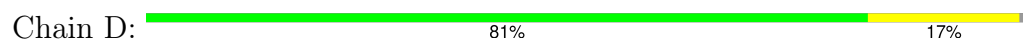
- Molecule 1: Isoform 4 of Epidermal growth factor receptor







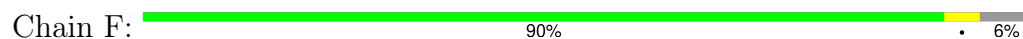
- Molecule 1: Isoform 4 of Epidermal growth factor receptor



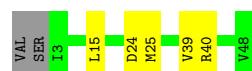
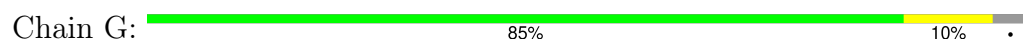
- Molecule 2: Proepiregulin



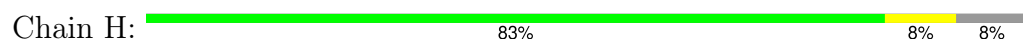
- Molecule 2: Proepiregulin



- Molecule 2: Proepiregulin



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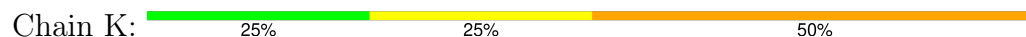


- Molecule 3: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 3: 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



- Molecule 3: 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



- Molecule 4: 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



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



## 4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.97Å 201.33Å 92.15Å 90.00° 99.04° 90.00°	Depositor
Resolution (Å)	44.38 – 3.50	Depositor
% Data completeness (in resolution range)	99.2 (44.38-3.50)	Depositor
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.33 (at 3.48Å)	Xtriage
Refinement program	PHENIX dev_3915	Depositor
R, $R_{free}$	0.266 , 0.315	Depositor
Wilson B-factor (Å <sup>2</sup> )	139.5	Xtriage
Anisotropy	0.268	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	16443	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	141.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, 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	0.25	0/3708	0.46	0/5051
1	B	0.25	0/3867	0.44	0/5253
1	C	0.27	0/3731	0.49	0/5082
1	D	0.25	0/3690	0.45	0/5032
2	E	0.24	0/377	0.42	0/508
2	F	0.25	0/354	0.42	0/476
2	G	0.24	0/358	0.41	0/484
2	H	0.24	0/339	0.42	0/458
All	All	0.25	0/16424	0.46	0/22344

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [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	3643	0	3335	36	0
1	B	3793	0	3550	47	0
1	C	3663	0	3362	59	0
1	D	3620	0	3297	60	0
2	E	371	0	345	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	347	0	302	1	0
2	G	351	0	309	3	0
2	H	332	0	280	2	0
3	I	50	0	43	0	0
3	K	50	0	43	2	0
3	M	50	0	43	1	0
4	J	61	0	52	0	0
5	L	28	0	25	0	0
6	A	14	0	13	2	0
6	B	28	0	26	0	0
6	C	14	0	13	2	0
6	D	28	0	26	4	0
All	All	16443	0	15064	206	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 7.

All (206) 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:D:33:ASN:HD21	6:D:601:NAG:H61	1.16	1.03
1:D:33:ASN:ND2	6:D:601:NAG:H61	1.81	0.96
1:C:323:ASP:HB3	3:K:2:NAG:H62	1.59	0.84
1:D:413:SER:HB3	1:D:435:GLY:HA3	1.60	0.82
1:C:420:ASN:OD1	1:C:444:ASN:ND2	2.16	0.78
1:C:328:ASN:HA	1:C:363:LEU:HD21	1.68	0.74
1:C:287:CYS:HB3	1:C:291:SER:HB3	1.71	0.72
1:D:300:ARG:HH11	1:D:405:ARG:HH22	1.33	0.72
1:B:187:THR:OG1	1:B:198:ARG:NH1	2.22	0.72
1:C:142:ASP:O	1:C:198:ARG:NH2	2.23	0.70
1:D:396:PHE:HA	1:D:399:LEU:HD13	1.73	0.70
1:D:416:VAL:HB	1:D:439:ILE:HG12	1.74	0.69
1:C:386:TRP:HD1	1:C:421:ILE:HG21	1.58	0.69
1:B:371:LEU:HB2	1:B:395:ALA:HB1	1.75	0.68
1:A:371:LEU:HB2	1:A:395:ALA:HB1	1.76	0.67
1:A:316:ILE:HG13	1:A:327:ILE:HG22	1.76	0.67
1:D:12:ASN:HD22	1:D:17:LEU:HD11	1.61	0.65
1:B:46:VAL:HG11	1:B:52:LEU:HD11	1.79	0.64
1:B:241:PRO:HB2	1:B:259:GLY:HA2	1.79	0.63
1:B:447:TYR:HE2	1:B:494:PRO:HD3	1.64	0.63
1:C:276:VAL:HG13	1:C:284:VAL:HG13	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:314:ASN:HD22	1:D:319:GLY:HA3	1.64	0.62
1:C:213:ALA:HB3	1:C:226:VAL:HG13	1.81	0.61
1:A:241:PRO:HB2	1:A:259:GLY:HA2	1.81	0.60
1:B:393:LEU:HD12	1:B:426:LEU:HD11	1.81	0.60
1:D:162:SER:O	1:D:163:CYS:HB3	2.02	0.60
1:D:33:ASN:ND2	6:D:601:NAG:C6	2.62	0.60
1:D:321:PHE:CE1	1:D:331:ASN:HB2	2.37	0.60
6:C:601:NAG:O7	6:C:601:NAG:O3	2.13	0.60
1:B:355:ASP:HB3	1:B:358:THR:HG22	1.84	0.60
1:A:444:ASN:HA	1:A:470:ARG:HD2	1.84	0.59
1:C:278:THR:HG22	1:C:280:HIS:H	1.66	0.59
1:A:194:GLN:HG3	1:D:219:PRO:HB3	1.84	0.59
1:B:98:LEU:HD21	1:B:125:ARG:HG2	1.84	0.58
1:B:349:PRO:HG3	1:B:385:ALA:HB2	1.84	0.58
1:C:118:GLU:OE1	1:C:198:ARG:NH2	2.30	0.58
1:D:327:ILE:HD11	1:D:345:LEU:HD22	1.86	0.57
1:A:177:GLY:H	1:A:182:ASN:HB3	1.68	0.57
2:H:33:GLU:HB3	2:H:36:TYR:HD2	1.69	0.57
1:C:211:GLN:HB3	1:C:236:CYS:HB2	1.85	0.57
1:C:276:VAL:HG22	1:C:302:CYS:HB2	1.88	0.56
1:C:491:CYS:HB2	1:C:499:CYS:HA	1.87	0.56
1:C:416:VAL:HG12	1:C:442:ASN:HD21	1.71	0.56
1:C:416:VAL:HG13	1:C:419:LEU:HD21	1.88	0.56
1:A:476:LYS:HD2	1:A:481:VAL:HG21	1.87	0.56
1:D:77:LEU:HB2	1:D:113:MET:HG2	1.89	0.55
1:B:252:GLN:HA	1:C:284:VAL:HG23	1.88	0.55
1:D:33:ASN:HD21	6:D:601:NAG:C6	2.05	0.55
1:B:327:ILE:HD11	1:B:345:LEU:HD22	1.88	0.55
1:D:371:LEU:HB2	1:D:395:ALA:HB1	1.89	0.55
1:D:360:THR:OG1	3:M:1:NAG:O7	2.25	0.54
1:B:413:SER:HB2	1:B:435:GLY:HA3	1.88	0.54
1:A:77:LEU:HD12	1:A:80:LEU:HB2	1.88	0.54
1:B:82:ILE:HG21	1:B:226:VAL:HG11	1.88	0.54
1:D:82:ILE:HG21	1:D:226:VAL:HG11	1.89	0.54
1:D:6:VAL:HG12	1:D:36:VAL:HB	1.89	0.53
1:D:190:ILE:HG13	1:D:202:LYS:HE3	1.90	0.53
1:A:390:ARG:HG3	1:A:394:HIS:CE1	2.44	0.53
1:A:55:LEU:HB3	1:A:77:LEU:HD22	1.90	0.53
1:A:440:SER:HA	1:A:467:ILE:O	2.09	0.53
1:B:488:PRO:HG2	1:B:501:SER:HB3	1.89	0.53
1:B:123:ALA:HB1	1:B:152:MET:HG2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:447:TYR:OH	1:C:480:GLN:HB3	2.08	0.53
1:D:435:GLY:O	1:D:462:GLN:HG2	2.09	0.52
1:B:417:VAL:HA	1:B:440:SER:O	2.09	0.52
1:D:142:ASP:OD1	1:D:188:LYS:N	2.42	0.52
1:D:316:ILE:HD13	1:D:346:HIS:H	1.75	0.52
1:A:175:CYS:HA	1:A:183:CYS:HA	1.92	0.52
2:G:39:VAL:HG13	2:G:40:ARG:HG3	1.92	0.52
1:C:64:TYR:CD1	1:C:89:TYR:HB2	2.44	0.51
1:A:291:SER:HB3	1:A:302:CYS:HB3	1.92	0.51
1:B:316:ILE:HD12	1:B:346:HIS:H	1.75	0.51
1:D:243:LEU:HD21	1:D:261:TYR:CE1	2.45	0.51
1:C:360:THR:OG1	3:K:1:NAG:O7	2.18	0.51
1:A:487:SER:HB2	1:A:488:PRO:HD3	1.92	0.50
1:C:422:THR:HG22	1:C:444:ASN:HB3	1.93	0.50
1:D:355:ASP:H	1:D:360:THR:HB	1.76	0.50
1:C:326:SER:HB2	1:C:348:LEU:HD12	1.94	0.50
1:B:75:ILE:HG13	1:B:108:LEU:HD13	1.94	0.50
1:C:349:PRO:HG3	1:C:385:ALA:HB2	1.94	0.49
1:C:419:LEU:HG	1:C:442:ASN:OD1	2.12	0.49
1:C:416:VAL:HB	1:C:439:ILE:HG12	1.94	0.49
1:D:284:VAL:HG12	1:D:286:ALA:H	1.78	0.49
1:C:200:ARG:NH1	1:C:218:GLY:HA2	2.28	0.49
1:C:52:LEU:HB3	1:C:55:LEU:HD12	1.94	0.49
1:D:434:ASP:HA	1:D:462:GLN:HE21	1.77	0.49
2:E:4:THR:HB	2:E:21:TYR:HB3	1.96	0.48
1:A:416:VAL:HG13	1:A:419:LEU:HD22	1.96	0.48
1:C:28:GLN:O	1:C:32:ASN:HB2	2.14	0.48
1:B:380:PHE:HB2	1:B:413:SER:O	2.14	0.47
1:C:41:LEU:O	1:C:65:VAL:HA	2.14	0.47
1:D:454:LYS:HD3	1:D:457:PHE:HD2	1.79	0.47
1:C:136:GLU:OE1	1:C:136:GLU:N	2.41	0.47
1:C:200:ARG:HG3	1:C:206:ASP:HB3	1.96	0.47
1:D:261:TYR:HE2	1:D:270:LYS:HA	1.80	0.47
1:D:12:ASN:ND2	1:D:17:LEU:HD11	2.29	0.47
1:B:200:ARG:HB2	1:B:206:ASP:HB3	1.96	0.47
1:C:114:ARG:HA	1:C:176:TRP:CD1	2.50	0.47
1:C:230:PHE:CE1	1:C:265:VAL:HA	2.50	0.47
1:C:472:GLU:O	1:C:476:LYS:HG3	2.14	0.47
1:D:411:GLN:O	1:D:436:ASP:HB2	2.15	0.47
1:B:412:PHE:HB3	1:B:438:ILE:HG22	1.96	0.47
1:D:426:LEU:HB2	1:D:492:TRP:CD1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:PHE:HD2	1:A:154:MET:HB3	1.79	0.46
1:D:163:CYS:SG	1:D:164:GLN:N	2.88	0.46
1:D:291:SER:HA	1:D:305:CYS:N	2.30	0.46
1:D:382:LEU:HD22	1:D:408:GLN:HE21	1.80	0.46
1:A:99:SER:HA	1:A:128:ASN:HB3	1.97	0.46
1:B:222:SER:HB3	1:B:235:THR:HG22	1.98	0.46
1:D:321:PHE:HE1	1:D:331:ASN:HB2	1.79	0.46
1:B:439:ILE:HB	1:B:466:ILE:HG23	1.98	0.46
1:D:403:ARG:O	1:D:433:SER:HB2	2.14	0.46
1:B:413:SER:N	1:B:436:ASP:O	2.49	0.46
1:C:200:ARG:HG2	1:C:206:ASP:O	2.16	0.46
1:D:321:PHE:CD1	1:D:331:ASN:HB2	2.51	0.46
1:D:421:ILE:HG13	1:D:445:LEU:HD13	1.98	0.46
1:C:229:LYS:CD	1:C:238:ASP:HA	2.46	0.45
1:D:271:CYS:SG	1:D:277:VAL:HG22	2.57	0.45
1:A:198:ARG:NH1	1:A:214:ALA:O	2.50	0.45
1:D:291:SER:HA	1:D:305:CYS:H	1.80	0.45
1:B:427:ARG:NH1	1:B:498:ASP:OD1	2.49	0.45
1:B:489:GLU:O	1:B:489:GLU:HG2	2.17	0.45
1:C:350:VAL:HG11	2:G:15:LEU:HD13	1.97	0.45
1:D:434:ASP:HA	1:D:462:GLN:NE2	2.32	0.45
1:A:44:THR:HA	1:A:68:ALA:O	2.16	0.45
1:A:7:CYS:SG	1:A:37:VAL:HG22	2.57	0.45
1:A:33:ASN:ND2	6:A:3301:NAG:O5	2.43	0.45
1:A:196:SER:HB2	1:D:209:HIS:NE2	2.32	0.45
1:B:82:ILE:HD11	1:B:120:LEU:HG	1.98	0.45
1:B:326:SER:HB2	1:B:348:LEU:HD13	1.98	0.45
1:B:101:TYR:HB3	1:B:130:PRO:HD2	1.98	0.44
1:B:170:CYS:SG	1:B:175:CYS:HB3	2.57	0.44
1:B:447:TYR:HD2	1:B:493:GLY:HA2	1.81	0.44
1:C:56:LYS:HG3	1:C:76:PRO:HB2	1.98	0.44
1:A:136:GLU:HG3	1:A:156:PHE:CG	2.53	0.44
1:D:19:THR:HG23	1:D:22:ASP:H	1.81	0.44
1:B:440:SER:HA	1:B:467:ILE:O	2.18	0.44
1:C:200:ARG:HE	1:C:206:ASP:HA	1.82	0.44
1:B:95:LEU:HB3	1:B:124:VAL:HG22	1.99	0.44
1:C:341:ILE:HG13	1:C:345:LEU:HD11	1.99	0.44
1:B:386:TRP:CG	1:B:387:PRO:HD2	2.53	0.44
1:D:81:GLN:HA	1:D:115:ASN:O	2.18	0.44
1:B:177:GLY:HA3	1:B:182:ASN:HD22	1.83	0.44
1:C:10:THR:O	1:C:40:ASN:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:447:TYR:CD2	1:B:493:GLY:HA2	2.52	0.43
1:C:93:TYR:CE1	1:C:125:ARG:HB2	2.52	0.43
1:B:160:LEU:HD12	1:B:162:SER:H	1.82	0.43
1:A:98:LEU:HD21	1:A:125:ARG:HG2	2.00	0.43
1:B:194:GLN:HG3	1:C:219:PRO:HB3	2.01	0.43
1:D:242:PRO:O	1:D:256:ASN:ND2	2.45	0.43
1:C:403:ARG:O	1:C:433:SER:HB2	2.18	0.43
1:C:413:SER:HB3	1:C:435:GLY:HA3	2.01	0.43
1:A:441:GLY:H	1:A:468:SER:HB2	1.82	0.43
1:B:386:TRP:CD2	1:B:387:PRO:HD2	2.54	0.43
1:B:263:PHE:HB2	1:B:283:CYS:SG	2.59	0.43
1:B:464:THR:HG22	1:B:466:ILE:HG13	1.99	0.43
1:D:408:GLN:HG3	1:D:409:HIS:CD2	2.54	0.43
1:A:439:ILE:HG22	1:A:469:ASN:HD21	1.84	0.43
1:D:291:SER:OG	1:D:304:LYS:HA	2.19	0.43
1:C:230:PHE:CE2	1:C:240:CYS:HB2	2.54	0.42
1:A:419:LEU:HB3	1:A:421:ILE:HG12	2.01	0.42
1:C:90:GLU:HB3	1:C:91:ASN:H	1.67	0.42
6:C:601:NAG:HO3	6:C:601:NAG:C7	2.19	0.42
1:B:487:SER:OG	1:B:488:PRO:HD3	2.19	0.42
1:C:380:PHE:CE1	1:C:408:GLN:HB2	2.55	0.42
1:A:327:ILE:HG23	1:A:346:HIS:O	2.19	0.42
1:C:328:ASN:OD1	1:C:331:ASN:ND2	2.52	0.42
1:A:414:LEU:HB3	1:A:437:VAL:HG22	2.01	0.42
1:B:194:GLN:HG3	1:C:219:PRO:CB	2.49	0.42
1:D:28:GLN:O	1:D:32:ASN:HB2	2.19	0.42
1:D:43:ILE:HD12	1:D:67:ILE:HG12	2.00	0.42
1:A:483:HIS:CD2	1:A:496:PRO:HG3	2.55	0.42
1:C:44:THR:HA	1:C:68:ALA:O	2.20	0.42
2:E:22:LEU:O	2:E:26:SER:N	2.52	0.42
2:H:42:GLU:HG2	2:H:43:HIS:N	2.34	0.42
1:A:33:ASN:HD21	6:A:3301:NAG:H62	1.84	0.42
1:C:243:LEU:HD11	1:C:261:TYR:CE1	2.55	0.42
1:C:386:TRP:CD1	1:C:421:ILE:HG21	2.47	0.42
1:A:47:GLN:O	1:A:72:VAL:HG22	2.19	0.41
1:A:118:GLU:HG3	1:A:198:ARG:NH2	2.35	0.41
1:A:271:CYS:HB2	1:A:277:VAL:HG22	2.02	0.41
1:D:83:ILE:HB	1:D:119:ILE:HG12	2.02	0.41
1:B:492:TRP:NE1	1:B:500:VAL:HG12	2.34	0.41
1:C:274:ASN:HD21	1:C:405:ARG:HA	1.85	0.41
2:F:15:LEU:HD12	2:F:42:GLU:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:ALA:HB3	1:A:361:PRO:HG2	2.02	0.41
1:C:385:ALA:HA	1:C:419:LEU:HB3	2.02	0.41
1:C:40:ASN:OD1	1:C:63:GLY:HA3	2.21	0.41
1:D:44:THR:HA	1:D:68:ALA:O	2.20	0.41
1:D:489:GLU:OE2	1:D:489:GLU:N	2.54	0.41
1:D:380:PHE:HB2	1:D:413:SER:O	2.20	0.41
1:D:347:ILE:HD13	1:D:386:TRP:HZ3	1.86	0.41
1:D:481:VAL:HG12	1:D:482:CYS:N	2.36	0.41
1:D:218:GLY:H	1:D:223:ASP:HB3	1.85	0.41
1:A:335:PHE:HA	1:A:338:CYS:SG	2.61	0.41
1:B:443:LYS:HB3	1:B:443:LYS:HE2	1.79	0.41
1:C:186:LEU:HD13	1:C:189:ILE:HD11	2.02	0.41
1:B:36:VAL:HG13	1:B:60:GLU:HG3	2.02	0.40
1:C:25:LEU:O	1:C:29:ARG:HG3	2.22	0.40
1:D:426:LEU:HB2	1:D:492:TRP:NE1	2.36	0.40
1:B:114:ARG:HA	1:B:176:TRP:CD1	2.55	0.40
1:C:101:TYR:HB3	1:C:130:PRO:HG2	2.03	0.40
1:C:378:THR:HA	1:C:403:ARG:HB2	2.04	0.40
1:D:142:ASP:OD1	1:D:187:THR:N	2.52	0.40
1:D:485:LEU:HD23	1:D:485:LEU:H	1.87	0.40
2:G:24:ASP:OD1	2:G:25:MET:N	2.54	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	498/507 (98%)	466 (94%)	31 (6%)	1 (0%)	44	75
1	B	501/507 (99%)	473 (94%)	28 (6%)	0	100	100
1	C	498/507 (98%)	463 (93%)	35 (7%)	0	100	100
1	D	496/507 (98%)	462 (93%)	31 (6%)	3 (1%)	22	56

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	46/48 (96%)	45 (98%)	1 (2%)	0	100	100
2	F	43/48 (90%)	42 (98%)	1 (2%)	0	100	100
2	G	44/48 (92%)	41 (93%)	3 (7%)	0	100	100
2	H	42/48 (88%)	41 (98%)	1 (2%)	0	100	100
All	All	2168/2220 (98%)	2033 (94%)	131 (6%)	4 (0%)	44	75

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	162	SER
1	D	163	CYS
1	D	305	CYS
1	A	488	PRO

### 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	385/447 (86%)	383 (100%)	2 (0%)	86	93
1	B	418/447 (94%)	417 (100%)	1 (0%)	92	97
1	C	391/447 (88%)	389 (100%)	2 (0%)	86	93
1	D	382/447 (86%)	381 (100%)	1 (0%)	91	96
2	E	43/44 (98%)	43 (100%)	0	100	100
2	F	38/44 (86%)	38 (100%)	0	100	100
2	G	39/44 (89%)	39 (100%)	0	100	100
2	H	37/44 (84%)	37 (100%)	0	100	100
All	All	1733/1964 (88%)	1727 (100%)	6 (0%)	91	96

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	212	CYS
1	A	290	ASP
1	B	133	CYS
1	C	90	GLU
1	C	305	CYS
1	D	163	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	12	ASN
1	D	33	ASN
1	D	409	HIS

### 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 ⓘ

19 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	$\# Z  > 2$
3	NAG	I	1	3,1	14,14,15	0.37	0	17,19,21	0.49	0
3	NAG	I	2	3	14,14,15	0.18	0	17,19,21	0.54	0
3	BMA	I	3	3	11,11,12	0.71	0	15,15,17	0.72	0
3	MAN	I	4	3	11,11,12	0.74	0	15,15,17	0.89	1 (6%)
4	NAG	J	1	4,1	14,14,15	0.47	0	17,19,21	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NAG	J	2	4	14,14,15	0.21	0	17,19,21	0.54	0
4	BMA	J	3	4	11,11,12	0.64	0	15,15,17	0.85	0
4	MAN	J	4	4	11,11,12	1.02	1 (9%)	15,15,17	1.87	2 (13%)
4	MAN	J	5	4	11,11,12	0.79	0	15,15,17	1.15	2 (13%)
3	NAG	K	1	3,1	14,14,15	0.63	0	17,19,21	1.10	3 (17%)
3	NAG	K	2	3	14,14,15	0.34	0	17,19,21	0.87	1 (5%)
3	BMA	K	3	3	11,11,12	0.64	0	15,15,17	0.72	0
3	MAN	K	4	3	11,11,12	0.67	0	15,15,17	0.95	2 (13%)
5	NAG	L	1	5,1	14,14,15	0.91	1 (7%)	17,19,21	1.05	2 (11%)
5	NAG	L	2	5	14,14,15	0.30	0	17,19,21	0.47	0
3	NAG	M	1	3,1	14,14,15	0.58	0	17,19,21	0.79	0
3	NAG	M	2	3	14,14,15	0.22	0	17,19,21	0.78	0
3	BMA	M	3	3	11,11,12	0.75	0	15,15,17	1.16	2 (13%)
3	MAN	M	4	3	11,11,12	0.66	0	15,15,17	0.98	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
3	NAG	I	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1
3	BMA	I	3	3	-	1/2/19/22	0/1/1/1
3	MAN	I	4	3	-	0/2/19/22	0/1/1/1
4	NAG	J	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	J	2	4	-	2/6/23/26	0/1/1/1
4	BMA	J	3	4	-	0/2/19/22	0/1/1/1
4	MAN	J	4	4	-	0/2/19/22	0/1/1/1
4	MAN	J	5	4	-	1/2/19/22	0/1/1/1
3	NAG	K	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	K	2	3	-	2/6/23/26	0/1/1/1
3	BMA	K	3	3	-	0/2/19/22	0/1/1/1
3	MAN	K	4	3	-	1/2/19/22	0/1/1/1
5	NAG	L	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	L	2	5	-	0/6/23/26	0/1/1/1
3	NAG	M	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	M	2	3	-	2/6/23/26	0/1/1/1
3	BMA	M	3	3	-	0/2/19/22	0/1/1/1
3	MAN	M	4	3	-	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)
5	L	1	NAG	O5-C1	-2.70	1.39	1.43
4	J	4	MAN	O5-C5	2.38	1.48	1.43

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	4	MAN	C1-O5-C5	6.15	120.43	112.19
4	J	5	MAN	C1-O5-C5	3.32	116.64	112.19
3	M	3	BMA	C1-O5-C5	2.69	115.80	112.19
5	L	1	NAG	C2-N2-C7	2.40	126.12	122.90
3	K	2	NAG	C1-O5-C5	2.39	115.39	112.19
5	L	1	NAG	C4-C3-C2	2.27	114.34	111.02
3	M	4	MAN	O2-C2-C3	-2.26	105.48	110.15
3	I	4	MAN	O2-C2-C3	-2.23	105.52	110.15
4	J	4	MAN	O2-C2-C3	-2.23	105.53	110.15
3	M	4	MAN	C1-O5-C5	2.22	115.16	112.19
3	K	1	NAG	C3-C4-C5	2.20	114.22	110.23
3	K	4	MAN	C1-O5-C5	2.19	115.13	112.19
3	K	1	NAG	C2-N2-C7	2.16	125.80	122.90
3	K	1	NAG	C1-C2-N2	2.15	113.82	110.43
3	K	4	MAN	O2-C2-C3	-2.12	105.77	110.15
3	M	3	BMA	C1-C2-C3	2.02	112.59	109.64
4	J	5	MAN	O2-C2-C3	-2.01	105.98	110.15

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	K	1	NAG	C1-C2-N2-C7
5	L	1	NAG	C1-C2-N2-C7
4	J	2	NAG	O5-C5-C6-O6
4	J	2	NAG	C4-C5-C6-O6
3	K	2	NAG	O5-C5-C6-O6
3	K	2	NAG	C4-C5-C6-O6
3	K	4	MAN	O5-C5-C6-O6
3	I	3	BMA	O5-C5-C6-O6
3	K	1	NAG	O5-C5-C6-O6
3	M	1	NAG	C1-C2-N2-C7
3	M	1	NAG	C3-C2-N2-C7
3	M	2	NAG	C3-C2-N2-C7
3	M	2	NAG	C1-C2-N2-C7

*Continued on next page...*

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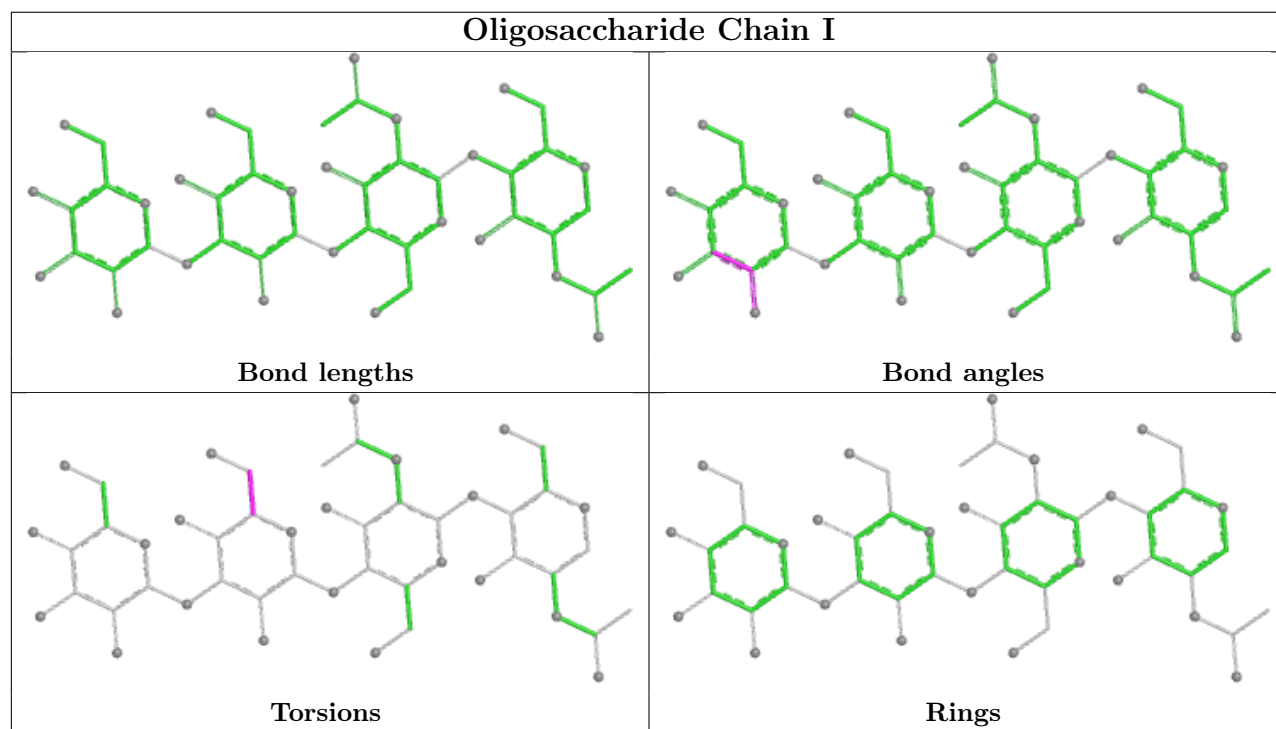
Mol	Chain	Res	Type	Atoms
4	J	5	MAN	C4-C5-C6-O6

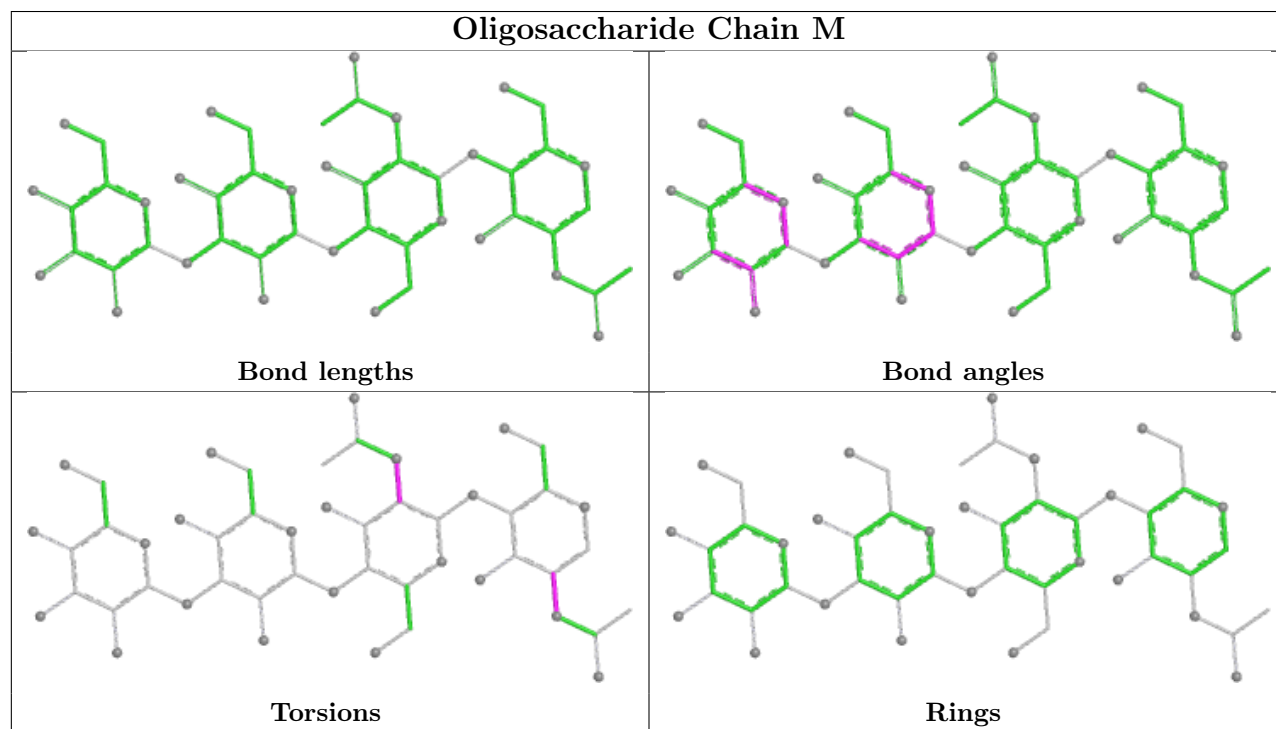
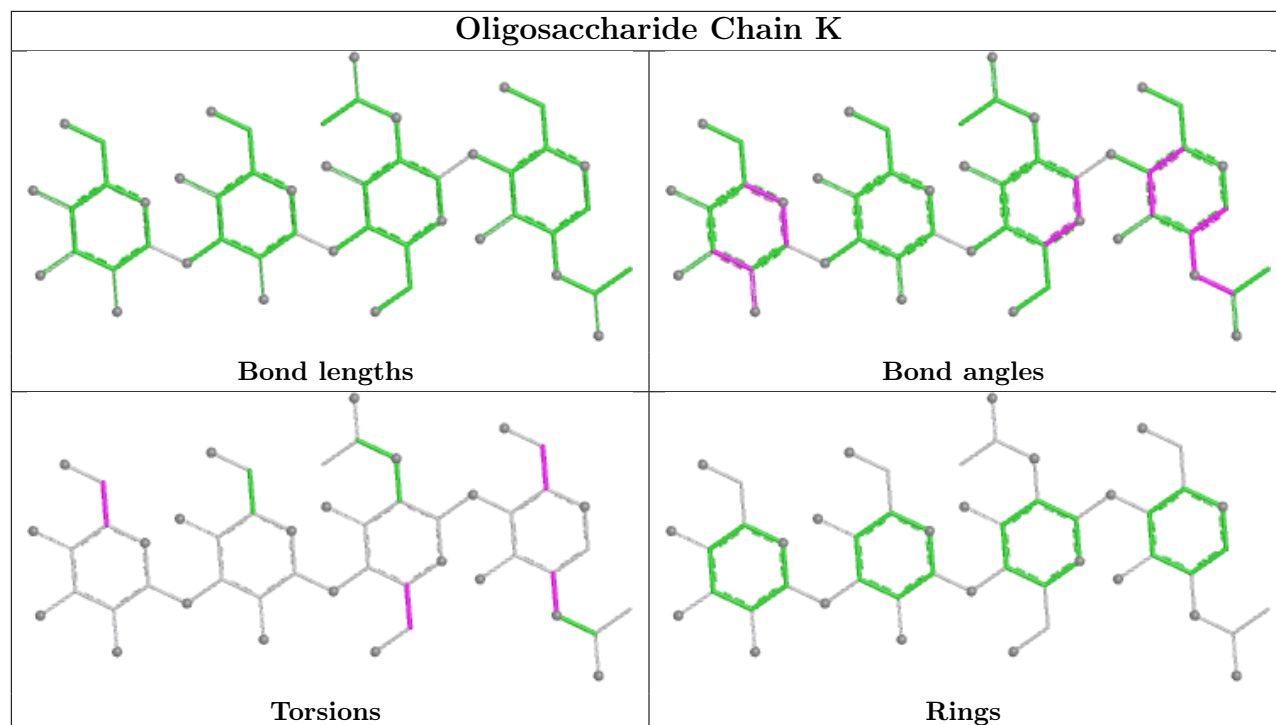
There are no ring outliers.

3 monomers are involved in 3 short contacts:

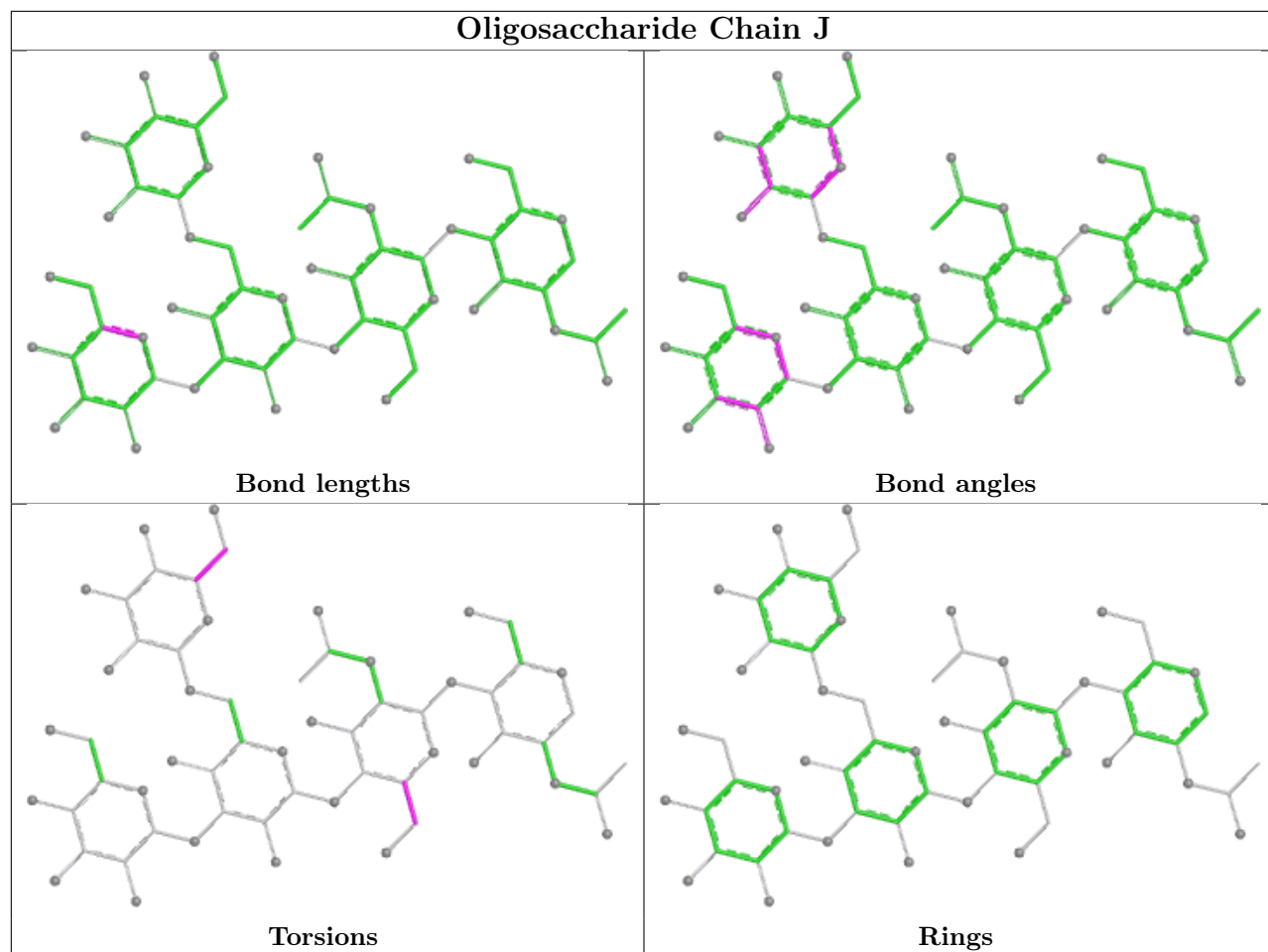
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	M	1	NAG	1	0
3	K	1	NAG	1	0
3	K	2	NAG	1	0

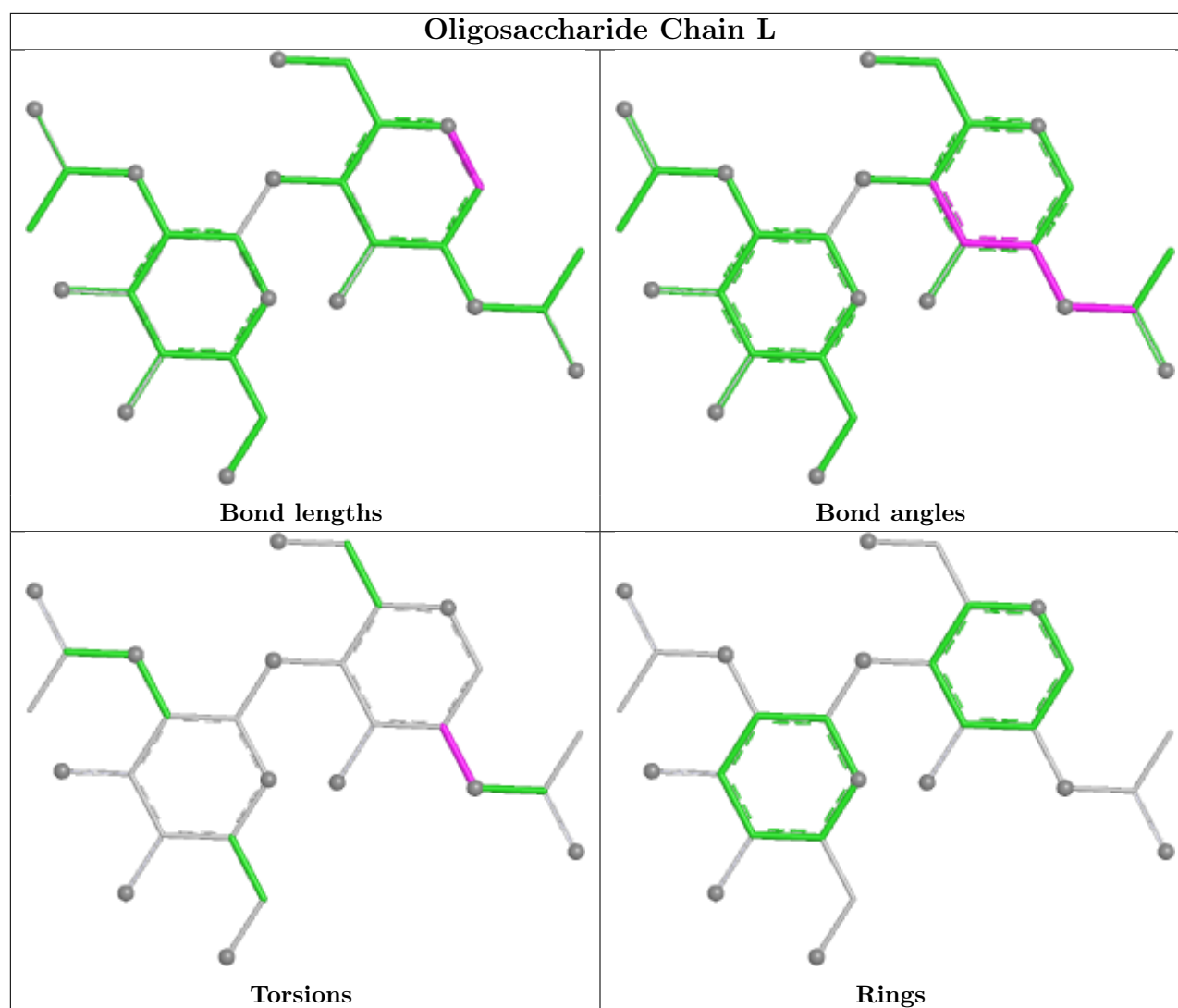
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











## 5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
6	NAG	B	601	1	14,14,15	0.23	0	17,19,21	0.59	0
6	NAG	A	3301	1	14,14,15	0.35	0	17,19,21	0.45	0
6	NAG	B	602	1	14,14,15	0.19	0	17,19,21	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	NAG	C	601	1	14,14,15	0.27	0	17,19,21	0.46	0
6	NAG	D	602	1	14,14,15	0.29	0	17,19,21	0.42	0
6	NAG	D	601	1	14,14,15	0.24	0	17,19,21	0.51	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
6	NAG	B	601	1	-	0/6/23/26	0/1/1/1
6	NAG	A	3301	1	-	0/6/23/26	0/1/1/1
6	NAG	B	602	1	-	0/6/23/26	0/1/1/1
6	NAG	C	601	1	-	3/6/23/26	0/1/1/1
6	NAG	D	602	1	-	0/6/23/26	0/1/1/1
6	NAG	D	601	1	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	601	NAG	O5-C5-C6-O6
6	D	601	NAG	C4-C5-C6-O6
6	C	601	NAG	O5-C5-C6-O6
6	C	601	NAG	C1-C2-N2-C7
6	C	601	NAG	C3-C2-N2-C7
6	D	601	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	3301	NAG	2	0
6	C	601	NAG	2	0
6	D	601	NAG	4	0

## 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](#)

EDS failed to run properly - this section is therefore empty.

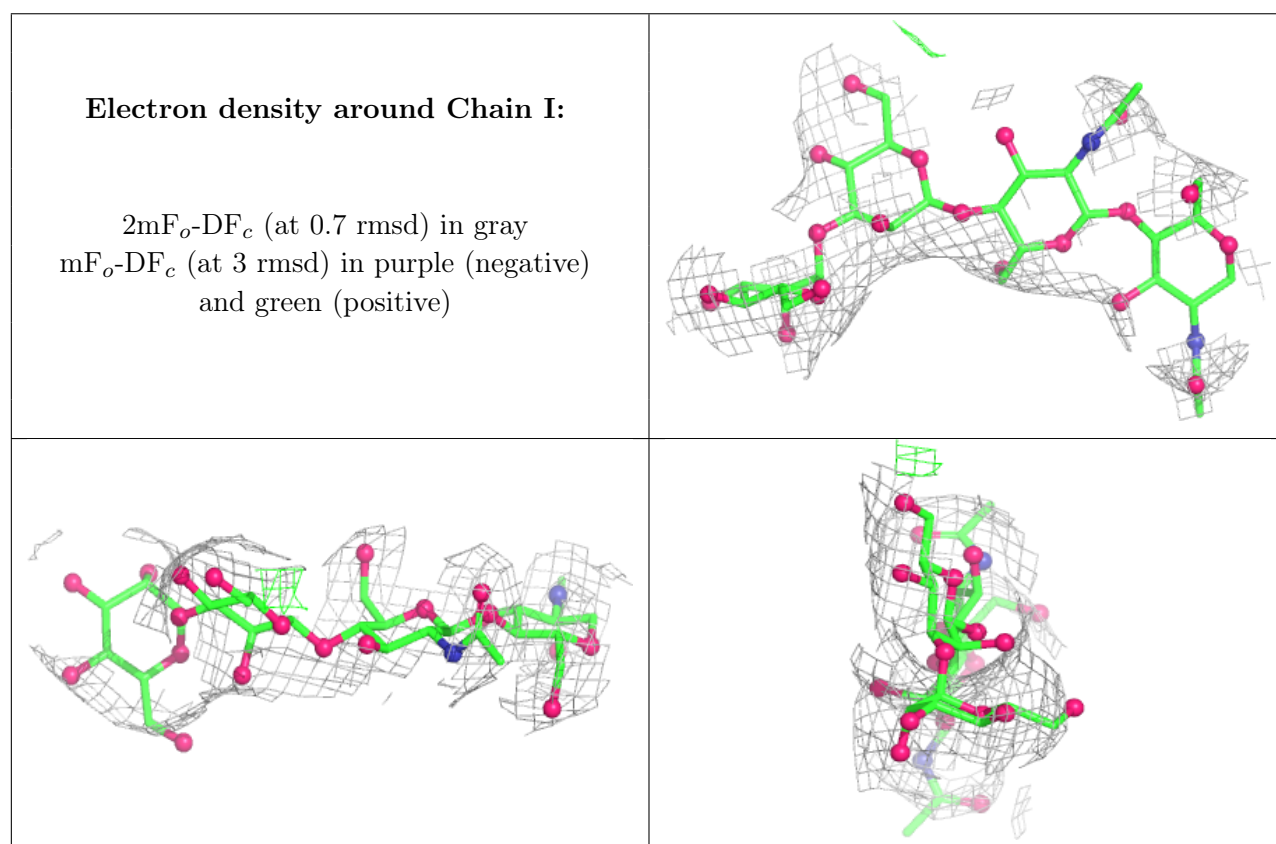
### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS failed to run properly - this section is therefore empty.

### 6.3 Carbohydrates [i](#)

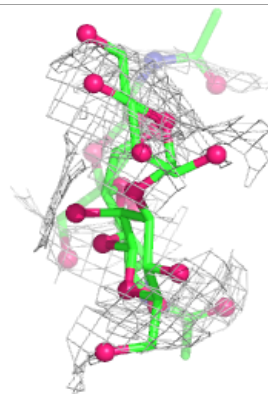
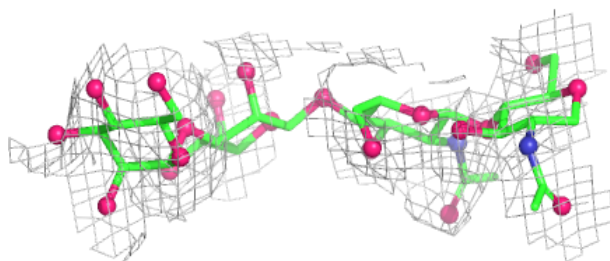
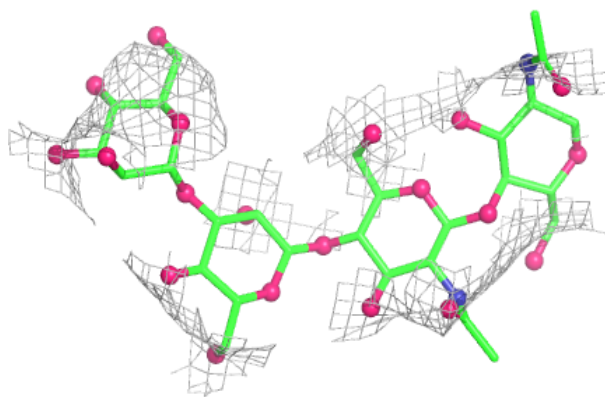
EDS failed to run properly - this section is therefore empty.

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.

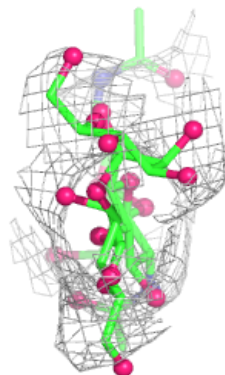
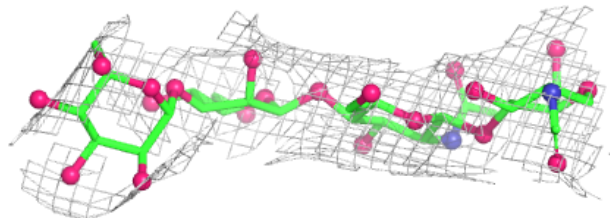
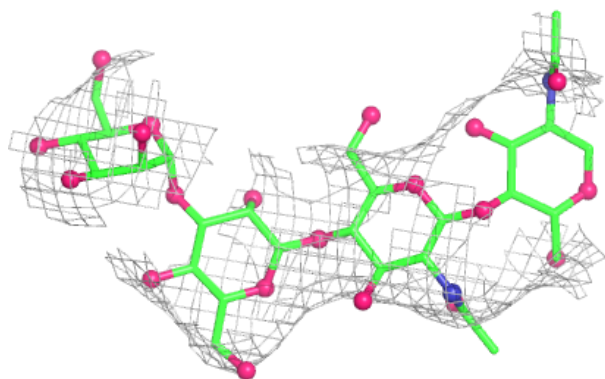


**Electron density around Chain K:**

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

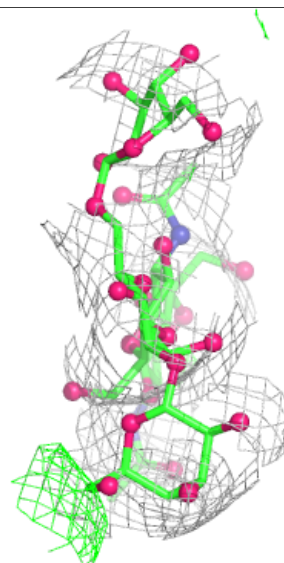
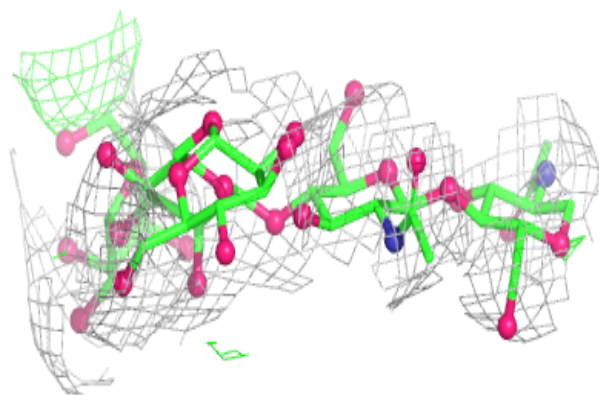
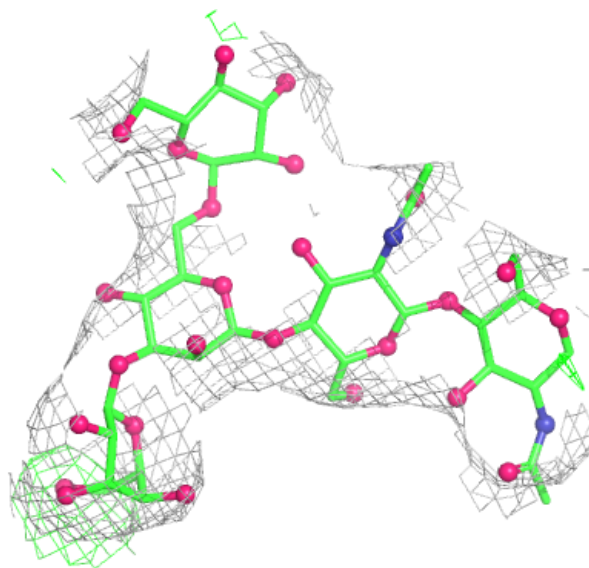
**Electron density around Chain M:**

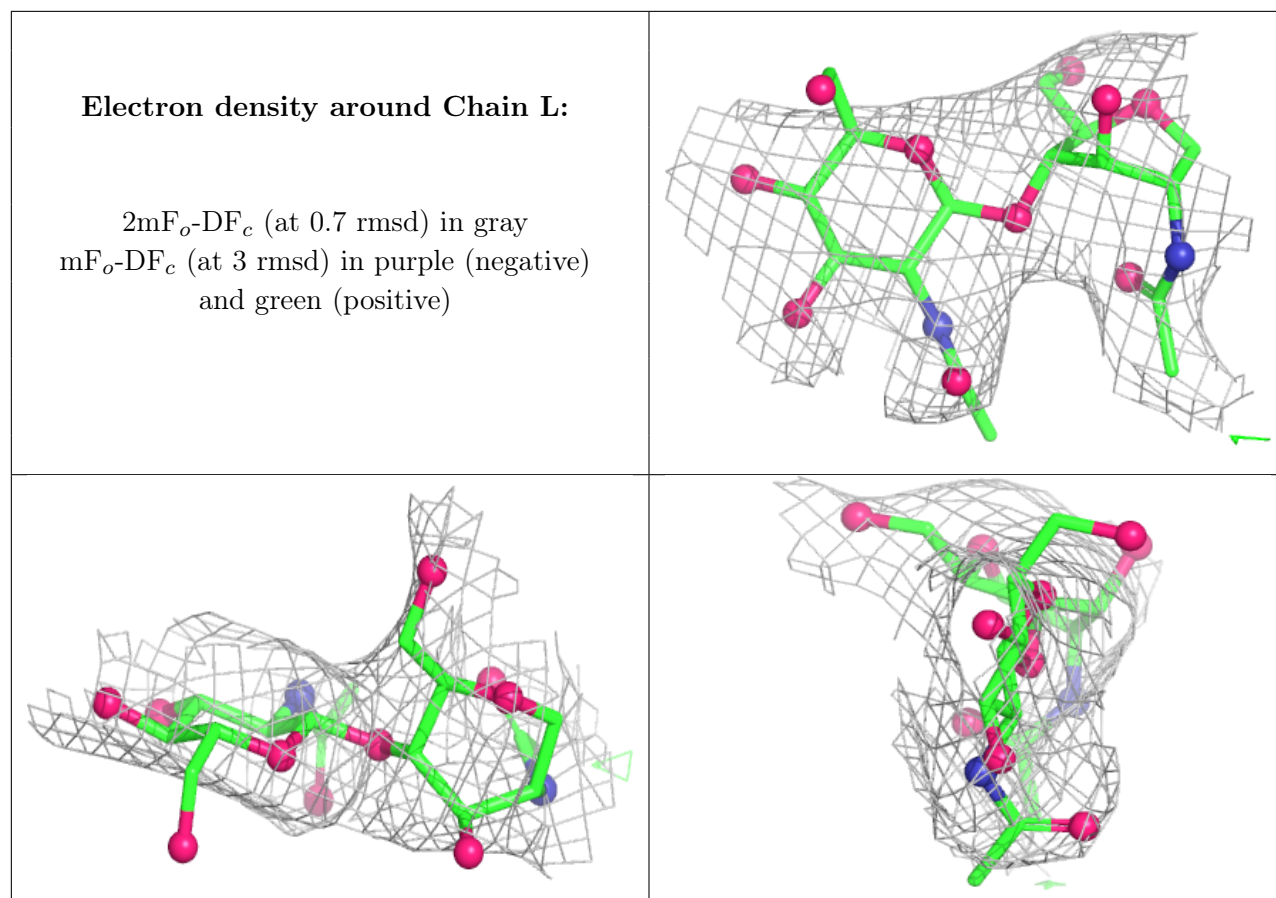
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around Chain J:**

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





## 6.4 Ligands [i](#)

EDS failed to run properly - this section is therefore empty.

## 6.5 Other polymers [i](#)

EDS failed to run properly - this section is therefore empty.