



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:29 PM UTC

PDB ID : 2PI8 / pdb_00002pi8
Title : Crystal structure of E. coli MltA with bound chitohexaose
Authors : van Straaten, K.E.; Barends, T.R.M.; Dijkstra, B.W.; Thunnissen, A.M.W.H.
Deposited on : 2007-04-13
Resolution : 2.25 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

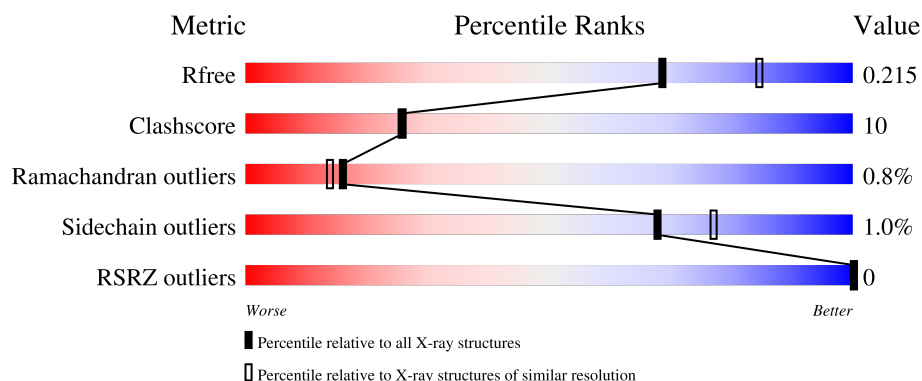
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	345	
1	B	345	
1	C	345	
1	D	345	
2	E	6	

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Mol	Chain	Length	Quality of chain
2	F	6	50% 50%
2	G	6	33% 67%
2	H	6	50% 50%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11043 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 Membrane-bound lytic murein transglycosylase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	Se	0	0	0
			2599	1645	466	480	8			
1	B	335	Total	C	N	O	Se	0	0	0
			2634	1666	472	488	8			
1	C	332	Total	C	N	O	Se	0	0	0
			2608	1651	468	481	8			
1	D	335	Total	C	N	O	Se	0	0	0
			2634	1666	472	488	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	modified residue	UNP P0A935
A	77	MSE	MET	modified residue	UNP P0A935
A	87	MSE	MET	modified residue	UNP P0A935
A	123	MSE	MET	modified residue	UNP P0A935
A	131	SER	PRO	conflict	UNP P0A935
A	154	MSE	MET	modified residue	UNP P0A935
A	159	MSE	MET	modified residue	UNP P0A935
A	206	MSE	MET	modified residue	UNP P0A935
A	208	MSE	MET	modified residue	UNP P0A935
A	272	ILE	LEU	conflict	UNP P0A935
A	293	MSE	MET	modified residue	UNP P0A935
A	308	ALA	ASP	engineered mutation	UNP P0A935
B	1	MSE	-	modified residue	UNP P0A935
B	77	MSE	MET	modified residue	UNP P0A935
B	87	MSE	MET	modified residue	UNP P0A935
B	123	MSE	MET	modified residue	UNP P0A935
B	131	SER	PRO	conflict	UNP P0A935
B	154	MSE	MET	modified residue	UNP P0A935
B	159	MSE	MET	modified residue	UNP P0A935
B	206	MSE	MET	modified residue	UNP P0A935
B	208	MSE	MET	modified residue	UNP P0A935

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Chain	Residue	Modelled	Actual	Comment	Reference
B	272	ILE	LEU	conflict	UNP P0A935
B	293	MSE	MET	modified residue	UNP P0A935
B	308	ALA	ASP	engineered mutation	UNP P0A935
C	1	MSE	-	modified residue	UNP P0A935
C	77	MSE	MET	modified residue	UNP P0A935
C	87	MSE	MET	modified residue	UNP P0A935
C	123	MSE	MET	modified residue	UNP P0A935
C	131	SER	PRO	conflict	UNP P0A935
C	154	MSE	MET	modified residue	UNP P0A935
C	159	MSE	MET	modified residue	UNP P0A935
C	206	MSE	MET	modified residue	UNP P0A935
C	208	MSE	MET	modified residue	UNP P0A935
C	272	ILE	LEU	conflict	UNP P0A935
C	293	MSE	MET	modified residue	UNP P0A935
C	308	ALA	ASP	engineered mutation	UNP P0A935
D	1	MSE	-	modified residue	UNP P0A935
D	77	MSE	MET	modified residue	UNP P0A935
D	87	MSE	MET	modified residue	UNP P0A935
D	123	MSE	MET	modified residue	UNP P0A935
D	131	SER	PRO	conflict	UNP P0A935
D	154	MSE	MET	modified residue	UNP P0A935
D	159	MSE	MET	modified residue	UNP P0A935
D	206	MSE	MET	modified residue	UNP P0A935
D	208	MSE	MET	modified residue	UNP P0A935
D	272	ILE	LEU	conflict	UNP P0A935
D	293	MSE	MET	modified residue	UNP P0A935
D	308	ALA	ASP	engineered mutation	UNP P0A935

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



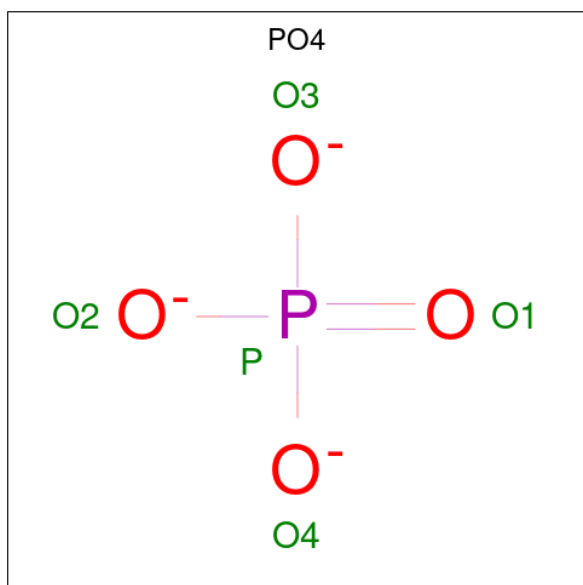
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	6	Total	C	N	O	0	0	0
			85	48	6	31			
2	F	6	Total	C	N	O	0	0	0
			85	48	6	31			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	6	Total	C	N	O	0	0	0
			85	48	6	31			
2	H	6	Total	C	N	O	0	0	0
			85	48	6	31			

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		

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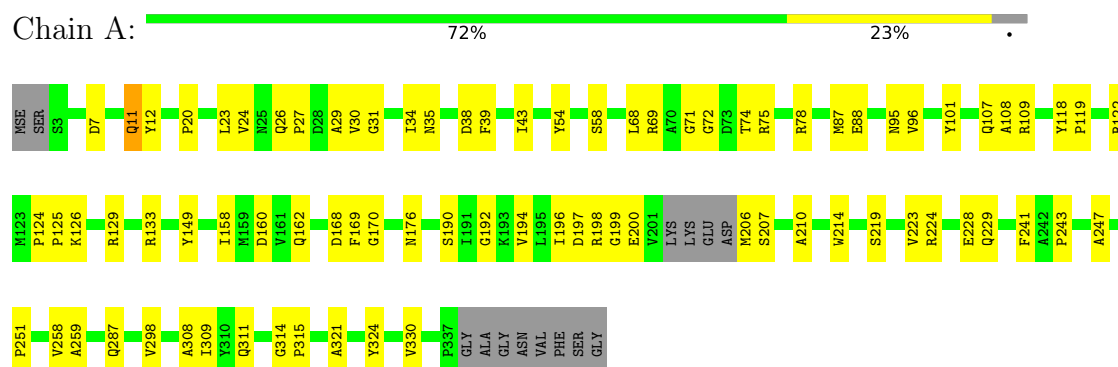
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	61	Total 61	O 61	0	0
4	C	46	Total 46	O 46	0	0
4	D	45	Total 45	O 45	0	0

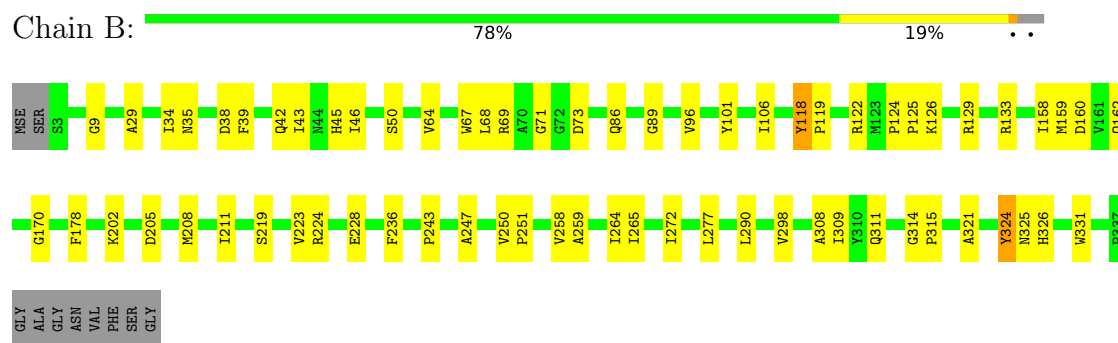
3 Residue-property plots

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

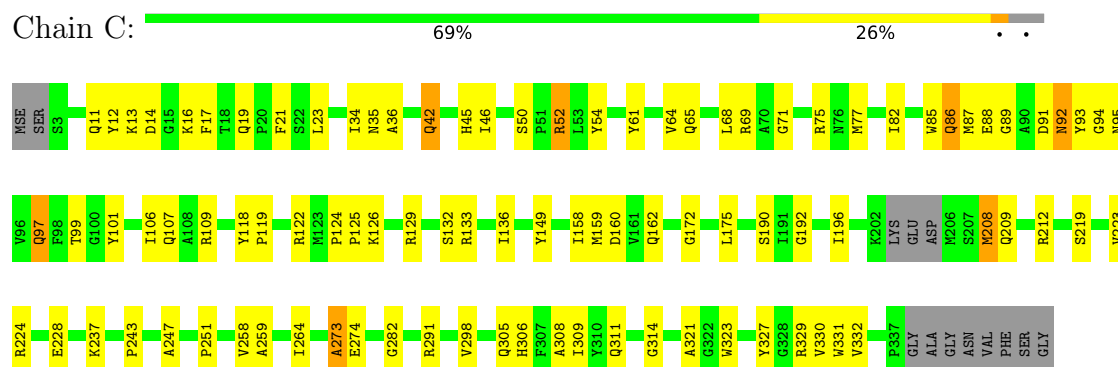
• Molecule 1: Membrane-bound lytic murein transglycosylase A



• Molecule 1: Membrane-bound lytic murein transglycosylase A

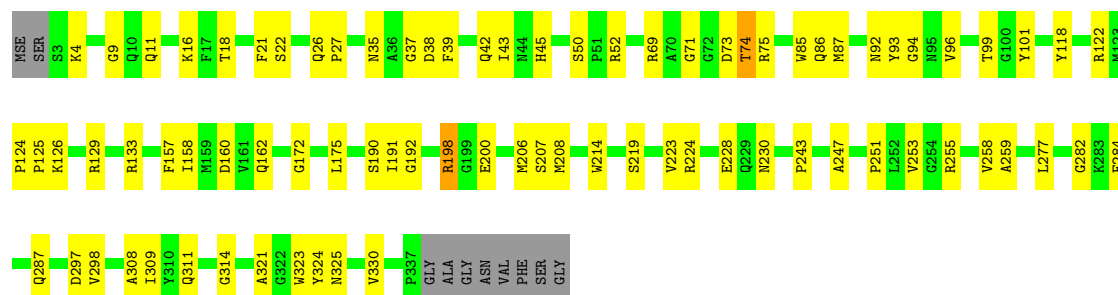


• Molecule 1: Membrane-bound lytic murein transglycosylase A



- Molecule 1: Membrane-bound lytic murein transglycosylase A

Chain D:  74% 23% ..



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

Chain E:  50% 50%



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

Chain F:  50% 50%



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

Chain G:  33% 67%



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

Chain H:  50% 50%

MAG1
MAG2
MAG3
MAG4
MAG5
MAG6

4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	91.00Å 91.00Å 187.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.00 – 2.25 36.00 – 2.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (36.00-2.25) 99.9 (36.00-2.25)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.16Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.182 , 0.226 0.173 , 0.215	Depositor DCC
R_{free} test set	6610 reflections (7.21%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 24.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.499 for -h,-k,l 0.499 for h,-h-k,-l 0.499 for -k,-h,-l	Xtriage
Reported twinning fraction	0.500 for &STRIP%twin_oper	Depositor
Outliers	0 of 91662 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	11043	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, PO4

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	0.34	1/2661 (0.0%)	0.86	6/3592 (0.2%)
1	B	0.31	0/2696	0.83	8/3637 (0.2%)
1	C	0.42	2/2670 (0.1%)	0.84	5/3603 (0.1%)
1	D	0.34	1/2696 (0.0%)	0.83	9/3637 (0.2%)
All	All	0.36	4/10723 (0.0%)	0.84	28/14469 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	19	GLN	CD-OE1	5.20	1.33	1.23
1	C	97	GLN	CD-OE1	5.13	1.33	1.23
1	A	11	GLN	CD-OE1	5.09	1.33	1.23
1	D	11	GLN	CD-OE1	5.08	1.33	1.23

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	314	GLY	CA-C-N	6.76	126.00	118.97
1	B	314	GLY	C-N-CA	6.76	126.00	118.97
1	C	118	TYR	CA-C-N	6.69	126.65	119.76
1	C	118	TYR	C-N-CA	6.69	126.65	119.76
1	C	314	GLY	CA-C-N	6.68	125.92	118.97
1	C	314	GLY	C-N-CA	6.68	125.92	118.97
1	B	118	TYR	CA-C-N	6.66	126.62	119.76
1	B	118	TYR	C-N-CA	6.66	126.62	119.76
1	A	314	GLY	CA-C-N	6.62	125.86	118.97
1	A	314	GLY	C-N-CA	6.62	125.86	118.97
1	D	314	GLY	CA-C-N	6.53	125.77	118.97
1	D	314	GLY	C-N-CA	6.53	125.77	118.97
1	A	118	TYR	CA-C-N	6.47	126.43	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	118	TYR	C-N-CA	6.47	126.43	119.76
1	D	118	TYR	CA-C-N	6.40	126.35	119.76
1	D	118	TYR	C-N-CA	6.40	126.35	119.76
1	A	324	TYR	N-CA-C	6.19	118.53	108.63
1	B	325	ASN	N-CA-C	-6.03	103.63	111.71
1	A	12	TYR	N-CA-C	-5.82	105.80	114.64
1	B	50	SER	CA-C-N	5.66	125.77	119.32
1	B	50	SER	C-N-CA	5.66	125.77	119.32
1	D	26	GLN	CA-C-N	5.65	126.90	119.84
1	D	26	GLN	C-N-CA	5.65	126.90	119.84
1	D	325	ASN	N-CA-C	-5.60	104.69	112.03
1	D	50	SER	CA-C-N	5.30	124.96	119.56
1	D	50	SER	C-N-CA	5.30	124.96	119.56
1	C	273	ALA	N-CA-C	5.19	117.07	109.24
1	B	324	TYR	N-CA-C	5.18	121.83	110.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	2599	0	2497	51	0
1	B	2634	0	2534	47	0
1	C	2608	0	2510	66	0
1	D	2634	0	2534	54	0
2	E	85	0	75	2	0
2	F	85	0	75	3	0
2	G	85	0	75	7	0
2	H	85	0	75	7	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
4	A	46	0	0	0	0
4	B	61	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	46	0	0	0	0
4	D	45	0	0	0	0
All	All	11043	0	10375	202	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 (202) 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:200:GLU:HG2	1:A:214:TRP:HE1	1.27	0.98
1:D:190:SER:HB3	2:H:1:NAG:H81	1.47	0.95
1:A:259:ALA:HB3	1:A:308:ALA:HB3	1.62	0.80
1:D:259:ALA:HB3	1:D:308:ALA:HB3	1.62	0.79
1:A:27:PRO:HG2	1:A:74:THR:HB	1.65	0.79
1:A:96:VAL:HG21	1:A:330:VAL:HG12	1.65	0.78
1:B:259:ALA:HB3	1:B:308:ALA:HB3	1.66	0.78
1:C:259:ALA:HB3	1:C:308:ALA:HB3	1.66	0.78
1:C:208:MSE:SE	2:G:1:NAG:H3	2.36	0.76
1:B:101:TYR:CE2	2:F:3:NAG:H62	2.24	0.73
1:A:29:ALA:HB3	1:A:74:THR:HG21	1.70	0.72
1:D:157:PHE:CZ	1:D:191:ILE:HG13	2.26	0.71
1:B:202:LYS:HB2	1:B:205:ASP:HB2	1.74	0.70
1:C:23:LEU:H	1:C:23:LEU:HD23	1.59	0.68
1:A:190:SER:OG	2:E:1:NAG:H81	1.96	0.65
1:D:27:PRO:HG2	1:D:74:THR:HB	1.77	0.65
1:D:96:VAL:HG21	1:D:330:VAL:HG12	1.77	0.65
1:C:86:GLN:HG2	1:C:331:TRP:NE1	2.12	0.64
1:C:101:TYR:CE2	2:G:3:NAG:H62	2.33	0.63
1:D:42:GLN:HG2	1:D:253:VAL:HB	1.80	0.63
1:A:258:VAL:HG12	1:A:309:ILE:HA	1.80	0.63
1:C:258:VAL:HG12	1:C:309:ILE:HA	1.80	0.62
1:C:89:GLY:HA2	1:C:305:GLN:HB3	1.82	0.61
1:B:89:GLY:HA2	1:B:96:VAL:HG22	1.81	0.61
1:C:52:ARG:H	1:C:52:ARG:HD3	1.65	0.61
1:D:258:VAL:HG12	1:D:309:ILE:HA	1.82	0.61
1:C:237:LYS:HD3	2:G:6:NAG:H83	1.83	0.60
1:A:30:VAL:HG13	1:A:31:GLY:H	1.67	0.59
1:C:50:SER:HA	1:C:52:ARG:HE	1.68	0.58
1:B:277:LEU:HD11	1:D:69:ARG:HD2	1.85	0.58
1:C:158:ILE:O	1:C:162:GLN:HG2	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:274:GLU:HG3	1:C:329:ARG:HH21	1.69	0.57
1:C:107:GLN:HG2	1:C:109:ARG:HH12	1.68	0.57
1:B:29:ALA:HB1	1:B:290:LEU:HD22	1.86	0.56
1:B:69:ARG:HH12	1:D:255:ARG:HH12	1.51	0.56
1:B:35:ASN:HB2	1:B:38:ASP:HB2	1.87	0.56
1:D:35:ASN:HB2	1:D:38:ASP:HB2	1.87	0.56
1:A:11:GLN:H	1:A:95:ASN:HD21	1.55	0.55
1:B:158:ILE:O	1:B:162:GLN:HG2	2.07	0.55
1:D:45:HIS:HB3	1:D:251:PRO:HG2	1.89	0.54
1:C:158:ILE:HD12	1:C:212:ARG:NH2	2.23	0.54
1:C:46:ILE:HG22	1:C:54:TYR:HB2	1.89	0.54
1:C:75:ARG:HG2	1:C:75:ARG:HH11	1.73	0.54
1:B:258:VAL:HG12	1:B:309:ILE:HA	1.90	0.53
1:C:45:HIS:HB3	1:C:251:PRO:HG2	1.90	0.53
1:B:219:SER:O	1:B:223:VAL:HG23	2.09	0.53
1:D:158:ILE:O	1:D:162:GLN:HG2	2.08	0.53
1:C:77:MSE:HE2	1:C:82:ILE:HG21	1.91	0.52
1:C:95:ASN:HB3	1:C:327:TYR:HB3	1.91	0.52
1:C:99:THR:OG1	2:G:1:NAG:H61	2.10	0.52
1:B:106:ILE:HD12	1:B:159:MSE:SE	2.60	0.52
1:C:126:LYS:HD2	1:C:126:LYS:N	2.25	0.52
1:D:124:PRO:HG2	1:D:129:ARG:HG3	1.91	0.51
1:B:67:TRP:HE1	1:B:73:ASP:C	2.19	0.51
1:C:101:TYR:CZ	2:G:3:NAG:H62	2.45	0.51
1:D:190:SER:CB	2:H:1:NAG:H81	2.31	0.51
1:A:158:ILE:O	1:A:162:GLN:HG2	2.10	0.51
1:C:42:GLN:HE21	1:C:42:GLN:HA	1.76	0.51
1:B:86:GLN:HB2	1:B:331:TRP:CD1	2.45	0.51
1:D:192:GLY:H	2:H:1:NAG:H83	1.76	0.51
1:A:311:GLN:HG3	1:A:321:ALA:HB2	1.94	0.50
1:B:272:ILE:CG2	1:B:331:TRP:HB2	2.41	0.50
1:A:126:LYS:HD2	1:A:126:LYS:N	2.26	0.50
1:D:157:PHE:HZ	1:D:191:ILE:HG13	1.71	0.50
1:D:219:SER:O	1:D:223:VAL:HG23	2.11	0.50
1:C:88:GLU:HB2	1:C:92:ASN:HA	1.94	0.49
1:C:162:GLN:HE22	2:G:2:NAG:C7	2.25	0.49
1:A:26:GLN:HG2	1:A:75:ARG:O	2.11	0.49
1:A:287:GLN:OE1	1:C:69:ARG:HB3	2.12	0.49
1:C:219:SER:O	1:C:223:VAL:HG23	2.12	0.49
1:A:39:PHE:O	1:A:43:ILE:HG12	2.12	0.49
1:D:4:LYS:H	1:D:4:LYS:HD2	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:200:GLU:HB3	1:D:214:TRP:HE1	1.76	0.49
1:A:30:VAL:HG21	1:C:71:GLY:HA2	1.93	0.49
1:D:18:THR:HG23	1:D:92:ASN:HB3	1.94	0.49
1:A:109:ARG:HH21	1:B:315:PRO:HG2	1.77	0.49
1:A:200:GLU:HG2	1:A:214:TRP:NE1	2.10	0.49
1:C:124:PRO:HG2	1:C:129:ARG:HG3	1.93	0.49
1:C:106:ILE:HD12	1:C:159:MSE:SE	2.63	0.49
1:D:198:ARG:HD2	1:D:198:ARG:N	2.28	0.49
1:A:87:MSE:SE	1:A:330:VAL:HG13	2.64	0.48
1:B:224:ARG:O	1:B:228:GLU:HG3	2.13	0.48
1:B:126:LYS:HD2	1:B:126:LYS:N	2.28	0.48
1:D:133:ARG:HH22	1:D:160:ASP:CG	2.21	0.48
1:B:272:ILE:HG23	1:B:331:TRP:HB2	1.96	0.48
1:D:247:ALA:HB2	1:D:298:VAL:HG22	1.96	0.48
1:A:96:VAL:HG21	1:A:330:VAL:CG1	2.40	0.48
1:A:315:PRO:HD3	1:B:170:GLY:O	2.14	0.48
1:D:21:PHE:HB3	1:D:85:TRP:CE3	2.49	0.47
1:A:30:VAL:HG13	1:A:31:GLY:N	2.27	0.47
1:B:45:HIS:HB3	1:B:251:PRO:HG2	1.96	0.47
1:C:61:TYR:HA	1:C:64:VAL:HG22	1.95	0.47
1:C:175:LEU:HD22	1:D:172:GLY:HA2	1.96	0.47
1:A:124:PRO:HB2	1:A:125:PRO:HD2	1.97	0.47
1:A:219:SER:O	1:A:223:VAL:HG23	2.15	0.47
1:C:133:ARG:HH22	1:C:160:ASP:CG	2.23	0.47
1:C:190:SER:OG	2:G:1:NAG:H83	2.15	0.47
1:D:52:ARG:HG2	1:D:52:ARG:HH11	1.80	0.47
1:B:39:PHE:O	1:B:43:ILE:HG12	2.15	0.47
1:C:224:ARG:O	1:C:228:GLU:HG3	2.15	0.47
1:A:243:PRO:HB2	1:A:251:PRO:HB3	1.97	0.47
1:B:208:MSE:SE	1:B:211:ILE:HD12	2.65	0.47
1:A:247:ALA:HB2	1:A:298:VAL:HG22	1.97	0.46
1:D:277:LEU:HD12	1:D:287:GLN:HG3	1.98	0.46
1:A:11:GLN:H	1:A:95:ASN:ND2	2.13	0.46
1:D:282:GLY:HA3	1:D:323:TRP:CD2	2.50	0.46
1:D:73:ASP:C	1:D:75:ARG:H	2.23	0.46
1:A:194:VAL:HG11	1:A:229:GLN:HG3	1.97	0.46
1:B:69:ARG:NH1	1:D:255:ARG:HH12	2.14	0.46
1:C:192:GLY:O	1:C:196:ILE:HG13	2.16	0.46
1:D:101:TYR:CE2	2:H:3:NAG:H62	2.50	0.46
1:B:124:PRO:HG2	1:B:129:ARG:HG3	1.97	0.46
1:C:36:ALA:HB1	1:C:68:LEU:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:21:PHE:HB3	1:D:85:TRP:HE3	1.81	0.46
1:D:126:LYS:HD2	1:D:126:LYS:N	2.30	0.46
1:C:172:GLY:HA2	1:D:175:LEU:HD22	1.96	0.46
1:B:122:ARG:O	1:B:124:PRO:HD3	2.15	0.46
1:A:107:GLN:HA	1:A:168:ASP:O	2.16	0.46
1:C:122:ARG:O	1:C:124:PRO:HD3	2.16	0.46
1:D:122:ARG:O	1:D:124:PRO:HD3	2.16	0.46
1:A:122:ARG:O	1:A:124:PRO:HD3	2.15	0.45
1:A:124:PRO:HG2	1:A:129:ARG:HG3	1.97	0.45
1:A:170:GLY:O	1:B:315:PRO:HD3	2.16	0.45
1:D:206:MSE:HE3	1:D:207:SER:H	1.81	0.45
1:B:64:VAL:O	1:B:68:LEU:HD13	2.17	0.45
1:D:224:ARG:O	1:D:228:GLU:HG3	2.17	0.45
1:B:133:ARG:HH22	1:B:160:ASP:CG	2.24	0.45
1:D:243:PRO:HB2	1:D:251:PRO:HB3	1.98	0.45
1:B:247:ALA:HB2	1:B:298:VAL:HG22	1.98	0.45
1:A:109:ARG:HE	1:B:315:PRO:HG2	1.81	0.45
1:C:247:ALA:HB2	1:C:298:VAL:HG22	1.99	0.45
1:C:42:GLN:HE21	1:C:42:GLN:CA	2.29	0.44
1:A:54:TYR:O	1:A:58:SER:HB2	2.17	0.44
1:B:69:ARG:C	1:B:71:GLY:H	2.26	0.44
1:D:22:SER:HB2	1:D:86:GLN:HB3	1.98	0.44
1:A:78:ARG:HD2	1:A:78:ARG:N	2.32	0.44
1:B:124:PRO:HB2	1:B:125:PRO:HD2	2.00	0.44
1:B:69:ARG:HH12	1:D:255:ARG:NH1	2.15	0.44
1:B:243:PRO:HB2	1:B:251:PRO:HB3	2.00	0.44
1:C:12:TYR:C	1:C:14:ASP:H	2.24	0.44
1:C:311:GLN:HG3	1:C:321:ALA:HB2	2.00	0.44
1:C:282:GLY:HA3	1:C:323:TRP:CD2	2.53	0.44
1:C:243:PRO:HB2	1:C:251:PRO:HB3	1.99	0.44
1:A:133:ARG:HH22	1:A:160:ASP:CG	2.26	0.43
1:D:16:LYS:HE2	1:D:93:TYR:CE2	2.53	0.43
1:D:311:GLN:HG3	1:D:321:ALA:HB2	1.99	0.43
1:B:208:MSE:HE3	2:F:2:NAG:O7	2.18	0.43
1:D:87:MSE:SE	1:D:330:VAL:HG13	2.68	0.43
1:C:88:GLU:HG3	1:C:89:GLY:N	2.32	0.43
1:A:108:ALA:O	1:A:169:PHE:HA	2.18	0.43
1:C:11:GLN:HE22	1:C:329:ARG:NH1	2.17	0.43
1:A:34:ILE:HG13	1:A:35:ASN:N	2.32	0.43
1:A:206:MSE:SE	1:A:206:MSE:C	3.12	0.43
1:C:11:GLN:NE2	1:C:329:ARG:NH1	2.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:191:ILE:HD11	1:D:230:ASN:HB2	2.01	0.43
1:A:71:GLY:O	1:C:34:ILE:HG21	2.18	0.43
1:A:101:TYR:CE2	2:E:3:NAG:H62	2.54	0.43
1:B:264:ILE:HB	1:B:265:ILE:HD12	2.01	0.43
1:B:311:GLN:HG3	1:B:321:ALA:HB2	2.01	0.43
1:C:23:LEU:H	1:C:23:LEU:CD2	2.30	0.43
1:A:176:ASN:HA	1:A:241:PHE:H	1.84	0.42
1:A:224:ARG:O	1:A:228:GLU:HG3	2.19	0.42
1:C:42:GLN:O	1:C:46:ILE:HG13	2.18	0.42
1:C:35:ASN:HB3	1:C:291:ARG:NH1	2.34	0.42
1:B:46:ILE:HG12	1:B:250:VAL:HG11	2.01	0.42
1:C:21:PHE:HZ	1:C:264:ILE:HG22	1.84	0.42
1:A:20:PRO:HA	1:A:88:GLU:HG2	2.01	0.42
1:A:192:GLY:O	1:A:196:ILE:HG13	2.19	0.42
1:D:124:PRO:HB2	1:D:125:PRO:HD2	2.01	0.42
1:A:68:LEU:HD12	1:A:69:ARG:N	2.33	0.42
1:B:101:TYR:CZ	2:F:3:NAG:H62	2.54	0.42
1:C:17:PHE:CZ	1:C:86:GLN:HB3	2.54	0.42
1:C:85:TRP:HB2	1:C:332:VAL:HB	2.01	0.42
1:C:87:MSE:O	1:C:94:GLY:HA2	2.19	0.42
1:C:273:ALA:HB2	1:C:330:VAL:HG23	2.00	0.42
1:C:88:GLU:HG3	1:C:89:GLY:H	1.85	0.42
1:C:132:SER:O	1:C:136:ILE:HG13	2.20	0.42
1:B:34:ILE:HB	1:D:37:GLY:H	1.85	0.42
1:C:175:LEU:CD2	1:D:172:GLY:HA2	2.50	0.42
1:A:11:GLN:N	1:A:95:ASN:HD21	2.16	0.42
1:A:207:SER:H	1:A:210:ALA:HB3	1.84	0.41
1:B:42:GLN:O	1:B:46:ILE:HG13	2.20	0.41
1:B:118:TYR:HA	1:B:119:PRO:HD3	1.87	0.41
1:C:65:GLN:O	1:C:69:ARG:HD3	2.20	0.41
1:D:9:GLY:HA2	1:D:284:PHE:CE2	2.56	0.41
1:C:16:LYS:HA	1:C:92:ASN:O	2.20	0.41
1:C:119:PRO:HG3	1:C:149:TYR:CE2	2.55	0.41
1:D:39:PHE:O	1:D:43:ILE:HG12	2.20	0.41
1:A:119:PRO:HG3	1:A:149:TYR:CE2	2.56	0.41
1:C:124:PRO:HB2	1:C:125:PRO:HD2	2.02	0.41
1:A:109:ARG:NH2	1:B:315:PRO:HG2	2.36	0.41
1:C:91:ASP:C	1:C:93:TYR:H	2.29	0.41
1:D:99:THR:OG1	2:H:1:NAG:H62	2.20	0.41
1:D:297:ASP:OD1	2:H:4:NAG:H3	2.20	0.41
1:D:99:THR:HA	1:D:324:TYR:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:ASN:HA	4:B:1142:HOH:O	2.20	0.40
1:B:326:HIS:ND1	1:B:326:HIS:N	2.69	0.40
1:C:97:GLN:HG3	1:C:306:HIS:CE1	2.56	0.40
1:D:101:TYR:CZ	2:H:3:NAG:H62	2.55	0.40
1:A:35:ASN:HB2	1:A:38:ASP:HB2	2.04	0.40
1:B:178:PHE:HB3	1:B:236:PHE:HB3	2.03	0.40
1:D:4:LYS:HD2	1:D:4:LYS:N	2.36	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	327/345 (95%)	302 (92%)	22 (7%)	3 (1%)	14	12
1	B	333/345 (96%)	315 (95%)	16 (5%)	2 (1%)	21	21
1	C	328/345 (95%)	311 (95%)	15 (5%)	2 (1%)	21	21
1	D	333/345 (96%)	310 (93%)	20 (6%)	3 (1%)	14	12
All	All	1321/1380 (96%)	1238 (94%)	73 (6%)	10 (1%)	16	14

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	GLY
1	D	71	GLY
1	B	9	GLY
1	C	13	LYS
1	C	92	ASN
1	D	74	THR
1	B	324	TYR
1	A	72	GLY

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Mol	Chain	Res	Type
1	A	24	VAL
1	D	94	GLY

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	266/267 (100%)	262 (98%)	4 (2%)	57	68
1	B	270/267 (101%)	270 (100%)	0	100	100
1	C	267/267 (100%)	262 (98%)	5 (2%)	50	61
1	D	270/267 (101%)	268 (99%)	2 (1%)	76	82
All	All	1073/1068 (100%)	1062 (99%)	11 (1%)	68	77

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	23	LEU
1	A	197	ASP
1	A	198	ARG
1	C	42	GLN
1	C	52	ARG
1	C	86	GLN
1	C	208	MSE
1	C	209	GLN
1	D	198	ARG
1	D	208	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	92	ASN
1	A	110	HIS

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Mol	Chain	Res	Type
1	A	162	GLN
1	A	281	ASN
1	A	305	GLN
1	B	11	GLN
1	B	57	GLN
1	B	92	ASN
1	B	110	HIS
1	C	76	ASN
1	C	281	ASN
1	C	287	GLN
1	D	76	ASN
1	D	79	GLN
1	D	162	GLN
1	D	280	ASN
1	D	287	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	$\# Z > 2$
2	NAG	E	1	2	15,15,15	0.38	0	21,21,21	0.56	0
2	NAG	E	2	2	14,14,15	0.47	0	17,19,21	0.65	0
2	NAG	E	3	2	14,14,15	0.40	0	17,19,21	0.68	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	E	4	2	14,14,15	0.40	0	17,19,21	0.69	1 (5%)
2	NAG	E	5	2	14,14,15	0.43	0	17,19,21	0.66	0
2	NAG	E	6	2	14,14,15	0.43	0	17,19,21	0.62	0
2	NAG	F	1	2	15,15,15	0.40	0	21,21,21	0.62	0
2	NAG	F	2	2	14,14,15	0.46	0	17,19,21	0.74	0
2	NAG	F	3	2	14,14,15	0.42	0	17,19,21	0.69	0
2	NAG	F	4	2	14,14,15	0.40	0	17,19,21	0.67	1 (5%)
2	NAG	F	5	2	14,14,15	0.44	0	17,19,21	0.68	0
2	NAG	F	6	2	14,14,15	0.42	0	17,19,21	0.59	0
2	NAG	G	1	2	15,15,15	0.39	0	21,21,21	0.58	0
2	NAG	G	2	2	14,14,15	0.44	0	17,19,21	0.69	0
2	NAG	G	3	2	14,14,15	0.42	0	17,19,21	0.69	0
2	NAG	G	4	2	14,14,15	0.41	0	17,19,21	0.66	0
2	NAG	G	5	2	14,14,15	0.41	0	17,19,21	0.64	0
2	NAG	G	6	2	14,14,15	0.40	0	17,19,21	0.64	0
2	NAG	H	1	2	15,15,15	0.39	0	21,21,21	0.65	0
2	NAG	H	2	2	14,14,15	0.46	0	17,19,21	0.67	0
2	NAG	H	3	2	14,14,15	0.41	0	17,19,21	0.69	0
2	NAG	H	4	2	14,14,15	0.38	0	17,19,21	0.66	0
2	NAG	H	5	2	14,14,15	0.39	0	17,19,21	0.65	0
2	NAG	H	6	2	14,14,15	0.43	0	17,19,21	0.63	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	1	2	-	2/6/26/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	E	3	2	-	1/6/23/26	0/1/1/1
2	NAG	E	4	2	-	0/6/23/26	0/1/1/1
2	NAG	E	5	2	-	0/6/23/26	0/1/1/1
2	NAG	E	6	2	-	0/6/23/26	0/1/1/1
2	NAG	F	1	2	-	1/6/26/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	NAG	F	3	2	-	2/6/23/26	0/1/1/1
2	NAG	F	4	2	-	0/6/23/26	0/1/1/1
2	NAG	F	5	2	-	0/6/23/26	0/1/1/1
2	NAG	F	6	2	-	2/6/23/26	0/1/1/1
2	NAG	G	1	2	-	0/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
2	NAG	G	3	2	-	2/6/23/26	0/1/1/1
2	NAG	G	4	2	-	0/6/23/26	0/1/1/1
2	NAG	G	5	2	-	0/6/23/26	0/1/1/1
2	NAG	G	6	2	-	0/6/23/26	0/1/1/1
2	NAG	H	1	2	-	2/6/26/26	0/1/1/1
2	NAG	H	2	2	-	2/6/23/26	0/1/1/1
2	NAG	H	3	2	-	2/6/23/26	0/1/1/1
2	NAG	H	4	2	-	0/6/23/26	0/1/1/1
2	NAG	H	5	2	-	0/6/23/26	0/1/1/1
2	NAG	H	6	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	4	NAG	C2-N2-C7	-2.11	120.08	122.90
2	E	4	NAG	C2-N2-C7	-2.11	120.08	122.90

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	1	NAG	O5-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
2	H	1	NAG	C4-C5-C6-O6
2	F	3	NAG	C4-C5-C6-O6
2	F	3	NAG	O5-C5-C6-O6
2	G	3	NAG	C4-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
2	H	3	NAG	C4-C5-C6-O6
2	G	3	NAG	O5-C5-C6-O6
2	H	3	NAG	O5-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
2	F	6	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6

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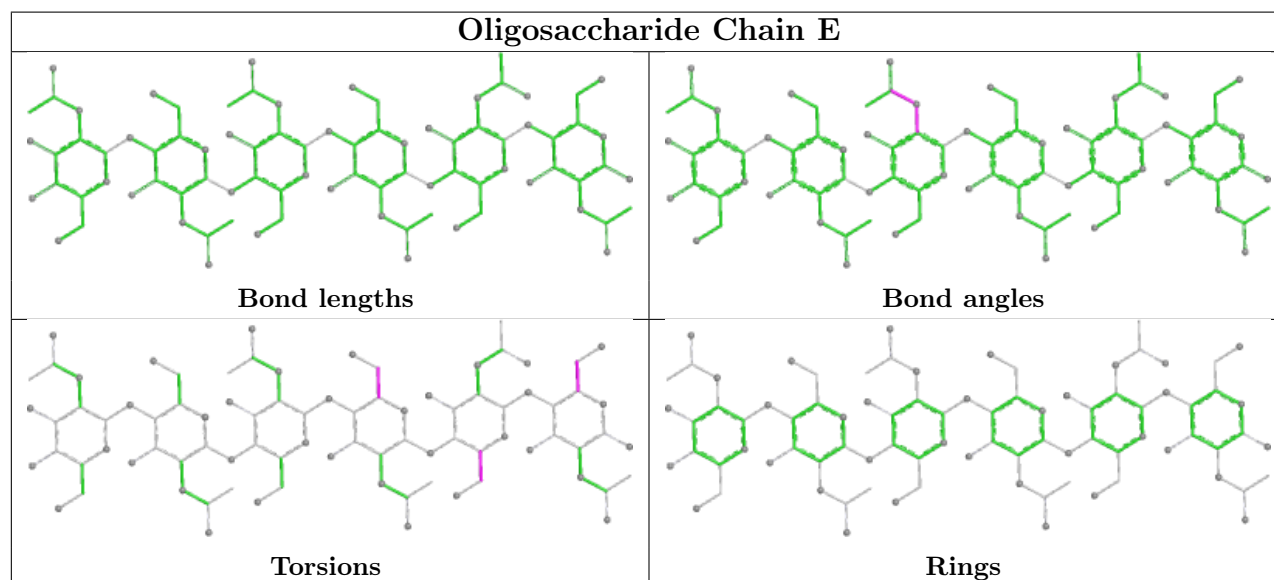
Mol	Chain	Res	Type	Atoms
2	F	1	NAG	O5-C5-C6-O6
2	E	3	NAG	C4-C5-C6-O6
2	F	6	NAG	O5-C5-C6-O6

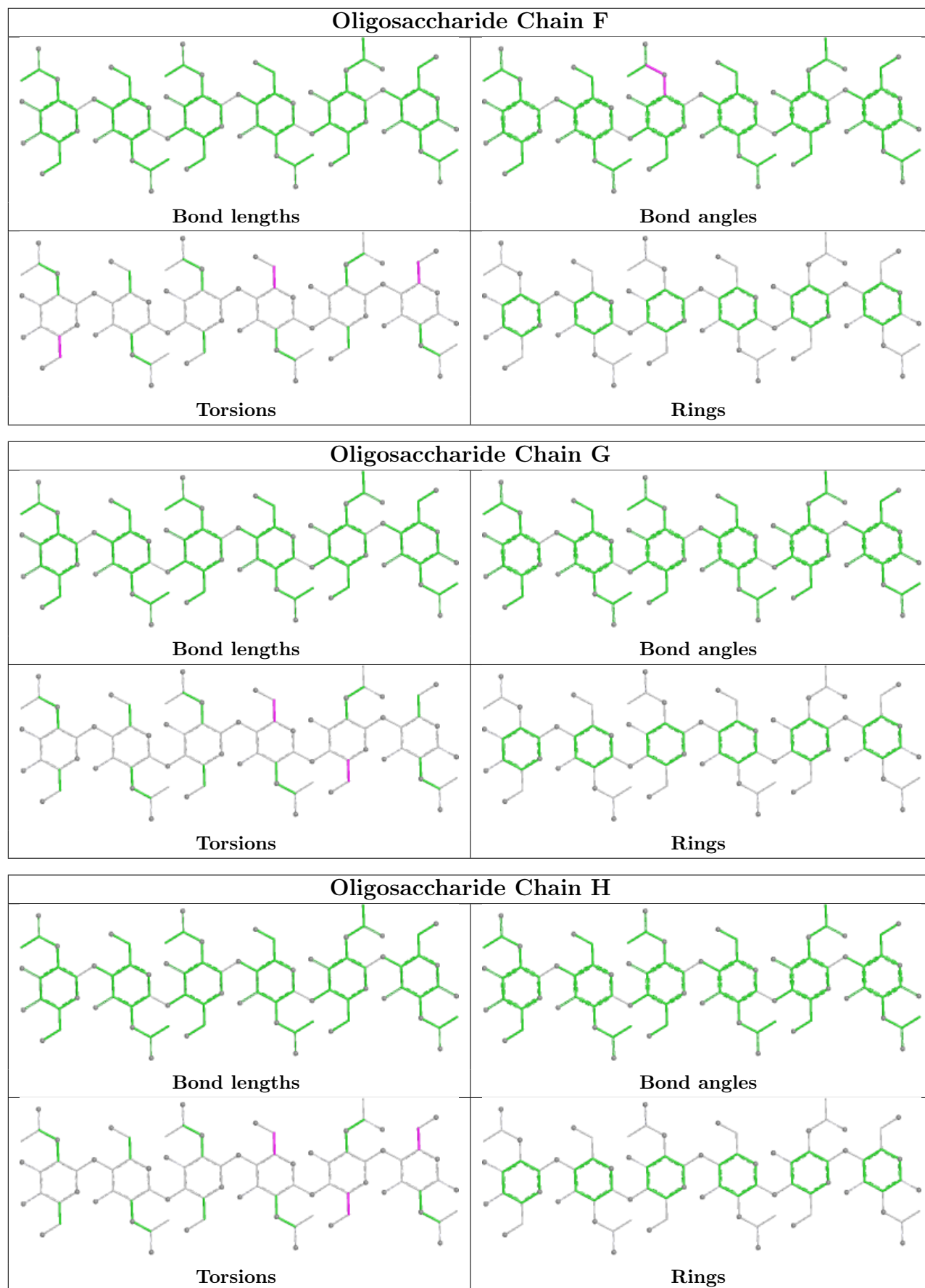
There are no ring outliers.

11 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	3	NAG	2	0
2	E	1	NAG	1	0
2	F	2	NAG	1	0
2	H	3	NAG	2	0
2	G	1	NAG	3	0
2	H	1	NAG	4	0
2	G	2	NAG	1	0
2	G	6	NAG	1	0
2	G	3	NAG	2	0
2	E	3	NAG	1	0
2	H	4	NAG	1	0

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





5.6 Ligand geometry

6 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PO4	C	701	-	4,4,4	1.72	0	6,6,6	0.46	0
3	PO4	C	697	-	4,4,4	1.74	0	6,6,6	0.46	0
3	PO4	B	696	-	4,4,4	1.73	0	6,6,6	0.46	0
3	PO4	A	699	-	4,4,4	1.78	2 (50%)	6,6,6	0.46	0
3	PO4	D	698	-	4,4,4	1.76	0	6,6,6	0.47	0
3	PO4	D	700	-	4,4,4	1.74	0	6,6,6	0.46	0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	699	PO4	P-O4	-2.04	1.48	1.54
3	A	699	PO4	P-O2	-2.03	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	323/345 (93%)	-1.49	0 100 100	20, 37, 54, 72	0
1	B	327/345 (94%)	-1.53	0 100 100	16, 36, 52, 66	0
1	C	324/345 (93%)	-1.53	0 100 100	17, 35, 52, 63	0
1	D	327/345 (94%)	-1.50	0 100 100	17, 38, 57, 65	0
All	All	1301/1380 (94%)	-1.51	0 100 100	16, 36, 54, 72	0

There are no RSRZ outliers to report.

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.

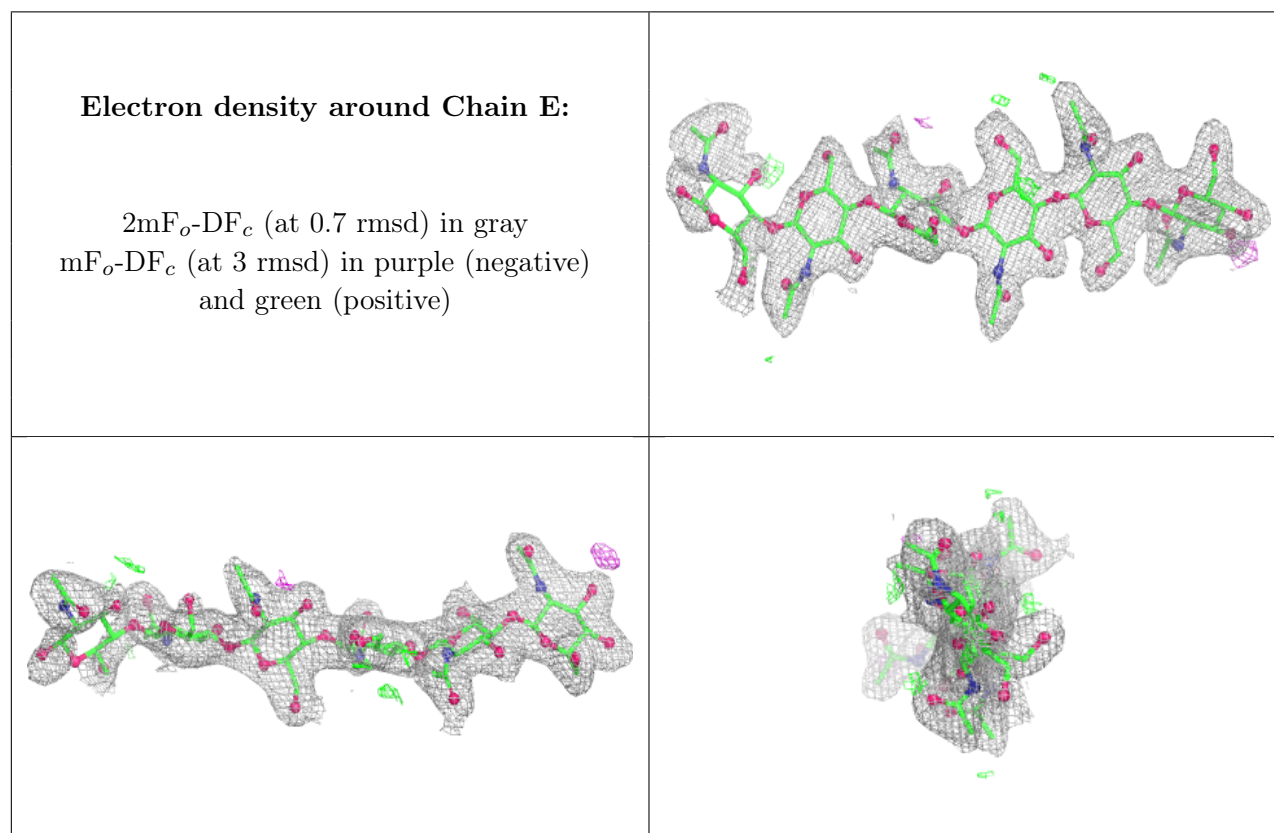
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	E	1	15/15	0.98	0.05	54,61,62,62	0
2	NAG	E	2	14/15	0.99	0.03	39,45,48,49	0
2	NAG	E	5	14/15	0.99	0.02	22,25,27,29	0
2	NAG	E	6	14/15	0.99	0.03	31,33,35,36	0
2	NAG	G	1	15/15	0.99	0.04	50,55,57,58	0
2	NAG	G	2	14/15	0.99	0.03	37,41,44,46	0
2	NAG	G	3	14/15	0.99	0.03	31,37,39,40	0
2	NAG	H	1	15/15	0.99	0.04	38,44,47,48	0
2	NAG	H	5	14/15	0.99	0.02	23,25,27,28	0

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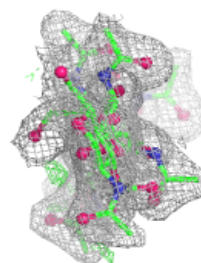
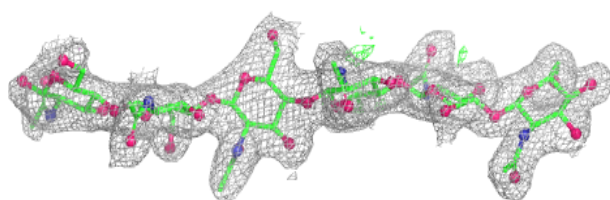
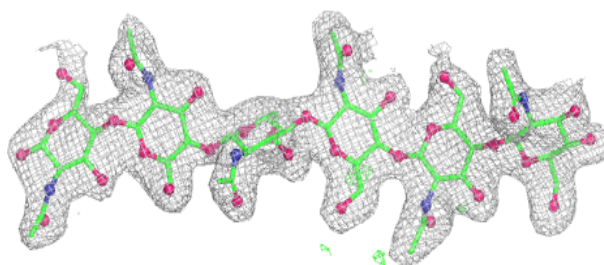
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	H	6	14/15	0.99	0.03	24,31,35,37	0
2	NAG	F	5	14/15	1.00	0.02	19,21,25,26	0
2	NAG	F	6	14/15	1.00	0.02	26,28,30,32	0
2	NAG	E	3	14/15	1.00	0.02	27,30,35,38	0
2	NAG	E	4	14/15	1.00	0.02	24,26,27,27	0
2	NAG	F	1	15/15	1.00	0.03	42,49,51,53	0
2	NAG	G	4	14/15	1.00	0.02	24,29,32,34	0
2	NAG	G	5	14/15	1.00	0.02	22,26,31,31	0
2	NAG	G	6	14/15	1.00	0.02	23,29,32,36	0
2	NAG	F	2	14/15	1.00	0.02	30,36,39,39	0
2	NAG	H	2	14/15	1.00	0.02	32,35,38,40	0
2	NAG	H	3	14/15	1.00	0.03	24,28,31,33	0
2	NAG	H	4	14/15	1.00	0.02	19,22,24,26	0
2	NAG	F	3	14/15	1.00	0.02	23,26,29,32	0
2	NAG	F	4	14/15	1.00	0.02	16,22,22,24	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.

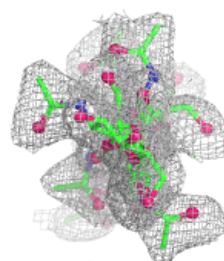
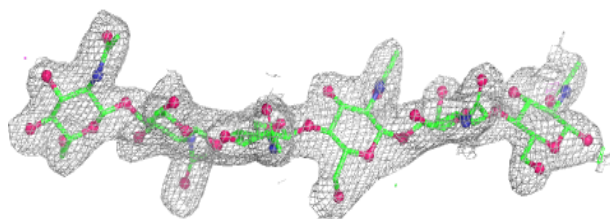
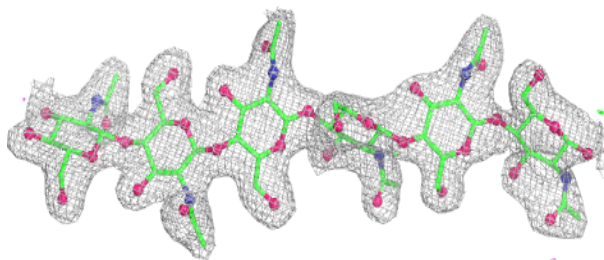


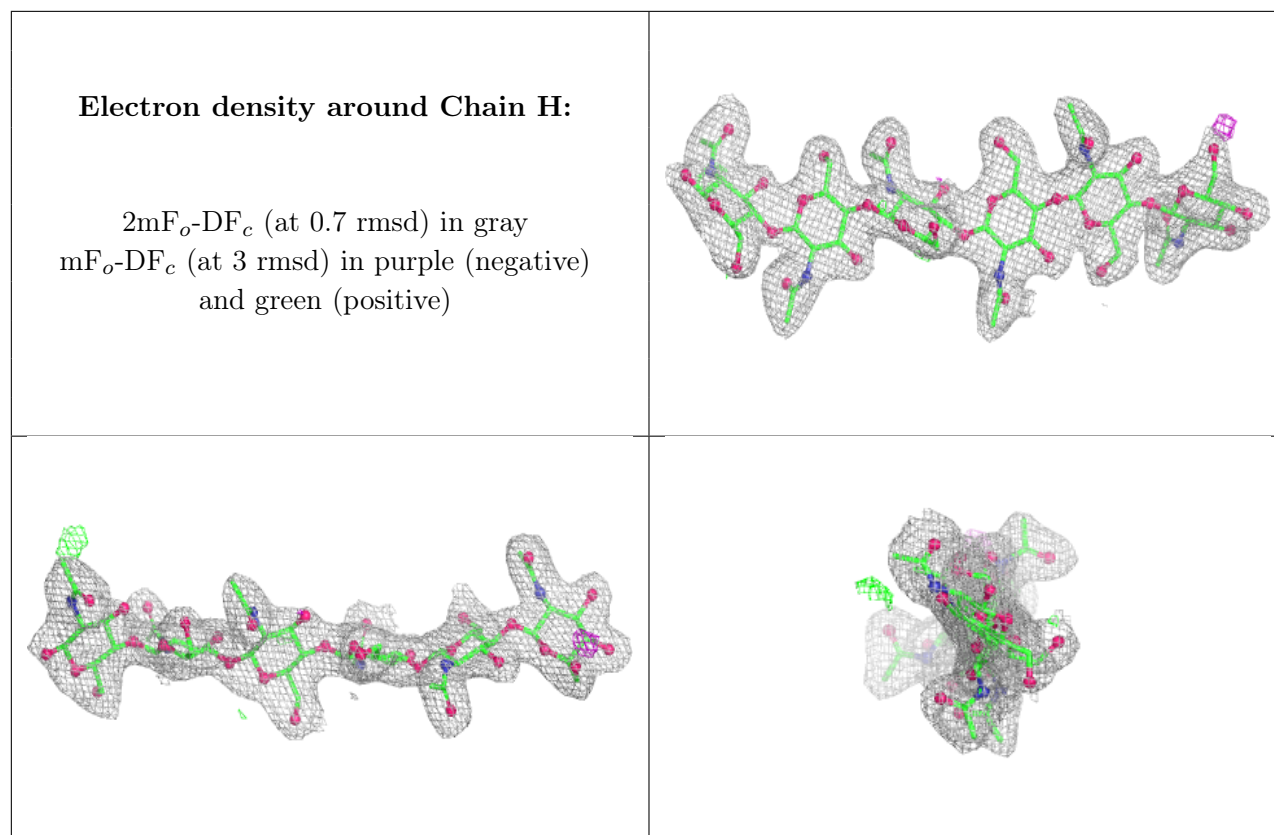
Electron density around Chain F:

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

**Electron density around Chain G:**

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





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($\text{\AA}^2$)	Q<0.9
3	PO4	C	701	5/5	0.99	0.02	44,44,46,46	0
3	PO4	B	696	5/5	1.00	0.02	27,28,28,29	0
3	PO4	C	697	5/5	1.00	0.02	23,24,25,25	0
3	PO4	A	699	5/5	1.00	0.02	20,21,22,25	0
3	PO4	D	698	5/5	1.00	0.03	20,22,24,25	0
3	PO4	D	700	5/5	1.00	0.02	42,42,42,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.