Full wwPDB X-ray Structure Validation Report (i)

Nov 14, 2023 – 01:05 PM JST

PDB ID	:	5ZD4
Title	:	Crystal structure of MBP-fused BIL1/BZR1 in complex with double-stranded DNA
Authors	:	Nosaki, S.; Miyakawa, T.; Xu, Y.; Nakamura, A.; Hirabayashi, K.; Tanokura, M.
Deposited on	:	2018-02-22
Resolution	:	2.17 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

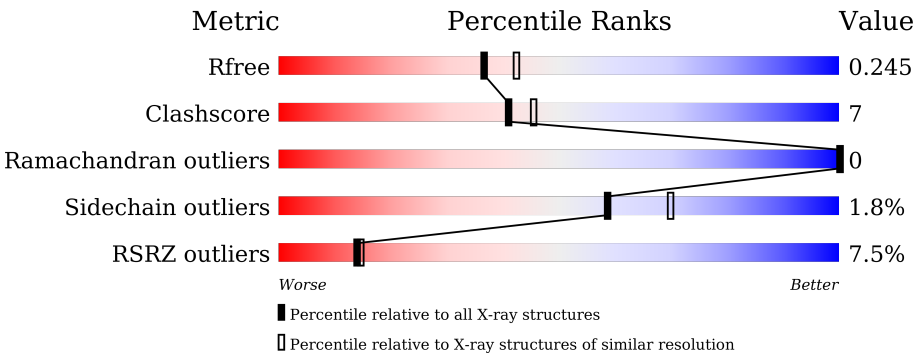
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	453	7%, 79%, 16%, • •
1	B	453	9%, 83%, 13%, •
1	C	453	5%, 86%, 9%, •
1	D	453	9%, 72%, 18%, 9%
2	E	15	60%, 33%, 7%
2	F	15	80%, 20%

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Mol	Chain	Length	Quality of chain
2	G	15	80% 20%
2	H	15	87% 13%
3	I	2	100%
3	J	2	50% 50%
3	K	2	100%
3	L	2	50% 50%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 15004 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 Maltose-binding periplasmic protein,Protein BRASSINAZOLE-RESISTANT 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	C	436	3394	2166	581	639	8	0	0	0
1	D	411	3208	2047	552	601	8	0	0	0
1	A	436	3394	2166	581	639	8	0	0	0
1	B	435	3385	2160	579	638	8	0	0	0

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-368	MET	-	expression tag	UNP P0AEY0
C	-286	ALA	ASP	engineered mutation	UNP P0AEY0
C	-285	ALA	LYS	engineered mutation	UNP P0AEY0
C	-196	ALA	GLU	engineered mutation	UNP P0AEY0
C	-195	ALA	ASN	engineered mutation	UNP P0AEY0
C	-129	ALA	LYS	engineered mutation	UNP P0AEY0
C	-9	ALA	GLU	engineered mutation	UNP P0AEY0
C	-6	ALA	LYS	engineered mutation	UNP P0AEY0
C	-5	ALA	ASP	engineered mutation	UNP P0AEY0
C	-1	ASN	-	linker	UNP P0AEY0
C	0	ALA	-	linker	UNP P0AEY0
D	-368	MET	-	expression tag	UNP P0AEY0
D	-286	ALA	ASP	engineered mutation	UNP P0AEY0
D	-285	ALA	LYS	engineered mutation	UNP P0AEY0
D	-196	ALA	GLU	engineered mutation	UNP P0AEY0
D	-195	ALA	ASN	engineered mutation	UNP P0AEY0
D	-129	ALA	LYS	engineered mutation	UNP P0AEY0
D	-9	ALA	GLU	engineered mutation	UNP P0AEY0
D	-6	ALA	LYS	engineered mutation	UNP P0AEY0
D	-5	ALA	ASP	engineered mutation	UNP P0AEY0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	ASN	-	linker	UNP P0AEY0
D	0	ALA	-	linker	UNP P0AEY0
A	-368	MET	-	expression tag	UNP P0AEY0
A	-286	ALA	ASP	engineered mutation	UNP P0AEY0
A	-285	ALA	LYS	engineered mutation	UNP P0AEY0
A	-196	ALA	GLU	engineered mutation	UNP P0AEY0
A	-195	ALA	ASN	engineered mutation	UNP P0AEY0
A	-129	ALA	LYS	engineered mutation	UNP P0AEY0
A	-9	ALA	GLU	engineered mutation	UNP P0AEY0
A	-6	ALA	LYS	engineered mutation	UNP P0AEY0
A	-5	ALA	ASP	engineered mutation	UNP P0AEY0
A	-1	ASN	-	linker	UNP P0AEY0
A	0	ALA	-	linker	UNP P0AEY0
B	-368	MET	-	expression tag	UNP P0AEY0
B	-286	ALA	ASP	engineered mutation	UNP P0AEY0
B	-285	ALA	LYS	engineered mutation	UNP P0AEY0
B	-196	ALA	GLU	engineered mutation	UNP P0AEY0
B	-195	ALA	ASN	engineered mutation	UNP P0AEY0
B	-129	ALA	LYS	engineered mutation	UNP P0AEY0
B	-9	ALA	GLU	engineered mutation	UNP P0AEY0
B	-6	ALA	LYS	engineered mutation	UNP P0AEY0
B	-5	ALA	ASP	engineered mutation	UNP P0AEY0
B	-1	ASN	-	linker	UNP P0AEY0
B	0	ALA	-	linker	UNP P0AEY0

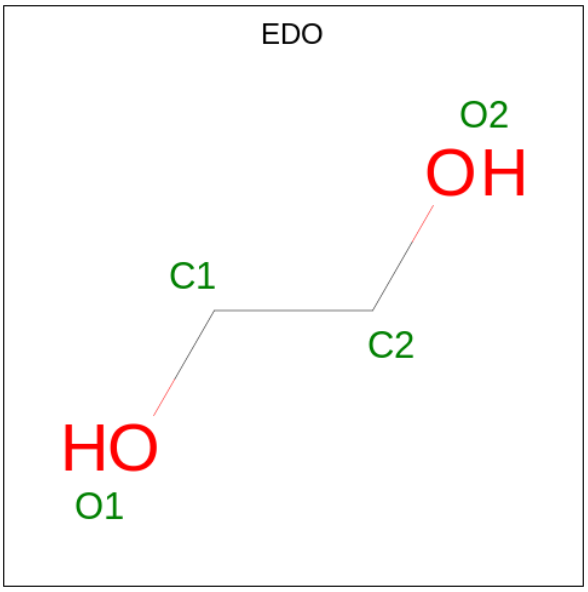
- Molecule 2 is a DNA chain called DNA (5'-D(*TP*TP*CP*AP*CP*AP*CP*GP*TP*GP*TP*GP*AP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	G	15	305	147	57	87	14	0	0	0
2	H	15	305	147	57	87	14	0	0	0
2	E	15	305	147	57	87	14	0	0	0
2	F	15	305	147	57	87	14	0	0	0

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
3	I	2	23	12	11	0	0	0
3	J	2	23	12	11	0	0	0
3	K	2	23	12	11	0	0	0
3	L	2	23	12	11	0	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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

- Molecule 5 is water.

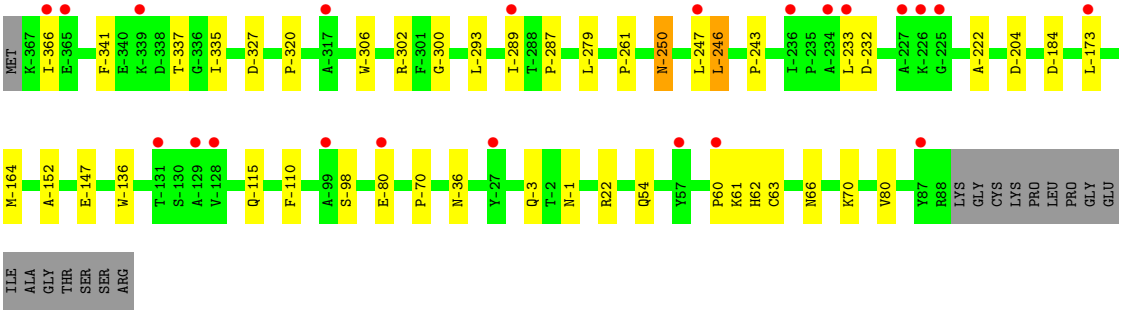
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	C	55	55	55	0	0
			Total	O		
5	D	30	30	30	0	0
			Total	O		
5	G	13	13	13	0	0
			Total	O		
5	H	13	13	13	0	0
			Total	O		
5	A	62	62	62	0	0
			Total	O		
5	B	94	94	94	0	0
			Total	O		
5	E	7	7	7	0	0
			Total	O		
5	F	9	9	9	0	0

3 Residue-property plots (i)

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

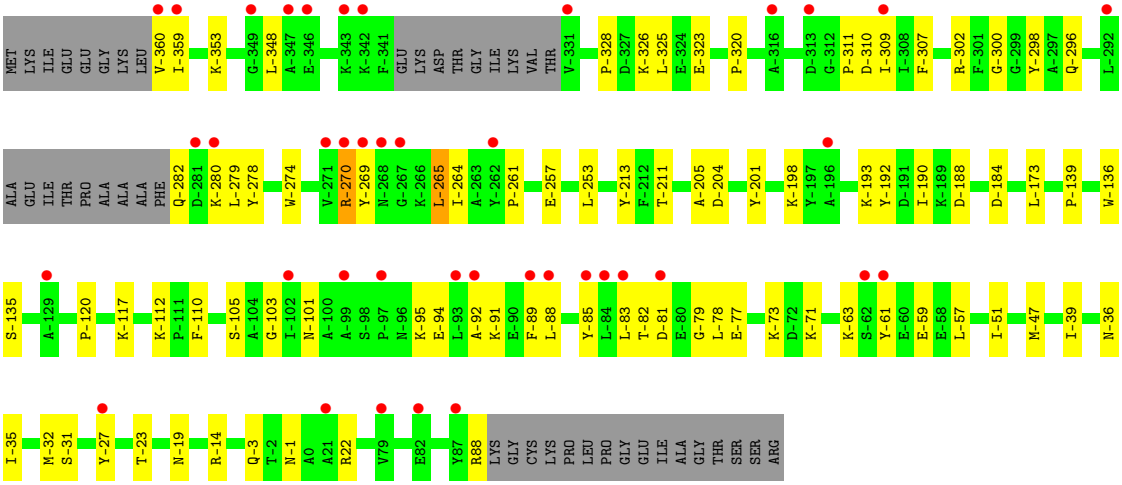
• Molecule 1: Maltose-binding periplasmic protein,Protein BRASSINAZOLE-RESISTANT 1

Chain C:



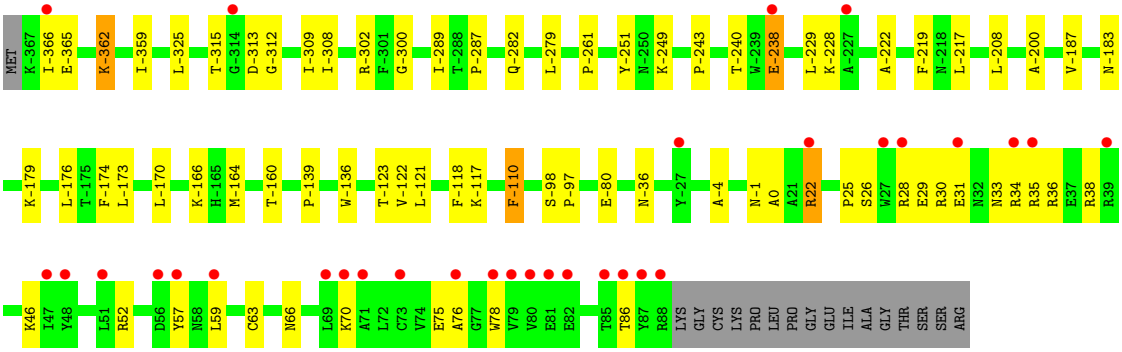
• Molecule 1: Maltose-binding periplasmic protein,Protein BRASSINAZOLE-RESISTANT 1

Chain D:



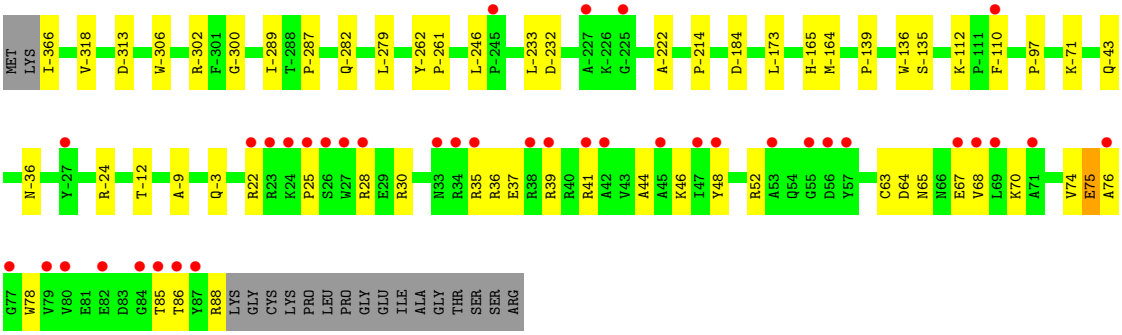
• Molecule 1: Maltose-binding periplasmic protein,Protein BRASSINAZOLE-RESISTANT 1

Chain A:



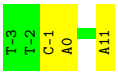
• Molecule 1: Maltose-binding periplasmic protein,Protein BRASSINAZOLE-RESISTANT 1

Chain B:



• Molecule 2: DNA (5'-D(*TP*TP*CP*AP*CP*AP*CP*GP*TP*GP*TP*GP*AP*AP*A)-3')

Chain G:

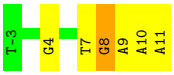


• Molecule 2: DNA (5'-D(*TP*TP*CP*AP*CP*AP*CP*GP*TP*GP*TP*GP*AP*AP*A)-3')

Chain H:

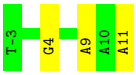
• Molecule 2: DNA (5'-D(*TP*TP*CP*AP*CP*AP*CP*GP*TP*GP*TP*GP*AP*AP*A)-3')

Chain E:



• Molecule 2: DNA (5'-D(*TP*TP*CP*AP*CP*AP*CP*GP*TP*GP*TP*GP*AP*AP*A)-3')

Chain F: 80% 20%



• Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain I: 100%

• Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain J: 50% 50%

• Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain K: 100%

• Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain L: 50% 50%

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	102.70Å 92.60Å 111.97Å 90.00° 100.42° 90.00°	Depositor
Resolution (Å)	43.60 – 2.17 43.60 – 2.17	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.60-2.17) 99.8 (43.60-2.17)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.58 (at 2.18Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.217 , 0.244 0.218 , 0.245	Depositor DCC
R_{free} test set	5449 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	35.4	Xtriage
Anisotropy	0.509	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 36.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	15004	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.28 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3527e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: GLC, EDO

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	0.37	0/3473	0.50	0/4713
1	B	0.31	0/3464	0.49	0/4702
1	C	0.27	0/3473	0.45	0/4713
1	D	0.39	0/3283	0.51	0/4454
2	E	0.99	3/342 (0.9%)	1.02	2/526 (0.4%)
2	F	0.94	1/342 (0.3%)	1.00	1/526 (0.2%)
2	G	0.85	0/342	1.01	1/526 (0.2%)
2	H	0.97	0/342	0.98	1/526 (0.2%)
All	All	0.43	4/15061 (0.0%)	0.56	5/20686 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	9	DA	O3'-P	-6.37	1.53	1.61
2	F	9	DA	O3'-P	-6.31	1.53	1.61
2	E	8	DG	O3'-P	-5.33	1.54	1.61
2	E	10	DA	O3'-P	-5.03	1.55	1.61

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	G	11	DA	C1'-O4'-C4'	-7.63	102.47	110.10
2	E	11	DA	C1'-O4'-C4'	-7.41	102.69	110.10
2	F	11	DA	C1'-O4'-C4'	-7.04	103.06	110.10
2	H	11	DA	C1'-O4'-C4'	-6.84	103.25	110.10
2	E	8	DG	C1'-O4'-C4'	-6.19	103.91	110.10

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	3394	0	3365	67	0
1	B	3385	0	3352	40	0
1	C	3394	0	3365	27	0
1	D	3208	0	3173	72	0
2	E	305	0	171	2	0
2	F	305	0	171	1	0
2	G	305	0	171	1	0
2	H	305	0	171	1	0
3	I	23	0	21	0	0
3	J	23	0	21	1	0
3	K	23	0	21	0	0
3	L	23	0	21	0	0
4	A	8	0	12	1	0
4	B	8	0	12	1	0
4	C	8	0	12	1	0
4	D	4	0	6	0	0
5	A	62	0	0	0	0
5	B	94	0	0	1	0
5	C	55	0	0	0	0
5	D	30	0	0	0	0
5	E	7	0	0	0	0
5	F	9	0	0	0	0
5	G	13	0	0	0	0
5	H	13	0	0	0	0
All	All	15004	0	14065	204	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 (204) 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:-240:THR:HG22	1:A:-238:GLU:H	1.10	1.11
1:C:-366:ILE:HD12	1:C:-98:SER:HA	1.38	1.04
1:A:-240:THR:HG22	1:A:-238:GLU:N	1.79	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-77:GLU:HG2	1:D:-73:LYS:HE3	1.48	0.95
1:B:-12:THR:HG23	1:B:-9:ALA:H	1.32	0.94
1:D:-82:THR:HG23	1:D:-79:GLY:H	1.33	0.93
1:D:-204:ASP:OD2	1:D:-117:LYS:HE2	1.69	0.92
1:B:64:ASP:HB2	1:B:67:GLU:HG2	1.53	0.89
1:C:-250:ASN:ND2	1:C:-247:LEU:H	1.71	0.88
1:C:-289:ILE:HG22	1:C:-287:PRO:HD3	1.57	0.87
1:A:-365:GLU:H	1:A:-362:LYS:HE3	1.39	0.86
1:D:-204:ASP:CG	1:D:-117:LYS:HE2	1.99	0.83
1:A:-240:THR:CG2	1:A:-238:GLU:H	1.94	0.80
1:A:31:GLU:HG2	1:A:34:ARG:NH2	1.97	0.79
1:D:-193:LYS:NZ	1:D:-192:TYR:O	2.16	0.78
1:B:36:ARG:CD	1:B:39:ARG:HH21	1.97	0.77
1:D:-77:GLU:HG2	1:D:-73:LYS:CE	2.15	0.75
1:C:-232:ASP:OD1	1:C:-222:ALA:N	2.20	0.75
1:A:-289:ILE:HG22	1:A:-287:PRO:HD3	1.70	0.73
1:C:-246:LEU:HD21	1:C:-243:PRO:HA	1.72	0.71
1:C:60:PRO:HG2	1:C:63:CYS:HB3	1.74	0.70
1:A:-325:LEU:HD13	1:A:-308:ILE:HD11	1.74	0.70
1:C:-250:ASN:HD22	1:C:-247:LEU:H	1.40	0.70
1:A:31:GLU:HG2	1:A:34:ARG:HH21	1.57	0.69
1:D:-269:TYR:HB3	1:D:-264:ILE:HD11	1.73	0.69
1:D:-95:LYS:HZ2	1:D:-91:LYS:HB2	1.58	0.69
1:D:-360:VAL:O	1:D:-359:ILE:HD13	1.93	0.68
1:A:-366:ILE:HD12	1:A:-98:SER:HA	1.75	0.68
1:D:-309:ILE:HD11	1:D:-92:ALA:HB1	1.75	0.68
1:C:-1:ASN:OD1	1:C:22:ARG:NH2	2.27	0.68
1:D:-326:LYS:HD2	1:D:-323:GLU:OE1	1.94	0.67
1:A:-240:THR:HG21	1:A:-238:GLU:HB2	1.77	0.67
1:B:-289:ILE:HG22	1:B:-287:PRO:HD3	1.77	0.66
1:D:-81:ASP:O	1:D:-77:GLU:HB2	1.96	0.65
1:B:36:ARG:HD3	1:B:39:ARG:HH21	1.61	0.65
1:D:-193:LYS:HD2	1:D:-192:TYR:H	1.60	0.65
1:D:-359:ILE:HD11	1:D:-310:ASP:HB2	1.79	0.65
1:B:-246:LEU:HD21	1:B:-233:LEU:HD21	1.79	0.65
1:D:-359:ILE:CD1	1:D:-309:ILE:HB	2.27	0.65
1:D:-348:LEU:HD11	1:D:-89:PHE:HZ	1.61	0.65
1:D:-83:LEU:HA	1:D:-78:LEU:HD11	1.79	0.65
1:A:-240:THR:CG2	1:A:-238:GLU:HB2	2.27	0.64
1:B:41:ARG:NH2	2:F:4:DG:N7	2.46	0.64
1:A:-362:LYS:C	1:A:-362:LYS:HD2	2.17	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:-306:TRP:CD1	1:B:-302:ARG:HG3	2.33	0.64
1:D:-269:TYR:HB3	1:D:-264:ILE:CD1	2.28	0.64
1:C:-279:LEU:HD23	1:C:-261:PRO:HG2	1.80	0.63
1:A:26:SER:OG	1:A:28:ARG:NH2	2.32	0.62
4:C:203:EDO:H12	2:H:1:DC:H5"	1.81	0.62
1:A:-365:GLU:N	1:A:-362:LYS:HE3	2.12	0.62
1:A:-251:TYR:OH	1:A:-249:LYS:NZ	2.34	0.60
1:A:-1:ASN:HA	1:A:22:ARG:HD2	1.83	0.60
1:A:-365:GLU:H	1:A:-362:LYS:CE	2.10	0.60
1:A:-279:LEU:HD23	1:A:-261:PRO:HG2	1.83	0.60
1:D:-19:ASN:OD1	1:D:-14:ARG:NH1	2.34	0.60
1:B:-279:LEU:HD23	1:B:-261:PRO:HG2	1.84	0.60
1:D:-184:ASP:HB2	1:D:-3:GLN:HB2	1.84	0.60
1:D:-83:LEU:HA	1:D:-78:LEU:CD1	2.31	0.59
1:C:-184:ASP:HB2	1:C:-3:GLN:HB2	1.84	0.59
1:D:-353:LYS:NZ	1:D:-257:GLU:OE2	2.35	0.59
1:D:-326:LYS:HG2	1:D:-323:GLU:HB2	1.82	0.59
1:B:37:GLU:O	1:B:41:ARG:HG3	2.03	0.59
1:D:-1:ASN:OD1	1:D:22:ARG:NH1	2.35	0.59
1:D:-135:SER:HB3	1:D:-71:LYS:HD3	1.85	0.59
1:D:-326:LYS:CD	1:D:-323:GLU:OE1	2.51	0.58
1:A:46:LYS:HE2	1:B:85:THR:HA	1.84	0.58
1:A:34:ARG:O	1:A:38:ARG:HG3	2.03	0.58
1:D:-326:LYS:CG	1:D:-323:GLU:HB2	2.33	0.58
1:D:-296:GLN:NE2	1:D:-269:TYR:OH	2.37	0.58
1:B:25:PRO:HG2	1:B:30:ARG:NE	2.19	0.58
1:D:-61:TYR:CZ	1:D:-57:LEU:HD21	2.39	0.58
1:D:-300:GLY:HA3	1:D:-36:ASN:O	2.04	0.57
1:C:-341:PHE:O	1:C:-337:THR:HG22	2.05	0.57
1:B:36:ARG:CD	1:B:39:ARG:NH2	2.67	0.57
1:D:-359:ILE:HD12	1:D:-309:ILE:HB	1.85	0.56
1:A:76:ALA:HB1	1:B:75:GLU:OE2	2.05	0.56
1:A:25:PRO:HG2	1:A:30:ARG:NE	2.21	0.55
1:A:22:ARG:HD3	1:A:22:ARG:N	2.21	0.55
1:D:-264:ILE:HG22	1:D:-264:ILE:O	2.07	0.54
1:A:-4:ALA:O	1:A:0:ALA:N	2.39	0.54
1:D:-61:TYR:CE2	1:D:-57:LEU:HD11	2.42	0.54
1:A:28:ARG:NH1	1:A:29:GLU:HG3	2.21	0.54
1:B:-318:VAL:HB	1:B:-313:ASP:HB2	1.88	0.54
1:A:-122:VAL:HG12	1:A:-121:LEU:O	2.07	0.54
1:A:22:ARG:HD3	1:A:22:ARG:H	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-279:LEU:HD23	1:D:-261:PRO:HG2	1.89	0.54
1:C:61:LYS:HG2	1:C:62:HIS:CD2	2.42	0.54
1:A:75:GLU:OE2	1:B:88:ARG:NH2	2.41	0.54
1:C:-246:LEU:HD11	1:C:-233:LEU:HD11	1.90	0.53
1:D:-326:LYS:HD2	1:D:-323:GLU:CD	2.28	0.53
1:D:-77:GLU:O	1:D:-73:LYS:HD2	2.09	0.53
1:D:-309:ILE:HA	1:D:-103:GLY:O	2.09	0.52
1:B:36:ARG:HD2	1:B:39:ARG:NH2	2.24	0.52
1:B:65:ASN:HA	1:B:68:VAL:HG12	1.91	0.52
1:D:-270:ARG:HB2	1:D:-265:LEU:HD22	1.91	0.52
1:D:-205:ALA:O	1:D:-112:LYS:NZ	2.43	0.52
1:D:-39:ILE:HD12	1:D:-39:ILE:O	2.10	0.52
1:A:-366:ILE:CD1	1:A:-98:SER:HA	2.40	0.52
1:D:-201:TYR:OH	1:D:-188:ASP:OD1	2.25	0.52
1:A:-183:ASN:O	1:A:-179:LYS:HG3	2.10	0.52
1:D:-359:ILE:HD12	1:D:-309:ILE:H	1.75	0.51
1:A:-217:LEU:HD23	1:A:-160:THR:HB	1.91	0.51
1:B:52:ARG:NH2	1:B:63:CYS:O	2.40	0.51
1:D:-51:ILE:O	1:D:-47:MET:HG2	2.12	0.50
1:A:-366:ILE:HD12	1:A:-97:PRO:CD	2.41	0.50
1:C:-152:ALA:HA	1:C:-147:GLU:OE1	2.11	0.50
1:D:-95:LYS:NZ	1:D:-91:LYS:HD2	2.27	0.50
1:A:-366:ILE:HD12	1:A:-97:PRO:HD3	1.93	0.49
1:B:-232:ASP:OD2	1:B:-165:HIS:HD2	1.94	0.49
1:D:-302:ARG:NH1	1:D:-31:SER:HA	2.28	0.49
1:C:-306:TRP:CD1	1:C:-302:ARG:HG3	2.48	0.49
1:A:-229:LEU:HD13	1:A:-222:ALA:HA	1.95	0.49
1:A:-219:PHE:O	1:A:-217:LEU:HD22	2.12	0.49
1:C:-300:GLY:HA3	1:C:-36:ASN:O	2.13	0.48
1:D:-190:ILE:HD11	1:D:-35:ILE:HD12	1.95	0.48
1:B:78:TRP:CE3	1:B:86:THR:HG22	2.48	0.48
1:A:78:TRP:CE3	1:A:86:THR:HG22	2.48	0.48
1:B:-287:PRO:HG2	1:B:-282:GLN:HG3	1.96	0.48
4:B:203:EDO:H21	5:B:303:HOH:O	2.13	0.48
1:C:-204:ASP:HA	1:C:-115:GLN:OE1	2.14	0.48
1:B:-246:LEU:HD11	1:B:-233:LEU:HD11	1.96	0.48
1:D:-95:LYS:HZ2	1:D:-91:LYS:HD2	1.78	0.48
1:D:-253:LEU:HD22	1:D:-120:PRO:HD3	1.95	0.48
1:A:-110:PHE:CE1	4:A:203:EDO:H11	2.49	0.47
1:D:-32:MET:HE2	1:D:-32:MET:HA	1.95	0.47
1:B:-12:THR:CG2	1:B:-9:ALA:H	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:ARG:NH2	1:A:63:CYS:O	2.46	0.47
1:D:-311:PRO:O	1:D:-101:ASN:HB2	2.14	0.47
1:C:54:GLN:O	1:D:88:ARG:NH1	2.48	0.46
1:A:33:ASN:HA	1:A:36:ARG:HH12	1.80	0.46
1:B:-300:GLY:HA3	1:B:-36:ASN:O	2.15	0.46
1:A:-118:PHE:CE2	1:A:-117:LYS:HD2	2.50	0.46
1:B:70:LYS:O	1:B:74:VAL:HG13	2.15	0.46
1:C:-366:ILE:CD1	1:C:-98:SER:HA	2.26	0.46
1:A:-251:TYR:CD2	1:A:-243:PRO:HG3	2.50	0.46
2:G:-1:DC:H2"	2:G:0:DA:C8	2.51	0.46
1:A:-217:LEU:CD2	1:A:-160:THR:HB	2.45	0.46
1:A:57:TYR:HB2	1:A:59:LEU:CD2	2.46	0.46
1:D:-27:TYR:O	1:D:-23:THR:HG23	2.16	0.46
1:B:-139:PRO:HA	1:B:-136:TRP:CE2	2.52	0.45
1:A:-139:PRO:HA	1:A:-136:TRP:CE2	2.51	0.45
1:A:34:ARG:HD3	2:E:4:DG:H3'	1.97	0.45
1:A:34:ARG:HB3	1:A:38:ARG:NH1	2.32	0.45
1:A:66:ASN:O	1:A:70:LYS:HG3	2.16	0.45
1:B:-173:LEU:HD12	1:B:-164:MET:HE1	1.97	0.45
1:C:70:LYS:HG2	1:C:80:VAL:HG11	1.99	0.45
1:D:-320:PRO:HG3	1:D:-298:TYR:CE1	2.52	0.45
1:D:-278:TYR:HA	1:D:-63:LYS:HE3	1.99	0.45
1:A:-300:GLY:HA3	1:A:-36:ASN:O	2.17	0.45
1:B:44:ALA:O	1:B:48:TYR:HD1	1.99	0.45
1:D:-328:PRO:HG2	1:D:-325:LEU:HB3	1.99	0.45
1:D:-326:LYS:HE2	1:D:-326:LYS:HB2	1.70	0.45
1:D:-274:TRP:CZ3	1:D:-261:PRO:HD2	2.52	0.44
1:C:-136:TRP:HB2	1:C:-70:PRO:HG2	1.99	0.44
1:D:-211:THR:HG21	1:D:-173:LEU:HD22	1.99	0.44
1:C:-246:LEU:CD2	1:C:-243:PRO:HA	2.42	0.44
1:D:-309:ILE:HD13	1:D:-88:LEU:HD11	1.99	0.44
1:D:-307:PHE:HA	1:D:-105:SER:O	2.17	0.44
1:D:-77:GLU:CG	1:D:-73:LYS:HE3	2.33	0.44
1:A:-228:LYS:NZ	1:A:-166:LYS:O	2.50	0.44
1:A:28:ARG:HH12	1:A:29:GLU:HG3	1.83	0.44
1:A:-123:THR:OG1	1:A:-122:VAL:N	2.51	0.43
1:D:-63:LYS:O	1:D:-59:GLU:HG3	2.18	0.43
1:A:-176:LEU:CD1	1:A:-173:LEU:HD23	2.49	0.43
1:A:46:LYS:NZ	1:B:85:THR:HG22	2.34	0.43
1:D:-139:PRO:HA	1:D:-136:TRP:CE2	2.53	0.43
1:A:-173:LEU:HD12	1:A:-164:MET:HE1	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-200:ALA:O	1:A:-187:VAL:HA	2.19	0.43
1:C:-337:THR:HG23	1:C:-335:ILE:HG12	2.01	0.43
1:B:-135:SER:OG	1:B:-71:LYS:HD3	2.18	0.43
1:A:-325:LEU:CD1	1:A:-308:ILE:HD11	2.46	0.42
1:D:-193:LYS:HD2	1:D:-192:TYR:N	2.32	0.42
1:A:-315:THR:HG22	1:A:-313:ASP:OD1	2.19	0.42
1:B:41:ARG:HG3	1:B:41:ARG:H	1.71	0.42
1:C:-320:PRO:HA	1:C:-293:LEU:CD1	2.50	0.42
1:D:-359:ILE:HD12	1:D:-309:ILE:N	2.35	0.42
1:D:-77:GLU:HG3	1:D:-61:TYR:OH	2.20	0.42
1:A:-251:TYR:CE2	1:A:-243:PRO:HG3	2.55	0.42
1:B:-214:PRO:HD3	1:B:-24:ARG:HG3	2.02	0.42
2:E:7:DT:H2"	2:E:8:DG:O4'	2.20	0.41
1:A:33:ASN:HA	1:A:36:ARG:NH1	2.36	0.41
1:B:64:ASP:CB	1:B:67:GLU:HG2	2.37	0.41
1:C:-173:LEU:HD12	1:C:-164:MET:HE1	2.02	0.41
1:D:-280:LYS:O	1:D:-63:LYS:HD2	2.20	0.41
1:A:-174:PHE:CE2	1:A:-170:LEU:HD11	2.55	0.41
1:B:-366:ILE:HG21	1:B:-97:PRO:HD2	2.02	0.41
1:A:22:ARG:NH2	1:A:22:ARG:O	2.52	0.41
1:C:-366:ILE:H	1:C:-366:ILE:HG12	1.57	0.41
1:D:-95:LYS:HE3	1:D:-94:GLU:HG2	2.02	0.41
1:D:-85:TYR:O	1:D:-82:THR:HG22	2.21	0.41
1:A:-366:ILE:HG22	1:A:-362:LYS:HZ3	1.86	0.41
1:A:-359:ILE:HG12	1:A:-309:ILE:HB	2.01	0.41
1:A:-287:PRO:HG2	1:A:-282:GLN:CG	2.50	0.41
1:B:-289:ILE:HD12	1:B:-262:TYR:CE2	2.56	0.41
1:D:-359:ILE:CD1	1:D:-309:ILE:H	2.33	0.40
1:D:-61:TYR:CD2	1:D:-57:LEU:HD11	2.57	0.40
1:B:-232:ASP:HA	1:B:-222:ALA:HB2	2.03	0.40
1:B:-112:LYS:HB2	1:B:-112:LYS:HE3	1.92	0.40
1:A:-366:ILE:HG23	1:A:-312:GLY:O	2.21	0.40
1:A:76:ALA:HB2	1:B:76:ALA:HB2	2.03	0.40
1:C:66:ASN:O	1:C:70:LYS:HG3	2.21	0.40
1:A:-187:VAL:HG11	1:A:0:ALA:HB3	2.03	0.40
1:B:-184:ASP:HB2	1:B:-3:GLN:HB2	2.04	0.40
1:D:-213:TYR:HB2	3:J:2:GLC:O5	2.20	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	434/453 (96%)	428 (99%)	6 (1%)	0	100	100
1	B	433/453 (96%)	428 (99%)	5 (1%)	0	100	100
1	C	434/453 (96%)	427 (98%)	7 (2%)	0	100	100
1	D	405/453 (89%)	397 (98%)	8 (2%)	0	100	100
All	All	1706/1812 (94%)	1680 (98%)	26 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	343/356 (96%)	335 (98%)	8 (2%)	50	60
1	B	342/356 (96%)	335 (98%)	7 (2%)	55	66
1	C	343/356 (96%)	338 (98%)	5 (2%)	65	76
1	D	324/356 (91%)	319 (98%)	5 (2%)	65	76
All	All	1352/1424 (95%)	1327 (98%)	25 (2%)	59	70

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	-327	ASP
1	C	-250	ASN

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Mol	Chain	Res	Type
1	C	-246	LEU
1	C	-110	PHE
1	C	-80	GLU
1	D	-282	GLN
1	D	-270	ARG
1	D	-265	LEU
1	D	-198	LYS
1	D	-110	PHE
1	A	-362	LYS
1	A	-302	ARG
1	A	-238	GLU
1	A	-208	LEU
1	A	-110	PHE
1	A	-80	GLU
1	A	22	ARG
1	A	35	ARG
1	B	-110	PHE
1	B	-43	GLN
1	B	22	ARG
1	B	28	ARG
1	B	35	ARG
1	B	46	LYS
1	B	75	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	-250	ASN
1	C	54	GLN
1	D	-296	GLN
1	D	-43	GLN
1	A	-3	GLN
1	B	-296	GLN
1	B	-165	HIS
1	B	-43	GLN
1	B	54	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

8 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	GLC	I	1	3	12,12,12	0.52	0	17,17,17	0.59	0
3	GLC	I	2	3	11,11,12	0.59	0	15,15,17	0.77	0
3	GLC	J	1	3	12,12,12	0.54	0	17,17,17	0.52	0
3	GLC	J	2	3	11,11,12	0.60	0	15,15,17	0.77	0
3	GLC	K	1	3	12,12,12	0.48	0	17,17,17	0.63	0
3	GLC	K	2	3	11,11,12	0.57	0	15,15,17	0.84	0
3	GLC	L	1	3	12,12,12	0.52	0	17,17,17	0.75	0
3	GLC	L	2	3	11,11,12	0.52	0	15,15,17	1.01	1 (6%)

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	GLC	I	1	3	-	1/2/22/22	0/1/1/1
3	GLC	I	2	3	-	0/2/19/22	0/1/1/1
3	GLC	J	1	3	-	0/2/22/22	0/1/1/1
3	GLC	J	2	3	-	0/2/19/22	0/1/1/1
3	GLC	K	1	3	-	0/2/22/22	0/1/1/1
3	GLC	K	2	3	-	0/2/19/22	0/1/1/1
3	GLC	L	1	3	-	0/2/22/22	0/1/1/1
3	GLC	L	2	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	L	2	GLC	C1-O5-C5	2.67	115.81	112.19

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	I	1	GLC	C4-C5-C6-O6

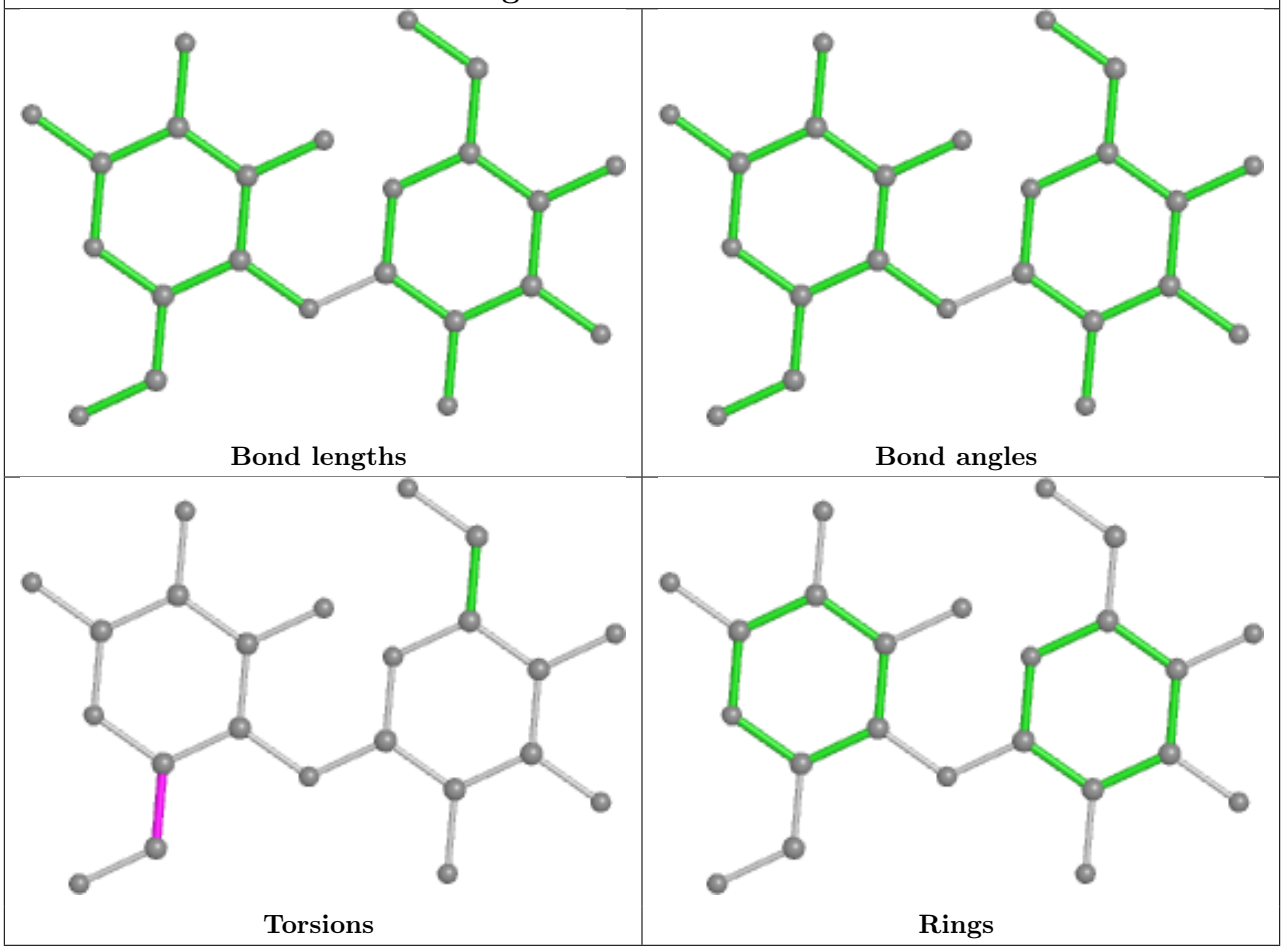
There are no ring outliers.

1 monomer is involved in 1 short contact:

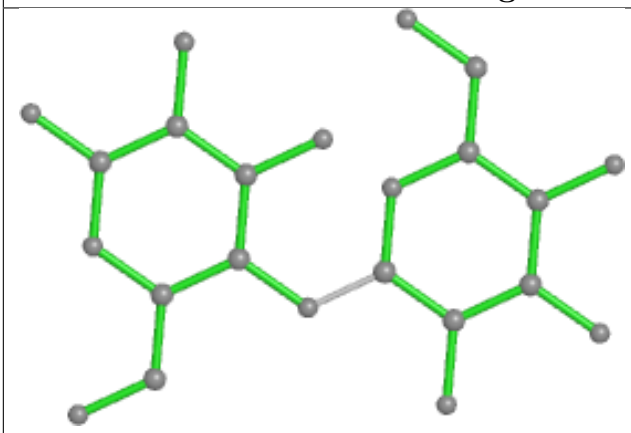
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	2	GLC	1	0

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

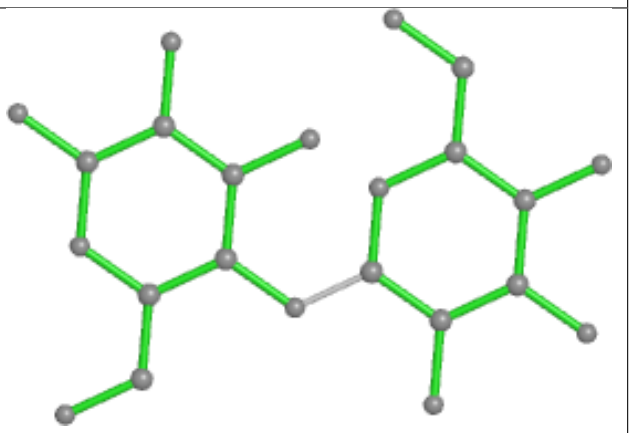
Oligosaccharide Chain I



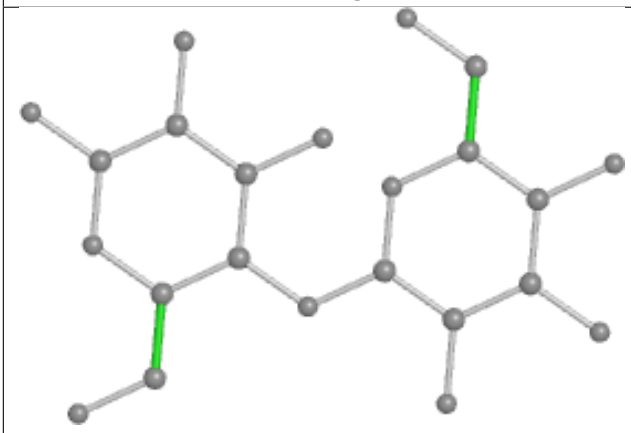
Oligosaccharide Chain J



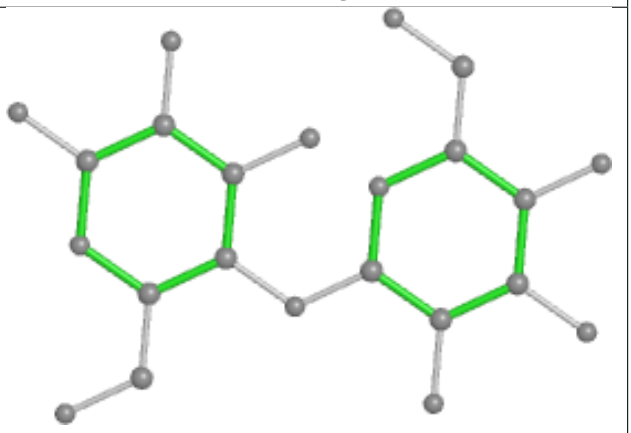
Bond lengths



Bond angles

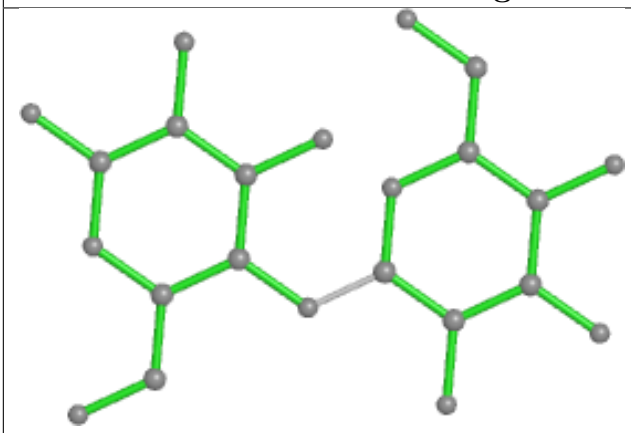


Torsions

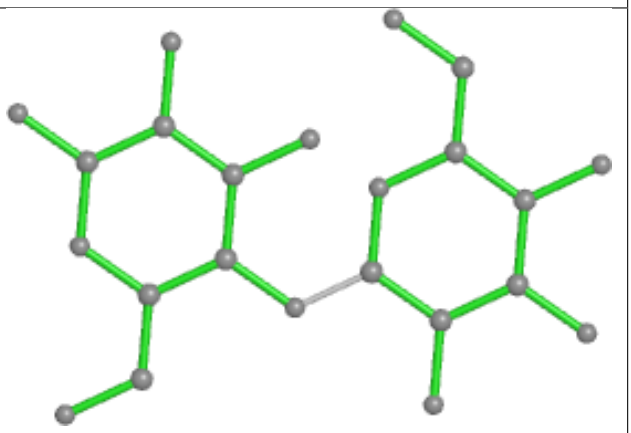


Rings

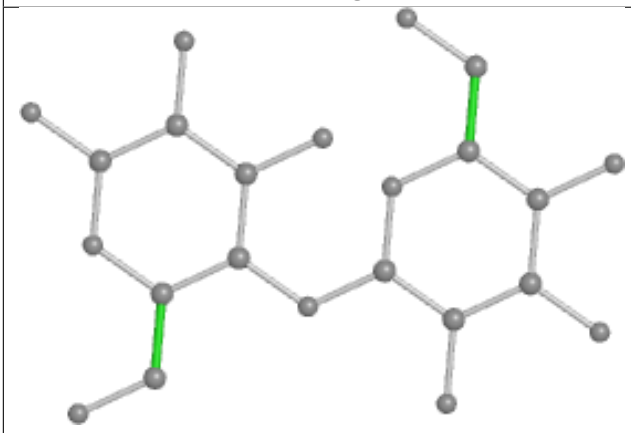
Oligosaccharide Chain K



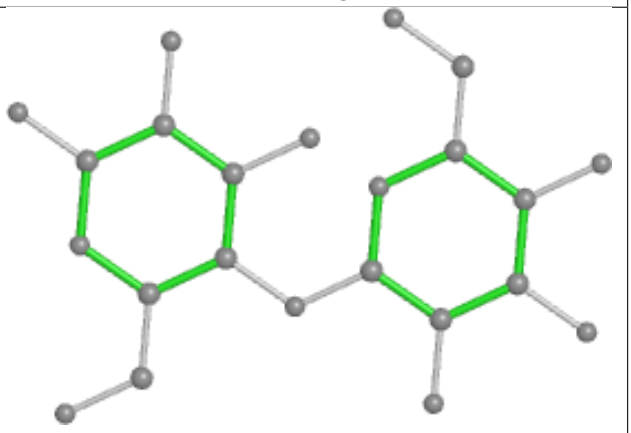
Bond lengths



Bond angles

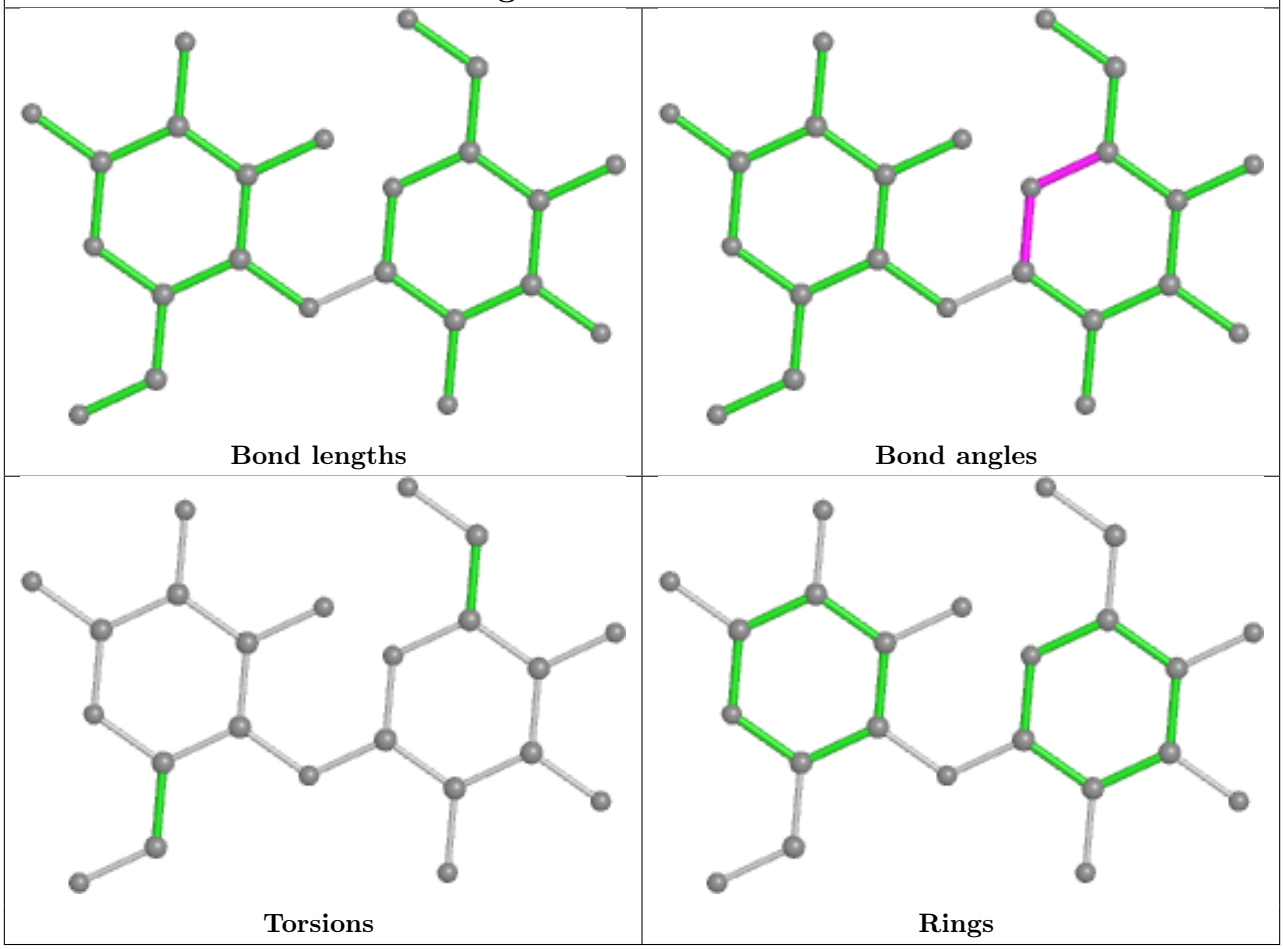


Torsions



Rings

Oligosaccharide Chain L



5.6 Ligand geometry (i)

7 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	$\#\|Z\| > 2$
4	EDO	B	202	-	3,3,3	0.47	0	2,2,2	0.30	0
4	EDO	C	203	-	3,3,3	0.46	0	2,2,2	0.33	0
4	EDO	A	202	-	3,3,3	0.47	0	2,2,2	0.30	0
4	EDO	C	202	-	3,3,3	0.46	0	2,2,2	0.32	0
4	EDO	A	203	-	3,3,3	0.46	0	2,2,2	0.31	0
4	EDO	D	202	-	3,3,3	0.46	0	2,2,2	0.36	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	EDO	B	203	-	3,3,3	0.46	0	2,2,2	0.45	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	202	-	-	0/1/1/1	-
4	EDO	C	203	-	-	0/1/1/1	-
4	EDO	A	202	-	-	1/1/1/1	-
4	EDO	C	202	-	-	0/1/1/1	-
4	EDO	A	203	-	-	1/1/1/1	-
4	EDO	D	202	-	-	0/1/1/1	-
4	EDO	B	203	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	203	EDO	O1-C1-C2-O2
4	A	202	EDO	O1-C1-C2-O2
4	A	203	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	203	EDO	1	0
4	A	203	EDO	1	0
4	B	203	EDO	1	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)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	436/453 (96%)	0.53	32 (7%)	15	15	26, 43, 79, 89	0
1	B	435/453 (96%)	0.54	39 (8%)	9	9	21, 40, 81, 95	0
1	C	436/453 (96%)	0.44	22 (5%)	28	30	27, 49, 78, 118	0
1	D	411/453 (90%)	0.67	40 (9%)	7	8	29, 56, 91, 101	0
2	E	15/15 (100%)	-0.22	0	100	100	46, 58, 76, 78	0
2	F	15/15 (100%)	-0.13	0	100	100	41, 64, 77, 80	0
2	G	15/15 (100%)	-0.09	0	100	100	30, 37, 62, 64	0
2	H	15/15 (100%)	-0.25	0	100	100	30, 42, 54, 55	0
All	All	1778/1872 (94%)	0.52	133 (7%)	14	15	21, 48, 83, 118	0

All (133) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	27	TRP	8.2
1	C	-129	ALA	7.6
1	B	27	TRP	6.8
1	D	-316	ALA	6.6
1	B	42	ALA	5.9
1	B	80	VAL	5.4
1	B	76	ALA	4.8
1	B	85	THR	4.8
1	A	87	TYR	4.7
1	A	80	VAL	4.7
1	A	47	ILE	4.5
1	D	-292	LEU	4.4
1	A	57	TYR	4.4
1	A	22	ARG	4.4
1	D	-129	ALA	4.4
1	D	-81	ASP	4.4

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Mol	Chain	Res	Type	RSRZ
1	D	-196	ALA	4.2
1	D	-102	ILE	4.2
1	C	-227	ALA	4.2
1	B	84	GLY	4.1
1	B	82	GLU	4.1
1	A	85	THR	4.0
1	A	79	VAL	3.9
1	B	24	LYS	3.9
1	C	-339	LYS	3.8
1	B	34	ARG	3.7
1	D	-92	ALA	3.7
1	B	-27	TYR	3.7
1	D	-97	PRO	3.7
1	C	-234	ALA	3.6
1	C	-225	GLY	3.6
1	B	87	TYR	3.6
1	D	-89	PHE	3.5
1	B	57	TYR	3.5
1	B	69	LEU	3.5
1	D	-342	LYS	3.5
1	B	26	SER	3.4
1	C	87	TYR	3.4
1	B	79	VAL	3.4
1	B	39	ARG	3.3
1	D	-93	LEU	3.3
1	B	53	ALA	3.2
1	B	33	ASN	3.2
1	A	35	ARG	3.2
1	B	25	PRO	3.2
1	B	77	GLY	3.2
1	A	48	TYR	3.1
1	A	76	ALA	3.0
1	D	-281	ASP	3.0
1	B	35	ARG	3.0
1	D	79	VAL	3.0
1	A	82	GLU	3.0
1	A	-366	ILE	3.0
1	A	70	LYS	3.0
1	A	56	ASP	3.0
1	C	-226	LYS	3.0
1	A	-27	TYR	3.0
1	B	23	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	39	ARG	2.9
1	D	-309	ILE	2.9
1	B	48	TYR	2.9
1	C	-128	VAL	2.9
1	A	81	GLU	2.9
1	B	28	ARG	2.9
1	B	86	THR	2.9
1	D	-267	GLY	2.9
1	A	69	LEU	2.8
1	D	-269	TYR	2.8
1	A	71	ALA	2.7
1	C	-173	LEU	2.7
1	A	-238	GLU	2.7
1	D	-27	TYR	2.7
1	A	28	ARG	2.7
1	B	56	ASP	2.7
1	B	45	ALA	2.7
1	B	38	ARG	2.7
1	C	-366	ILE	2.6
1	B	71	ALA	2.6
1	A	-227	ALA	2.6
1	D	-346	GLU	2.6
1	D	-262	TYR	2.6
1	C	-99	ALA	2.6
1	B	41	ARG	2.6
1	D	-84	LEU	2.6
1	C	-317	ALA	2.6
1	D	-331	VAL	2.6
1	A	86	THR	2.6
1	D	-349	GLY	2.6
1	D	-347	ALA	2.5
1	C	-27	TYR	2.5
1	A	73	CYS	2.5
1	D	-359	ILE	2.5
1	D	-268	ASN	2.5
1	D	-88	LEU	2.5
1	A	59	LEU	2.5
1	B	22	ARG	2.5
1	D	-61	TYR	2.5
1	B	55	GLY	2.5
1	D	-343	LYS	2.5
1	D	-270	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	87	TYR	2.4
1	A	78	TRP	2.4
1	D	-85	TYR	2.4
1	D	-99	ALA	2.4
1	B	47	ILE	2.4
1	C	-247	LEU	2.4
1	C	60	PRO	2.4
1	D	-83	LEU	2.3
1	C	-236	ILE	2.3
1	C	-233	LEU	2.3
1	A	51	LEU	2.3
1	D	-62	SER	2.3
1	A	31	GLU	2.3
1	A	34	ARG	2.3
1	D	82	GLU	2.3
1	B	-225	GLY	2.3
1	C	-289	ILE	2.3
1	C	-365	GLU	2.2
1	B	67	GLU	2.2
1	D	-360	VAL	2.2
1	C	-131	THR	2.2
1	B	-110	PHE	2.2
1	B	68	VAL	2.2
1	C	-80	GLU	2.2
1	B	-227	ALA	2.2
1	B	-245	PRO	2.2
1	A	88	ARG	2.1
1	D	-280	LYS	2.1
1	C	57	TYR	2.1
1	D	-313	ASP	2.1
1	D	21	ALA	2.0
1	D	-271	VAL	2.0
1	A	-314	GLY	2.0

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

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

6.3 Carbohydrates (i)

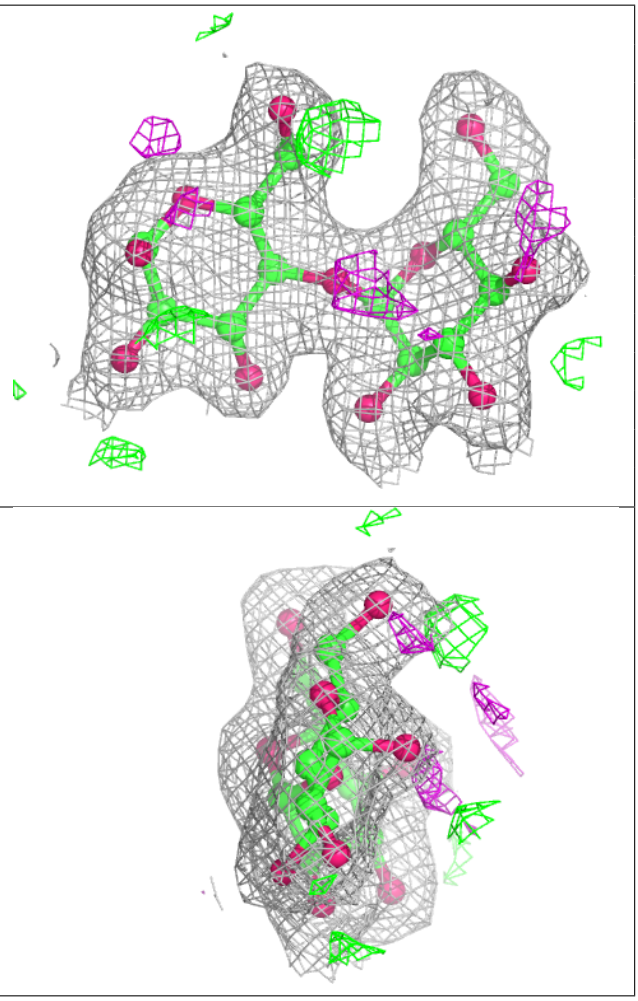
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GLC	J	1	12/12	0.94	0.17	40,48,53,54	0
3	GLC	J	2	11/12	0.94	0.13	41,43,47,47	0
3	GLC	K	1	12/12	0.95	0.19	29,32,36,36	0
3	GLC	I	1	12/12	0.96	0.14	32,34,39,41	0
3	GLC	K	2	11/12	0.96	0.13	27,29,31,34	0
3	GLC	I	2	11/12	0.97	0.13	27,30,33,33	0
3	GLC	L	1	12/12	0.97	0.17	26,29,32,34	0
3	GLC	L	2	11/12	0.97	0.18	21,25,27,30	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.

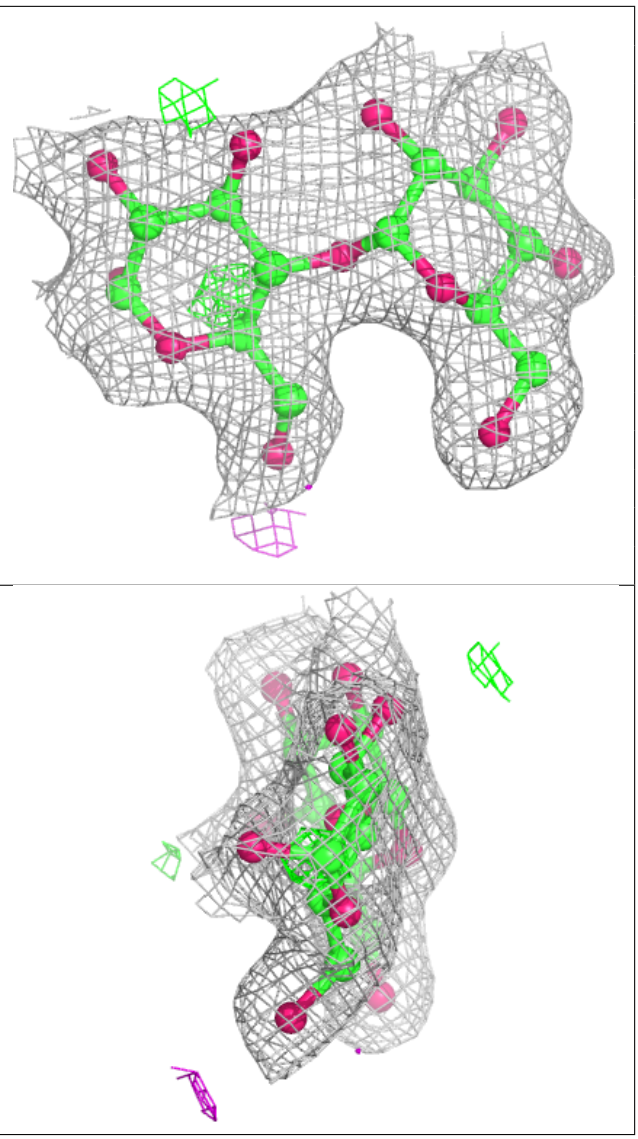
Electron density around Chain I:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)



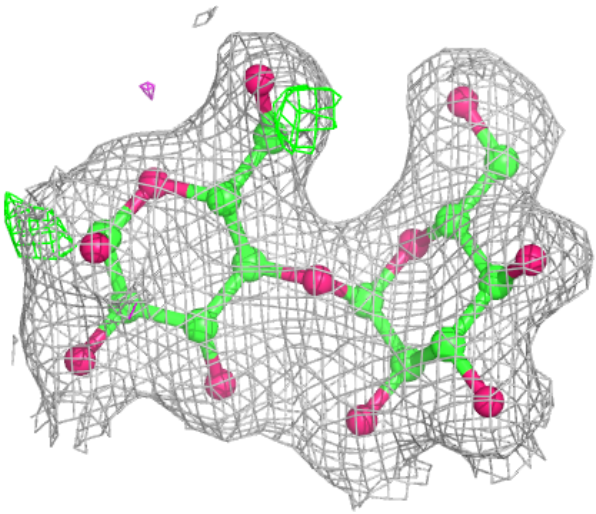
Electron density around Chain J:

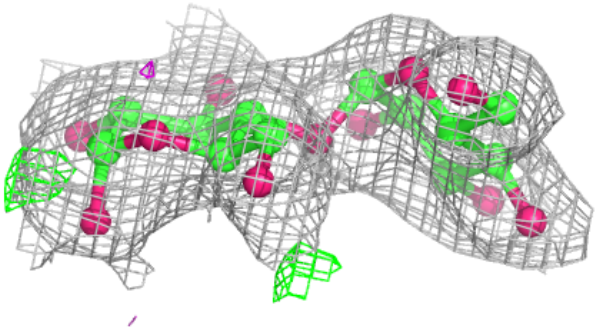
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

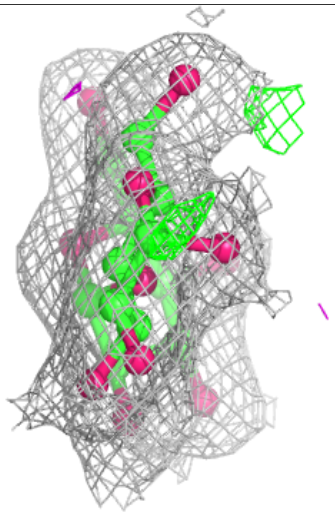


Electron density around Chain K:

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

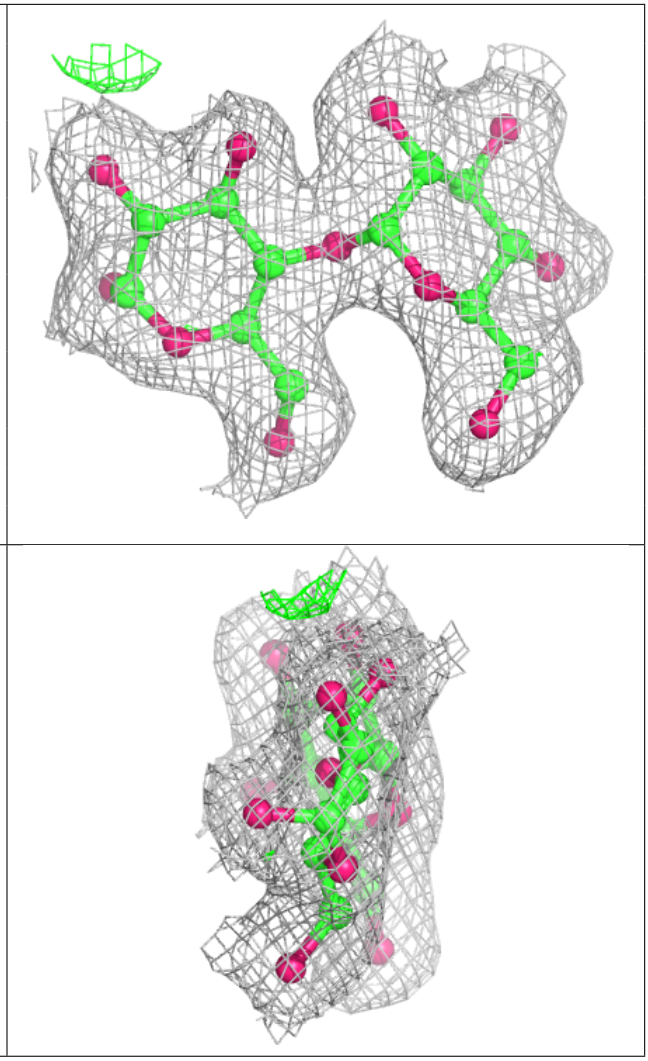






Electron density around Chain L:

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



6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	EDO	C	203	4/4	0.77	0.20	42,44,47,49	0
4	EDO	A	203	4/4	0.78	0.21	46,47,48,48	0
4	EDO	C	202	4/4	0.83	0.18	38,39,40,41	0
4	EDO	D	202	4/4	0.88	0.20	41,43,44,44	0
4	EDO	B	203	4/4	0.88	0.23	36,42,45,46	0
4	EDO	A	202	4/4	0.96	0.08	36,38,38,39	0
4	EDO	B	202	4/4	0.97	0.10	28,31,32,33	0

6.5 Other polymers (i)

There are no such residues in this entry.