



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:30 PM UTC

PDB ID : 7LFR / pdb_00007lfr
Title : Crystal structure of the epidermal growth factor receptor extracellular region with R84K mutation in complex with epiregulin crystallized with spermine
Authors : Hu, C.; Leche II, C.A.; Stayrook, S.E.; Ferguson, K.M.; Lemmon, M.A.
Deposited on : 2021-01-18
Resolution : 3.20 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

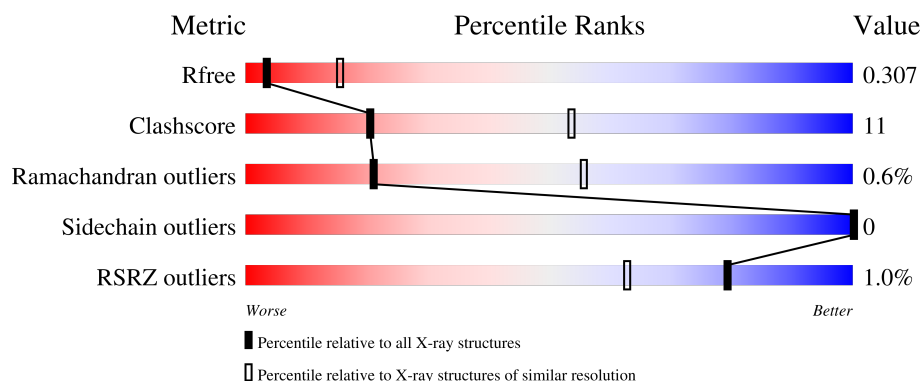
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	502	2% 72% 28%
1	B	502	83% 16% .
2	C	48	4% 75% 23% .
2	D	48	81% 17% .
3	F	2	100%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BMA	F	2	-	-	X	-
4	NAG	A	601	-	-	X	-
5	MAN	A	602	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8418 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 Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	500	Total	C	N	O	S	0	0	0
			3773	2338	664	729	42			
1	B	499	Total	C	N	O	S	0	0	0
			3785	2345	661	737	42			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	84	LYS	ARG	engineered mutation	UNP P00533
A	502	HIS	-	expression tag	UNP P00533
B	84	LYS	ARG	engineered mutation	UNP P00533
B	502	HIS	-	expression tag	UNP P00533

- Molecule 2 is a protein called Proepiregulin.

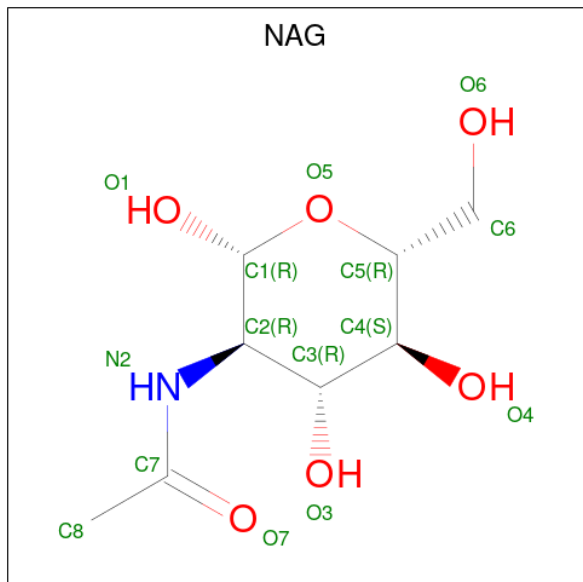
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	47	Total	C	N	O	S	0	0	0
			362	222	62	70	8			
2	D	47	Total	C	N	O	S	0	0	0
			370	230	62	70	8			

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



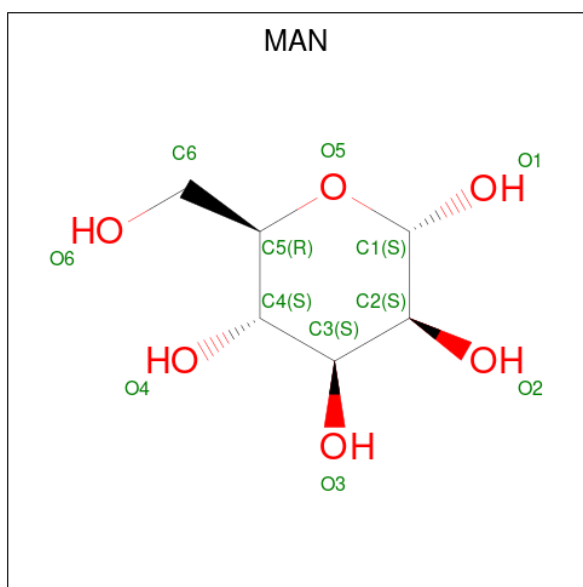
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	2	Total	C	N	O	0	0	0
			25	14	1	10			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



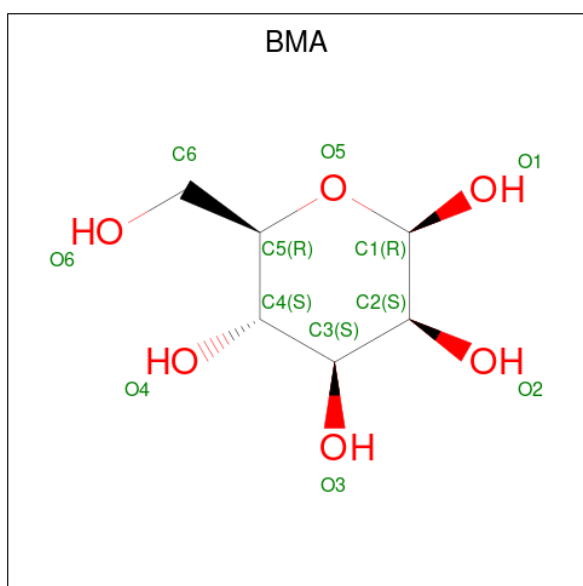
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is alpha-D-mannopyranose (CCD ID: MAN) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			11	6	5		
5	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 6 is beta-D-mannopyranose (CCD ID: BMA) (formula: $C_6H_{12}O_6$).

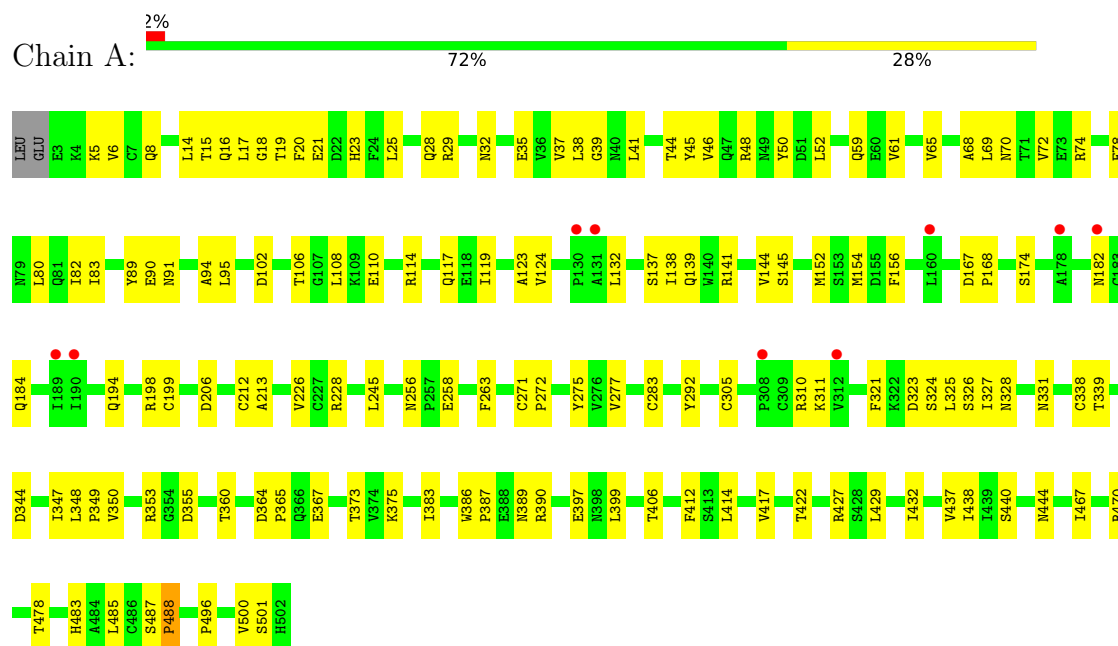


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			11	6	5		

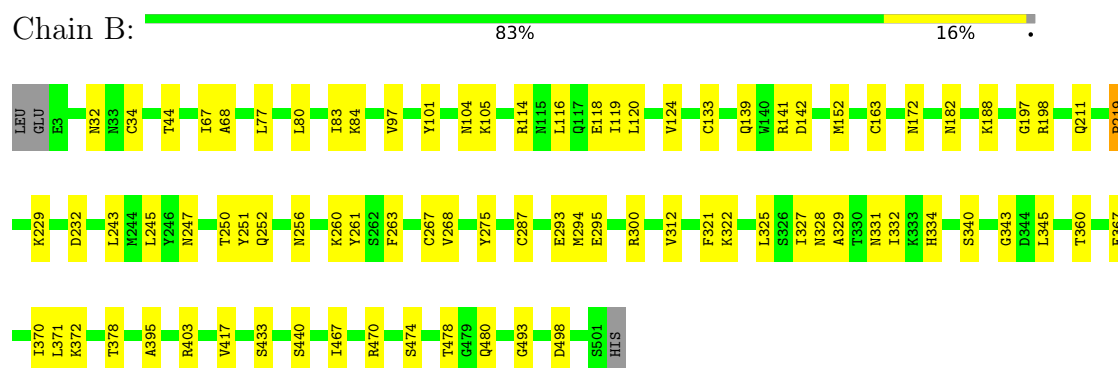
3 Residue-property plots [i](#)

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

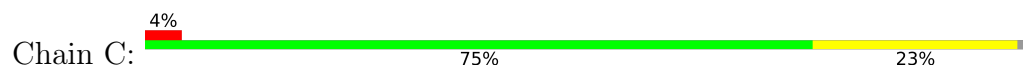
• Molecule 1: Epidermal growth factor receptor



• Molecule 1: Epidermal growth factor receptor

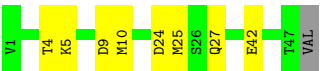
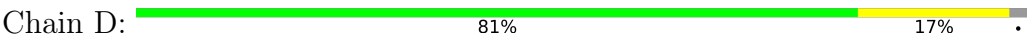


• Molecule 2: Proepiregulin

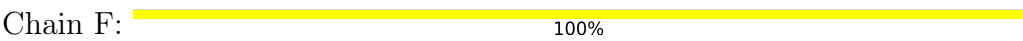




● Molecule 2: Proepiregulin



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.59Å 87.21Å 198.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.60 – 3.20 43.60 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.60-3.20) 99.8 (43.60-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.250 , 0.303 0.262 , 0.307	Depositor DCC
R_{free} test set	1115 reflections (4.01%)	wwPDB-VP
Wilson B-factor (Å ²)	96.5	Xtriage
Anisotropy	0.066	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8418	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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, MAN, BMA

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	0.16	0/3844	0.40	0/5216
1	B	0.14	0/3856	0.39	0/5229
2	C	0.21	0/368	0.46	0/495
2	D	0.14	0/377	0.41	0/507
All	All	0.15	0/8445	0.40	0/11447

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3773	0	3569	115	0
1	B	3785	0	3592	52	0
2	C	362	0	329	11	0
2	D	370	0	343	5	0
3	F	25	0	22	10	0
4	A	14	0	13	23	0
4	B	56	0	52	2	0
5	A	11	0	10	10	0
5	B	11	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	11	0	10	1	0
All	All	8418	0	7950	184	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 11.

All (184) 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:331:ASN:HD22	4:A:601:NAG:C6	1.04	1.60
1:A:331:ASN:ND2	4:A:601:NAG:C6	1.90	1.31
1:A:328:ASN:HD21	4:A:601:NAG:C1	1.46	1.29
1:A:331:ASN:ND2	4:A:601:NAG:O5	1.71	1.23
1:A:331:ASN:ND2	4:A:601:NAG:H62	1.48	1.21
1:A:328:ASN:ND2	4:A:601:NAG:C1	2.05	1.16
1:A:323:ASP:OD2	3:F:2:BMA:H62	1.46	1.15
1:A:328:ASN:OD1	4:A:601:NAG:C1	2.01	1.08
4:A:601:NAG:O4	3:F:1:NAG:C1	2.02	1.08
5:A:602:MAN:C1	3:F:2:BMA:O4	2.03	1.06
1:A:331:ASN:HD22	4:A:601:NAG:C5	1.72	1.03
1:A:328:ASN:CG	4:A:601:NAG:C1	2.31	1.03
1:A:331:ASN:ND2	4:A:601:NAG:C5	2.23	1.01
1:A:331:ASN:HD22	4:A:601:NAG:H62	0.81	0.97
5:A:602:MAN:C1	3:F:2:BMA:H61	1.95	0.95
5:A:602:MAN:C1	3:F:2:BMA:C4	2.56	0.84
1:A:91:ASN:HD21	5:A:602:MAN:H61	1.42	0.83
1:A:95:LEU:HB3	1:A:124:VAL:HG12	1.62	0.82
1:B:331:ASN:HB3	4:B:603:NAG:H62	1.63	0.80
1:A:16:GLN:NE2	1:A:20:PHE:CD1	2.51	0.77
1:A:331:ASN:CG	4:A:601:NAG:H62	2.09	0.77
1:A:16:GLN:HE22	1:A:20:PHE:HD1	1.30	0.76
1:A:328:ASN:HD21	4:A:601:NAG:C2	1.98	0.75
5:A:602:MAN:C1	3:F:2:BMA:C6	2.66	0.73
1:A:331:ASN:HB3	4:A:601:NAG:H62	1.68	0.73
1:A:414:LEU:HB3	1:A:437:VAL:HG22	1.70	0.72
1:A:331:ASN:CB	4:A:601:NAG:H62	2.19	0.71
2:C:20:ILE:HD11	2:C:31:ARG:HB2	1.72	0.70
1:A:331:ASN:ND2	4:A:601:NAG:O6	2.25	0.70
1:B:332:ILE:HG13	1:B:370:ILE:HD11	1.73	0.69
1:B:327:ILE:HD11	1:B:345:LEU:HD22	1.75	0.69
1:B:114:ARG:NH1	1:B:182:ASN:OD1	2.26	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:ASP:HB3	1:A:106:THR:H	1.60	0.67
1:A:91:ASN:ND2	5:A:602:MAN:H61	2.08	0.66
1:B:260:LYS:HG2	1:B:267:CYS:HB3	1.78	0.65
1:A:16:GLN:NE2	1:A:20:PHE:HD1	1.93	0.64
1:A:41:LEU:HB3	1:A:65:VAL:HG22	1.78	0.63
1:A:292:TYR:CD2	1:A:305:CYS:HB3	2.33	0.63
1:B:294:MET:O	1:B:300:ARG:HA	1.99	0.63
1:B:312:VAL:HG22	1:B:340:SER:HB2	1.79	0.63
1:A:326:SER:HB2	1:A:348:LEU:HB2	1.81	0.62
1:A:328:ASN:OD1	4:A:601:NAG:O5	2.17	0.62
1:A:91:ASN:ND2	5:A:602:MAN:C6	2.62	0.62
1:A:323:ASP:OD2	3:F:2:BMA:C6	2.36	0.62
1:B:141:ARG:HD2	1:B:141:ARG:H	1.62	0.62
1:B:232:ASP:HA	1:B:267:CYS:HB2	1.80	0.62
1:A:272:PRO:HB2	1:A:275:TYR:HD1	1.63	0.62
1:A:213:ALA:HB3	1:A:226:VAL:HB	1.81	0.61
1:B:478:THR:HG23	1:B:480:GLN:OE1	1.99	0.61
1:A:194:GLN:HG3	1:B:219:PRO:HB2	1.81	0.61
1:B:250:THR:O	1:B:252:GLN:N	2.31	0.61
1:B:77:LEU:HD22	1:B:80:LEU:HD12	1.82	0.60
1:A:83:ILE:HB	1:A:119:ILE:HG12	1.83	0.60
1:A:470:ARG:HH22	1:A:478:THR:HG21	1.66	0.60
1:B:372:LYS:HG2	1:B:395:ALA:HA	1.83	0.60
1:A:78:GLU:OE1	1:A:114:ARG:NH2	2.35	0.60
1:B:232:ASP:OD1	1:B:260:LYS:NZ	2.33	0.60
2:D:4:THR:HG22	2:D:5:LYS:H	1.67	0.59
1:A:485:LEU:HD11	1:A:496:PRO:HB3	1.85	0.58
1:A:117:GLN:O	1:A:198:ARG:NH1	2.37	0.58
1:B:295:GLU:HG2	1:B:300:ARG:HG2	1.86	0.57
1:A:108:LEU:HG	1:A:132:LEU:HD13	1.85	0.57
1:A:440:SER:HA	1:A:467:ILE:O	2.04	0.57
1:A:74:ARG:HG3	1:A:110:GLU:HB2	1.86	0.56
1:A:95:LEU:HB3	1:A:124:VAL:CG1	2.32	0.56
1:B:470:ARG:NH1	1:B:474:SER:OG	2.38	0.56
1:A:21:GLU:OE2	1:A:50:TYR:OH	2.20	0.56
1:A:69:LEU:HD11	2:C:22:LEU:HD21	1.86	0.56
1:A:16:GLN:NE2	1:A:18:GLY:O	2.38	0.56
4:A:601:NAG:HO4	3:F:1:NAG:C1	2.16	0.56
1:A:487:SER:N	1:A:488:PRO:HD3	2.21	0.56
1:A:364:ASP:HB3	1:A:367:GLU:HG3	1.86	0.56
1:A:344:ASP:OD1	1:A:406:THR:OG1	2.24	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:602:MAN:C1	3:F:2:BMA:H4	2.35	0.55
1:B:367:GLU:O	1:B:370:ILE:HG12	2.07	0.55
5:A:602:MAN:C1	3:F:2:BMA:HO4	2.20	0.54
1:B:378:THR:HA	1:B:403:ARG:HB2	1.89	0.54
1:B:141:ARG:HD2	1:B:141:ARG:N	2.23	0.54
1:A:375:LYS:HA	1:A:399:LEU:HA	1.89	0.53
1:B:440:SER:HA	1:B:467:ILE:O	2.08	0.53
1:A:347:ILE:HD12	1:A:383:ILE:HG12	1.91	0.53
1:B:118:GLU:OE2	1:B:198:ARG:NH1	2.41	0.53
1:A:17:LEU:HD22	2:C:36:TYR:O	2.11	0.51
1:A:321:PHE:HE1	1:A:331:ASN:HB2	1.76	0.50
1:B:211:GLN:HE22	1:B:229:LYS:HD2	1.75	0.50
1:B:44:THR:HA	1:B:68:ALA:O	2.12	0.49
1:A:19:THR:HG23	1:A:21:GLU:HB2	1.94	0.49
1:B:293:GLU:OE2	1:B:403:ARG:NH1	2.45	0.49
1:A:137:SER:HB2	1:A:174:SER:HA	1.94	0.49
1:A:5:LYS:N	1:A:35:GLU:OE2	2.40	0.48
1:A:327:ILE:HD11	1:A:347:ILE:HG12	1.96	0.48
2:C:20:ILE:HD11	2:C:31:ARG:CB	2.39	0.48
1:A:310:ARG:HG2	1:A:311:LYS:H	1.77	0.48
1:B:371:LEU:HB2	1:B:395:ALA:HB1	1.95	0.48
1:A:331:ASN:HB3	4:A:601:NAG:C6	2.42	0.48
2:C:16:HIS:H	2:C:41:CYS:HB2	1.78	0.48
2:C:4:THR:OG1	2:C:21:TYR:HB3	2.14	0.48
1:A:14:LEU:HD13	2:C:22:LEU:HD22	1.95	0.48
1:A:256:ASN:OD1	1:A:258:GLU:N	2.47	0.48
1:A:483:HIS:CD2	1:A:496:PRO:HD3	2.48	0.48
1:A:328:ASN:HD21	4:A:601:NAG:HN2	1.62	0.47
1:A:198:ARG:HD3	1:A:213:ALA:O	2.14	0.47
1:A:292:TYR:HD2	1:A:305:CYS:HB3	1.79	0.47
1:A:139:GLN:HG3	1:A:141:ARG:NH2	2.29	0.47
1:B:263:PHE:HE1	1:B:275:TYR:CE2	2.33	0.47
1:A:154:MET:O	1:A:154:MET:HG3	2.14	0.47
1:A:29:ARG:CZ	2:C:46:LEU:HD21	2.45	0.47
1:A:386:TRP:CG	1:A:387:PRO:HD2	2.50	0.46
1:A:6:VAL:HG13	1:A:38:LEU:HD23	1.97	0.46
1:B:325:LEU:HD22	2:D:42:GLU:HG2	1.96	0.46
1:B:328:ASN:ND2	1:B:360:THR:HG21	2.31	0.46
2:D:9:ASP:OD1	2:D:10:MET:HG2	2.15	0.46
1:A:91:ASN:ND2	5:A:602:MAN:O6	2.48	0.46
1:A:167:ASP:CG	1:A:168:PRO:HD2	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:ASN:O	1:B:105:LYS:HG3	2.15	0.46
1:B:142:ASP:HA	1:B:188:LYS:HD2	1.99	0.46
1:A:82:ILE:HG21	1:A:226:VAL:HG11	1.98	0.45
2:C:21:TYR:CE1	2:C:23:VAL:HA	2.51	0.45
1:A:412:PHE:CE1	1:A:438:ILE:HB	2.51	0.45
1:A:500:VAL:HG13	1:A:501:SER:N	2.31	0.45
1:B:84:LYS:HG2	1:B:120:LEU:HD12	1.98	0.45
1:B:245:LEU:HG	1:B:256:ASN:HB2	1.98	0.45
1:A:23:HIS:CD2	1:A:45:TYR:HB2	2.51	0.45
1:A:325:LEU:HB2	4:A:601:NAG:O7	2.17	0.45
1:A:37:VAL:HB	1:A:61:VAL:HG22	1.99	0.45
1:A:46:VAL:HG11	1:A:52:LEU:HD11	1.99	0.45
2:C:21:TYR:HE1	2:C:23:VAL:HA	1.82	0.44
1:B:417:VAL:HA	1:B:440:SER:O	2.18	0.44
1:A:212:CYS:O	1:A:228:ARG:NH2	2.51	0.44
1:A:94:ALA:N	1:A:123:ALA:O	2.45	0.44
1:B:83:ILE:HB	1:B:119:ILE:HG12	2.00	0.44
1:A:124:VAL:CG2	1:A:152:MET:HE1	2.47	0.44
1:B:287:CYS:SG	1:B:293:GLU:HG2	2.58	0.44
1:A:199:CYS:HB2	1:A:206:ASP:O	2.18	0.43
1:B:139:GLN:HE22	1:B:172:ASN:HD22	1.66	0.43
1:A:48:ARG:HA	1:A:72:VAL:HA	2.00	0.43
1:A:328:ASN:ND2	1:A:360:THR:HG21	2.33	0.43
1:A:28:GLN:O	1:A:32:ASN:HB2	2.18	0.43
1:B:67:ILE:O	1:B:97:VAL:HA	2.18	0.43
1:A:397:GLU:HB2	1:A:427:ARG:NH1	2.34	0.43
1:A:365:PRO:HB3	1:A:387:PRO:HB3	2.01	0.43
1:B:80:LEU:HD22	1:B:116:LEU:HD13	2.01	0.43
1:A:338:CYS:O	1:A:373:THR:HG23	2.19	0.43
1:B:403:ARG:O	1:B:433:SER:HB2	2.19	0.43
1:A:429:LEU:HD21	1:A:432:ILE:HD11	2.01	0.42
1:B:32:ASN:C	1:B:34:CYS:H	2.27	0.42
1:A:144:VAL:HG12	1:A:145:SER:O	2.19	0.42
1:A:389:ASN:OD1	1:A:389:ASN:N	2.51	0.42
1:A:349:PRO:O	1:A:353:ARG:HG2	2.20	0.42
1:A:386:TRP:CD2	1:A:387:PRO:HD2	2.54	0.42
1:A:422:THR:HA	1:A:444:ASN:O	2.19	0.42
1:B:101:TYR:OH	2:D:24:ASP:OD2	2.32	0.42
1:B:343:GLY:N	1:B:378:THR:OG1	2.52	0.42
1:A:70:ASN:HB3	1:A:72:VAL:HG12	2.01	0.42
1:B:260:LYS:HE2	1:B:268:VAL:C	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:LEU:HD11	1:A:256:ASN:HD22	1.84	0.42
1:A:59:GLN:HA	1:A:80:LEU:HA	2.02	0.42
1:A:387:PRO:HG2	1:A:390:ARG:HG3	2.02	0.42
1:A:15:THR:HG21	2:C:38:GLY:O	2.20	0.41
1:A:44:THR:HA	1:A:68:ALA:O	2.20	0.41
2:D:25:MET:O	2:D:27:GLN:HG3	2.20	0.41
1:A:21:GLU:O	1:A:25:LEU:HD23	2.20	0.41
1:A:95:LEU:O	1:A:124:VAL:HA	2.20	0.41
1:A:8:GLN:O	1:A:39:GLY:N	2.52	0.41
1:A:263:PHE:HB2	1:A:283:CYS:SG	2.60	0.41
1:B:124:VAL:HB	1:B:152:MET:HE3	2.02	0.41
1:B:243:LEU:HD21	1:B:261:TYR:CE1	2.55	0.41
1:A:310:ARG:HG2	1:A:311:LYS:N	2.36	0.41
1:A:310:ARG:HD2	1:A:339:THR:H	1.85	0.41
1:B:133:CYS:HB3	1:B:163:CYS:HB3	1.83	0.41
1:B:247:ASN:HB3	1:B:250:THR:O	2.20	0.41
1:A:114:ARG:HB3	1:A:182:ASN:HA	2.03	0.41
1:A:321:PHE:CE1	1:A:331:ASN:HB2	2.56	0.41
1:A:324:SER:OG	4:A:601:NAG:H4	2.21	0.41
4:B:604:NAG:O4	6:B:605:BMA:O5	2.27	0.41
1:A:271:CYS:SG	1:A:277:VAL:HG22	2.60	0.40
1:A:350:VAL:CG2	1:A:355:ASP:HB2	2.51	0.40
1:A:138:ILE:HA	1:A:184:GLN:OE1	2.22	0.40
1:A:89:TYR:CD2	1:A:90:GLU:HG2	2.55	0.40
1:B:142:ASP:OD1	1:B:198:ARG:NH2	2.54	0.40
1:B:321:PHE:HZ	1:B:334:HIS:HB2	1.86	0.40
1:B:493:GLY:N	1:B:498:ASP:HB3	2.36	0.40
1:B:328:ASN:OD1	1:B:329:ALA:N	2.50	0.40
1:A:417:VAL:HA	1:A:440:SER:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	498/502 (99%)	457 (92%)	39 (8%)	2 (0%)	30	62
1	B	497/502 (99%)	461 (93%)	32 (6%)	4 (1%)	16	50
2	C	45/48 (94%)	42 (93%)	3 (7%)	0	100	100
2	D	45/48 (94%)	42 (93%)	3 (7%)	0	100	100
All	All	1085/1100 (99%)	1002 (92%)	77 (7%)	6 (1%)	21	56

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	251	TYR
1	A	156	PHE
1	B	219	PRO
1	B	322	LYS
1	B	197	GLY
1	A	488	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	415/441 (94%)	415 (100%)	0	100	100
1	B	421/441 (96%)	421 (100%)	0	100	100
2	C	41/44 (93%)	41 (100%)	0	100	100
2	D	43/44 (98%)	43 (100%)	0	100	100
All	All	920/970 (95%)	920 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	79	ASN

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Mol	Chain	Res	Type
1	A	91	ASN
1	A	134	ASN
1	A	211	GLN
1	A	252	GLN
1	A	328	ASN
1	A	331	ASN
1	A	409	HIS
1	B	172	ASN
1	B	211	GLN
1	B	247	ASN
1	B	398	ASN
1	B	409	HIS
2	C	16	HIS
2	C	27	GLN
2	D	27	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 ⓘ

2 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	F	1	3	14,14,15	0.41	0	17,19,21	0.62	0
3	BMA	F	2	3	11,11,12	0.21	0	15,15,17	0.69	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	F	1	3	-	1/6/23/26	0/1/1/1
3	BMA	F	2	3	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

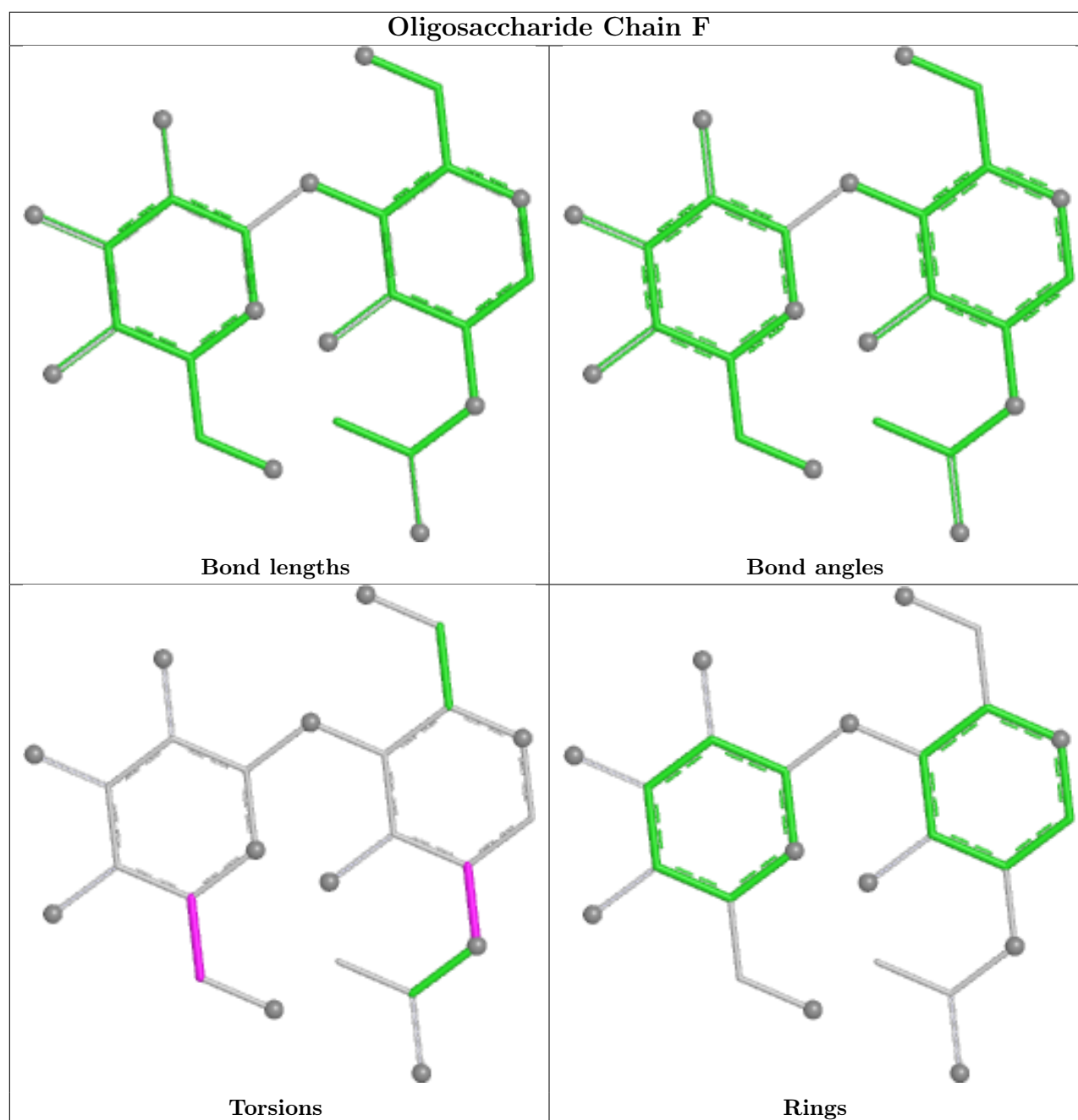
Mol	Chain	Res	Type	Atoms
3	F	2	BMA	O5-C5-C6-O6
3	F	2	BMA	C4-C5-C6-O6
3	F	1	NAG	C1-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1	NAG	2	0
3	F	2	BMA	8	0

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



5.6 Ligand geometry [i](#)

8 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	NAG	B	601	1	14,14,15	0.23	0	17,19,21	0.53	0
5	MAN	B	606	-	11,11,12	0.66	0	15,15,17	1.05	2 (13%)
6	BMA	B	605	-	11,11,12	0.59	0	15,15,17	0.70	0
4	NAG	B	602	1	14,14,15	0.27	0	17,19,21	0.43	0
4	NAG	B	604	-	14,14,15	0.28	0	17,19,21	0.40	0
4	NAG	A	601	-	14,14,15	0.54	0	17,19,21	0.53	0
5	MAN	A	602	-	11,11,12	0.70	0	15,15,17	0.96	2 (13%)
4	NAG	B	603	-	14,14,15	0.23	0	17,19,21	0.47	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	NAG	B	601	1	-	2/6/23/26	0/1/1/1
5	MAN	B	606	-	-	0/2/19/22	0/1/1/1
6	BMA	B	605	-	-	1/2/19/22	0/1/1/1
4	NAG	B	602	1	-	1/6/23/26	0/1/1/1
4	NAG	B	604	-	-	0/6/23/26	0/1/1/1
4	NAG	A	601	-	-	0/6/23/26	0/1/1/1
5	MAN	A	602	-	-	1/2/19/22	0/1/1/1
4	NAG	B	603	-	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	606	MAN	C1-O5-C5	2.49	115.53	112.19
5	A	602	MAN	C1-O5-C5	2.42	115.43	112.19
5	A	602	MAN	O2-C2-C3	-2.18	105.63	110.15
5	B	606	MAN	O2-C2-C3	-2.11	105.78	110.15

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	601	NAG	O5-C5-C6-O6
4	B	601	NAG	C4-C5-C6-O6
4	B	602	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	A	602	MAN	O5-C5-C6-O6
4	B	603	NAG	C1-C2-N2-C7
6	B	605	BMA	O5-C5-C6-O6

There are no ring outliers.

5 monomers are involved in 35 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	605	BMA	1	0
4	B	604	NAG	1	0
4	A	601	NAG	23	0
5	A	602	MAN	10	0
4	B	603	NAG	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	500/502 (99%)	0.26	9 (1%) 67 48	57, 100, 138, 160	0
1	B	499/502 (99%)	0.04	0 100 100	54, 82, 114, 139	0
2	C	47/48 (97%)	0.46	2 (4%) 40 24	78, 99, 119, 128	0
2	D	47/48 (97%)	0.09	0 100 100	56, 71, 87, 102	0
All	All	1093/1100 (99%)	0.16	11 (1%) 79 63	54, 90, 130, 160	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	190	ILE	2.6
2	C	47	THR	2.6
1	A	130	PRO	2.4
1	A	160	LEU	2.4
1	A	308	PRO	2.3
1	A	189	ILE	2.3
1	A	178	ALA	2.2
1	A	312	VAL	2.1
1	A	182	ASN	2.1
1	A	131	ALA	2.0
2	C	45	PHE	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.

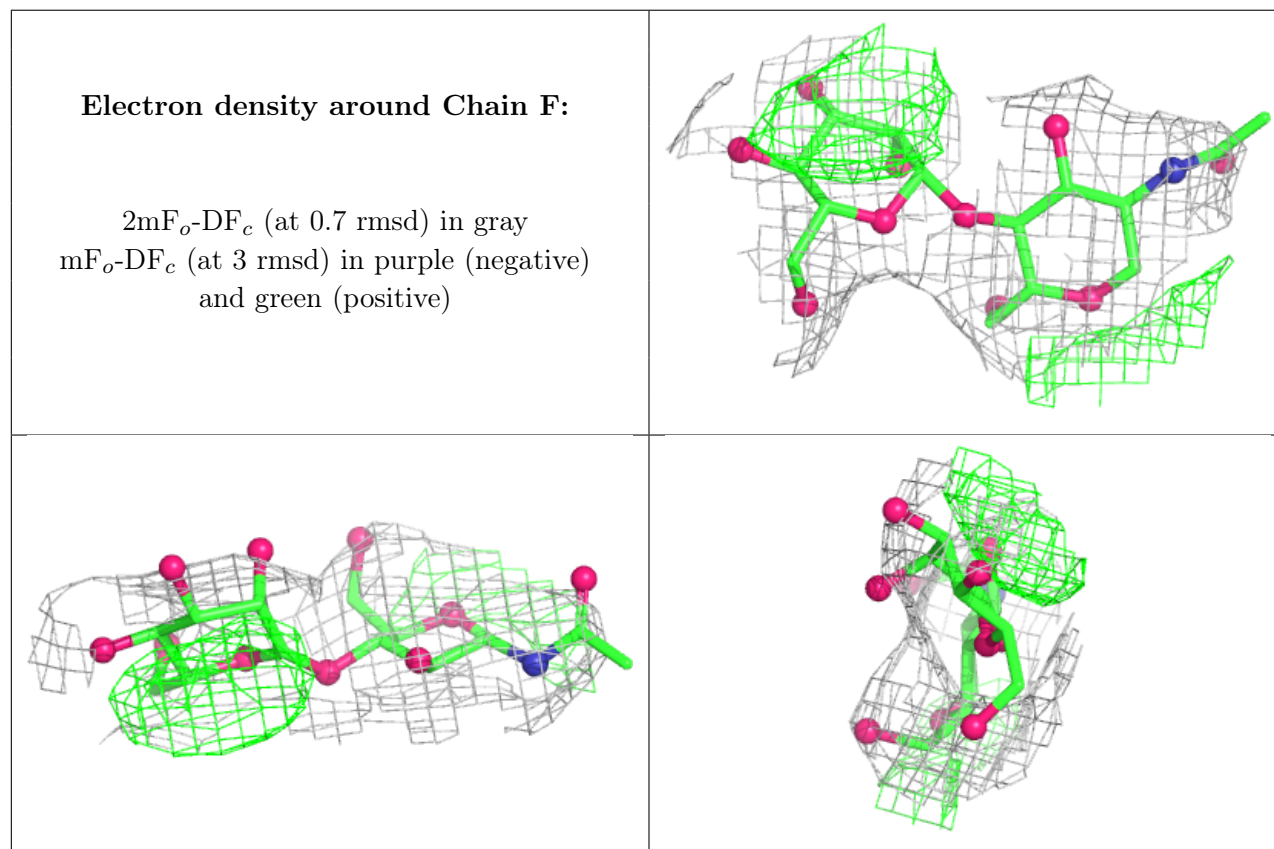
6.3 Carbohydrates [i](#)

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	NAG	F	1	14/15	-	-	90,104,115,121	0
3	BMA	F	2	11/12	-	-	117,124,130,149	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.



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
6	BMA	B	605	11/12	0.44	0.12	96,117,121,122	0
5	MAN	A	602	11/12	0.47	0.13	102,114,122,130	0
4	NAG	B	602	14/15	0.59	0.12	85,115,122,127	0
5	MAN	B	606	11/12	0.64	0.10	109,117,123,126	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	A	601	14/15	0.69	0.21	69,83,95,97	0
4	NAG	B	601	14/15	0.72	0.10	96,110,120,122	0
4	NAG	B	604	14/15	0.80	0.12	81,93,106,115	0
4	NAG	B	603	14/15	0.89	0.15	60,72,88,94	0

6.5 Other polymers [i](#)

There are no such residues in this entry.