



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:04 PM UTC

PDB ID : 5SWB / pdb_00005swb
Title : Crystal structure of N-glycan transport solute binding protein (NgtS) from Streptococcus pneumoniae in complex with Man5GlcNAc
Authors : Robb, M.; Boraston, A.B.
Deposited on : 2016-08-08
Resolution : 1.73 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

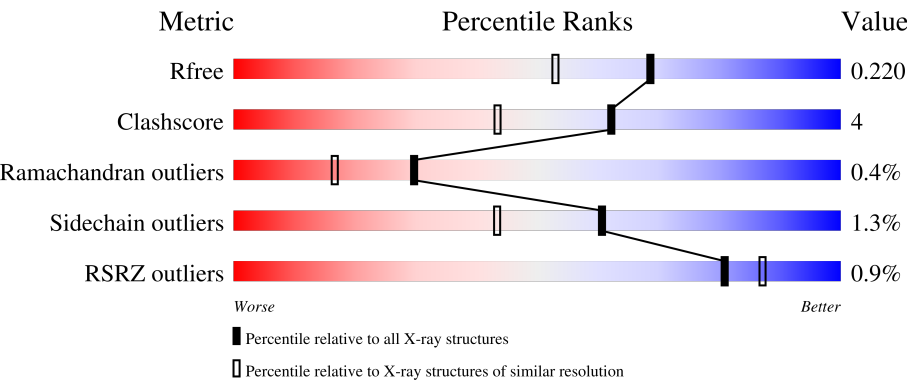
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.73 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.



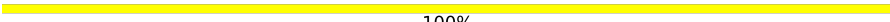
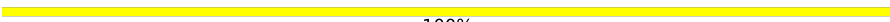
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1187 (1.74-1.74)
Clashscore	190562	1207 (1.74-1.74)
Ramachandran outliers	187476	1200 (1.74-1.74)
Sidechain outliers	187428	1200 (1.74-1.74)
RSRZ outliers	180081	1188 (1.74-1.74)

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	476	% 87%8%5%
1	C	476	% 89%6%5%
1	D	476	% 85%9%5%
1	G	476	% 83%11%5%
2	B	6	100%

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Mol	Chain	Length	Quality of chain
2	F	6	 100%
2	H	6	 100%
3	E	6	 100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16745 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 Extracellular solute-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	454	Total	C	N	O	S	0	8	0
			3583	2288	584	704	7			
1	C	451	Total	C	N	O	S	0	5	0
			3552	2263	581	701	7			
1	D	450	Total	C	N	O	S	0	5	0
			3548	2266	578	697	7			
1	G	454	Total	C	N	O	S	0	8	0
			3600	2300	587	706	7			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	initiating methionine	UNP A0A0T7JX40
A	492	LEU	-	expression tag	UNP A0A0T7JX40
A	493	GLU	-	expression tag	UNP A0A0T7JX40
A	494	HIS	-	expression tag	UNP A0A0T7JX40
A	495	HIS	-	expression tag	UNP A0A0T7JX40
A	496	HIS	-	expression tag	UNP A0A0T7JX40
A	497	HIS	-	expression tag	UNP A0A0T7JX40
A	498	HIS	-	expression tag	UNP A0A0T7JX40
A	499	HIS	-	expression tag	UNP A0A0T7JX40
C	24	MET	-	initiating methionine	UNP A0A0T7JX40
C	492	LEU	-	expression tag	UNP A0A0T7JX40
C	493	GLU	-	expression tag	UNP A0A0T7JX40
C	494	HIS	-	expression tag	UNP A0A0T7JX40
C	495	HIS	-	expression tag	UNP A0A0T7JX40
C	496	HIS	-	expression tag	UNP A0A0T7JX40
C	497	HIS	-	expression tag	UNP A0A0T7JX40
C	498	HIS	-	expression tag	UNP A0A0T7JX40
C	499	HIS	-	expression tag	UNP A0A0T7JX40
D	24	MET	-	initiating methionine	UNP A0A0T7JX40
D	492	LEU	-	expression tag	UNP A0A0T7JX40
D	493	GLU	-	expression tag	UNP A0A0T7JX40

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Chain	Residue	Modelled	Actual	Comment	Reference
D	494	HIS	-	expression tag	UNP A0A0T7JX40
D	495	HIS	-	expression tag	UNP A0A0T7JX40
D	496	HIS	-	expression tag	UNP A0A0T7JX40
D	497	HIS	-	expression tag	UNP A0A0T7JX40
D	498	HIS	-	expression tag	UNP A0A0T7JX40
D	499	HIS	-	expression tag	UNP A0A0T7JX40
G	24	MET	-	initiating methionine	UNP A0A0T7JX40
G	492	LEU	-	expression tag	UNP A0A0T7JX40
G	493	GLU	-	expression tag	UNP A0A0T7JX40
G	494	HIS	-	expression tag	UNP A0A0T7JX40
G	495	HIS	-	expression tag	UNP A0A0T7JX40
G	496	HIS	-	expression tag	UNP A0A0T7JX40
G	497	HIS	-	expression tag	UNP A0A0T7JX40
G	498	HIS	-	expression tag	UNP A0A0T7JX40
G	499	HIS	-	expression tag	UNP A0A0T7JX40

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	0	0	0
			70	38	1	31			
2	F	6	Total	C	N	O	0	0	0
			70	38	1	31			
2	H	6	Total	C	N	O	0	0	0
			70	38	1	31			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-5)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	6	Total	C	N	O	0	0	0
			69	37	1	31			

- Molecule 4 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	11	Total	Cd	0	0
			11	11		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	10	Total 10	Cd 10	0	0
4	D	6	Total 6	Cd 6	0	0
4	G	7	Total 7	Cd 7	0	0

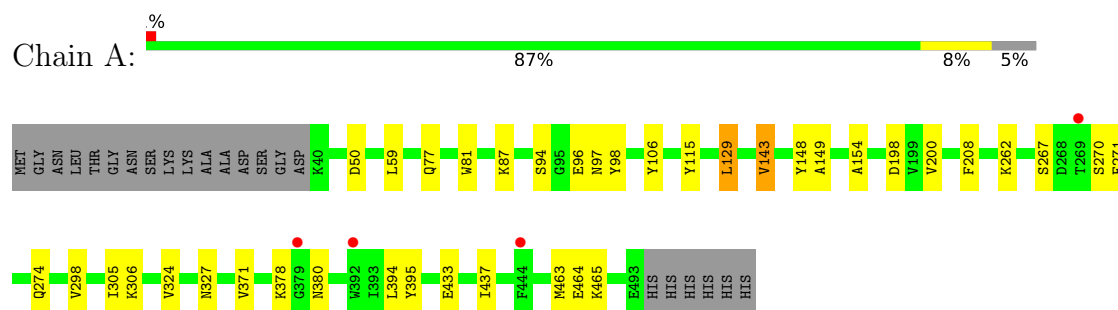
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	549	Total 549	O 549	0	0
5	C	500	Total 500	O 500	0	0
5	D	555	Total 555	O 555	0	0
5	G	545	Total 545	O 545	0	0

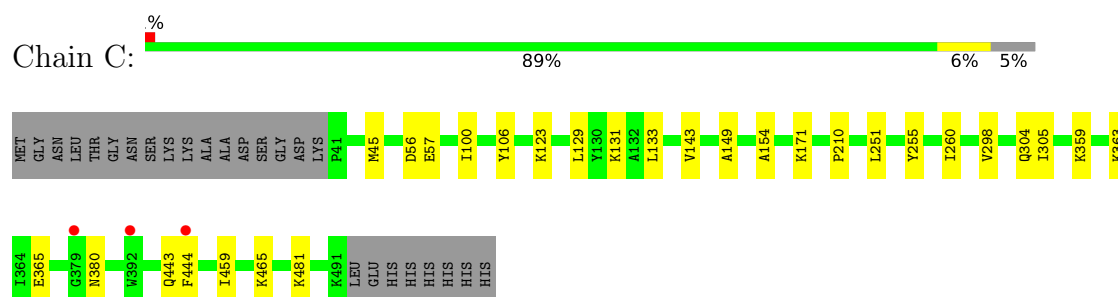
3 Residue-property plots [i](#)

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.

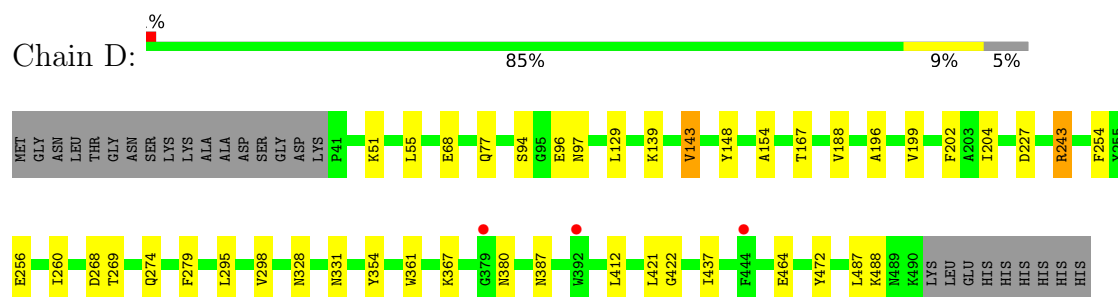
- Molecule 1: Extracellular solute-binding protein



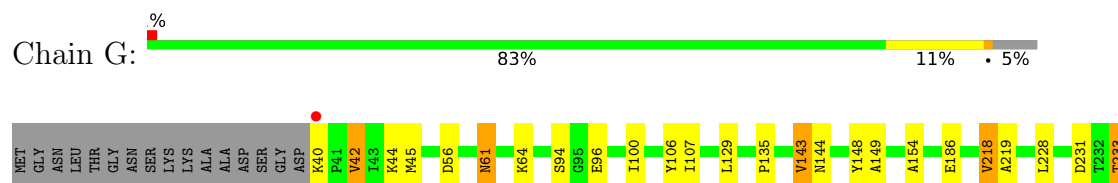
- Molecule 1: Extracellular solute-binding protein

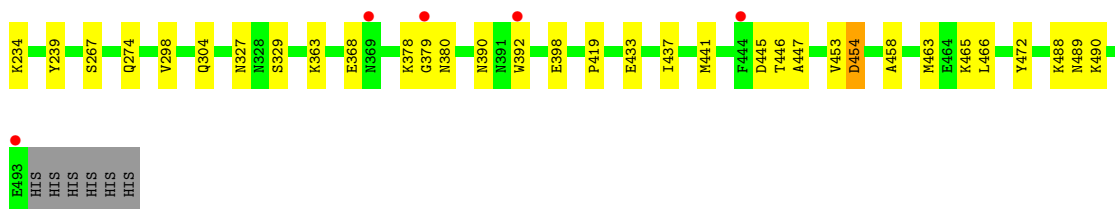


- Molecule 1: Extracellular solute-binding protein



- Molecule 1: Extracellular solute-binding protein





- Molecule 2: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain B:  100%



- Molecule 2: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%



- Molecule 2: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  100%



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

Chain E:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	104.18Å 110.60Å 105.84Å 90.00° 109.52° 90.00°	Depositor
Resolution (Å)	99.76 – 1.73 99.76 – 1.73	Depositor EDS
% Data completeness (in resolution range)	99.9 (99.76-1.73) 99.9 (99.76-1.73)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.28 (at 1.73Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.180 , 0.220 0.180 , 0.220	Depositor DCC
R_{free} test set	11777 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.019 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16745	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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 36.14 % 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 5.2350e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.19	3/3682 (0.1%)	1.15	4/4994 (0.1%)
1	C	1.17	0/3641	1.10	5/4938 (0.1%)
1	D	1.16	1/3635 (0.0%)	1.13	2/4930 (0.0%)
1	G	1.19	6/3701 (0.2%)	1.16	10/5019 (0.2%)
All	All	1.18	10/14659 (0.1%)	1.13	21/19881 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	454	ASP	CA-CB	6.66	1.63	1.53
1	G	107	ILE	N-CA	6.38	1.54	1.46
1	G	445	ASP	CA-CB	5.95	1.62	1.53
1	A	208	PHE	N-CA	5.73	1.53	1.46
1	G	329	SER	N-CA	5.41	1.52	1.46
1	G	379	GLY	C-O	5.35	1.29	1.23
1	D	387	ASN	CA-CB	5.33	1.60	1.53
1	A	81	TRP	N-CA	5.30	1.52	1.46
1	G	454	ASP	CG-OD1	-5.12	1.15	1.25
1	A	394	LEU	N-CA	5.08	1.52	1.46

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	40	LYS	CA-C-N	-10.26	109.49	119.85
1	G	40	LYS	C-N-CA	-10.26	109.49	119.85
1	D	243	ARG	NE-CZ-NH2	-7.78	112.20	119.20
1	G	454	ASP	CB-CG-OD2	7.42	135.46	118.40
1	G	454	ASP	OD1-CG-OD2	-6.19	108.05	122.90
1	D	243	ARG	NE-CZ-NH1	6.15	127.65	121.50
1	G	379	GLY	N-CA-C	6.09	122.53	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	305	ILE	N-CA-C	5.96	117.17	108.53
1	A	50	ASP	CA-C-N	5.77	128.50	120.65
1	A	50	ASP	C-N-CA	5.77	128.50	120.65
1	A	305	ILE	N-CA-C	5.44	117.23	108.85
1	C	133	LEU	N-CA-C	5.37	118.30	109.76
1	C	459	ILE	CA-C-N	-5.24	113.64	119.19
1	C	459	ILE	C-N-CA	-5.24	113.64	119.19
1	G	446	THR	N-CA-C	5.22	116.65	111.07
1	G	61	ASN	N-CA-C	5.18	116.93	111.28
1	C	365	GLU	N-CA-CB	5.17	117.64	109.83
1	A	98	TYR	CA-C-O	-5.15	115.94	121.45
1	G	458	ALA	N-CA-C	5.09	116.91	111.36
1	G	398	GLU	N-CA-C	5.06	117.53	111.71
1	G	234	LYS	N-CA-C	5.00	117.55	109.40

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	3583	0	3526	23	0
1	C	3552	0	3482	23	0
1	D	3548	0	3501	34	0
1	G	3600	0	3543	44	1
2	B	70	0	60	0	0
2	F	70	0	60	0	0
2	H	70	0	60	0	0
3	E	69	0	58	0	0
4	A	11	0	0	0	1
4	C	10	0	0	1	0
4	D	6	0	0	1	1
4	G	7	0	0	0	1
5	A	549	0	0	7	1
5	C	500	0	0	9	1
5	D	555	0	0	16	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	G	545	0	0	14	1
All	All	16745	0	14290	118	5

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 4.

All (118) 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:C:56[B]:ASP:OD1	1:G:56[B]:ASP:CG	1.91	1.13
1:D:367:LYS:HE2	5:D:686:HOH:O	1.57	1.04
1:C:56[B]:ASP:CG	1:G:56[B]:ASP:CG	2.24	1.03
1:C:56[B]:ASP:OD1	1:G:56[B]:ASP:OD2	1.77	1.02
1:G:463:MET:CE	1:G:466:LEU:HD12	1.96	0.96
4:C:512:CD:CD	5:C:644:HOH:O	1.36	0.94
1:G:94:SER:OG	1:G:96:GLU:HG3	1.71	0.91
1:C:56[B]:ASP:OD1	1:G:56[B]:ASP:OD1	1.90	0.90
1:D:380:ASN:HA	5:D:748:HOH:O	1.71	0.90
1:G:463:MET:HE1	1:G:466:LEU:HD12	1.59	0.83
1:D:96:GLU:HG2	5:D:707:HOH:O	1.77	0.83
1:A:77:GLN:HE21	1:D:77:GLN:HE21	1.26	0.82
1:G:463:MET:HE2	1:G:466:LEU:HD12	1.63	0.81
1:G:231:ASP:OD1	1:G:233:THR:HG22	1.81	0.80
1:C:304[A]:GLN:OE1	5:C:601:HOH:O	1.99	0.79
1:G:304[B]:GLN:NE2	5:G:601:HOH:O	2.04	0.79
1:C:56[B]:ASP:CG	1:G:56[B]:ASP:OD1	2.24	0.79
1:A:306:LYS:HD3	5:A:610:HOH:O	1.86	0.75
1:A:327:ASN:ND2	5:A:602:HOH:O	2.19	0.74
1:G:463:MET:HE2	1:G:463:MET:HA	1.71	0.72
1:G:363:LYS:NZ	5:G:605:HOH:O	2.22	0.72
1:D:331:ASN:ND2	5:D:602:HOH:O	2.24	0.71
1:D:139:LYS:HE3	1:D:421:LEU:O	1.90	0.70
1:D:77:GLN:NE2	5:D:603:HOH:O	2.25	0.68
1:G:144:ASN:ND2	5:G:608:HOH:O	2.26	0.66
1:D:167:THR:OG1	5:D:601:HOH:O	2.12	0.66
1:A:274:GLN:NE2	5:A:601:HOH:O	2.15	0.65
4:D:509:CD:CD	5:D:789:HOH:O	1.65	0.63
1:C:444:PHE:CE1	1:C:465:LYS:HE3	2.34	0.63
1:A:94:SER:OG	1:A:96[B]:GLU:HG3	1.99	0.62
1:C:56[B]:ASP:OD2	1:G:56[B]:ASP:OD1	2.17	0.62
1:G:267:SER:HB3	5:G:756:HOH:O	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:ASP:OD2	1:C:131:LYS:HG2	2.00	0.61
1:G:463:MET:HE1	1:G:466:LEU:CD1	2.28	0.61
1:G:489:ASN:ND2	5:G:610:HOH:O	2.33	0.61
1:C:298[A]:VAL:HG11	5:C:1065:HOH:O	2.00	0.61
1:G:368:GLU:HA	5:G:993:HOH:O	2.01	0.60
1:D:487:LEU:HD12	5:D:724:HOH:O	2.01	0.60
1:C:123:LYS:NZ	5:C:609:HOH:O	2.34	0.59
1:G:390:ASN:OD1	1:G:392[A]:TRP:HB2	2.02	0.59
1:D:188[A]:VAL:HG12	1:D:279:PHE:HE2	1.65	0.59
1:D:202:PHE:CD2	1:D:260[A]:ILE:HD11	2.39	0.57
1:A:96[A]:GLU:HG2	1:A:97:ASN:N	2.19	0.57
1:G:433:GLU:OE2	5:G:602:HOH:O	2.17	0.57
1:A:298[A]:VAL:HG11	5:A:1119:HOH:O	2.05	0.57
1:G:218:VAL:O	1:G:219:ALA:HB3	2.05	0.56
1:A:433:GLU:O	1:A:437:ILE:HG12	2.05	0.56
1:D:274:GLN:NE2	5:D:605:HOH:O	2.31	0.55
1:A:200:VAL:HG21	1:A:271:PHE:CE1	2.40	0.55
1:D:139:LYS:HG2	1:D:422:GLY:HA3	1.89	0.55
1:D:202:PHE:HD2	1:D:260[A]:ILE:HD11	1.71	0.54
1:C:171:LYS:NZ	5:C:610:HOH:O	2.37	0.53
1:D:227:ASP:OD2	1:D:243:ARG:NH2	2.37	0.53
1:G:363:LYS:NZ	5:G:603:HOH:O	2.20	0.53
1:D:94:SER:OG	1:D:96:GLU:HG3	2.09	0.52
1:G:45:MET:HE2	1:G:100:ILE:HG21	1.92	0.52
1:G:61:ASN:HA	1:G:64:LYS:HE2	1.92	0.51
1:G:298[A]:VAL:HG11	5:G:1101:HOH:O	2.09	0.51
1:D:254:PHE:HB3	1:D:260[A]:ILE:HG12	1.92	0.51
1:G:231:ASP:CG	1:G:233:THR:HG22	2.36	0.50
1:C:444:PHE:CD1	1:C:465:LYS:HE3	2.47	0.50
1:D:328:ASN:HB2	5:D:935:HOH:O	2.10	0.50
1:D:204:ILE:HG12	1:D:260[A]:ILE:HD12	1.94	0.49
1:D:464:GLU:HG3	5:D:625:HOH:O	2.12	0.49
1:D:97:ASN:HA	5:D:935:HOH:O	2.12	0.49
1:A:267:SER:HB3	5:A:763:HOH:O	2.13	0.49
1:G:239:TYR:HE2	1:G:441:MET:HE1	1.77	0.48
1:D:202:PHE:HB3	1:D:260[A]:ILE:HD13	1.95	0.48
1:G:437:ILE:HD11	1:G:472:TYR:HA	1.96	0.48
1:C:363:LYS:NZ	5:C:607:HOH:O	2.31	0.48
1:G:143[A]:VAL:CG1	1:G:148:TYR:HD2	2.26	0.48
1:C:359:LYS:NZ	5:C:617:HOH:O	2.47	0.47
1:A:106:TYR:CE1	1:A:149:ALA:HB2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:LYS:NZ	5:A:610:HOH:O	2.39	0.46
1:D:488:LYS:CB	5:D:1083:HOH:O	2.62	0.46
1:A:465:LYS:HE2	1:A:465:LYS:HA	1.98	0.46
1:D:243:ARG:HB2	5:D:1005:HOH:O	2.16	0.46
1:C:210:PRO:HG3	1:C:251:LEU:CD1	2.47	0.45
1:D:51:LYS:HE3	1:D:55:LEU:HD23	1.97	0.45
1:G:490:LYS:NZ	5:G:624:HOH:O	2.49	0.45
1:D:227:ASP:CG	1:D:243:ARG:HH22	2.25	0.45
1:A:143:VAL:HG22	1:A:148:TYR:HE2	1.81	0.45
1:D:196:ALA:O	1:D:199:VAL:HG22	2.17	0.44
1:C:380:ASN:HA	5:C:866:HOH:O	2.17	0.44
1:D:256:GLU:HA	1:D:256:GLU:OE1	2.17	0.44
1:G:143[A]:VAL:HG13	1:G:148:TYR:CD2	2.53	0.44
1:G:228:LEU:HD22	1:G:419:PRO:HD2	1.98	0.44
1:A:306:LYS:CD	5:A:610:HOH:O	2.56	0.44
1:A:129:LEU:C	1:A:129:LEU:HD12	2.42	0.44
1:G:44:LYS:CD	5:G:684:HOH:O	2.65	0.44
1:D:295:LEU:HA	1:D:298:VAL:HG22	2.00	0.44
1:C:444:PHE:CZ	1:C:465:LYS:HE3	2.52	0.44
1:A:96[A]:GLU:HG2	1:A:97:ASN:O	2.18	0.43
1:A:115:TYR:CE1	1:A:324:VAL:HG21	2.54	0.43
1:A:143:VAL:HG22	1:A:148:TYR:CE2	2.53	0.43
1:G:447:ALA:C	1:G:453:VAL:HG22	2.44	0.43
1:C:255:TYR:HB2	1:C:260:ILE:HD11	2.00	0.42
1:G:106:TYR:CE1	1:G:149:ALA:HB2	2.53	0.42
1:G:274:GLN:NE2	1:G:378:LYS:O	2.49	0.42
1:A:371:VAL:HG13	1:A:395:TYR:HB3	2.01	0.42
1:C:57:GLU:OE1	5:C:602:HOH:O	2.21	0.42
1:D:268:ASP:O	1:D:269:THR:C	2.62	0.42
1:G:488:LYS:NZ	5:G:633:HOH:O	2.53	0.42
1:G:186:GLU:OE2	5:G:604:HOH:O	2.22	0.42
1:A:59:LEU:HD23	1:A:59:LEU:HA	1.92	0.42
1:C:106:TYR:CE1	1:C:149:ALA:HB2	2.55	0.42
1:D:354:TYR:HB3	1:D:361:TRP:CD1	2.55	0.41
1:D:412:LEU:HD12	5:D:945:HOH:O	2.20	0.41
1:G:143[A]:VAL:CG1	1:G:148:TYR:CD2	3.03	0.41
1:G:42:VAL:HG13	5:G:978:HOH:O	2.20	0.41
1:G:463:MET:HE2	1:G:463:MET:CA	2.47	0.41
1:C:443:GLN:HB2	1:C:444:PHE:CD1	2.56	0.41
1:A:87:LYS:HG2	5:D:991:HOH:O	2.20	0.41
1:D:143:VAL:HG22	1:D:148:TYR:CE2	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:218:VAL:O	1:G:219:ALA:CB	2.67	0.40
1:D:437:ILE:HD11	1:D:472:TYR:HA	2.03	0.40
1:C:45:MET:HA	1:C:100:ILE:O	2.22	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:454:ASP:OD2	4:G:511:CD:CD[2_555]	1.42	0.78
4:D:510:CD:CD	5:D:694:HOH:O[2_544]	1.53	0.67
4:A:511:CD:CD	5:C:612:HOH:O[2_454]	1.74	0.46
5:D:615:HOH:O	5:D:976:HOH:O[2_554]	2.00	0.20
5:A:823:HOH:O	5:G:1053:HOH:O[1_554]	2.09	0.11

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	460/476 (97%)	449 (98%)	8 (2%)	3 (1%)	18	6
1	C	454/476 (95%)	442 (97%)	11 (2%)	1 (0%)	43	29
1	D	453/476 (95%)	441 (97%)	11 (2%)	1 (0%)	43	29
1	G	460/476 (97%)	445 (97%)	13 (3%)	2 (0%)	30	16
All	All	1827/1904 (96%)	1777 (97%)	43 (2%)	7 (0%)	30	16

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	ALA
1	C	154	ALA
1	D	154	ALA

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Mol	Chain	Res	Type
1	G	154	ALA
1	G	380	ASN
1	A	378	LYS
1	A	380	ASN

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	389/408 (95%)	383 (98%)	6 (2%)	57	38
1	C	385/408 (94%)	382 (99%)	3 (1%)	73	64
1	D	385/408 (94%)	382 (99%)	3 (1%)	73	64
1	G	391/408 (96%)	381 (97%)	10 (3%)	40	17
All	All	1550/1632 (95%)	1528 (99%)	22 (1%)	61	41

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

Mol	Chain	Res	Type
1	A	129	LEU
1	A	143	VAL
1	A	262	LYS
1	A	270	SER
1	A	463	MET
1	A	464	GLU
1	C	129	LEU
1	C	143	VAL
1	C	481	LYS
1	D	68	GLU
1	D	129	LEU
1	D	143	VAL
1	G	42	VAL
1	G	129	LEU
1	G	135	PRO
1	G	143[A]	VAL
1	G	143[B]	VAL

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Mol	Chain	Res	Type
1	G	218	VAL
1	G	233	THR
1	G	327[A]	ASN
1	G	327[B]	ASN
1	G	465	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	160	GLN
1	A	391	ASN
1	C	160	GLN
1	C	369	ASN
1	D	160	GLN
1	G	160	GLN
1	G	360	ASN
1	G	443	GLN
1	G	489	ASN

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](#)

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	B	1	2	15,15,15	1.06	1 (6%)	21,21,21	0.98	0
2	BMA	B	2	2	11,11,12	1.23	1 (9%)	15,15,17	1.20	2 (13%)
2	MAN	B	3	2	11,11,12	0.83	0	15,15,17	1.47	2 (13%)
2	MAN	B	4	2	11,11,12	1.19	2 (18%)	15,15,17	1.42	2 (13%)
2	MAN	B	5	2	11,11,12	0.75	0	15,15,17	1.45	4 (26%)
2	MAN	B	6	2	11,11,12	0.87	0	15,15,17	0.98	1 (6%)
3	NAG	E	1	3	15,15,15	0.95	0	21,21,21	1.17	2 (9%)
3	BMA	E	2	3	11,11,12	1.25	1 (9%)	15,15,17	1.09	1 (6%)
3	MAN	E	3	3	9,9,12	1.99	2 (22%)	10,12,17	4.63	5 (50%)
3	MAN	E	4	3	11,11,12	1.39	2 (18%)	15,15,17	1.09	0
3	MAN	E	5	3	12,12,12	1.25	1 (8%)	17,17,17	1.78	4 (23%)
3	MAN	E	6	3	11,11,12	1.60	3 (27%)	15,15,17	1.22	2 (13%)
2	NAG	F	1	2	15,15,15	0.66	0	21,21,21	1.52	3 (14%)
2	BMA	F	2	2	11,11,12	1.08	1 (9%)	15,15,17	1.47	3 (20%)
2	MAN	F	3	2	11,11,12	1.24	1 (9%)	15,15,17	1.43	3 (20%)
2	MAN	F	4	2	11,11,12	1.23	2 (18%)	15,15,17	1.15	2 (13%)
2	MAN	F	5	2	11,11,12	0.77	0	15,15,17	1.61	3 (20%)
2	MAN	F	6	2	11,11,12	1.40	2 (18%)	15,15,17	1.21	1 (6%)
2	NAG	H	1	2	15,15,15	1.09	1 (6%)	21,21,21	1.37	2 (9%)
2	BMA	H	2	2	11,11,12	1.26	2 (18%)	15,15,17	1.01	1 (6%)
2	MAN	H	3	2	11,11,12	1.19	1 (9%)	15,15,17	1.97	3 (20%)
2	MAN	H	4	2	11,11,12	1.05	0	15,15,17	1.74	3 (20%)
2	MAN	H	5	2	11,11,12	0.88	0	15,15,17	1.32	1 (6%)
2	MAN	H	6	2	11,11,12	1.39	2 (18%)	15,15,17	1.24	2 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	2	-	0/6/26/26	0/1/1/1
2	BMA	B	2	2	-	0/2/19/22	0/1/1/1
2	MAN	B	3	2	-	0/2/19/22	0/1/1/1
2	MAN	B	4	2	-	0/2/19/22	0/1/1/1
2	MAN	B	5	2	-	0/2/19/22	0/1/1/1
2	MAN	B	6	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	E	1	3	-	0/6/26/26	0/1/1/1
3	BMA	E	2	3	-	0/2/19/22	0/1/1/1
3	MAN	E	3	3	-	-	0/1/1/1
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
3	MAN	E	5	3	-	0/2/22/22	0/1/1/1
3	MAN	E	6	3	-	0/2/19/22	0/1/1/1
2	NAG	F	1	2	-	0/6/26/26	0/1/1/1
2	BMA	F	2	2	-	0/2/19/22	0/1/1/1
2	MAN	F	3	2	-	0/2/19/22	0/1/1/1
2	MAN	F	4	2	-	0/2/19/22	0/1/1/1
2	MAN	F	5	2	-	1/2/19/22	0/1/1/1
2	MAN	F	6	2	-	0/2/19/22	0/1/1/1
2	NAG	H	1	2	-	0/6/26/26	0/1/1/1
2	BMA	H	2	2	-	0/2/19/22	0/1/1/1
2	MAN	H	3	2	-	0/2/19/22	0/1/1/1
2	MAN	H	4	2	-	0/2/19/22	0/1/1/1
2	MAN	H	5	2	-	0/2/19/22	0/1/1/1
2	MAN	H	6	2	-	0/2/19/22	0/1/1/1

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	3	MAN	C4-C3	3.95	1.58	1.52
3	E	3	MAN	C5-C4	3.45	1.60	1.52
2	H	3	MAN	O5-C5	-3.32	1.37	1.43
3	E	6	MAN	C2-C3	3.24	1.57	1.52
3	E	5	MAN	O1-C1	3.23	1.49	1.39
2	F	6	MAN	C4-C5	2.95	1.59	1.53
3	E	2	BMA	C2-C3	2.92	1.57	1.52
2	B	4	MAN	O2-C2	-2.79	1.37	1.43
2	F	4	MAN	O5-C1	-2.77	1.39	1.43
3	E	4	MAN	C2-C3	2.76	1.56	1.52
2	F	4	MAN	C2-C3	2.72	1.56	1.52
2	H	6	MAN	C1-C2	2.55	1.58	1.52
3	E	4	MAN	O5-C1	-2.47	1.39	1.43
2	F	2	BMA	C2-C3	2.38	1.56	1.52
3	E	6	MAN	O5-C1	2.33	1.47	1.43
2	F	6	MAN	C2-C3	2.31	1.56	1.52
2	B	4	MAN	O5-C1	-2.29	1.39	1.43
2	H	6	MAN	C2-C3	2.19	1.55	1.52
2	H	2	BMA	O3-C3	2.18	1.48	1.43
2	B	1	NAG	C1-C2	2.18	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	3	MAN	C2-C3	2.15	1.55	1.52
2	B	2	BMA	C1-C2	2.14	1.57	1.52
3	E	6	MAN	C4-C5	2.14	1.57	1.53
2	H	1	NAG	O7-C7	2.08	1.27	1.23
2	H	2	BMA	C1-C2	2.04	1.57	1.52

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	3	MAN	C5-O5-C1	10.94	129.07	111.42
3	E	3	MAN	C5-C4-C3	7.63	120.75	109.64
2	H	3	MAN	O3-C3-C2	-4.33	101.22	110.05
3	E	5	MAN	O1-C1-C2	4.17	121.08	108.98
2	F	5	MAN	C1-O5-C5	4.07	117.64	112.19
2	H	3	MAN	C1-O5-C5	4.05	117.61	112.19
3	E	3	MAN	C1-C2-C3	-3.49	104.56	109.64
2	B	3	MAN	O3-C3-C2	-3.47	102.97	110.05
3	E	3	MAN	O3-C3-C2	-3.25	103.43	110.05
2	H	4	MAN	C1-C2-C3	-3.19	105.00	109.64
2	H	3	MAN	C1-C2-C3	-3.19	105.01	109.64
3	E	3	MAN	C4-C3-C2	-3.17	107.16	110.92
2	H	5	MAN	C1-O5-C5	3.17	116.43	112.19
3	E	5	MAN	O5-C1-C2	-3.15	104.75	110.30
2	F	6	MAN	C3-C4-C5	-3.13	104.55	110.23
3	E	5	MAN	O5-C5-C6	3.10	114.12	106.44
2	H	4	MAN	O2-C2-C3	3.09	116.56	110.15
2	B	4	MAN	C3-C4-C5	-3.08	104.65	110.23
2	B	4	MAN	C1-O5-C5	3.03	116.24	112.19
2	F	2	BMA	C1-C2-C3	-3.02	105.24	109.64
2	F	3	MAN	C1-C2-C3	-3.00	105.28	109.64
2	F	5	MAN	O6-C6-C5	-2.97	101.24	111.33
2	H	1	NAG	C1-C2-N2	2.86	114.04	110.73
2	B	3	MAN	C1-C2-C3	-2.76	105.62	109.64
2	F	2	BMA	O2-C2-C3	-2.76	104.44	110.15
2	F	5	MAN	C6-C5-C4	-2.75	106.26	113.02
2	B	5	MAN	O6-C6-C5	-2.69	102.17	111.33
2	H	1	NAG	C3-C2-N2	-2.65	105.73	110.62
2	B	5	MAN	C3-C4-C5	2.65	115.03	110.23
2	F	1	NAG	C3-C4-C5	2.61	114.97	110.23
2	F	4	MAN	C3-C4-C5	-2.61	105.49	110.23
2	F	1	NAG	O1-C1-O5	-2.49	103.02	110.41
2	H	4	MAN	C1-O5-C5	2.48	115.51	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	1	NAG	O3-C3-C2	-2.47	104.67	109.58
2	F	1	NAG	O4-C4-C5	-2.46	103.26	109.32
3	E	5	MAN	C6-C5-C4	-2.39	107.16	113.02
2	F	4	MAN	O2-C2-C3	-2.32	105.34	110.15
2	B	2	BMA	O3-C3-C4	-2.28	105.00	110.38
3	E	6	MAN	O6-C6-C5	-2.28	103.58	111.33
2	H	2	BMA	O5-C5-C4	-2.27	105.31	110.83
2	F	3	MAN	O3-C3-C2	-2.26	105.43	110.05
2	B	5	MAN	C2-C3-C4	-2.21	106.97	110.86
2	B	5	MAN	O3-C3-C2	-2.21	105.54	110.05
2	F	2	BMA	O5-C5-C4	-2.13	105.65	110.83
2	H	6	MAN	O4-C4-C3	-2.08	105.47	110.38
3	E	6	MAN	O4-C4-C3	-2.08	105.48	110.38
2	H	6	MAN	C1-C2-C3	-2.07	106.64	109.64
2	B	6	MAN	O6-C6-C5	-2.06	104.31	111.33
2	B	2	BMA	C1-C2-C3	-2.06	106.65	109.64
3	E	2	BMA	C6-C5-C4	2.03	118.00	113.02
3	E	1	NAG	O5-C5-C6	2.02	111.46	106.44
2	F	3	MAN	O4-C4-C5	2.02	114.31	109.32

There are no chirality outliers.

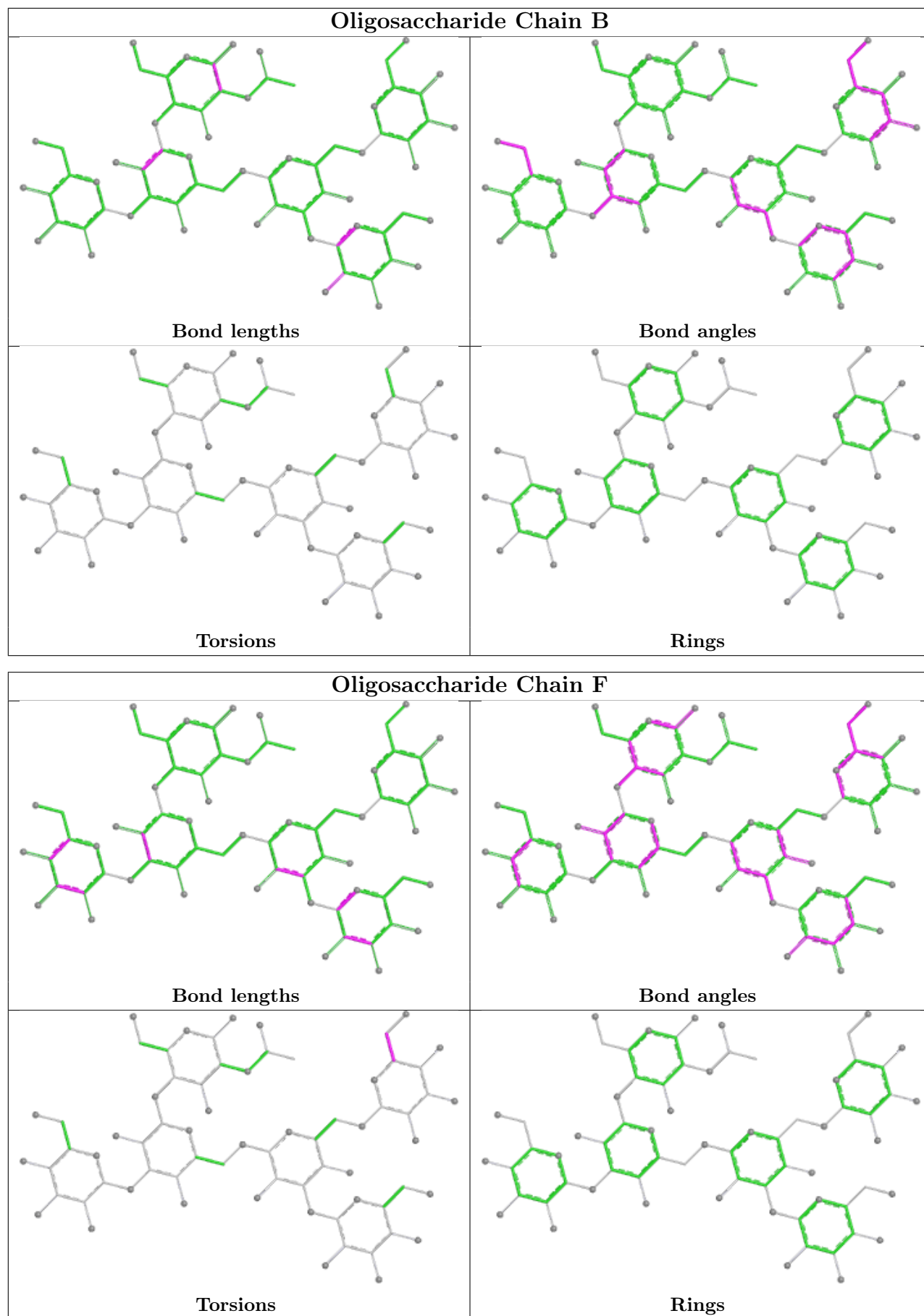
All (1) torsion outliers are listed below:

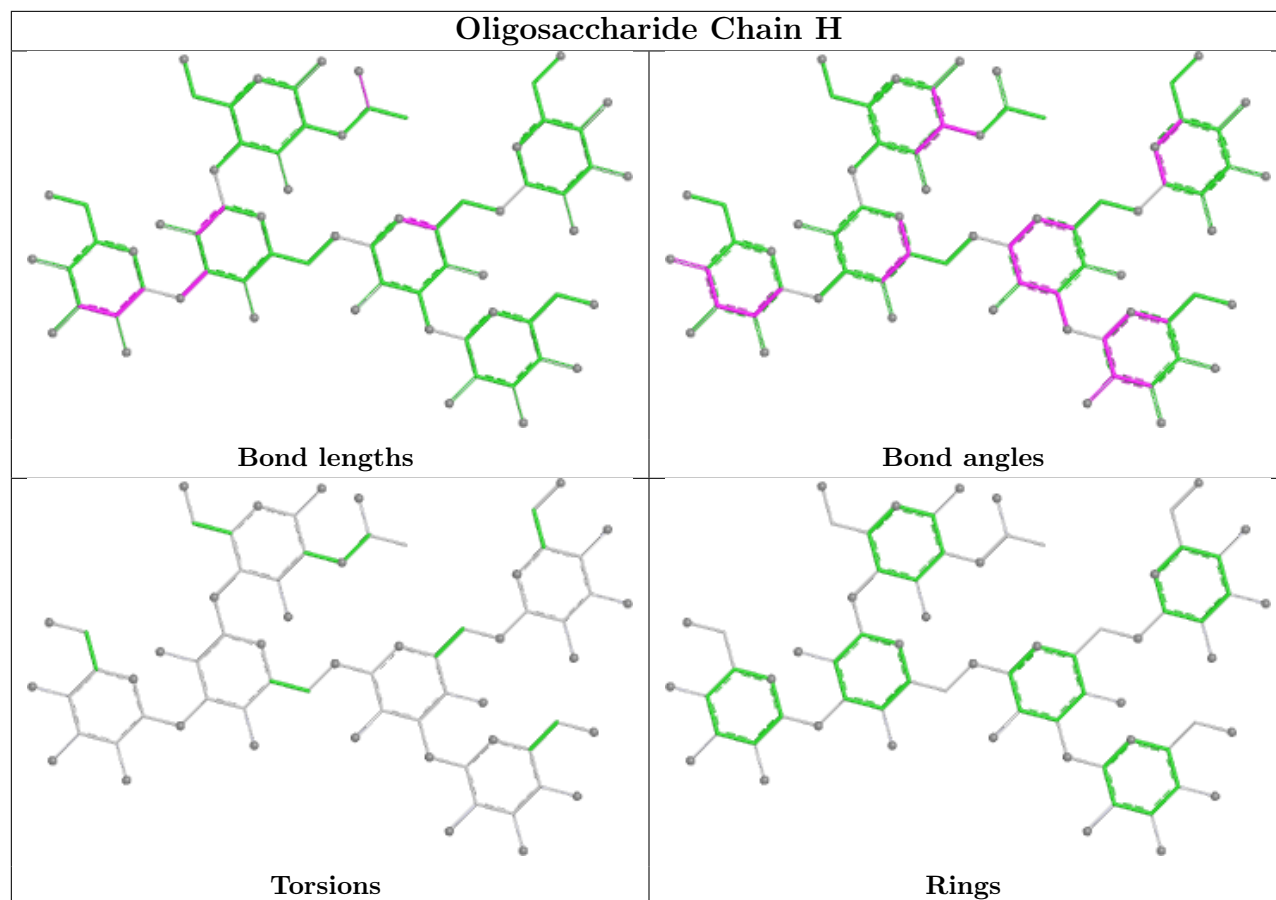
Mol	Chain	Res	Type	Atoms
2	F	5	MAN	C4-C5-C6-O6

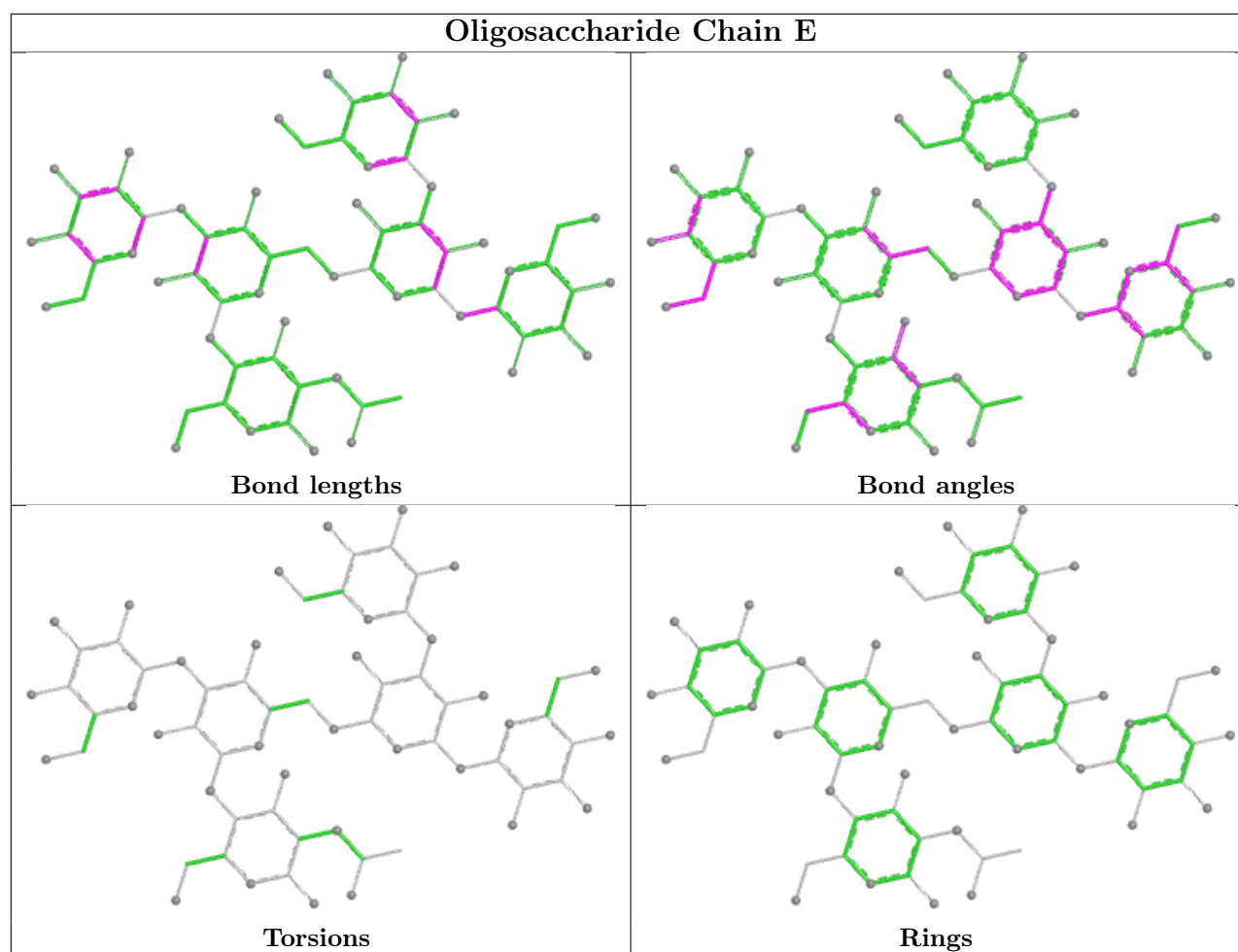
There are no ring outliers.

No monomer is involved in short contacts.

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







5.6 Ligand geometry [i](#)

Of 34 ligands modelled in this entry, 34 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

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 [i](#)

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	454/476 (95%)	-0.09	4 (0%) 81 87	15, 26, 42, 55	8 (1%)
1	C	451/476 (94%)	0.06	3 (0%) 84 89	15, 29, 46, 61	5 (1%)
1	D	450/476 (94%)	-0.03	3 (0%) 84 89	14, 27, 44, 61	5 (1%)
1	G	454/476 (95%)	0.06	6 (1%) 75 82	15, 27, 44, 63	9 (1%)
All	All	1809/1904 (95%)	-0.00	16 (0%) 81 87	14, 27, 44, 63	27 (1%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	392[A]	TRP	7.1
1	G	493	GLU	4.4
1	A	379	GLY	3.4
1	D	379	GLY	3.0
1	D	444	PHE	2.9
1	C	379	GLY	2.9
1	G	379	GLY	2.8
1	C	392	TRP	2.7
1	A	392	TRP	2.6
1	A	444	PHE	2.5
1	G	40	LYS	2.4
1	D	392	TRP	2.4
1	G	444	PHE	2.3
1	A	269	THR	2.2
1	G	369	ASN	2.1
1	C	444	PHE	2.1

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 ⓘ

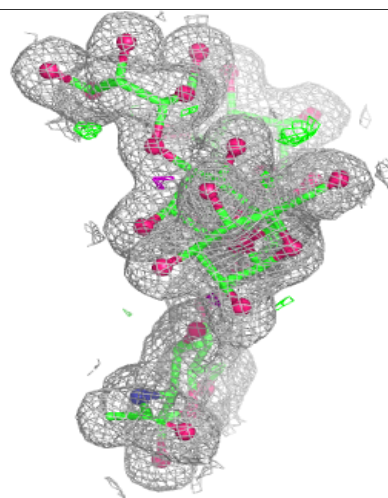
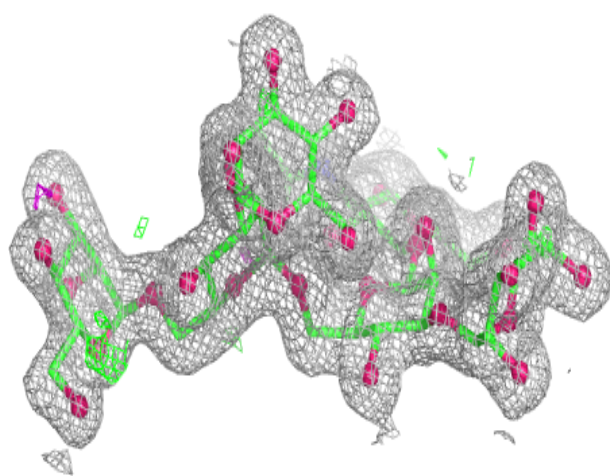
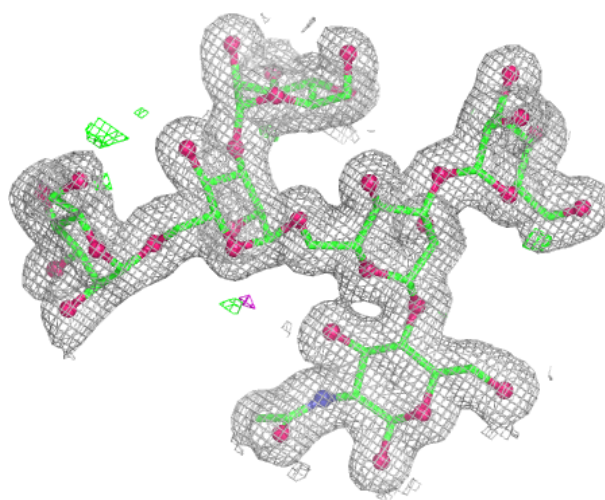
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	B	1	15/15	-	-	14,17,19,20	0
2	BMA	B	2	11/12	-	-	14,16,18,18	0
2	MAN	B	3	11/12	-	-	18,19,20,22	0
2	MAN	B	4	11/12	-	-	19,21,22,24	0
2	MAN	B	5	11/12	-	-	22,25,32,38	0
2	MAN	B	6	11/12	-	-	17,19,21,21	0
2	MAN	F	5	11/12	0.95	0.07	25,27,37,45	0
2	MAN	F	3	11/12	0.96	0.06	18,18,21,22	0
3	MAN	E	5	12/12	0.96	0.08	21,23,29,31	0
2	MAN	H	4	11/12	0.97	0.05	19,19,22,24	0
2	MAN	H	5	11/12	0.97	0.06	20,21,26,31	0
2	MAN	F	6	11/12	0.97	0.07	16,19,21,22	0
3	MAN	E	6	11/12	0.97	0.05	18,20,23,24	0
2	BMA	H	2	11/12	0.98	0.04	13,15,17,18	0
2	MAN	F	4	11/12	0.98	0.05	19,22,25,26	0
2	BMA	F	2	11/12	0.98	0.04	15,17,18,18	0
2	MAN	H	6	11/12	0.98	0.05	15,17,20,20	0
3	MAN	E	3	9/12	0.98	0.04	19,20,23,27	0
3	MAN	E	4	11/12	0.98	0.05	20,22,27,29	0
2	NAG	F	1	15/15	0.98	0.05	15,17,18,19	0
2	NAG	H	1	15/15	0.98	0.04	13,16,18,18	0
3	NAG	E	1	15/15	0.99	0.04	17,18,21,23	0
3	BMA	E	2	11/12	0.99	0.04	17,18,18,18	0
2	MAN	H	3	11/12	0.99	0.04	17,18,20,20	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.

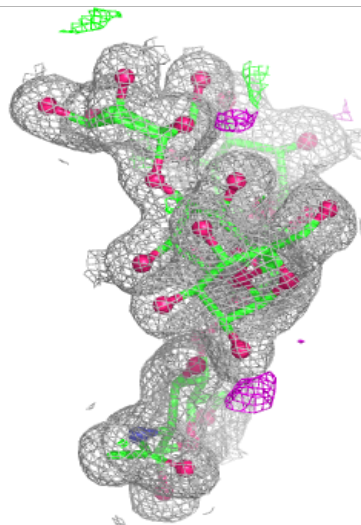
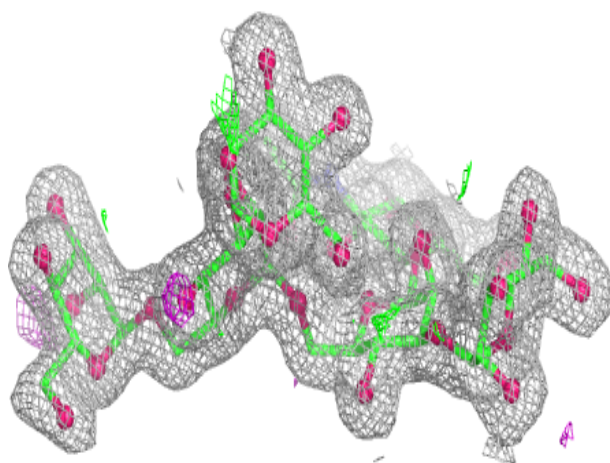
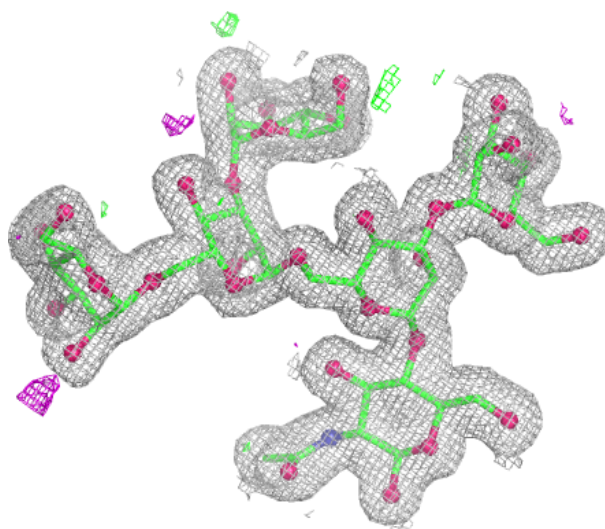
Electron density around Chain B:

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



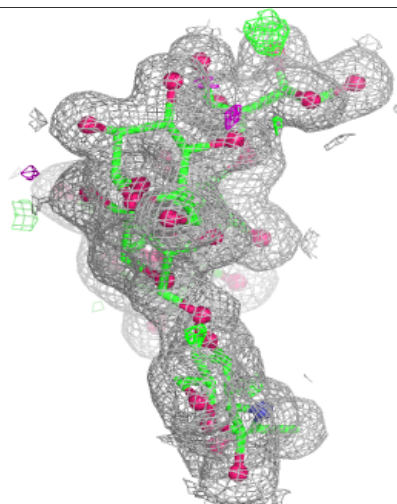
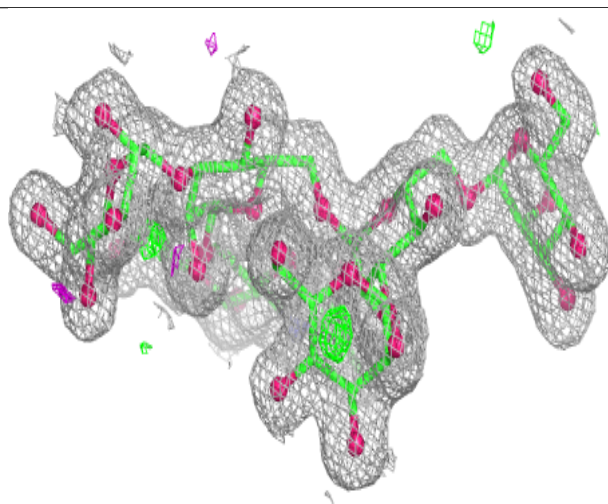
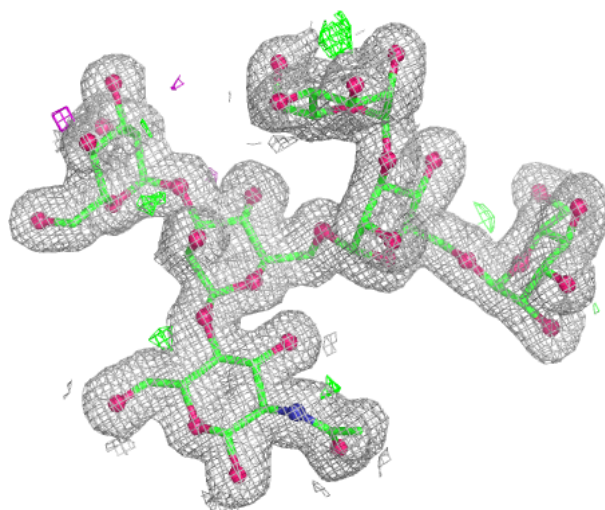
Electron density around Chain F:

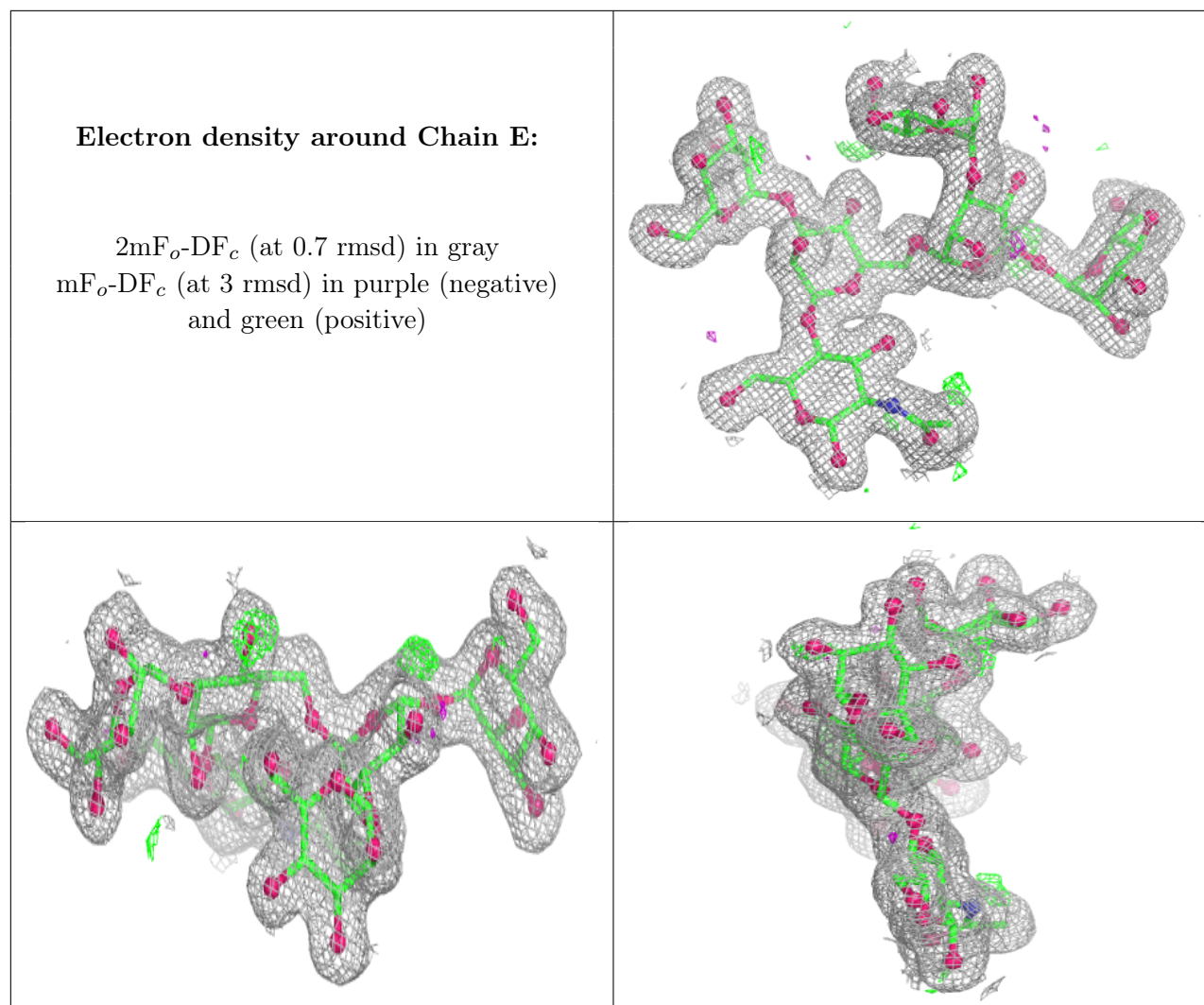
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain H:

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





6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CD	A	514	1/1	0.88	0.10	75,75,75,75	1
4	CD	G	512	1/1	0.88	0.17	136,136,136,136	0
4	CD	G	509	1/1	0.90	0.12	115,115,115,115	0
4	CD	G	513	1/1	0.91	0.07	23,23,23,23	1
4	CD	A	508	1/1	0.94	0.07	49,49,49,49	1
4	CD	A	517	1/1	0.95	0.10	47,47,47,47	1
4	CD	C	511	1/1	0.95	0.09	42,42,42,42	1
4	CD	G	508	1/1	0.95	0.06	27,27,27,27	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CD	D	511	1/1	0.96	0.06	31,31,31,31	1
4	CD	G	510	1/1	0.96	0.07	37,37,37,37	1
4	CD	D	512	1/1	0.96	0.06	41,41,41,41	1
4	CD	A	512	1/1	0.96	0.06	40,40,40,40	1
4	CD	C	515	1/1	0.97	0.09	42,42,42,42	1
4	CD	A	515	1/1	0.97	0.05	31,31,31,31	1
4	CD	C	509	1/1	0.97	0.05	40,40,40,40	1
4	CD	A	516	1/1	0.97	0.06	34,34,34,34	1
4	CD	C	510	1/1	0.98	0.03	25,25,25,25	1
4	CD	A	510	1/1	0.98	0.04	29,29,29,29	1
4	CD	C	512	1/1	0.98	0.06	35,35,35,35	1
4	CD	C	514	1/1	0.98	0.07	46,46,46,46	1
4	CD	A	511	1/1	0.98	0.06	43,43,43,43	1
4	CD	G	511	1/1	0.98	0.04	25,25,25,25	1
4	CD	C	516	1/1	0.98	0.06	48,48,48,48	1
4	CD	D	508	1/1	0.98	0.03	26,26,26,26	0
4	CD	C	507	1/1	0.99	0.03	24,24,24,24	1
4	CD	C	513	1/1	0.99	0.05	27,27,27,27	1
4	CD	C	508	1/1	0.99	0.03	29,29,29,29	0
4	CD	A	507	1/1	0.99	0.03	28,28,28,28	0
4	CD	A	513	1/1	0.99	0.05	32,32,32,32	1
4	CD	A	509	1/1	0.99	0.02	23,23,23,23	1
4	CD	D	509	1/1	0.99	0.05	31,31,31,31	1
4	CD	D	510	1/1	0.99	0.05	28,28,28,28	1
4	CD	D	501	1/1	1.00	0.04	24,24,24,24	1
4	CD	G	507	1/1	1.00	0.03	22,22,22,22	1

6.5 Other polymers [i](#)

There are no such residues in this entry.