



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 11:26 PM BST

PDB ID : 4LSX  
Title : Plant steroid receptor ectodomain bound to brassinolide and SERK1 co-receptor ectodomain  
Authors : Santiago, J.; Henzler, C.; Hothorn, M.  
Deposited on : 2013-07-23  
Resolution : 3.30 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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

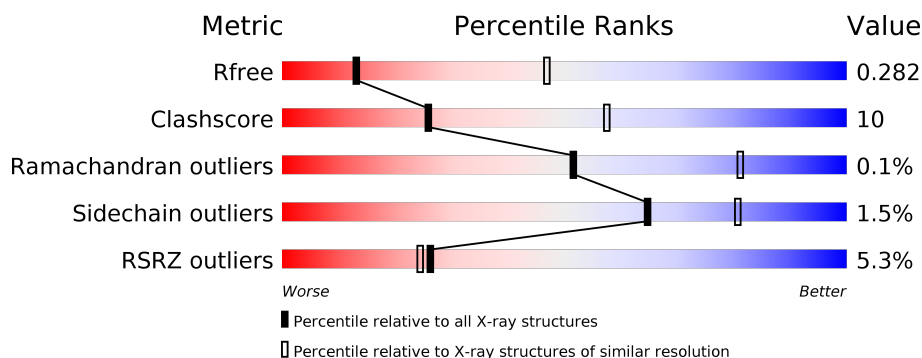
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.30 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.



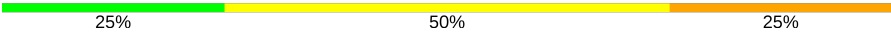
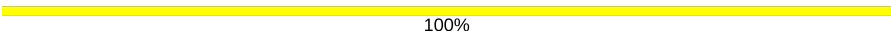


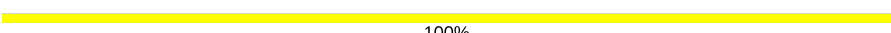
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 on 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  $\leq 5\%$ . 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	774	<div> <div>3%</div> <div> <div></div> <div>72%</div> <div>23%</div> <div>• •</div> </div> </div>
1	B	774	<div> <div>3%</div> <div> <div></div> <div>69%</div> <div>22%</div> <div>• 8%</div> </div> </div>
2	C	203	<div> <div>9%</div> <div> <div></div> <div>66%</div> <div>25%</div> <div>9%</div> </div> </div>
2	D	203	<div> <div>16%</div> <div> <div></div> <div>67%</div> <div>23%</div> <div>9%</div> </div> </div>
3	E	4	<div> <div></div> <div> <div>50%</div> <div>25%</div> <div>25%</div> </div> </div>
3	H	4	<div> <div></div> <div> <div>25%</div> <div>50%</div> <div>25%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	J	4	 25% 50% 25%
4	F	2	 100%
4	K	2	 100%
5	G	5	 40% 40% 20%
6	I	3	 100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
8	NAG	A	802	-	-	-	X

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 13920 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 Protein BRASSINOSTEROID INSENSITIVE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	740	Total	C	N	O	S	0	0	0
			5454	3421	914	1088	31			
1	B	710	Total	C	N	O	S	0	0	0
			5217	3283	868	1037	29			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	GLY	-	expression tag	UNP O22476
A	25	SER	-	expression tag	UNP O22476
A	26	SER	-	expression tag	UNP O22476
A	27	MET	-	expression tag	UNP O22476
A	28	GLY	-	expression tag	UNP O22476
A	643	GLU	GLY	engineered mutation	UNP O22476
A	789	LEU	-	expression tag	UNP O22476
A	790	GLU	-	expression tag	UNP O22476
A	791	ASN	-	expression tag	UNP O22476
A	792	LEU	-	expression tag	UNP O22476
A	793	TYR	-	expression tag	UNP O22476
A	794	PHE	-	expression tag	UNP O22476
A	795	GLN	-	expression tag	UNP O22476
A	796	GLY	-	expression tag	UNP O22476
A	797	ALA	-	expression tag	UNP O22476
B	24	GLY	-	expression tag	UNP O22476
B	25	SER	-	expression tag	UNP O22476
B	26	SER	-	expression tag	UNP O22476
B	27	MET	-	expression tag	UNP O22476
B	28	GLY	-	expression tag	UNP O22476
B	643	GLU	GLY	engineered mutation	UNP O22476
B	789	LEU	-	expression tag	UNP O22476
B	790	GLU	-	expression tag	UNP O22476
B	791	ASN	-	expression tag	UNP O22476
B	792	LEU	-	expression tag	UNP O22476

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Chain	Residue	Modelled	Actual	Comment	Reference
B	793	TYR	-	expression tag	UNP O22476
B	794	PHE	-	expression tag	UNP O22476
B	795	GLN	-	expression tag	UNP O22476
B	796	GLY	-	expression tag	UNP O22476
B	797	ALA	-	expression tag	UNP O22476

- Molecule 2 is a protein called Somatic embryogenesis receptor kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	185	Total	C	N	O	S	0	0	0
			1397	881	238	273	5			
2	D	185	Total	C	N	O	S	0	0	0
			1380	871	234	271	4			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	20	GLY	-	expression tag	UNP Q94AG2
C	21	SER	-	expression tag	UNP Q94AG2
C	22	SER	-	expression tag	UNP Q94AG2
C	23	MET	-	expression tag	UNP Q94AG2
C	214	LEU	-	expression tag	UNP Q94AG2
C	215	GLU	-	expression tag	UNP Q94AG2
C	216	ASN	-	expression tag	UNP Q94AG2
C	217	LEU	-	expression tag	UNP Q94AG2
C	218	TYR	-	expression tag	UNP Q94AG2
C	219	PHE	-	expression tag	UNP Q94AG2
C	220	GLN	-	expression tag	UNP Q94AG2
C	221	GLY	-	expression tag	UNP Q94AG2
C	222	ALA	-	expression tag	UNP Q94AG2
D	20	GLY	-	expression tag	UNP Q94AG2
D	21	SER	-	expression tag	UNP Q94AG2
D	22	SER	-	expression tag	UNP Q94AG2
D	23	MET	-	expression tag	UNP Q94AG2
D	214	LEU	-	expression tag	UNP Q94AG2
D	215	GLU	-	expression tag	UNP Q94AG2
D	216	ASN	-	expression tag	UNP Q94AG2
D	217	LEU	-	expression tag	UNP Q94AG2
D	218	TYR	-	expression tag	UNP Q94AG2
D	219	PHE	-	expression tag	UNP Q94AG2
D	220	GLN	-	expression tag	UNP Q94AG2
D	221	GLY	-	expression tag	UNP Q94AG2

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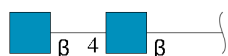
Chain	Residue	Modelled	Actual	Comment	Reference
D	222	ALA	-	expression tag	UNP Q94AG2

- 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.



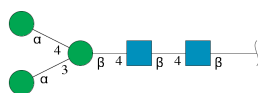
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	4	Total	C	N	O	0	0	0
			50	28	2	20			
3	H	4	Total	C	N	O	0	0	0
			50	28	2	20			
3	J	4	Total	C	N	O	0	0	0
			50	28	2	20			

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



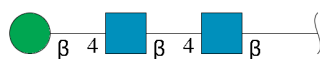
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	K	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



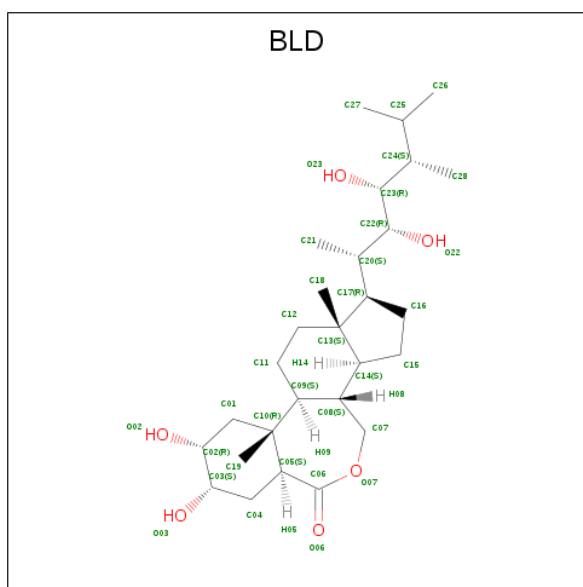
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	G	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 6 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.



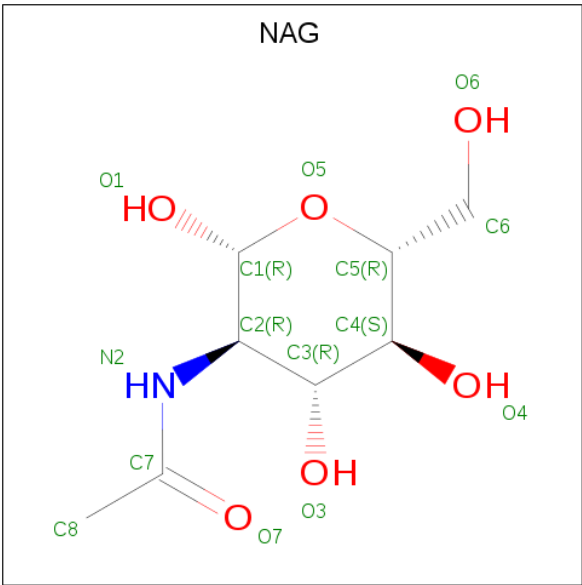
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	I	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 7 is Brassinolide (three-letter code: BLD) (formula:  $C_{28}H_{48}O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			34	28	6		
7	B	1	Total	C	O	0	0
			34	28	6		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



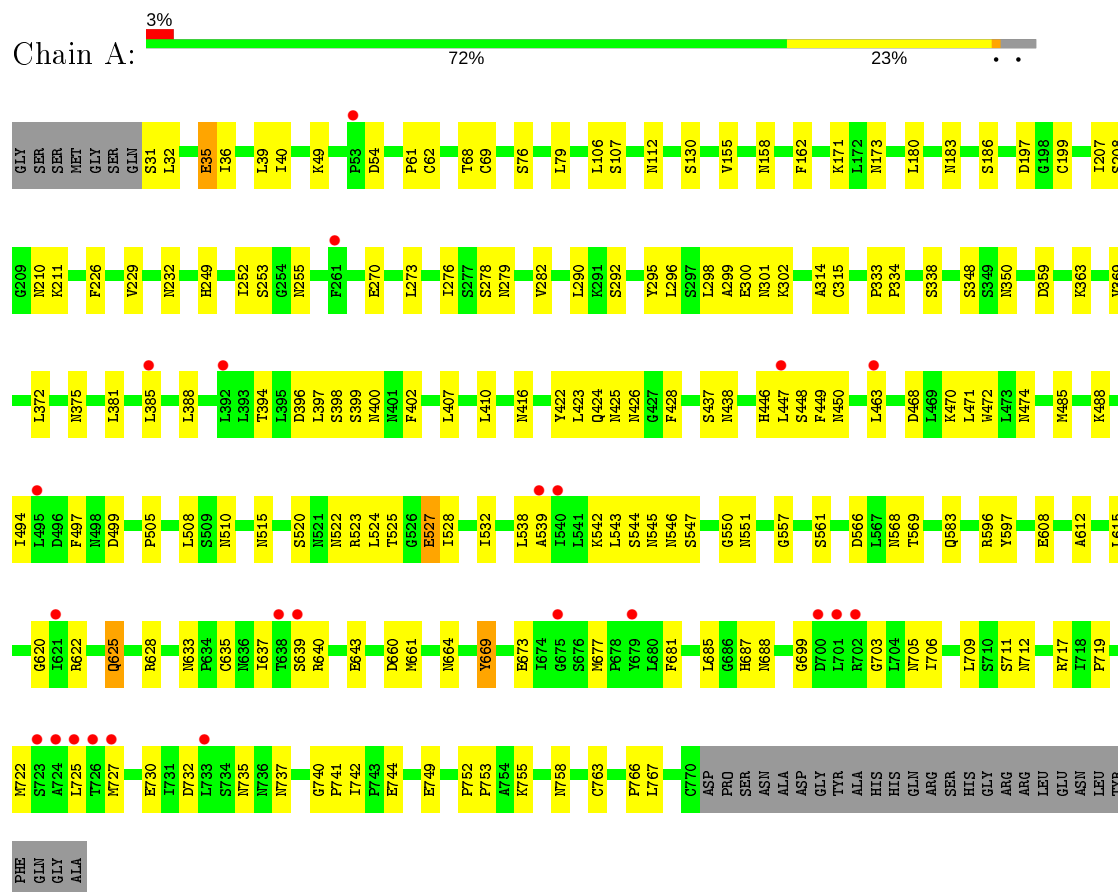
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	B	1	Total	C	N	O	0	0
			14	8	1	5		
8	C	1	Total	C	N	O	0	0
			14	8	1	5		
8	C	1	Total	C	N	O	0	0
			14	8	1	5		
8	D	1	Total	C	N	O	0	0
			14	8	1	5		
8	D	1	Total	C	N	O	0	0
			14	8	1	5		



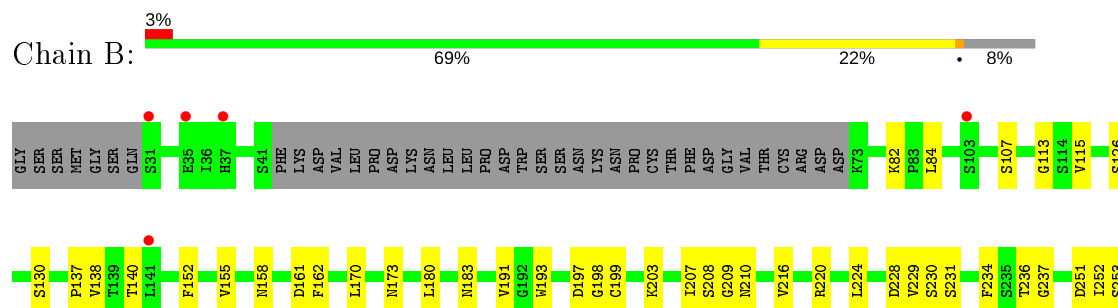
### 3 Residue-property plots [i](#)

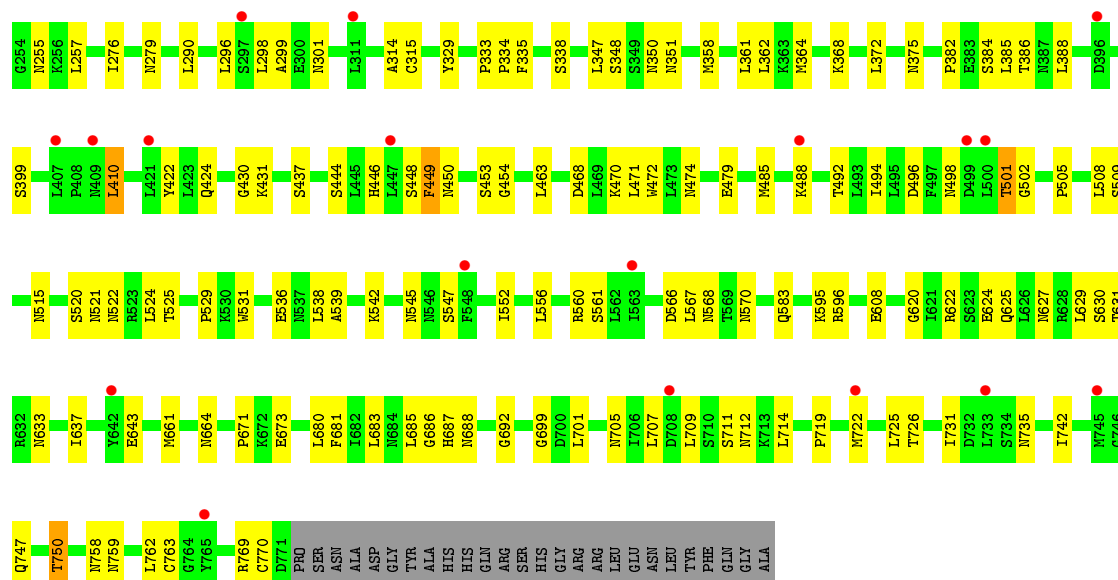
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

#### • Molecule 1: Protein BRASSINOSTEROID INSENSITIVE 1

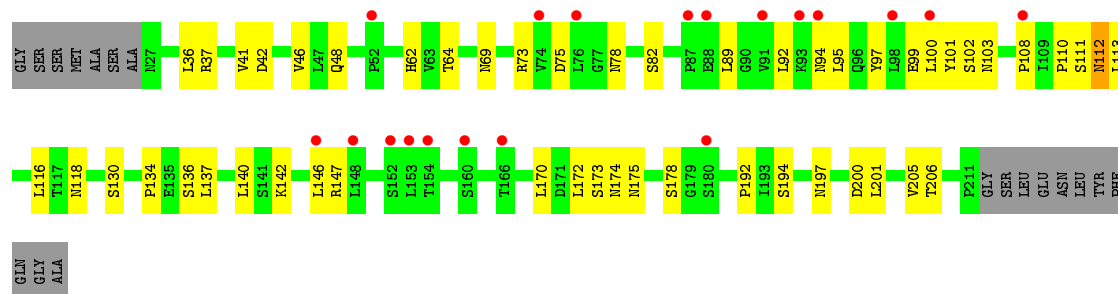


#### • Molecule 1: Protein BRASSINOSTEROID INSENSITIVE 1

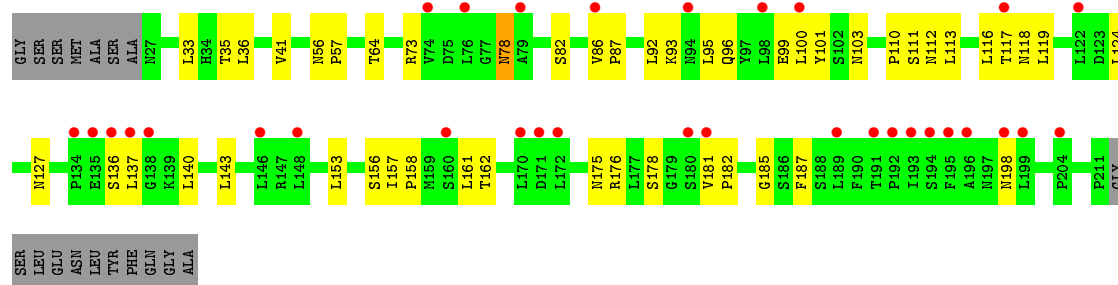




• Molecule 2: Somatic embryogenesis receptor kinase 1



• Molecule 2: Somatic embryogenesis receptor kinase 1



• 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



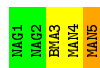
- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



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



- Molecule 6: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.90Å 69.90Å 873.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.53 – 3.30 48.53 – 3.30	Depositor EDS
% Data completeness (in resolution range)	98.2 (48.53-3.30) 96.4 (48.53-3.30)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.87 (at 3.33Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1334)	Depositor
R, $R_{free}$	0.246 , 0.285 0.243 , 0.282	Depositor DCC
$R_{free}$ test set	2010 reflections (5.70%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	86.0	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 75.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.458 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	13920	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	122.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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: BLD, BMA, NAG, 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	0.23	0/5553	0.45	0/7543
1	B	0.22	0/5311	0.44	0/7218
2	C	0.25	0/1427	0.47	0/1959
2	D	0.24	0/1410	0.47	0/1940
All	All	0.23	0/13701	0.45	0/18660

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	5454	0	5251	113	0
1	B	5217	0	5032	109	0
2	C	1397	0	1366	31	0
2	D	1380	0	1332	30	0
3	E	50	0	43	1	0
3	H	50	0	43	2	0
3	J	50	0	43	2	0
4	F	28	0	25	2	0
4	K	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	G	61	0	52	1	0
6	I	39	0	34	0	0
7	A	34	0	47	5	0
7	B	34	0	47	5	0
8	A	28	0	26	3	0
8	B	14	0	13	1	0
8	C	28	0	26	0	0
8	D	28	0	26	0	0
All	All	13920	0	13431	286	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 10.

All (286) 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:551:ASN:HD21	8:A:814:NAG:H82	1.39	0.86
1:A:290:LEU:HB3	1:A:314:ALA:HB2	1.71	0.73
2:C:64:THR:HB	2:C:73:ARG:HB2	1.72	0.72
2:D:140:LEU:HD12	2:D:143:LEU:HD22	1.72	0.71
1:B:198:GLY:HA2	1:B:220:ARG:HH12	1.56	0.71
2:C:94:ASN:HA	2:C:118:ASN:HD22	1.56	0.70
1:A:155:VAL:O	1:A:158:ASN:ND2	2.25	0.69
1:B:191:VAL:HG11	1:B:216:VAL:HG13	1.74	0.68
1:B:449:PHE:HE1	1:B:625:GLN:HB3	1.58	0.68
1:B:661:MET:O	1:B:664:ASN:ND2	2.26	0.67
1:A:685:LEU:O	1:A:688:ASN:ND2	2.27	0.67
1:B:347:LEU:O	1:B:350:ASN:ND2	2.28	0.67
1:B:290:LEU:HB3	1:B:314:ALA:HB2	1.77	0.67
1:A:633:ASN:ND2	1:A:635:CYS:SG	2.67	0.67
5:G:3:BMA:H62	5:G:5:MAN:H2	1.77	0.67
2:D:162:THR:HB	2:D:185:GLY:HA3	1.77	0.65
1:A:397:LEU:O	1:A:400:ASN:ND2	2.29	0.65
1:A:423:LEU:O	1:A:426:ASN:ND2	2.30	0.65
1:B:437:SER:HA	1:B:463:LEU:HD21	1.78	0.65
1:A:229:VAL:O	1:A:232:ASN:ND2	2.29	0.65
1:A:471:LEU:O	1:A:474:ASN:ND2	2.28	0.65
1:A:520:SER:O	1:A:522:ASN:ND2	2.30	0.64
1:A:687:HIS:HA	1:A:711:SER:HB2	1.79	0.64
1:A:753:PRO:HG3	1:A:766:PRO:HB2	1.78	0.64
1:A:468:ASP:HB3	1:A:470:LYS:HE3	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:685:LEU:O	1:B:688:ASN:ND2	2.32	0.63
1:A:709:LEU:O	1:A:712:ASN:ND2	2.31	0.63
2:C:172:LEU:O	2:C:175:ASN:ND2	2.31	0.63
1:A:276:ILE:O	1:A:279:ASN:ND2	2.31	0.62
1:B:155:VAL:O	1:B:158:ASN:ND2	2.28	0.62
1:B:180:LEU:O	1:B:183:ASN:ND2	2.33	0.62
1:A:35:GLU:O	1:A:39:LEU:HG	1.99	0.62
1:B:520:SER:O	1:B:522:ASN:ND2	2.31	0.62
1:B:515:ASN:HA	1:B:538:LEU:HA	1.81	0.62
2:D:100:LEU:O	2:D:103:ASN:ND2	2.32	0.61
1:A:207:ILE:HG22	1:A:229:VAL:HA	1.81	0.61
1:A:661:MET:O	1:A:664:ASN:ND2	2.34	0.61
2:D:110:PRO:HB2	2:D:113:LEU:HG	1.82	0.61
1:A:472:TRP:HB3	1:A:494:ILE:HG22	1.83	0.60
1:A:551:ASN:ND2	8:A:814:NAG:H82	2.15	0.60
1:B:431:LYS:HG3	1:B:454:GLY:HA3	1.82	0.60
1:B:567:LEU:O	1:B:570:ASN:ND2	2.34	0.60
1:A:180:LEU:O	1:A:183:ASN:ND2	2.33	0.60
1:B:747:GLN:HG2	2:D:101:TYR:OH	2.01	0.60
1:B:471:LEU:O	1:B:474:ASN:ND2	2.35	0.60
2:D:157:ILE:HD12	2:D:181:VAL:HG22	1.84	0.60
1:A:448:SER:O	1:A:450:ASN:ND2	2.34	0.60
1:A:545:ASN:HA	1:A:569:THR:HB	1.83	0.60
1:B:207:ILE:O	1:B:210:ASN:ND2	2.34	0.60
1:A:543:LEU:O	1:A:546:ASN:ND2	2.34	0.60
1:A:299:ALA:O	1:A:301:ASN:ND2	2.35	0.59
2:C:36:LEU:HD22	2:C:89:LEU:HD21	1.83	0.59
2:C:73:ARG:HG2	2:C:97:TYR:HB2	1.84	0.59
7:A:801:BLD:O02	2:C:62:HIS:ND1	2.35	0.59
1:A:437:SER:HA	1:A:463:LEU:HD21	1.83	0.59
2:C:100:LEU:O	2:C:103:ASN:ND2	2.36	0.59
2:D:92:LEU:HD12	2:D:95:LEU:HD22	1.83	0.59
1:A:719:PRO:HG2	1:A:722:MET:HG3	1.85	0.58
1:A:735:ASN:OD1	1:A:758:ASN:ND2	2.35	0.58
1:B:424:GLN:HE21	1:B:620:GLY:HA3	1.69	0.58
1:A:399:SER:O	1:A:622:ARG:NH2	2.37	0.58
2:C:92:LEU:HD12	2:C:95:LEU:HD22	1.86	0.57
1:B:372:LEU:O	1:B:375:ASN:ND2	2.34	0.57
2:D:87:PRO:O	2:D:112:ASN:ND2	2.35	0.57
2:C:174:ASN:OD1	2:C:197:ASN:ND2	2.28	0.57
3:E:2:NAG:H3	3:E:2:NAG:H83	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:732:ASP:HA	1:A:755:LYS:HB3	1.86	0.56
2:C:110:PRO:HB2	2:C:113:LEU:HG	1.87	0.56
1:B:735:ASN:OD1	1:B:758:ASN:ND2	2.37	0.56
1:A:527:GLU:OE1	1:A:528:ILE:N	2.34	0.56
1:B:399:SER:O	1:B:622:ARG:NH2	2.39	0.56
3:H:2:NAG:H3	3:H:2:NAG:H83	1.86	0.55
1:A:252:ILE:O	1:A:255:ASN:ND2	2.36	0.55
1:B:315:CYS:SG	1:B:338:SER:OG	2.64	0.55
1:A:706:ILE:HG12	1:A:730:GLU:HB3	1.88	0.55
7:A:801:BLD:O02	2:C:62:HIS:N	2.38	0.55
2:D:99:GLU:HB3	2:D:101:TYR:CE2	2.42	0.55
1:B:472:TRP:HB3	1:B:494:ILE:HG22	1.89	0.55
1:B:750:THR:HG23	2:D:73:ARG:HH12	1.72	0.55
2:D:153:LEU:HB2	2:D:175:ASN:HD21	1.71	0.55
1:A:253:SER:HB2	4:F:1:NAG:H62	1.89	0.55
1:A:424:GLN:HE21	1:A:620:GLY:HA3	1.71	0.54
2:D:95:LEU:HD23	2:D:116:LEU:HD13	1.90	0.54
1:A:525:THR:HG22	1:A:547:SER:HB2	1.89	0.54
2:D:64:THR:HB	2:D:73:ARG:HB2	1.88	0.54
1:A:68:THR:HB	1:A:76:SER:HB3	1.90	0.54
1:B:208:SER:O	1:B:210:ASN:ND2	2.41	0.54
1:A:717:ARG:HG2	1:A:740:GLY:HA3	1.90	0.54
1:A:515:ASN:HA	1:A:538:LEU:HA	1.90	0.54
1:A:542:LYS:NZ	1:A:566:ASP:OD1	2.41	0.54
1:B:536:GLU:O	1:B:561:SER:OG	2.20	0.54
1:A:596:ARG:NH2	1:A:643:GLU:OE2	2.39	0.53
1:A:640:ARG:NH2	2:C:69:ASN:OD1	2.40	0.53
1:A:597:TYR:HB2	1:A:615:LEU:HD11	1.90	0.52
1:B:203:LYS:HA	1:B:224:LEU:HA	1.91	0.52
1:B:207:ILE:HD11	1:B:229:VAL:HG12	1.91	0.52
1:B:560:ARG:NH1	1:B:583:GLN:OE1	2.41	0.52
1:A:375:ASN:O	1:A:622:ARG:NH2	2.43	0.52
1:A:568:ASN:OD1	1:A:569:THR:N	2.42	0.52
1:B:686:GLY:O	1:B:688:ASN:ND2	2.41	0.52
1:B:680:LEU:HD21	1:B:683:LEU:HB2	1.91	0.52
1:B:596:ARG:NH2	1:B:643:GLU:OE2	2.42	0.52
1:B:448:SER:O	1:B:450:ASN:ND2	2.43	0.52
1:A:717:ARG:HE	1:A:741:PRO:HD2	1.75	0.52
1:B:681:PHE:CD2	7:B:801:BLD:H112	2.45	0.52
1:B:252:ILE:O	1:B:255:ASN:ND2	2.43	0.51
1:A:270:GLU:OE1	1:A:292:SER:OG	2.23	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:MET:O	1:A:488:LYS:NZ	2.32	0.51
1:B:479:GLU:HG2	1:B:502:GLY:HA3	1.92	0.51
1:B:566:ASP:OD2	1:B:568:ASN:ND2	2.40	0.51
1:A:669:TYR:HD1	1:A:669:TYR:H	1.57	0.51
1:A:207:ILE:O	1:A:210:ASN:ND2	2.44	0.50
1:B:276:ILE:O	1:B:279:ASN:ND2	2.43	0.50
1:A:171:LYS:HA	1:A:197:ASP:HB2	1.93	0.50
1:A:542:LYS:HG3	1:A:566:ASP:HB3	1.92	0.50
2:D:93:LYS:O	2:D:118:ASN:ND2	2.44	0.50
1:B:348:SER:O	1:B:350:ASN:ND2	2.44	0.50
1:B:631:THR:O	3:J:2:NAG:O4	2.30	0.50
1:B:362:LEU:HD11	1:B:384:SER:HB2	1.92	0.50
1:B:505:PRO:HG2	1:B:508:LEU:HG	1.93	0.50
2:D:41:VAL:HB	2:D:82:SER:HB3	1.93	0.50
1:B:299:ALA:O	1:B:301:ASN:ND2	2.45	0.50
1:B:422:TYR:HE1	1:B:444:SER:HG	1.59	0.50
1:A:744:GLU:OE1	1:A:744:GLU:N	2.40	0.50
1:B:333:PRO:HB2	1:B:335:PHE:CD2	2.47	0.50
1:A:295:TYR:CZ	4:F:1:NAG:H82	2.47	0.50
1:A:226:PHE:HA	1:A:249:HIS:HB3	1.94	0.49
1:B:138:VAL:N	1:B:161:ASP:O	2.45	0.49
1:B:173:ASN:ND2	1:B:197:ASP:OD1	2.45	0.49
2:C:192:PRO:HG3	2:C:205:VAL:O	2.11	0.49
1:A:416:ASN:ND2	1:A:438:ASN:O	2.45	0.49
1:B:629:LEU:O	1:B:633:ASN:N	2.45	0.49
1:B:742:ILE:N	1:B:763:CYS:O	2.42	0.49
1:A:699:GLY:HA2	1:A:725:LEU:HD21	1.94	0.49
1:A:208:SER:O	1:A:210:ASN:ND2	2.45	0.49
1:A:488:LYS:HZ2	1:A:510:ASN:HB3	1.78	0.49
1:A:425:ASN:H	1:A:449:PHE:HB3	1.78	0.49
1:A:639:SER:HB3	1:A:640:ARG:HB2	1.94	0.49
1:A:673:GLU:N	1:A:673:GLU:OE1	2.46	0.49
1:A:278:SER:H	1:A:300:GLU:HB2	1.78	0.49
1:A:557:GLY:O	1:A:583:GLN:HB2	2.13	0.49
2:C:41:VAL:HB	2:C:82:SER:HB2	1.94	0.49
1:A:385:LEU:HD12	1:A:388:LEU:HD12	1.94	0.48
1:A:539:ALA:HA	1:A:561:SER:O	2.13	0.48
1:B:329:TYR:HB3	1:B:351:ASN:HB3	1.94	0.48
1:B:358:MET:HG3	1:B:382:PRO:HB2	1.95	0.48
2:C:118:ASN:HA	2:C:142:LYS:HD2	1.94	0.48
1:B:296:LEU:HD11	1:B:298:LEU:HD13	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:501:THR:HG22	1:B:502:GLY:H	1.78	0.48
1:B:542:LYS:NZ	1:B:566:ASP:OD1	2.47	0.48
2:C:146:LEU:O	2:C:170:LEU:HA	2.14	0.48
2:C:178:SER:HA	2:C:200:ASP:O	2.13	0.48
1:B:709:LEU:O	1:B:712:ASN:ND2	2.45	0.48
1:A:348:SER:O	1:A:350:ASN:ND2	2.47	0.48
1:B:335:PHE:O	1:B:338:SER:OG	2.22	0.48
1:A:173:ASN:ND2	1:A:197:ASP:OD1	2.47	0.47
1:A:742:ILE:N	1:A:763:CYS:O	2.45	0.47
2:C:78:ASN:HD22	2:C:102:SER:HB3	1.79	0.47
2:C:99:GLU:HB3	2:C:101:TYR:CE2	2.50	0.47
1:A:315:CYS:SG	1:A:338:SER:OG	2.71	0.47
1:A:372:LEU:O	1:A:375:ASN:ND2	2.41	0.47
1:A:527:GLU:HG2	1:A:550:GLY:HA3	1.97	0.47
1:B:209:GLY:H	1:B:231:SER:HB2	1.80	0.47
2:C:112:ASN:HD22	2:C:112:ASN:H	1.61	0.47
1:A:449:PHE:HE1	1:A:625:GLN:HB3	1.80	0.47
1:B:385:LEU:HD12	1:B:388:LEU:HD12	1.95	0.47
1:A:608:GLU:OE1	1:A:608:GLU:N	2.48	0.47
1:A:36:ILE:O	1:A:40:ILE:HG12	2.15	0.46
1:B:525:THR:HG22	1:B:547:SER:HB2	1.96	0.46
1:B:692:GLY:H	1:B:714:LEU:HA	1.80	0.46
1:B:386:THR:HG22	1:B:410:LEU:HA	1.97	0.46
1:B:687:HIS:HA	1:B:711:SER:HB2	1.96	0.46
1:A:62:CYS:HB3	1:A:69:CYS:HB3	1.91	0.46
1:A:664:ASN:N	1:A:688:ASN:OD1	2.46	0.46
1:A:752:PRO:HA	1:A:753:PRO:HD3	1.85	0.46
2:C:173:SER:N	2:C:194:SER:O	2.48	0.46
1:A:186:SER:HA	1:A:211:LYS:O	2.14	0.46
1:A:447:LEU:O	1:A:450:ASN:ND2	2.41	0.46
1:B:468:ASP:HB3	1:B:470:LYS:HE3	1.98	0.46
1:B:449:PHE:CE1	1:B:625:GLN:HB3	2.45	0.46
2:C:62:HIS:HB3	2:C:75:ASP:O	2.15	0.46
2:D:35:THR:HG22	2:D:86:VAL:HG11	1.98	0.46
1:B:759:ASN:HB3	1:B:762:LEU:HB2	1.99	0.45
1:B:333:PRO:HA	1:B:334:PRO:HD3	1.81	0.45
1:B:681:PHE:HD2	7:B:801:BLD:H112	1.81	0.45
1:B:236:THR:OG1	1:B:237:GLY:N	2.47	0.45
2:C:134:PRO:HB2	2:C:137:LEU:HG	1.98	0.45
2:C:42:ASP:OD2	2:C:46:VAL:N	2.50	0.45
1:A:722:MET:O	1:A:725:LEU:HG	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:182:PRO:HG2	2:D:187:PHE:CD2	2.52	0.45
1:A:333:PRO:HA	1:A:334:PRO:HD3	1.75	0.45
2:D:117:THR:OG1	2:D:118:ASN:ND2	2.50	0.45
2:D:124:LEU:O	2:D:127:ASN:ND2	2.50	0.45
1:A:249:HIS:NE2	1:A:273:LEU:HD22	2.32	0.45
1:A:402:PHE:HB3	1:A:428:PHE:CE1	2.52	0.45
1:B:529:PRO:HB3	1:B:531:TRP:CE2	2.53	0.44
1:B:552:ILE:HG23	1:B:556:LEU:HD12	1.99	0.44
2:D:111:SER:O	2:D:136:SER:OG	2.25	0.44
2:D:56:ASN:HB2	2:D:57:PRO:HD2	1.99	0.44
1:A:499:ASP:OD1	1:A:523:ARG:NH2	2.50	0.44
7:A:801:BLD:H328	7:A:801:BLD:H22	1.70	0.44
1:B:539:ALA:HA	1:B:561:SER:O	2.18	0.44
1:B:705:ASN:ND2	7:B:801:BLD:H101	2.32	0.44
2:D:137:LEU:O	2:D:140:LEU:HG	2.18	0.44
1:A:107:SER:HA	1:A:130:SER:O	2.18	0.44
2:D:175:ASN:O	2:D:198:ASN:HA	2.16	0.44
1:A:36:ILE:HB	1:A:61:PRO:HG3	2.00	0.44
1:B:673:GLU:OE1	1:B:673:GLU:N	2.49	0.44
2:C:116:LEU:O	2:C:140:LEU:HD22	2.18	0.44
1:A:396:ASP:OD1	1:A:398:SER:OG	2.23	0.44
1:B:430:GLY:O	1:B:453:SER:N	2.49	0.44
1:B:126:SER:HA	1:B:152:PHE:HB3	2.00	0.43
2:C:137:LEU:O	2:C:140:LEU:HG	2.17	0.43
1:A:505:PRO:HG2	1:A:508:LEU:HG	2.00	0.43
1:B:255:ASN:O	1:B:279:ASN:HA	2.18	0.43
1:B:113:GLY:HA2	8:B:802:NAG:H82	1.99	0.43
2:D:116:LEU:O	2:D:140:LEU:HD22	2.19	0.43
7:A:801:BLD:H121	7:A:801:BLD:H23	1.79	0.43
1:B:726:THR:OG1	2:D:78:ASN:HB2	2.18	0.43
1:B:276:ILE:HD11	1:B:298:LEU:HG	2.01	0.43
2:D:96:GLN:HA	2:D:119:LEU:HA	1.99	0.43
1:A:753:PRO:HB3	1:A:767:LEU:HD23	2.00	0.43
2:C:111:SER:O	2:C:136:SER:OG	2.27	0.43
1:B:361:LEU:HD23	1:B:364:MET:HE3	2.01	0.43
1:B:498:ASN:H	1:B:522:ASN:ND2	2.16	0.43
2:D:175:ASN:OD1	2:D:176:ARG:N	2.42	0.43
1:A:681:PHE:O	1:A:705:ASN:N	2.43	0.43
1:A:749:GLU:CD	2:C:147:ARG:HH12	2.22	0.43
1:B:624:GLU:O	1:B:627:ASN:ND2	2.52	0.42
1:B:130:SER:HB2	3:H:1:NAG:H62	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:VAL:HG13	1:A:394:THR:HB	2.01	0.42
1:A:705:ASN:HA	1:A:727:MET:O	2.20	0.42
1:A:296:LEU:HD11	1:A:298:LEU:HD13	2.00	0.42
1:A:449:PHE:CE1	1:A:625:GLN:HB3	2.54	0.42
1:B:699:GLY:HA2	1:B:725:LEU:HD21	2.02	0.42
1:B:630:SER:HB2	3:J:4:MAN:H62	2.00	0.42
1:A:612:ALA:N	1:A:660:ASP:OD2	2.30	0.42
1:B:251:ASP:OD1	1:B:253:SER:OG	2.23	0.42
2:C:108:PRO:HA	2:C:130:SER:O	2.20	0.42
1:B:498:ASN:H	1:B:522:ASN:HD21	1.67	0.42
1:B:707:LEU:O	1:B:731:ILE:HA	2.20	0.42
2:D:156:SER:HA	2:D:178:SER:O	2.20	0.42
1:A:677:MET:HE3	1:A:677:MET:HB2	1.94	0.42
7:B:801:BLD:H23	7:B:801:BLD:H121	1.69	0.42
1:B:170:LEU:N	1:B:193:TRP:O	2.50	0.42
1:B:769:ARG:HG2	1:B:770:CYS:N	2.35	0.42
1:A:422:TYR:HD1	1:A:446:HIS:HB2	1.85	0.41
1:A:79:LEU:HA	1:A:79:LEU:HD23	1.82	0.41
1:B:521:ASN:HA	1:B:545:ASN:HB3	2.01	0.41
1:A:449:PHE:CE1	1:A:628:ARG:HD2	2.55	0.41
1:B:229:VAL:HG23	1:B:252:ILE:HA	2.02	0.41
1:A:359:ASP:O	1:A:363:LYS:HG2	2.21	0.41
1:A:381:LEU:HB2	1:A:407:LEU:HD21	2.02	0.41
1:A:544:SER:O	1:A:546:ASN:ND2	2.53	0.41
1:B:485:MET:O	1:B:488:LYS:NZ	2.43	0.41
1:B:234:PHE:HB3	1:B:257:LEU:HD21	2.03	0.41
1:B:719:PRO:HG2	1:B:722:MET:HG3	2.02	0.41
1:B:137:PRO:HG2	1:B:140:THR:HG23	2.03	0.41
1:B:228:ASP:OD1	1:B:230:SER:OG	2.23	0.41
1:B:333:PRO:HB2	1:B:335:PHE:CE2	2.55	0.41
1:B:492:THR:HG21	1:B:595:LYS:HE3	2.03	0.41
1:B:671:PRO:HB2	1:B:673:GLU:OE1	2.21	0.41
1:A:688:ASN:HB2	1:A:712:ASN:OD1	2.21	0.41
1:B:496:ASP:HA	1:B:520:SER:H	1.86	0.41
1:A:282:VAL:HG12	1:A:302:LYS:HB2	2.02	0.41
1:A:31:SER:OG	1:A:32:LEU:N	2.54	0.41
1:A:375:ASN:O	1:A:400:ASN:HA	2.21	0.41
1:A:681:PHE:HA	1:A:703:GLY:O	2.21	0.41
1:A:737:ASN:O	1:A:737:ASN:ND2	2.54	0.41
1:B:115:VAL:HB	1:B:140:THR:HB	2.03	0.41
1:B:470:LYS:HG2	1:B:494:ILE:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:B:801:BLD:H115	7:B:801:BLD:H207	1.93	0.41
2:D:158:PRO:HB2	2:D:161:LEU:HG	2.03	0.41
7:A:801:BLD:H115	7:A:801:BLD:H207	1.89	0.40
2:C:201:LEU:HB3	2:C:206:THR:HG21	2.02	0.40
1:A:497:PHE:N	1:A:520:SER:O	2.44	0.40
1:B:368:LYS:HA	1:B:368:LYS:HD3	1.81	0.40
1:B:542:LYS:HG3	1:B:566:ASP:HB3	2.03	0.40
1:B:701:LEU:O	1:B:725:LEU:HD22	2.21	0.40
1:B:82:LYS:O	1:B:84:LEU:N	2.51	0.40
2:D:33:LEU:O	2:D:36:LEU:HB3	2.21	0.40
1:A:112:ASN:HD22	8:A:802:NAG:C7	2.33	0.40
1:A:106:LEU:HA	1:A:106:LEU:HD23	1.89	0.40
2:C:37:ARG:NH2	2:C:48:GLN:HA	2.36	0.40
1:B:509:SER:HB3	1:B:531:TRP:CD2	2.56	0.40
1:B:608:GLU:OE1	1:B:608:GLU:N	2.55	0.40

There are no symmetry-related clashes.

## 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	738/774 (95%)	710 (96%)	27 (4%)	1 (0%)	51	81
1	B	706/774 (91%)	680 (96%)	25 (4%)	1 (0%)	51	81
2	C	183/203 (90%)	176 (96%)	7 (4%)	0	100	100
2	D	183/203 (90%)	177 (97%)	6 (3%)	0	100	100
All	All	1810/1954 (93%)	1743 (96%)	65 (4%)	2 (0%)	51	81

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	637	ILE
1	B	637	ILE

### 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	611/680 (90%)	600 (98%)	11 (2%)	59	78
1	B	580/680 (85%)	571 (98%)	9 (2%)	62	79
2	C	165/182 (91%)	164 (99%)	1 (1%)	86	91
2	D	161/182 (88%)	160 (99%)	1 (1%)	86	91
All	All	1517/1724 (88%)	1495 (98%)	22 (2%)	65	81

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	GLU
1	A	49	LYS
1	A	54	ASP
1	A	162	PHE
1	A	199	CYS
1	A	410	LEU
1	A	524	LEU
1	A	527	GLU
1	A	532	ILE
1	A	625	GLN
1	A	669	TYR
1	B	107	SER
1	B	162	PHE
1	B	199	CYS
1	B	410	LEU
1	B	446	HIS
1	B	449	PHE
1	B	501	THR
1	B	524	LEU
1	B	750	THR

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Mol	Chain	Res	Type
2	C	112	ASN
2	D	78	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	ASN
1	A	551	ASN
1	B	705	ASN
2	C	78	ASN
2	C	112	ASN
2	D	94	ASN
2	D	118	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 ⓘ

24 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	E	1	1,3	14,14,15	0.26	0	17,19,21	0.46	0
3	NAG	E	2	3	14,14,15	0.47	0	17,19,21	1.26	1 (5%)
3	BMA	E	3	3	11,11,12	0.70	0	15,15,17	1.06	0
3	MAN	E	4	3	11,11,12	0.75	0	15,15,17	0.96	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NAG	F	1	1,4	14,14,15	0.35	0	17,19,21	0.51	0
4	NAG	F	2	4	14,14,15	0.40	0	17,19,21	0.63	1 (5%)
5	NAG	G	1	1,5	14,14,15	0.30	0	17,19,21	0.40	0
5	NAG	G	2	5	14,14,15	0.22	0	17,19,21	0.36	0
5	BMA	G	3	5	11,11,12	0.98	0	15,15,17	0.96	0
5	MAN	G	4	5	11,11,12	0.74	0	15,15,17	0.98	2 (13%)
5	MAN	G	5	5	11,11,12	0.77	0	15,15,17	1.68	4 (26%)
3	NAG	H	1	1,3	14,14,15	0.32	0	17,19,21	0.42	0
3	NAG	H	2	3	14,14,15	0.48	0	17,19,21	1.25	1 (5%)
3	BMA	H	3	3	11,11,12	0.63	0	15,15,17	0.97	0
3	MAN	H	4	3	11,11,12	0.79	1 (9%)	15,15,17	0.98	2 (13%)
6	NAG	I	1	1,6	14,14,15	0.64	1 (7%)	17,19,21	0.57	0
6	NAG	I	2	6	14,14,15	0.66	1 (7%)	17,19,21	0.59	0
6	BMA	I	3	6	11,11,12	0.65	0	15,15,17	1.05	2 (13%)
3	NAG	J	1	1,3	14,14,15	0.23	0	17,19,21	0.46	0
3	NAG	J	2	3	14,14,15	0.50	0	17,19,21	0.46	0
3	BMA	J	3	3	11,11,12	1.37	3 (27%)	15,15,17	1.49	2 (13%)
3	MAN	J	4	3	11,11,12	0.83	1 (9%)	15,15,17	1.04	2 (13%)
4	NAG	K	1	1,4	14,14,15	0.42	0	17,19,21	0.39	0
4	NAG	K	2	4	14,14,15	0.24	0	17,19,21	0.39	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
3	NAG	E	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	4/6/23/26	0/1/1/1
3	BMA	E	3	3	-	2/2/19/22	0/1/1/1
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
4	NAG	F	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	F	2	4	-	2/6/23/26	0/1/1/1
5	NAG	G	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	G	2	5	-	0/6/23/26	0/1/1/1
5	BMA	G	3	5	-	0/2/19/22	0/1/1/1
5	MAN	G	4	5	-	2/2/19/22	0/1/1/1
5	MAN	G	5	5	-	0/2/19/22	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1

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

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	3	BMA	O5-C1	-3.05	1.38	1.43
6	I	2	NAG	O5-C1	-2.35	1.40	1.43
6	I	1	NAG	O5-C1	-2.30	1.40	1.43
3	J	4	MAN	O5-C1	-2.27	1.40	1.43
3	J	3	BMA	C4-C3	2.24	1.58	1.52
3	J	3	BMA	C2-C3	2.13	1.55	1.52
3	H	4	MAN	O5-C1	-2.05	1.40	1.43

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	2	NAG	C2-N2-C7	4.38	129.13	122.90
3	H	2	NAG	C2-N2-C7	4.37	129.12	122.90
5	G	5	MAN	C1-O5-C5	3.94	117.54	112.19
3	J	3	BMA	C2-C3-C4	3.93	117.69	110.89
5	G	5	MAN	O5-C1-C2	2.93	115.30	110.77
3	J	3	BMA	C3-C4-C5	2.83	115.28	110.24
5	G	5	MAN	C1-C2-C3	2.81	113.12	109.67
6	I	3	BMA	O2-C2-C3	-2.44	105.25	110.14
4	F	2	NAG	C1-O5-C5	2.37	115.41	112.19
6	I	3	BMA	C1-C2-C3	-2.37	106.75	109.67
3	H	4	MAN	O2-C2-C3	-2.32	105.48	110.14
5	G	5	MAN	O2-C2-C3	-2.32	105.49	110.14
3	J	4	MAN	O2-C2-C3	-2.28	105.57	110.14
3	J	4	MAN	C1-O5-C5	2.20	115.17	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	4	MAN	O2-C2-C3	-2.18	105.78	110.14
3	E	4	MAN	O2-C2-C3	-2.17	105.79	110.14
3	H	4	MAN	C1-O5-C5	2.07	115.00	112.19
5	G	4	MAN	C1-O5-C5	2.07	115.00	112.19
3	E	4	MAN	C1-O5-C5	2.01	114.91	112.19

There are no chirality outliers.

All (26) torsion outliers are listed below:

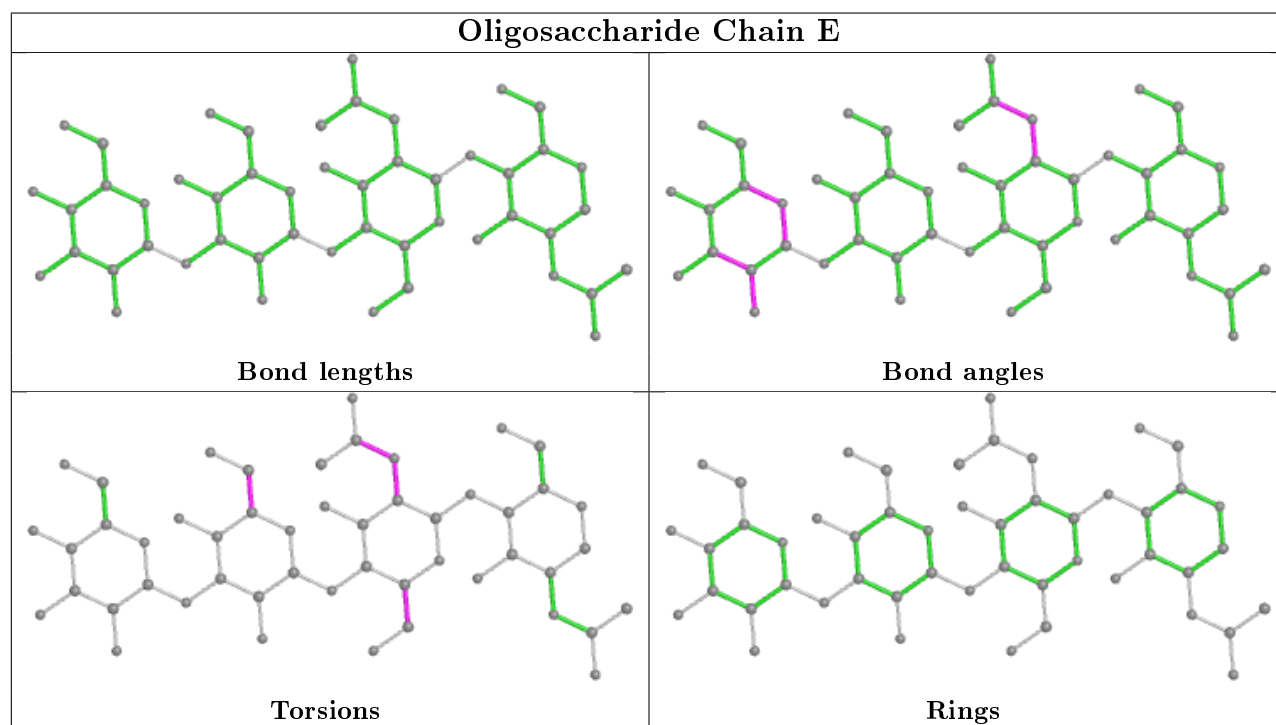
Mol	Chain	Res	Type	Atoms
4	F	1	NAG	C4-C5-C6-O6
4	F	2	NAG	O5-C5-C6-O6
4	F	1	NAG	O5-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
5	G	4	MAN	O5-C5-C6-O6
5	G	4	MAN	C4-C5-C6-O6
6	I	2	NAG	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
4	F	2	NAG	C4-C5-C6-O6
6	I	2	NAG	C4-C5-C6-O6
3	H	2	NAG	C8-C7-N2-C2
3	H	2	NAG	O7-C7-N2-C2
3	E	2	NAG	C8-C7-N2-C2
3	E	2	NAG	O7-C7-N2-C2
6	I	3	BMA	O5-C5-C6-O6
5	G	1	NAG	O5-C5-C6-O6
5	G	1	NAG	C4-C5-C6-O6
3	E	3	BMA	C4-C5-C6-O6
4	K	1	NAG	O5-C5-C6-O6
3	E	3	BMA	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
4	K	2	NAG	C4-C5-C6-O6
6	I	3	BMA	C4-C5-C6-O6
6	I	2	NAG	C1-C2-N2-C7
3	H	2	NAG	C3-C2-N2-C7
3	E	2	NAG	C3-C2-N2-C7

There are no ring outliers.

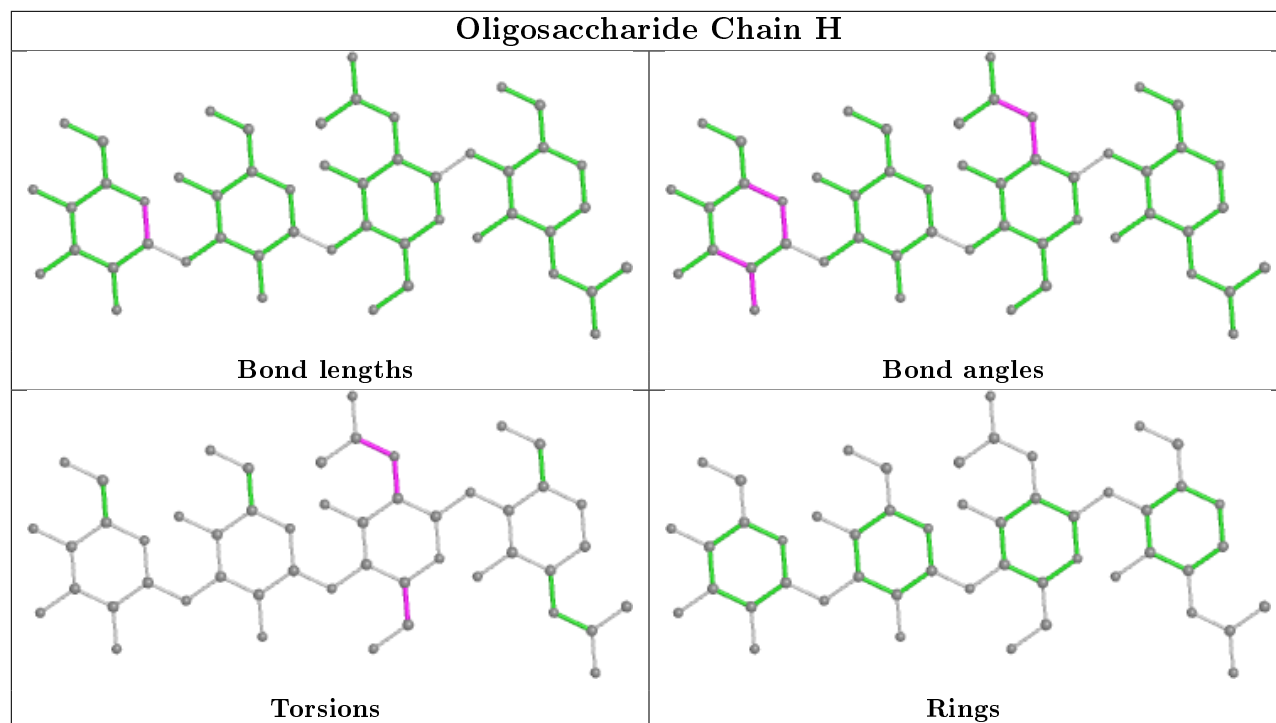
8 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	2	NAG	1	0
3	E	2	NAG	1	0
5	G	3	BMA	1	0
3	H	1	NAG	1	0
3	J	4	MAN	1	0
5	G	5	MAN	1	0
4	F	1	NAG	2	0
3	J	2	NAG	1	0

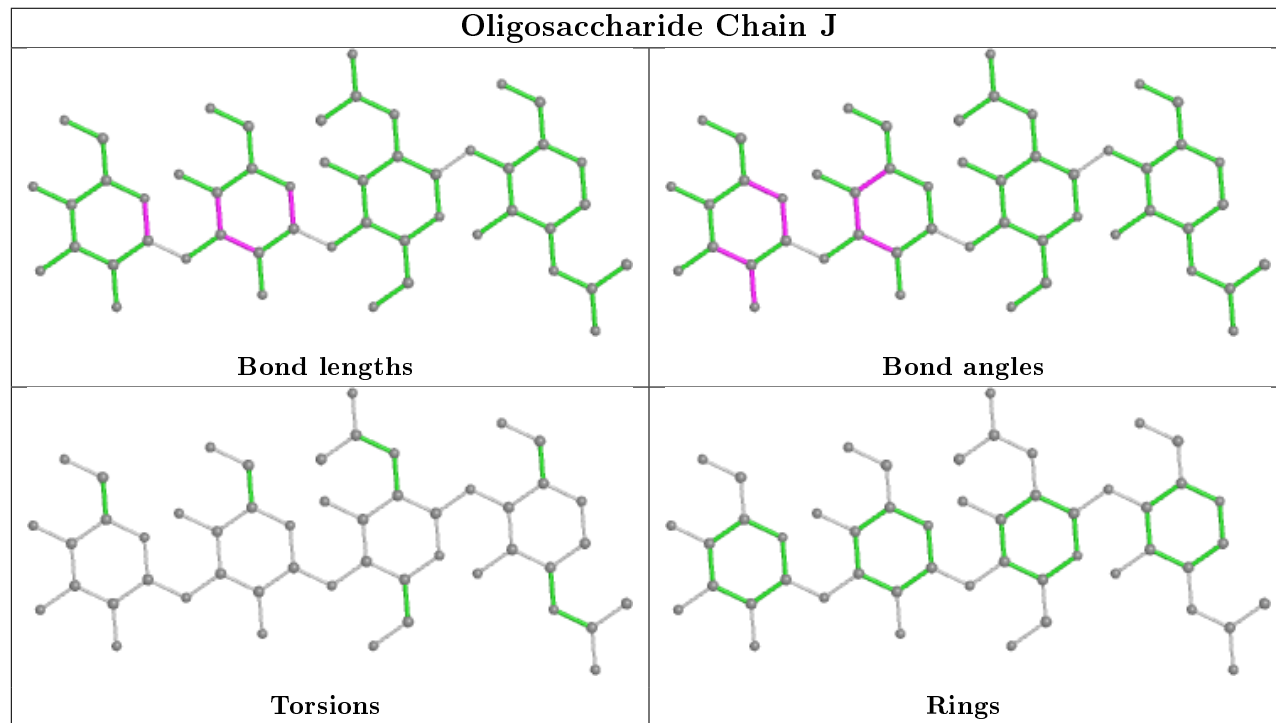
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

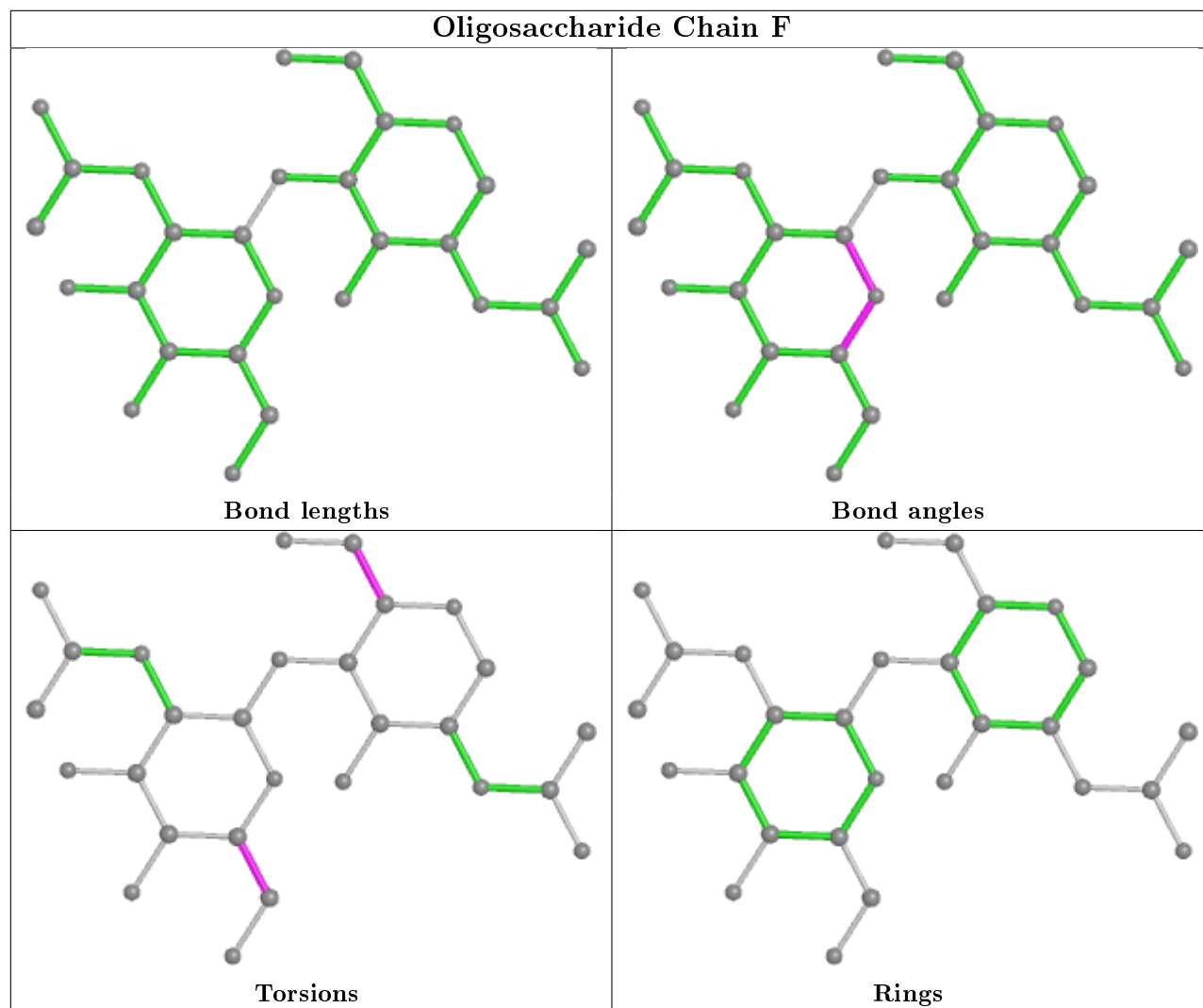


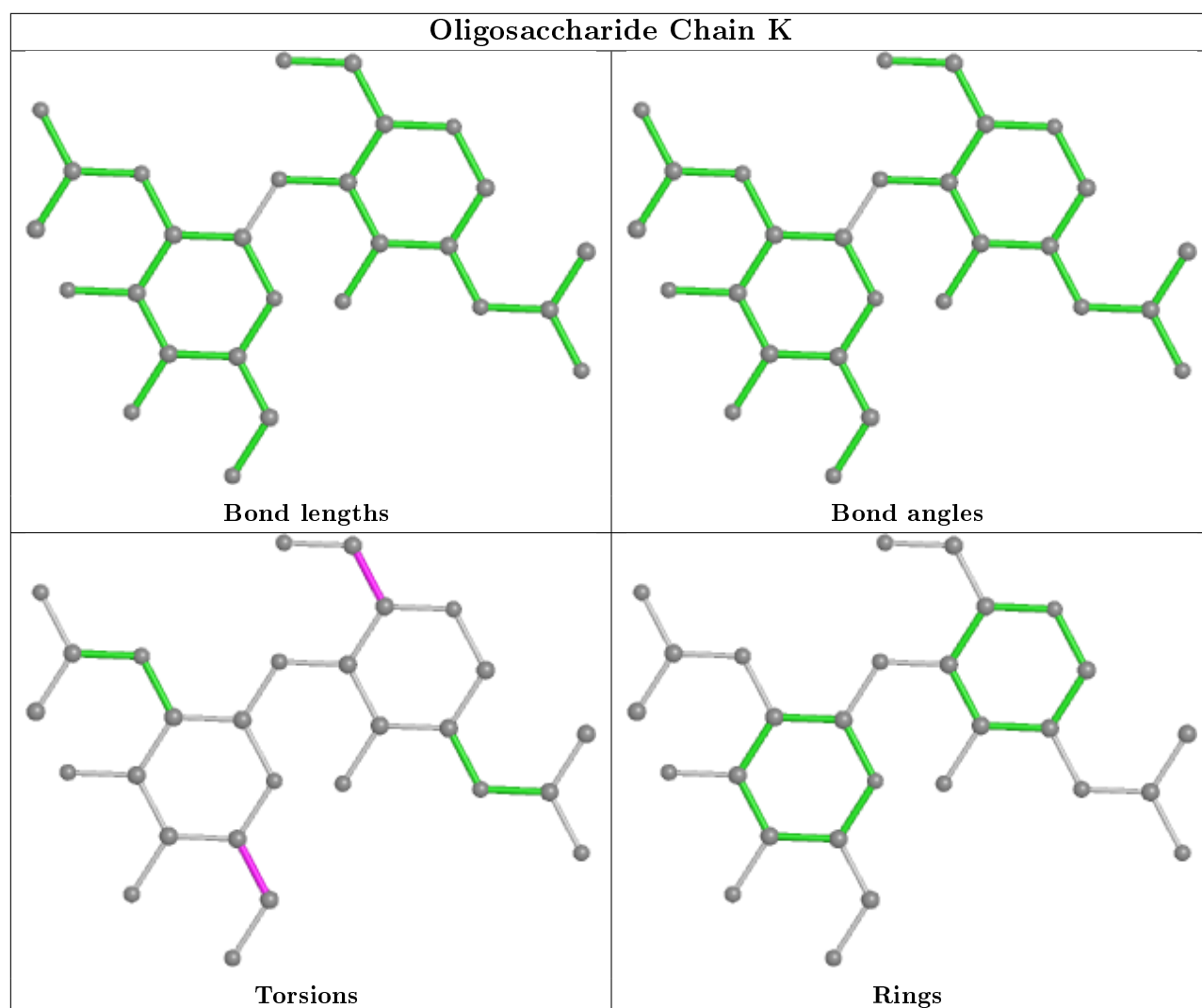
## Oligosaccharide Chain H

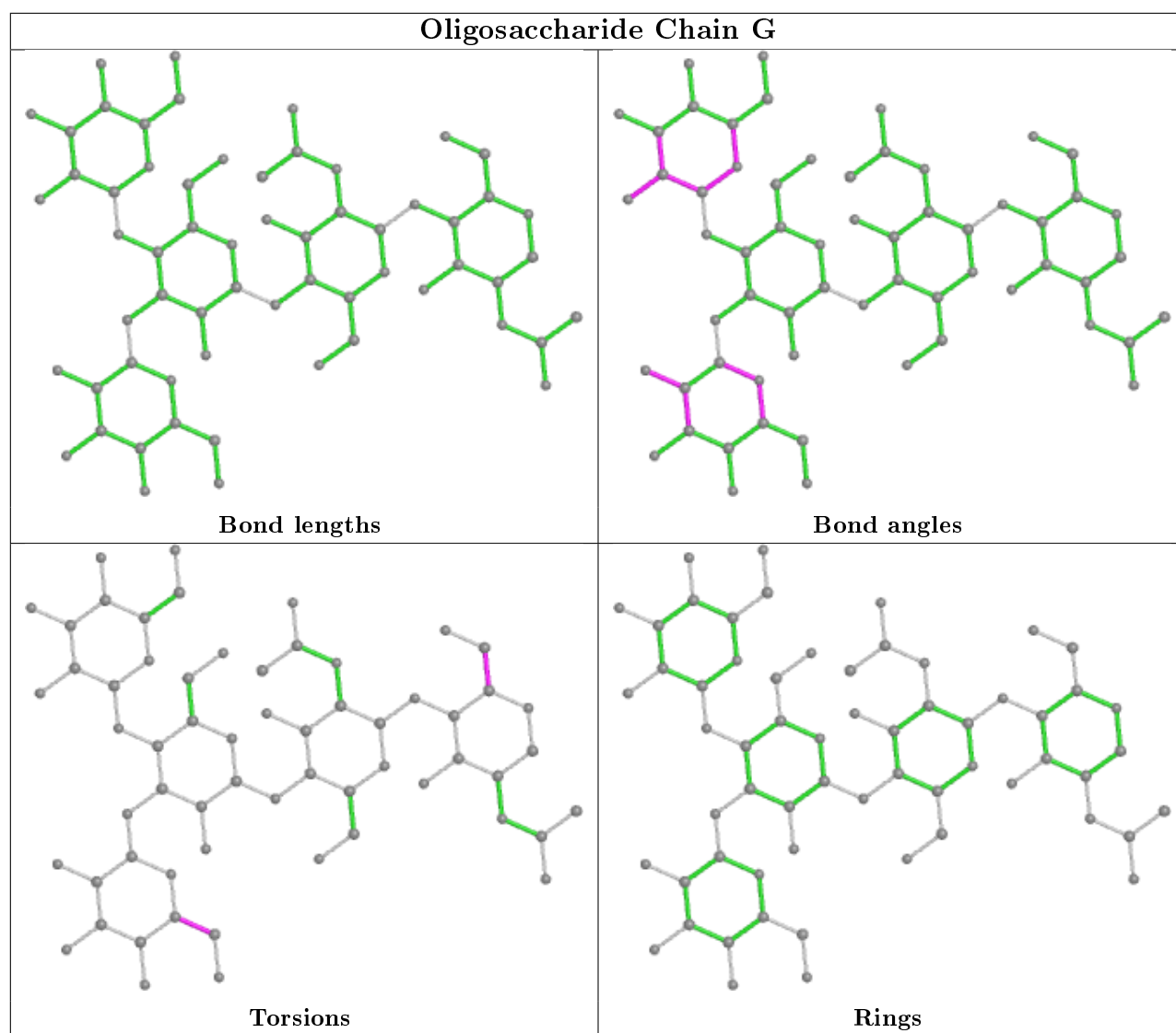


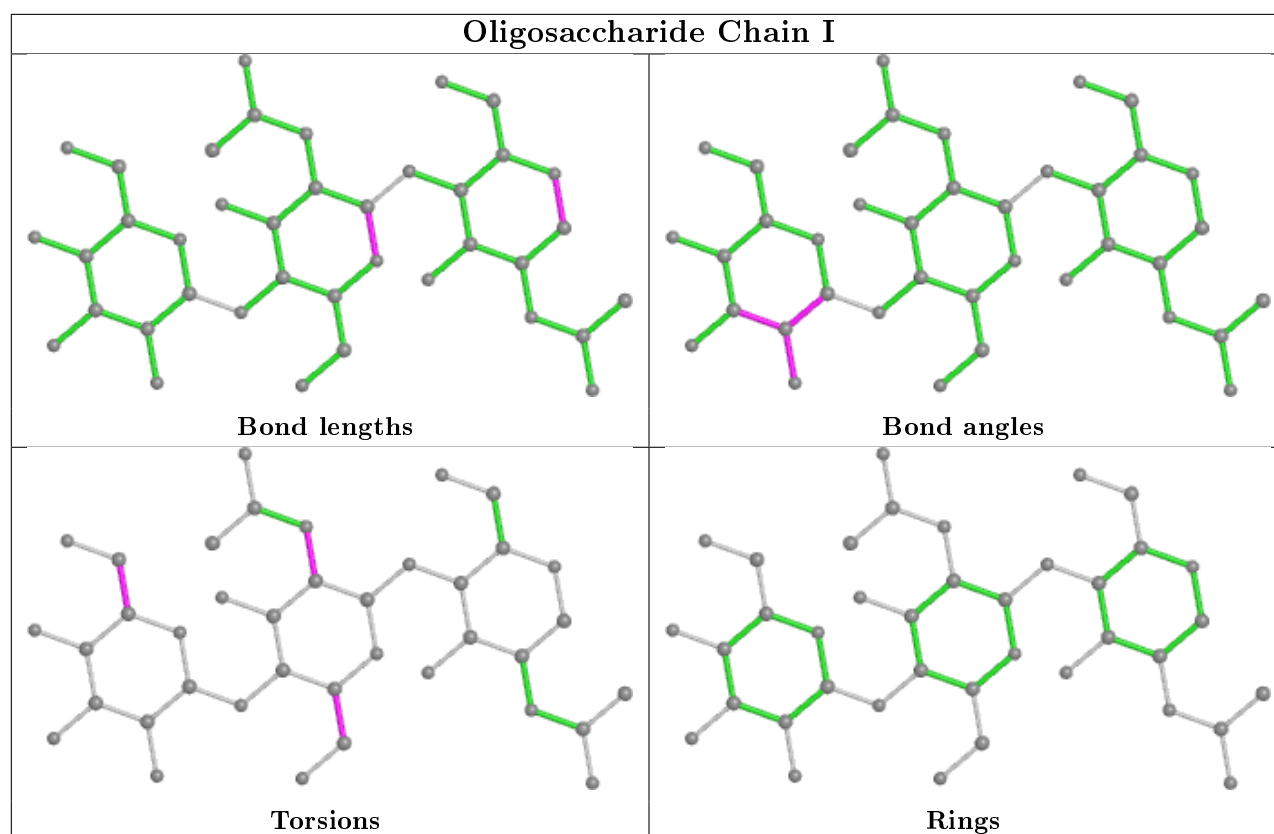
## Oligosaccharide Chain J











## 5.6 Ligand geometry [i](#)

9 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	$\# Z  > 2$
7	BLD	A	801	-	36,37,37	3.18	12 (33%)	46,59,59	1.98	14 (30%)
7	BLD	B	801	-	36,37,37	3.17	12 (33%)	46,59,59	1.86	12 (26%)
8	NAG	D	1002	2	14,14,15	1.67	1 (7%)	17,19,21	1.61	3 (17%)
8	NAG	B	802	1	14,14,15	0.24	0	17,19,21	0.43	0
8	NAG	A	802	1	14,14,15	0.36	0	17,19,21	0.59	0
8	NAG	C	1001	2	14,14,15	0.20	0	17,19,21	0.44	0
8	NAG	C	1000	2	14,14,15	1.12	1 (7%)	17,19,21	0.67	0
8	NAG	A	814	1	14,14,15	0.58	0	17,19,21	1.19	1 (5%)
8	NAG	D	1001	2	14,14,15	0.36	0	17,19,21	0.44	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
7	BLD	A	801	-	-	4/20/85/85	0/4/4/4
7	BLD	B	801	-	-	4/20/85/85	0/4/4/4
8	NAG	D	1002	2	-	3/6/23/26	0/1/1/1
8	NAG	B	802	1	-	2/6/23/26	0/1/1/1
8	NAG	A	802	1	-	2/6/23/26	0/1/1/1
8	NAG	C	1001	2	-	2/6/23/26	0/1/1/1
8	NAG	C	1000	2	-	2/6/23/26	0/1/1/1
8	NAG	A	814	1	-	0/6/23/26	0/1/1/1
8	NAG	D	1001	2	-	0/6/23/26	0/1/1/1

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	801	BLD	O07-C07	-11.47	1.30	1.45
7	B	801	BLD	O07-C07	-11.30	1.30	1.45
7	B	801	BLD	O06-C06	9.97	1.46	1.21
7	A	801	BLD	O06-C06	9.88	1.46	1.21
8	D	1002	NAG	C1-C2	5.80	1.61	1.52
7	B	801	BLD	C15-C14	-5.22	1.43	1.54
7	A	801	BLD	C15-C14	-5.19	1.43	1.54
7	A	801	BLD	C11-C09	4.94	1.62	1.53
7	B	801	BLD	C11-C09	4.79	1.61	1.53
8	C	1000	NAG	O5-C1	3.59	1.49	1.43
7	B	801	BLD	C19-C10	3.45	1.60	1.54
7	A	801	BLD	C19-C10	3.44	1.60	1.54
7	A	801	BLD	C20-C17	-3.33	1.49	1.54
7	B	801	BLD	O22-C22	-3.22	1.35	1.43
7	A	801	BLD	O22-C22	-3.17	1.35	1.43
7	B	801	BLD	C20-C17	-3.15	1.49	1.54
7	A	801	BLD	C28-C24	-2.93	1.47	1.53
7	B	801	BLD	C28-C24	-2.90	1.47	1.53
7	B	801	BLD	C16-C15	2.64	1.61	1.54
7	A	801	BLD	C16-C15	2.57	1.61	1.54
7	B	801	BLD	C08-C14	2.39	1.58	1.53
7	B	801	BLD	C13-C14	-2.35	1.50	1.55
7	A	801	BLD	C13-C14	-2.26	1.50	1.55
7	A	801	BLD	C08-C14	2.22	1.57	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	801	BLD	O07-C06	-2.07	1.32	1.34
7	B	801	BLD	O07-C06	-2.06	1.32	1.34

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	801	BLD	C15-C14-C08	-5.29	110.37	119.08
7	A	801	BLD	O07-C06-O06	-5.17	108.67	116.72
7	B	801	BLD	C01-C10-C05	5.12	115.22	107.06
7	B	801	BLD	C15-C14-C08	-5.06	110.74	119.08
7	B	801	BLD	O07-C06-O06	-4.91	109.07	116.72
7	A	801	BLD	C01-C10-C05	4.87	114.83	107.06
8	D	1002	NAG	C2-N2-C7	3.73	128.21	122.90
8	A	814	NAG	C1-O5-C5	3.52	116.97	112.19
8	D	1002	NAG	O5-C5-C4	-3.45	102.43	110.83
7	A	801	BLD	C17-C20-C22	-3.41	105.47	111.51
7	B	801	BLD	C11-C09-C08	-3.18	107.17	111.75
7	A	801	BLD	C11-C09-C08	-3.11	107.27	111.75
7	A	801	BLD	C12-C13-C17	2.89	120.89	116.57
7	A	801	BLD	C07-O07-C06	-2.73	117.16	121.01
7	B	801	BLD	C07-O07-C06	-2.72	117.16	121.01
7	B	801	BLD	C12-C13-C17	2.65	120.54	116.57
7	A	801	BLD	C16-C15-C14	-2.58	100.02	105.13
7	A	801	BLD	C16-C17-C20	-2.54	108.97	112.61
7	A	801	BLD	C12-C13-C14	2.45	111.08	107.27
7	B	801	BLD	C17-C20-C22	-2.39	107.29	111.51
7	A	801	BLD	C14-C08-C09	-2.38	105.90	109.09
7	B	801	BLD	C16-C15-C14	-2.33	100.51	105.13
7	A	801	BLD	C28-C24-C25	-2.25	108.76	112.47
7	B	801	BLD	C16-C17-C20	-2.23	109.42	112.61
7	A	801	BLD	C24-C23-C22	-2.18	109.93	114.85
7	B	801	BLD	C28-C24-C25	-2.14	108.94	112.47
7	B	801	BLD	C21-C20-C22	-2.13	107.09	111.39
8	D	1002	NAG	C4-C3-C2	2.12	114.13	111.02
7	A	801	BLD	C21-C20-C22	-2.11	107.13	111.39
7	B	801	BLD	C14-C08-C09	-2.08	106.31	109.09

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	801	BLD	C20-C22-C23-O23

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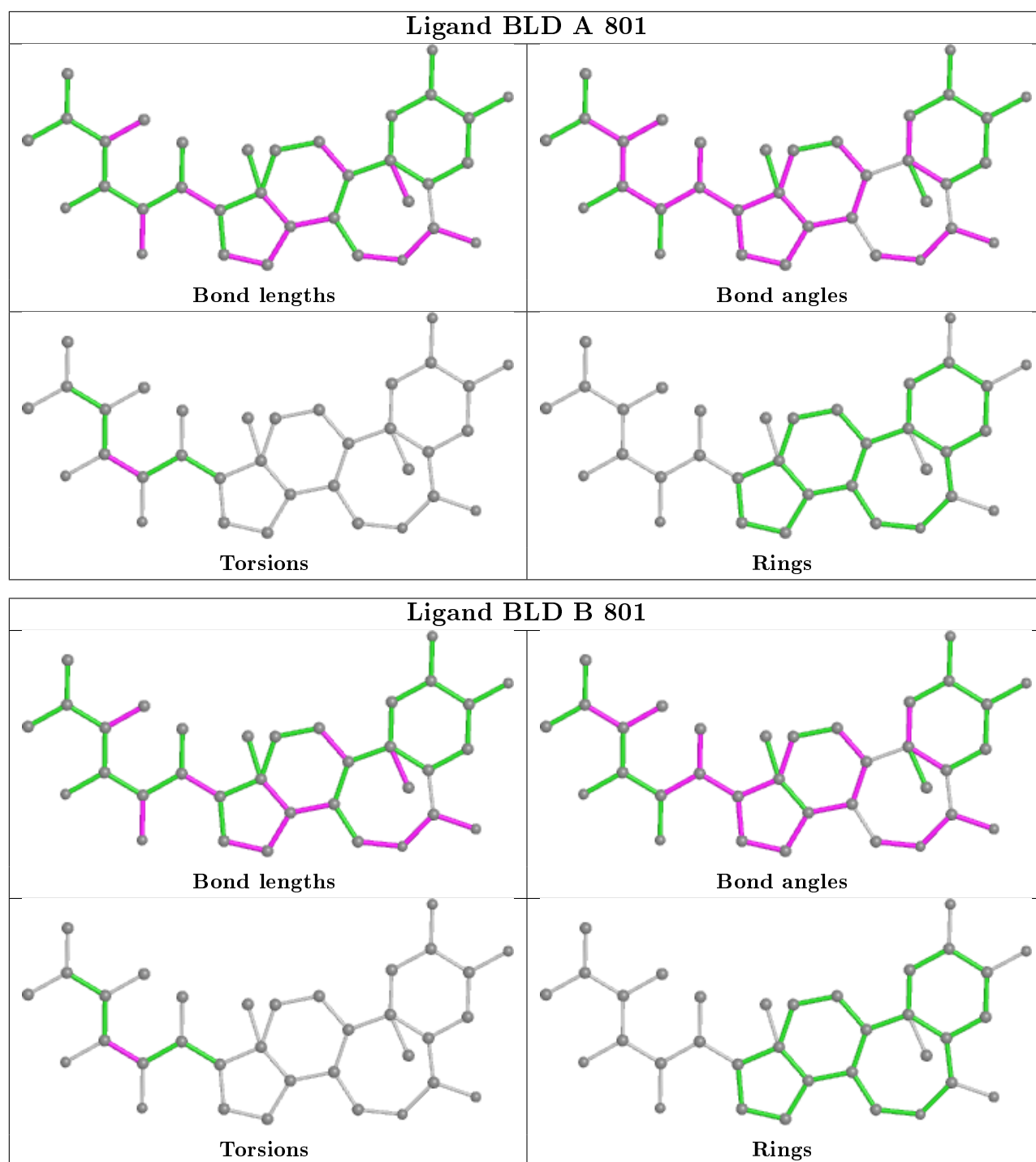
Mol	Chain	Res	Type	Atoms
7	A	801	BLD	C20-C22-C23-C24
7	A	801	BLD	O22-C22-C23-C24
7	B	801	BLD	C20-C22-C23-O23
7	B	801	BLD	O22-C22-C23-C24
8	B	802	NAG	C4-C5-C6-O6
8	B	802	NAG	O5-C5-C6-O6
8	A	802	NAG	C4-C5-C6-O6
8	A	802	NAG	O5-C5-C6-O6
8	C	1001	NAG	O5-C5-C6-O6
8	C	1000	NAG	O5-C5-C6-O6
8	C	1001	NAG	C4-C5-C6-O6
8	D	1002	NAG	C1-C2-N2-C7
8	C	1000	NAG	C4-C5-C6-O6
8	D	1002	NAG	O5-C5-C6-O6
7	B	801	BLD	O22-C22-C23-O23
7	B	801	BLD	C20-C22-C23-C24
8	D	1002	NAG	C3-C2-N2-C7
7	A	801	BLD	O22-C22-C23-O23

There are no ring outliers.

5 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	801	BLD	5	0
7	B	801	BLD	5	0
8	B	802	NAG	1	0
8	A	802	NAG	1	0
8	A	814	NAG	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

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	740/774 (95%)	0.03	23 (3%) 49 48	47, 114, 169, 224	0
1	B	710/774 (91%)	0.05	23 (3%) 47 46	57, 115, 166, 222	0
2	C	185/203 (91%)	0.41	19 (10%) 6 6	88, 145, 192, 210	0
2	D	185/203 (91%)	0.73	32 (17%) 1 1	105, 146, 225, 259	0
All	All	1820/1954 (93%)	0.15	97 (5%) 26 24	47, 121, 183, 259	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	726	THR	8.1
2	D	192	PRO	6.6
2	D	138	GLY	6.5
2	D	137	LEU	6.4
2	D	98	LEU	6.3
2	D	196	ALA	6.0
1	A	723	SER	5.8
2	D	146	LEU	5.8
2	D	195	PHE	5.3
1	B	31	SER	4.7
1	A	725	LEU	4.7
1	B	500	LEU	4.7
2	C	87	PRO	4.4
2	D	74	VAL	4.3
2	D	135	GLU	4.2
2	D	79	ALA	4.2
2	D	76	LEU	4.2
2	D	189	LEU	4.2
1	A	724	ALA	4.1
2	C	76	LEU	4.1
2	C	98	LEU	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	702	ARG	4.0
1	A	675	GLY	4.0
1	A	733	LEU	3.9
2	C	74	VAL	3.9
2	D	172	LEU	3.8
2	C	180	SER	3.8
2	D	100	LEU	3.7
2	D	148	LEU	3.7
2	D	94	ASN	3.6
2	D	198	ASN	3.6
2	D	194	SER	3.6
1	A	539	ALA	3.5
1	A	463	LEU	3.4
1	A	53	PRO	3.4
2	D	122	LEU	3.3
2	D	199	LEU	3.3
2	D	134	PRO	3.2
2	C	153	LEU	3.2
1	A	701	LEU	3.2
2	D	136	SER	3.2
2	D	171	ASP	3.1
2	C	146	LEU	3.1
1	B	103	SER	3.0
1	A	621	ILE	2.8
2	C	108	PRO	2.8
2	C	160	SER	2.8
2	C	152	SER	2.8
1	B	733	LEU	2.8
2	C	148	LEU	2.8
2	D	204	PRO	2.7
1	B	765	TYR	2.7
2	C	91	VAL	2.7
1	A	638	THR	2.7
1	B	35	GLU	2.6
1	A	679	TYR	2.6
1	A	540	ILE	2.6
2	C	154	THR	2.6
2	D	117	THR	2.6
2	D	191	THR	2.6
1	B	311	LEU	2.6
1	B	447	LEU	2.6
2	D	160	SER	2.6

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Mol	Chain	Res	Type	RSRZ
2	D	181	VAL	2.6
2	C	100	LEU	2.5
1	B	409	ASN	2.5
2	C	94	ASN	2.5
1	B	642	TYR	2.5
2	D	180	SER	2.5
1	A	392	LEU	2.4
2	C	88	GLU	2.4
1	B	708	ASP	2.4
2	C	93	LYS	2.4
1	B	722	MET	2.3
1	A	639	SER	2.3
1	A	727	MET	2.3
1	B	396	ASP	2.3
1	A	495	LEU	2.3
1	B	297	SER	2.3
2	C	52	PRO	2.3
1	B	488	LYS	2.3
1	A	261	PHE	2.2
1	B	37	HIS	2.2
2	D	86	VAL	2.2
1	B	499	ASP	2.2
2	D	170	LEU	2.1
1	A	700	ASP	2.1
1	B	745	MET	2.1
1	B	141	LEU	2.1
2	C	166	THR	2.1
1	B	548	PHE	2.1
1	B	563	ILE	2.1
1	A	447	LEU	2.1
1	B	407	LEU	2.1
1	A	385	LEU	2.0
1	B	421	LEU	2.0
2	D	193	ILE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> 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.

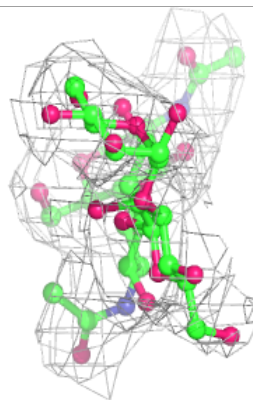
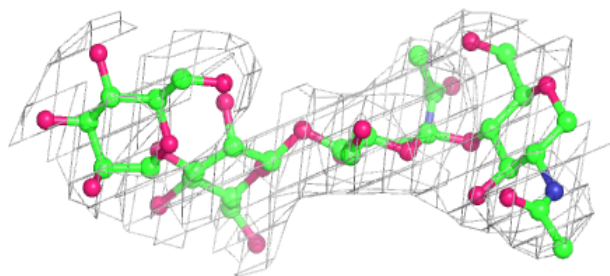
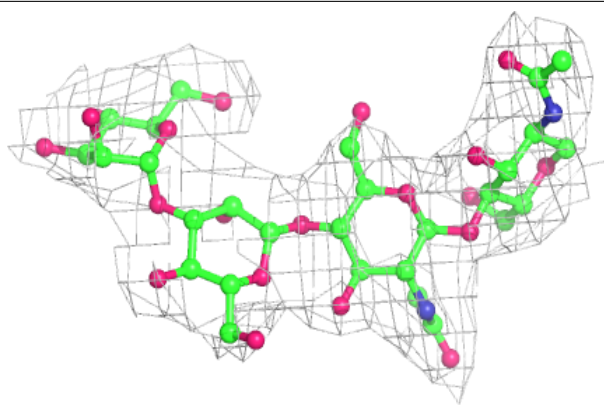
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	MAN	H	4	11/12	0.77	0.27	150,157,165,169	0
3	BMA	H	3	11/12	0.79	0.18	125,136,149,156	0
6	BMA	I	3	11/12	0.81	0.19	136,138,139,143	0
3	NAG	J	1	14/15	0.83	0.19	86,99,107,109	0
4	NAG	K	1	14/15	0.85	0.20	93,107,121,122	0
3	BMA	J	3	11/12	0.85	0.16	98,131,145,150	0
5	MAN	G	4	11/12	0.87	0.17	128,135,139,143	0
5	MAN	G	5	11/12	0.87	0.21	76,87,117,130	0
4	NAG	F	2	14/15	0.88	0.17	107,112,122,127	0
3	BMA	E	3	11/12	0.89	0.13	77,110,116,116	0
5	NAG	G	1	14/15	0.90	0.24	111,121,148,149	0
3	MAN	E	4	11/12	0.90	0.13	83,104,121,128	0
5	BMA	G	3	11/12	0.91	0.13	101,118,146,149	0
3	NAG	H	1	14/15	0.92	0.20	110,124,137,137	0
4	NAG	F	1	14/15	0.92	0.26	92,102,111,111	0
6	NAG	I	2	14/15	0.93	0.13	93,114,124,132	0
3	NAG	H	2	14/15	0.93	0.19	109,116,136,136	0
5	NAG	G	2	14/15	0.93	0.15	88,102,106,111	0
3	NAG	J	2	14/15	0.93	0.16	98,120,128,130	0
4	NAG	K	2	14/15	0.93	0.26	72,85,107,107	0
6	NAG	I	1	14/15	0.93	0.17	72,85,96,100	0
3	NAG	E	2	14/15	0.94	0.12	108,114,118,120	0
3	MAN	J	4	11/12	0.95	0.13	84,92,96,99	0
3	NAG	E	1	14/15	0.95	0.17	48,61,73,95	0

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

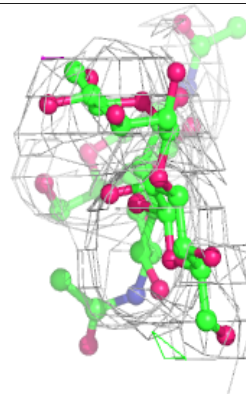
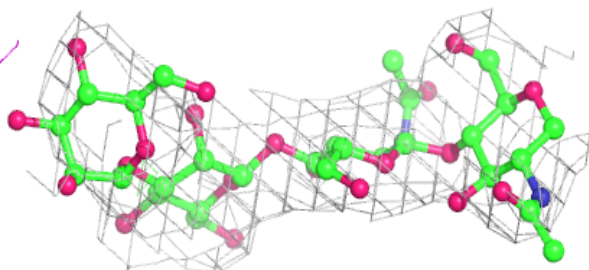
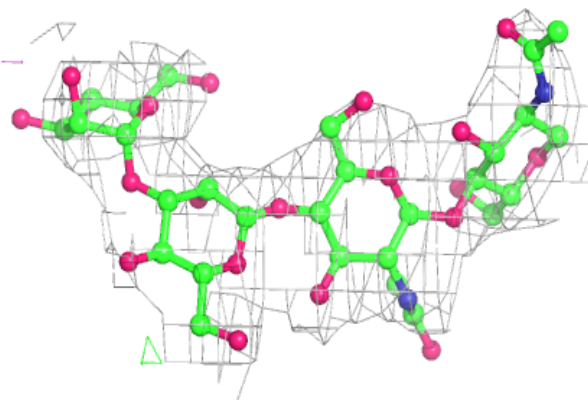


**Electron density around Chain E:**

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

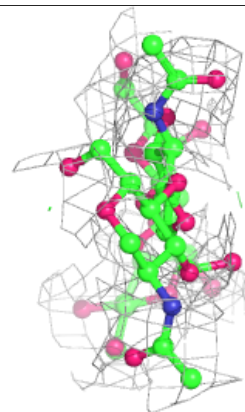
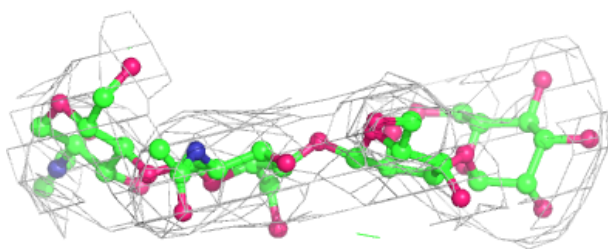
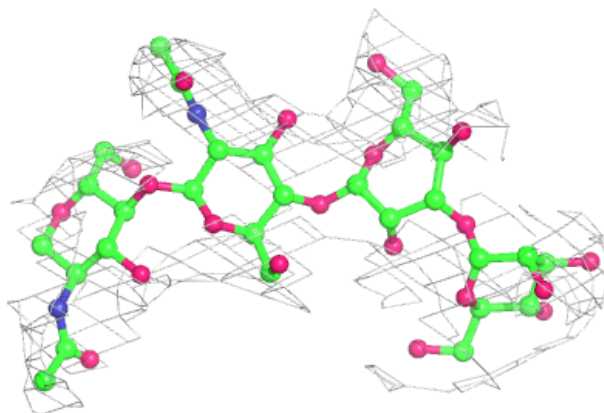
**Electron density around Chain H:**

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



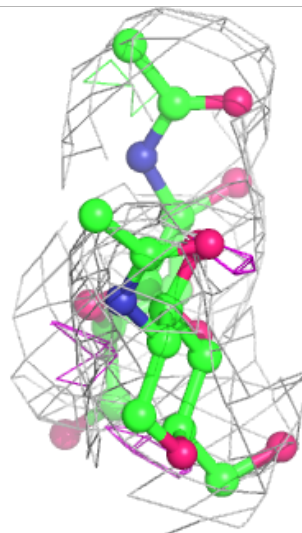
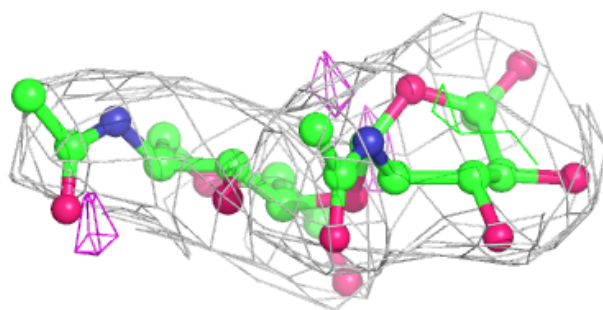
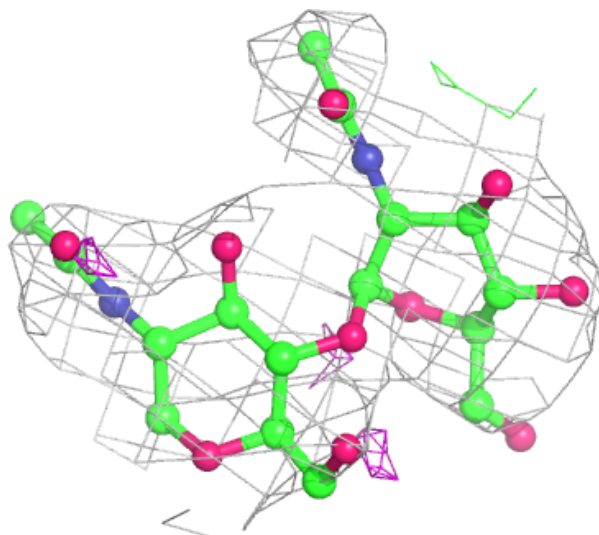
**Electron density around Chain J:**

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



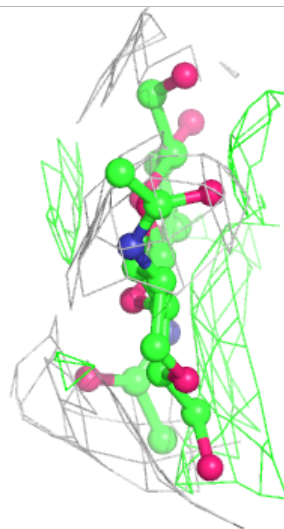
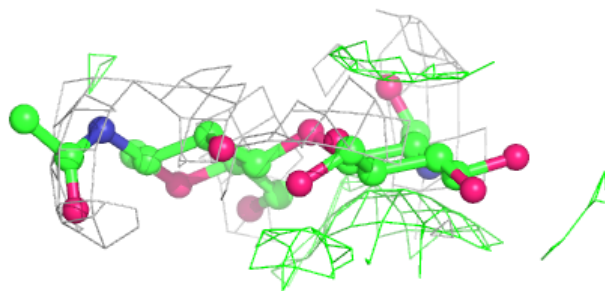
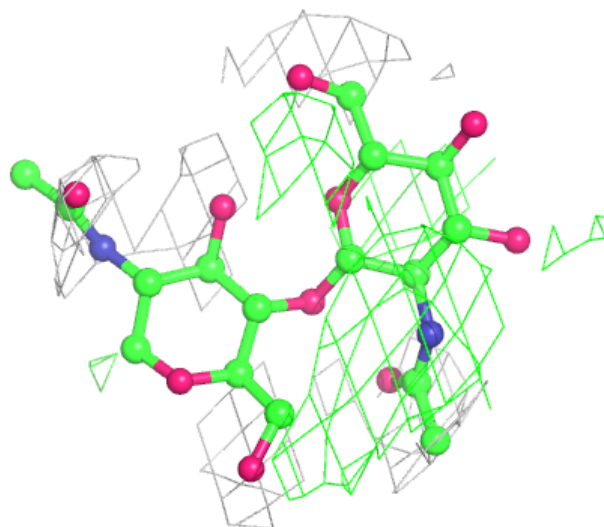
**Electron density around Chain F:**

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



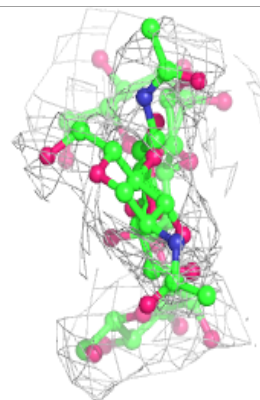
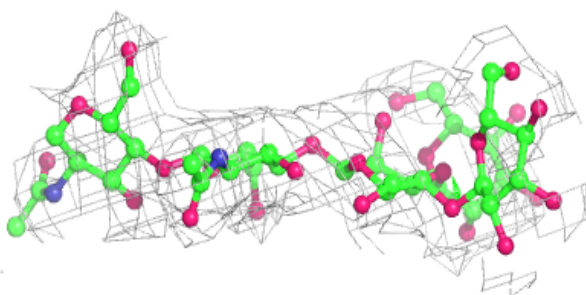
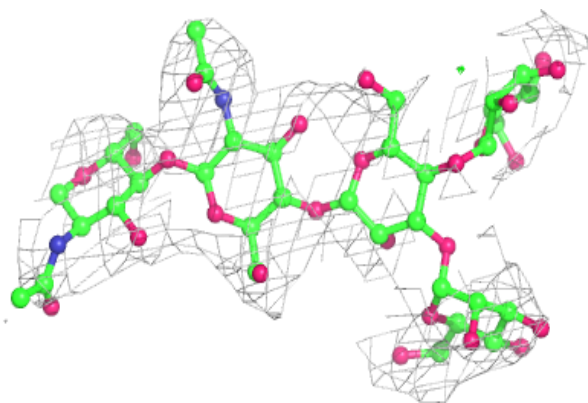
**Electron density around Chain K:**

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

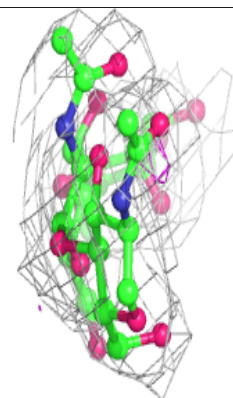
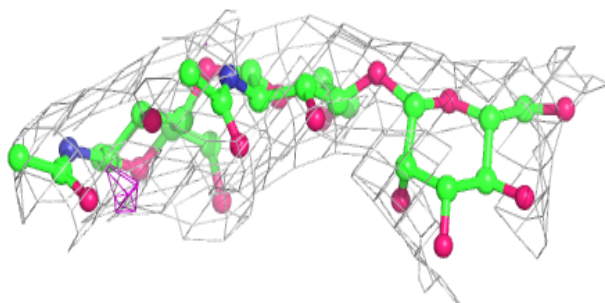
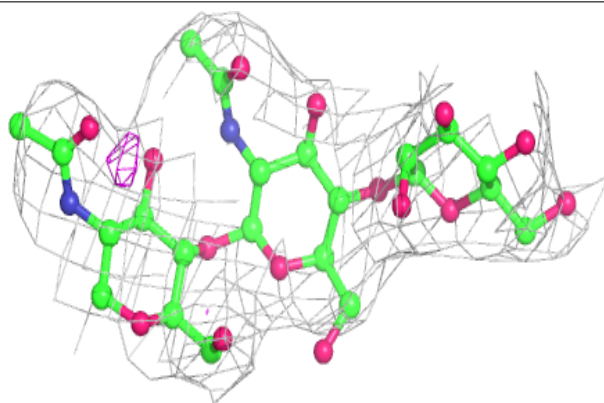


**Electron density around Chain G:**

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

**Electron density around Chain I:**

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

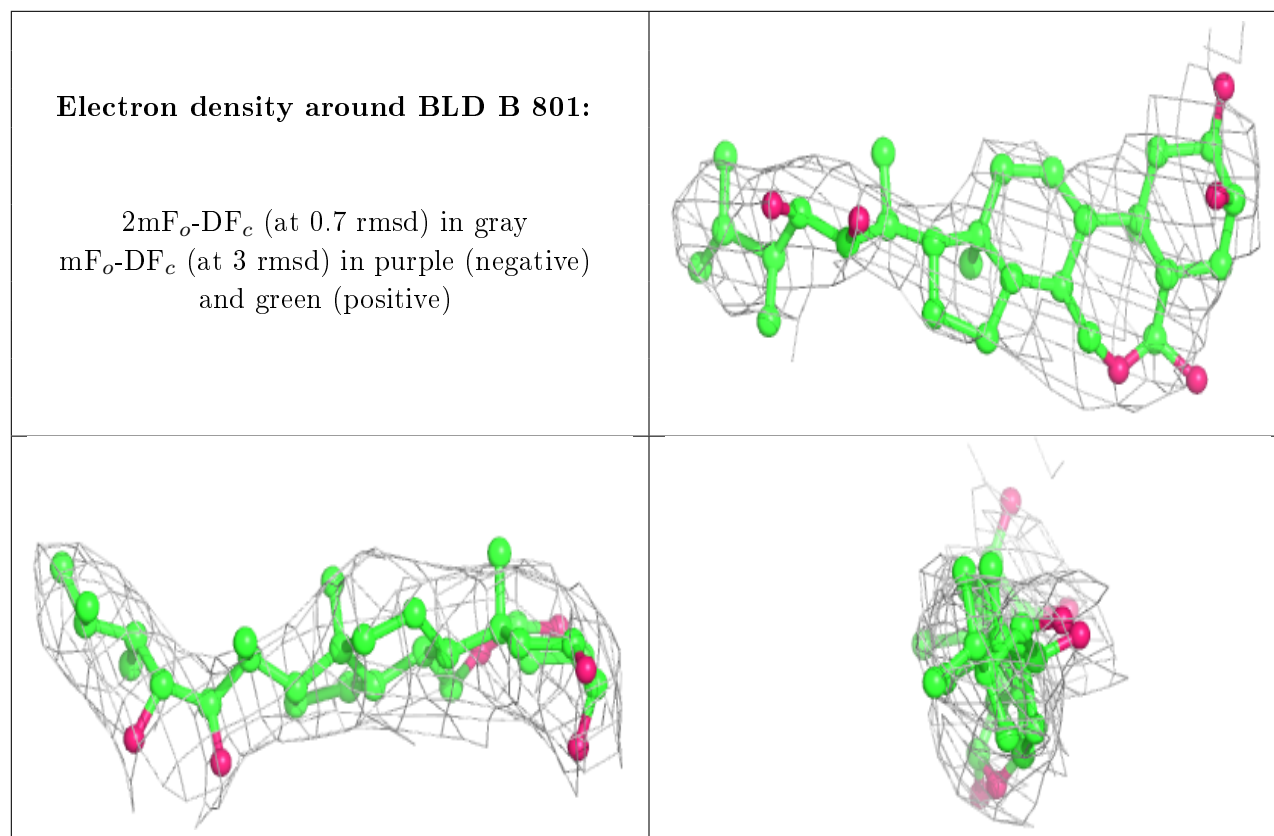


## 6.4 Ligands ⓘ

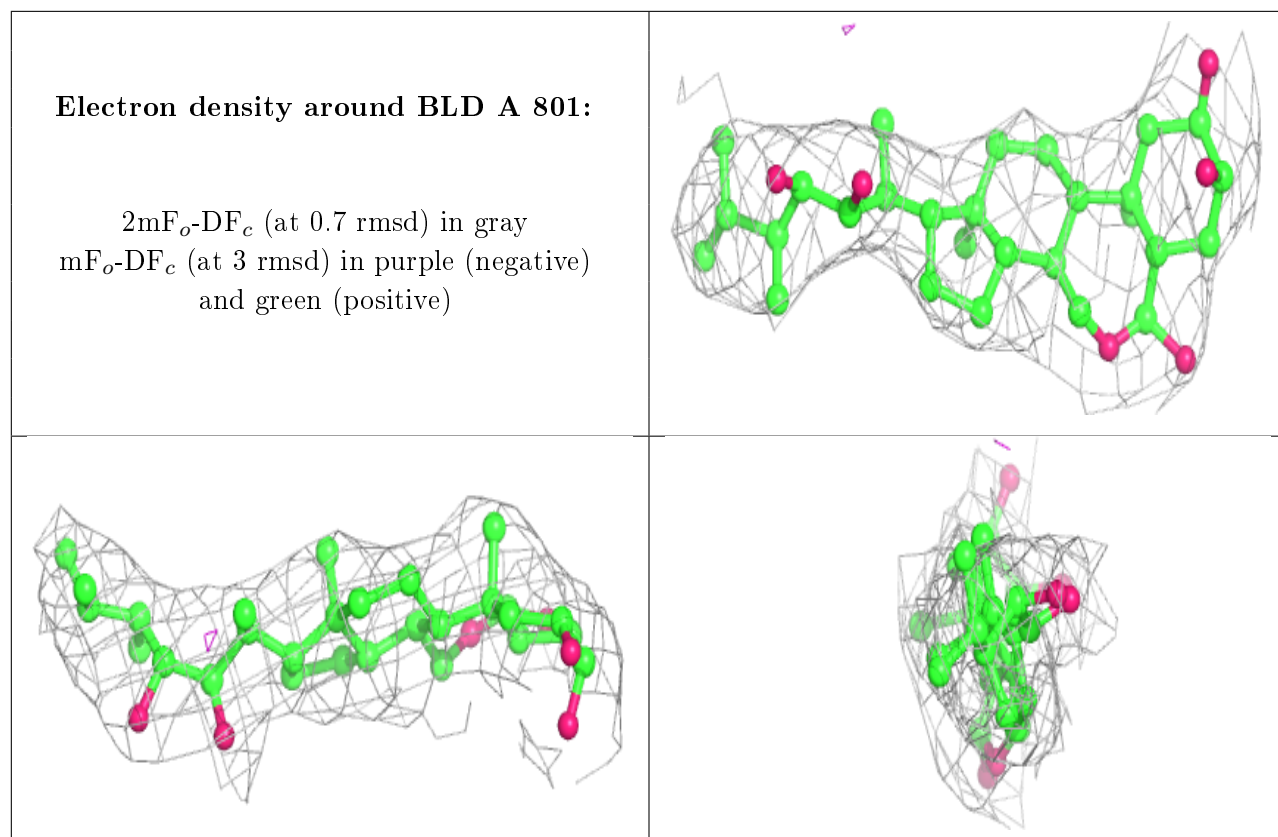
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
8	NAG	A	802	14/15	0.64	0.41	138,158,174,175	0
8	NAG	A	814	14/15	0.79	0.20	88,94,101,103	0
8	NAG	C	1001	14/15	0.84	0.18	77,91,108,109	0
8	NAG	D	1001	14/15	0.85	0.20	102,115,132,134	0
8	NAG	B	802	14/15	0.86	0.17	96,105,113,114	0
7	BLD	B	801	34/34	0.90	0.43	85,127,138,149	0
8	NAG	C	1000	14/15	0.90	0.19	96,99,105,106	0
8	NAG	D	1002	14/15	0.91	0.12	81,98,108,109	0
7	BLD	A	801	34/34	0.94	0.22	61,109,134,140	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.







## 6.5 Other polymers [i](#)

There are no such residues in this entry.