



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 25, 2026 – 05:30 PM EDT

PDB ID : 3U2S / pdb_00003u2s
Title : Crystal Structure of PG9 Fab in Complex with V1V2 Region from HIV-1 strain ZM109
Authors : McLellan, J.S.; Pancera, M.; Kwong, P.D.
Deposited on : 2011-10-04
Resolution : 1.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

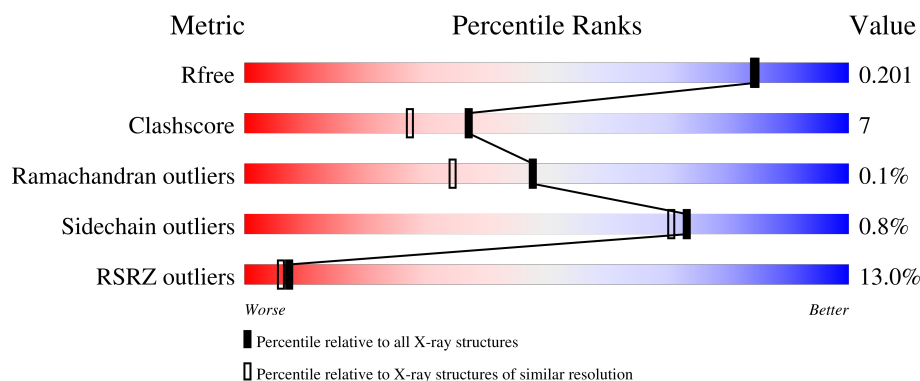
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	248	
1	H	248	
2	B	216	
2	L	216	
3	C	124	

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Mol	Chain	Length	Quality of chain
3	G	124	
4	D	7	
4	E	7	

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 16931 atoms, of which 7771 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 PG9 heavy chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	H	230	Total	C	H	N	O	S	0	0	0
			3447	1122	1667	298	351	9			
1	A	227	Total	C	H	N	O	S	0	0	0
			3415	1113	1652	295	346	9			

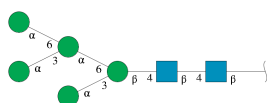
- Molecule 2 is a protein called PG9 light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	L	211	Total	C	H	N	O	S	0	0	0
			3103	974	1537	267	321	4			
2	B	211	Total	C	H	N	O	S	0	0	0
			3103	974	1537	267	321	4			

- Molecule 3 is a protein called Envelope glycoprotein gp120.

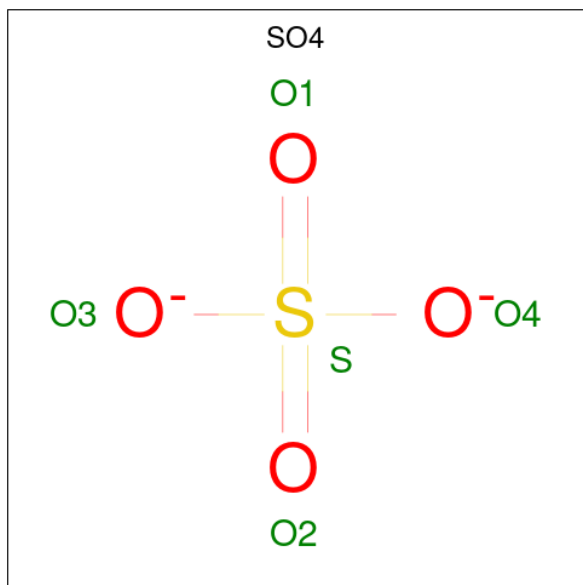
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	G	90	Total	C	H	N	O	S	0	0	0
			1379	440	677	117	141	4			
3	C	90	Total	C	H	N	O	S	0	0	0
			1386	442	681	117	142	4			

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



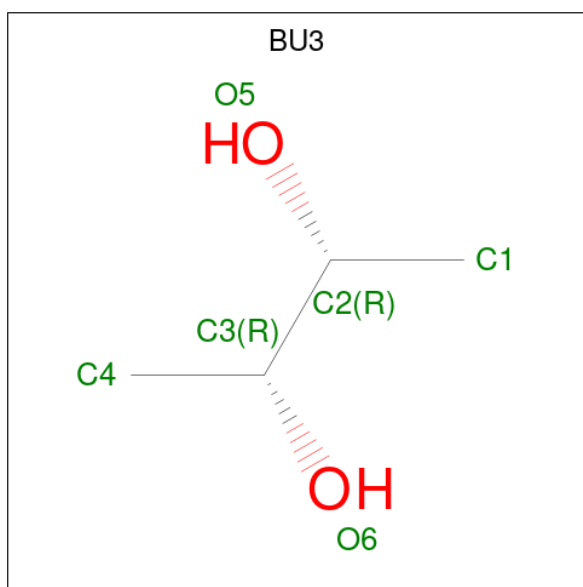
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	7	Total	C	N	O	0	0	0
			83	46	2	35			
4	E	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



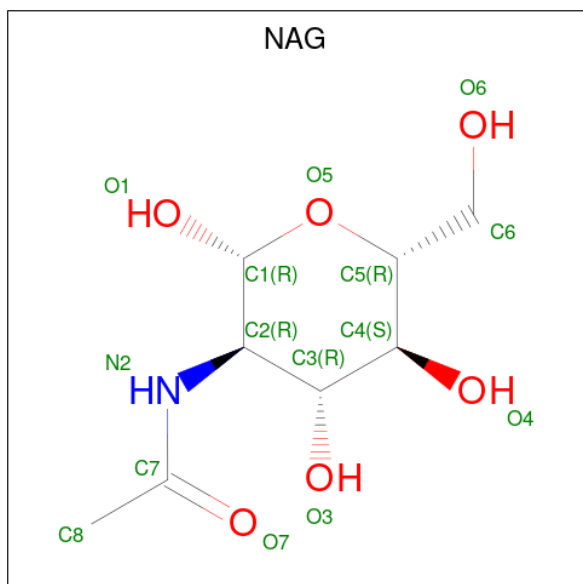
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	H	1	Total	O	S	0	0
			5	4	1		
5	L	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is (R,R)-2,3-BUTANEDIOL (CCD ID: BU3) (formula: C₄H₁₀O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	L	1	Total	C	H	O	0	0
			16	4	10	2		
6	B	1	Total	C	H	O	0	0
			16	4	10	2		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	G	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	1	Total	C	N	O	0	0
			14	8	1	5		

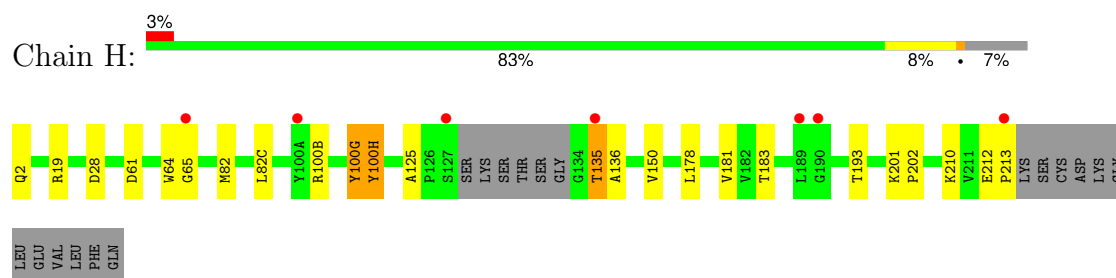
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	H	232	Total	O	0	0
			232	232		
8	L	186	Total	O	0	0
			186	186		
8	G	38	Total	O	0	0
			38	38		
8	A	203	Total	O	0	0
			203	203		
8	B	154	Total	O	0	0
			154	154		
8	C	34	Total	O	0	0
			34	34		

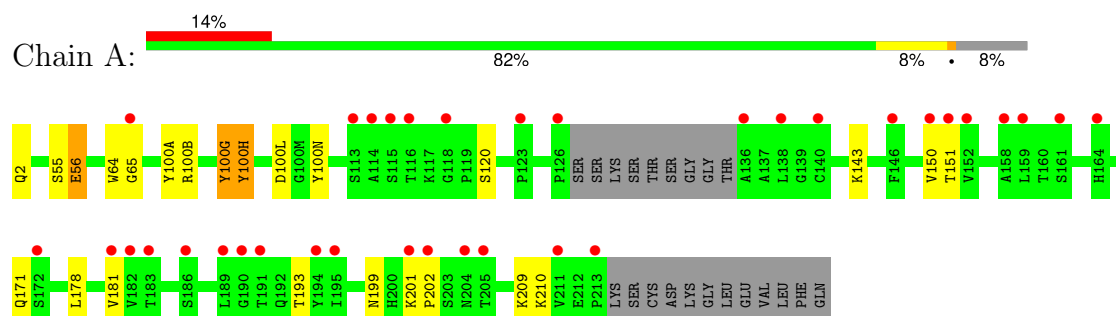
3 Residue-property plots [i](#)

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.

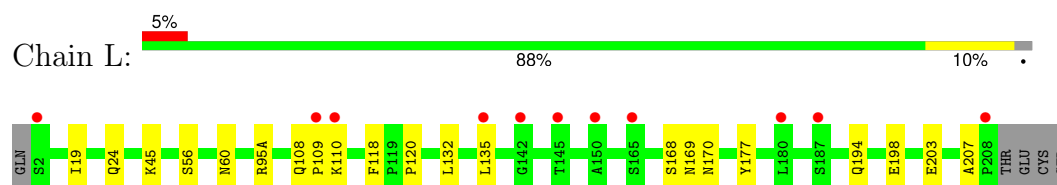
- Molecule 1: PG9 heavy chain



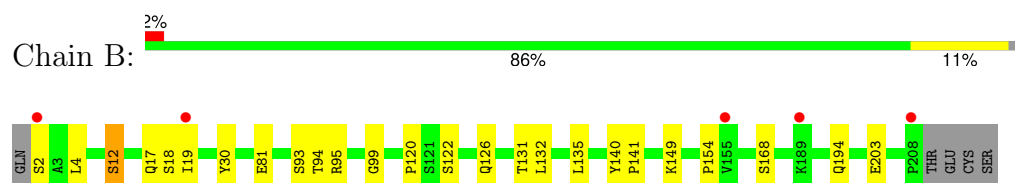
- Molecule 1: PG9 heavy chain



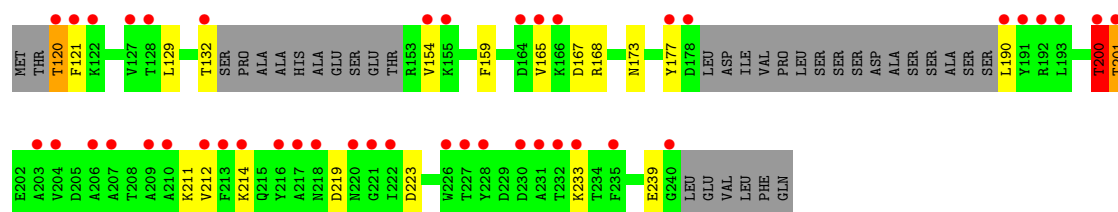
- Molecule 2: PG9 light chain



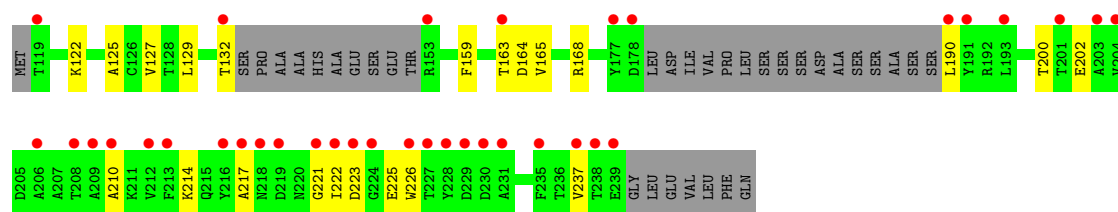
- Molecule 2: PG9 light chain



- Molecule 3: Envelope glycoprotein gp120



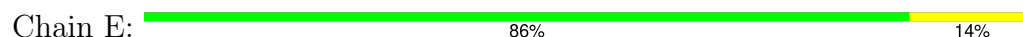
• Molecule 3: Envelope glycoprotein gp120



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.47Å 86.56Å 94.88Å 90.00° 92.08° 90.00°	Depositor
Resolution (Å)	41.58 – 1.80 41.58 – 1.80	Depositor EDS
% Data completeness (in resolution range)	91.0 (41.58-1.80) 90.9 (41.58-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 1.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_755)	Depositor
R, R_{free}	0.178 , 0.205 0.174 , 0.201	Depositor DCC
R_{free} test set	6169 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.533	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for k,h,-l 0.007 for -k,-h,-l 0.025 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16931	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, SO4, BU3, BMA, PCA, NAG, TYS

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	0.42	0/1770	0.71	0/2409
1	H	0.44	0/1787	0.70	0/2432
2	B	0.43	0/1601	0.70	0/2180
2	L	0.46	0/1601	0.70	0/2180
3	C	0.31	0/714	0.58	0/967
3	G	0.31	0/711	0.71	2/962 (0.2%)
All	All	0.42	0/8184	0.69	2/11130 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	G	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	201	THR	N-CA-C	5.85	117.75	108.79
3	G	200	THR	N-CA-C	5.57	118.63	109.72

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	G	200	THR	Peptide

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	1763	1652	1663	23	0
1	H	1780	1667	1678	22	0
2	B	1566	1537	1533	19	0
2	L	1566	1537	1533	24	0
3	C	705	681	680	20	0
3	G	702	677	676	17	0
4	D	83	0	70	0	0
4	E	83	0	70	0	0
5	B	15	0	0	0	0
5	H	5	0	0	0	0
5	L	5	0	0	1	0
6	B	6	10	10	0	0
6	L	6	10	10	0	0
7	C	14	0	13	0	0
7	G	14	0	13	0	0
8	A	203	0	0	8	0
8	B	154	0	0	5	0
8	C	34	0	0	2	0
8	G	38	0	0	6	0
8	H	232	0	0	5	1
8	L	186	0	0	11	1
All	All	9160	7771	7949	116	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

The worst 5 of 116 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:H:181:VAL:HG11	2:L:135:LEU:HD13	1.42	1.02
3:G:173:ASN:OD1	8:G:456:HOH:O	1.82	0.96
2:L:56:SER:HG	1:A:100(A):TYR:HH	1.15	0.92
1:A:171:GLN:NE2	8:A:377:HOH:O	2.10	0.84
2:L:110:LYS:NZ	8:L:239:HOH:O	2.10	0.83

All (1) 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)
8:H:338:HOH:O	8:L:277:HOH:O[2_645]	2.19	0.01

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	221/248 (89%)	217 (98%)	4 (2%)	0	100	100
1	H	224/248 (90%)	221 (99%)	3 (1%)	0	100	100
2	B	209/216 (97%)	204 (98%)	5 (2%)	0	100	100
2	L	209/216 (97%)	203 (97%)	6 (3%)	0	100	100
3	C	84/124 (68%)	82 (98%)	1 (1%)	1 (1%)	10	3
3	G	84/124 (68%)	82 (98%)	2 (2%)	0	100	100
All	All	1031/1176 (88%)	1009 (98%)	21 (2%)	1 (0%)	48	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	163	THR

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	189/207 (91%)	188 (100%)	1 (0%)	81	80
1	H	191/207 (92%)	190 (100%)	1 (0%)	81	80
2	B	178/183 (97%)	177 (99%)	1 (1%)	78	77
2	L	178/183 (97%)	178 (100%)	0	100	100
3	C	77/105 (73%)	76 (99%)	1 (1%)	61	54
3	G	76/105 (72%)	73 (96%)	3 (4%)	28	16
All	All	889/990 (90%)	882 (99%)	7 (1%)	73	70

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	G	214	LYS
1	A	56	GLU
3	C	127	VAL
2	B	12	SER
3	G	200	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	38	GLN
2	B	79	GLN
3	C	215	GLN
2	L	37	GLN
2	L	38	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	TYS	A	100(G)	-	15,16,17	1.12	1 (6%)	15,22,24	0.82	1 (6%)
1	TYS	H	100(H)	1	15,16,17	1.09	2 (13%)	15,22,24	0.49	0
1	PCA	A	2	1	7,8,9	2.19	2 (28%)	9,10,12	1.78	4 (44%)
1	TYS	A	100(H)	1	15,16,17	1.05	1 (6%)	15,22,24	0.62	0
1	PCA	H	2	1	7,8,9	2.14	2 (28%)	9,10,12	2.13	5 (55%)
1	TYS	H	100(G)	-	15,16,17	1.06	0	15,22,24	0.89	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TYS	A	100(G)	-	-	1/10/11/13	0/1/1/1
1	TYS	H	100(H)	1	-	3/10/11/13	0/1/1/1
1	PCA	A	2	1	-	0/0/11/13	0/1/1/1
1	TYS	A	100(H)	1	-	1/10/11/13	0/1/1/1
1	PCA	H	2	1	-	0/0/11/13	0/1/1/1
1	TYS	H	100(G)	-	-	2/10/11/13	0/1/1/1

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	2	PCA	CD-N	4.92	1.46	1.34
1	H	2	PCA	CD-N	4.61	1.45	1.34
1	H	2	PCA	CA-N	3.04	1.50	1.46
1	A	2	PCA	CA-N	2.84	1.49	1.46
1	A	100(H)	TYS	OH-S	2.30	1.62	1.58

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	2	PCA	CA-N-CD	-3.01	103.27	113.58
1	H	2	PCA	CG-CD-N	2.89	115.46	108.39
1	H	2	PCA	OE-CD-CG	-2.88	121.58	126.72
1	H	2	PCA	CB-CA-N	2.71	110.69	103.24
1	H	100(G)	TYS	OH-CZ-CE2	2.68	123.96	118.70

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	H	100(G)	TYS	CE1-CZ-OH-S
1	H	100(G)	TYS	CE2-CZ-OH-S
1	H	100(H)	TYS	CZ-OH-S-O1
1	H	100(H)	TYS	CZ-OH-S-O3
1	H	100(H)	TYS	CZ-OH-S-O2

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	100(G)	TYS	1	0
1	H	100(H)	TYS	2	0
1	A	100(H)	TYS	2	0
1	H	100(G)	TYS	1	0

5.5 Carbohydrates [i](#)

14 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	$\# Z > 2$
4	NAG	D	1	3,4	14,14,15	0.61	0	17,19,21	1.03	2 (11%)
4	NAG	D	2	4	14,14,15	0.65	0	17,19,21	0.89	0
4	BMA	D	3	4	11,11,12	0.76	0	15,15,17	0.95	1 (6%)
4	MAN	D	4	4	11,11,12	0.64	0	15,15,17	1.00	0
4	MAN	D	5	4	11,11,12	0.70	0	15,15,17	0.50	0
4	MAN	D	6	4	11,11,12	0.64	0	15,15,17	0.92	0
4	MAN	D	7	4	11,11,12	0.63	0	15,15,17	0.71	0
4	NAG	E	1	3,4	14,14,15	0.50	0	17,19,21	0.67	0
4	NAG	E	2	4	14,14,15	0.75	0	17,19,21	0.79	0
4	BMA	E	3	4	11,11,12	0.59	0	15,15,17	0.91	1 (6%)
4	MAN	E	4	4	11,11,12	0.76	0	15,15,17	0.92	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	E	5	4	11,11,12	0.69	0	15,15,17	0.63	0
4	MAN	E	6	4	11,11,12	0.62	0	15,15,17	0.95	0
4	MAN	E	7	4	11,11,12	0.63	0	15,15,17	0.61	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	D	1	3,4	-	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	0/6/23/26	0/1/1/1
4	BMA	D	3	4	-	0/2/19/22	0/1/1/1
4	MAN	D	4	4	-	0/2/19/22	0/1/1/1
4	MAN	D	5	4	-	0/2/19/22	0/1/1/1
4	MAN	D	6	4	-	0/2/19/22	0/1/1/1
4	MAN	D	7	4	-	2/2/19/22	0/1/1/1
4	NAG	E	1	3,4	-	0/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	0/2/19/22	0/1/1/1
4	MAN	E	4	4	-	0/2/19/22	0/1/1/1
4	MAN	E	5	4	-	0/2/19/22	0/1/1/1
4	MAN	E	6	4	-	0/2/19/22	0/1/1/1
4	MAN	E	7	4	-	0/2/19/22	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(°)
4	D	1	NAG	O5-C5-C6	2.35	112.23	107.66
4	D	1	NAG	C2-N2-C7	-2.24	119.90	122.90
4	D	3	BMA	C1-C2-C3	2.08	112.68	109.64
4	E	3	BMA	C1-C2-C3	2.03	112.60	109.64

There are no chirality outliers.

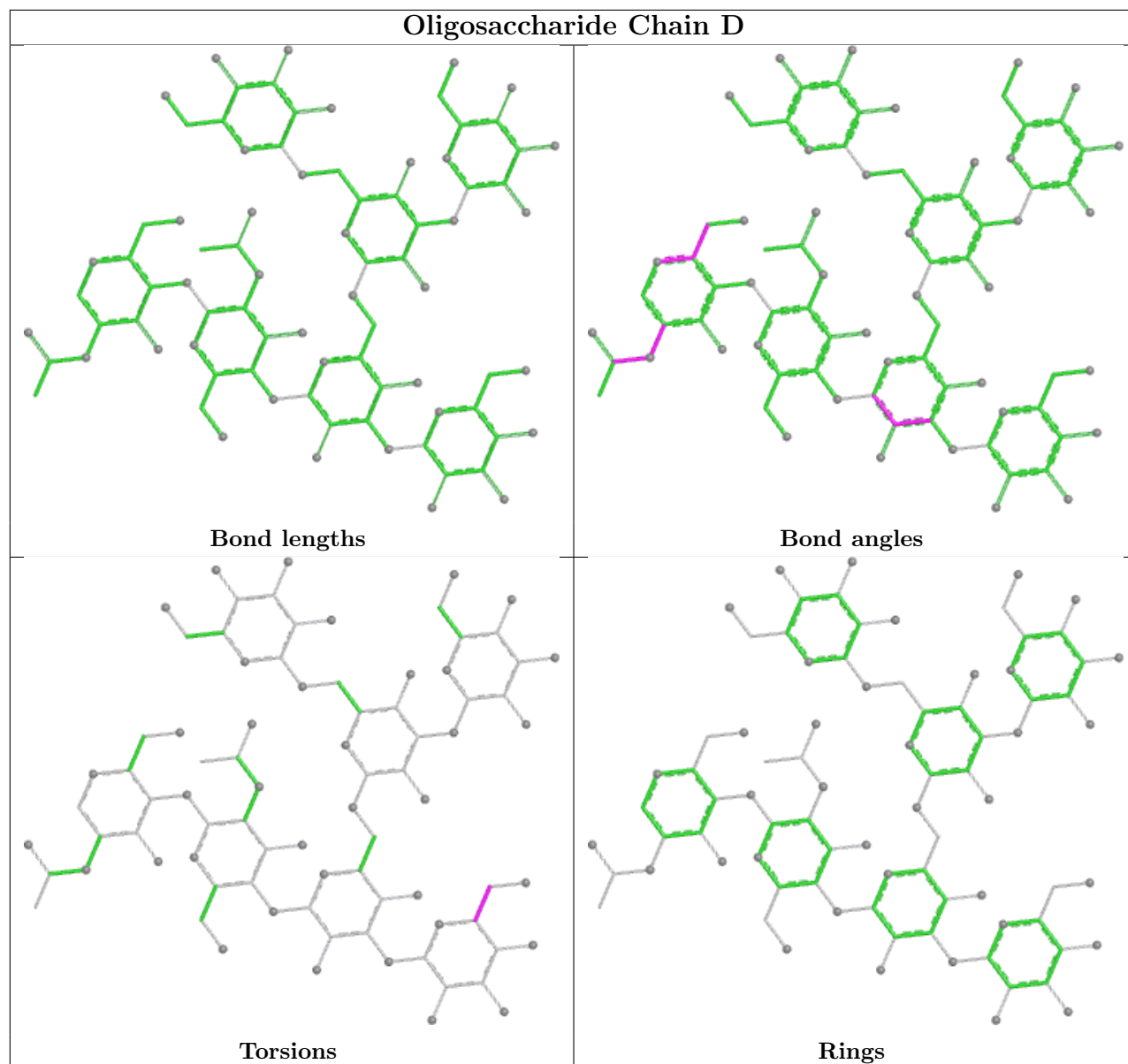
All (2) torsion outliers are listed below:

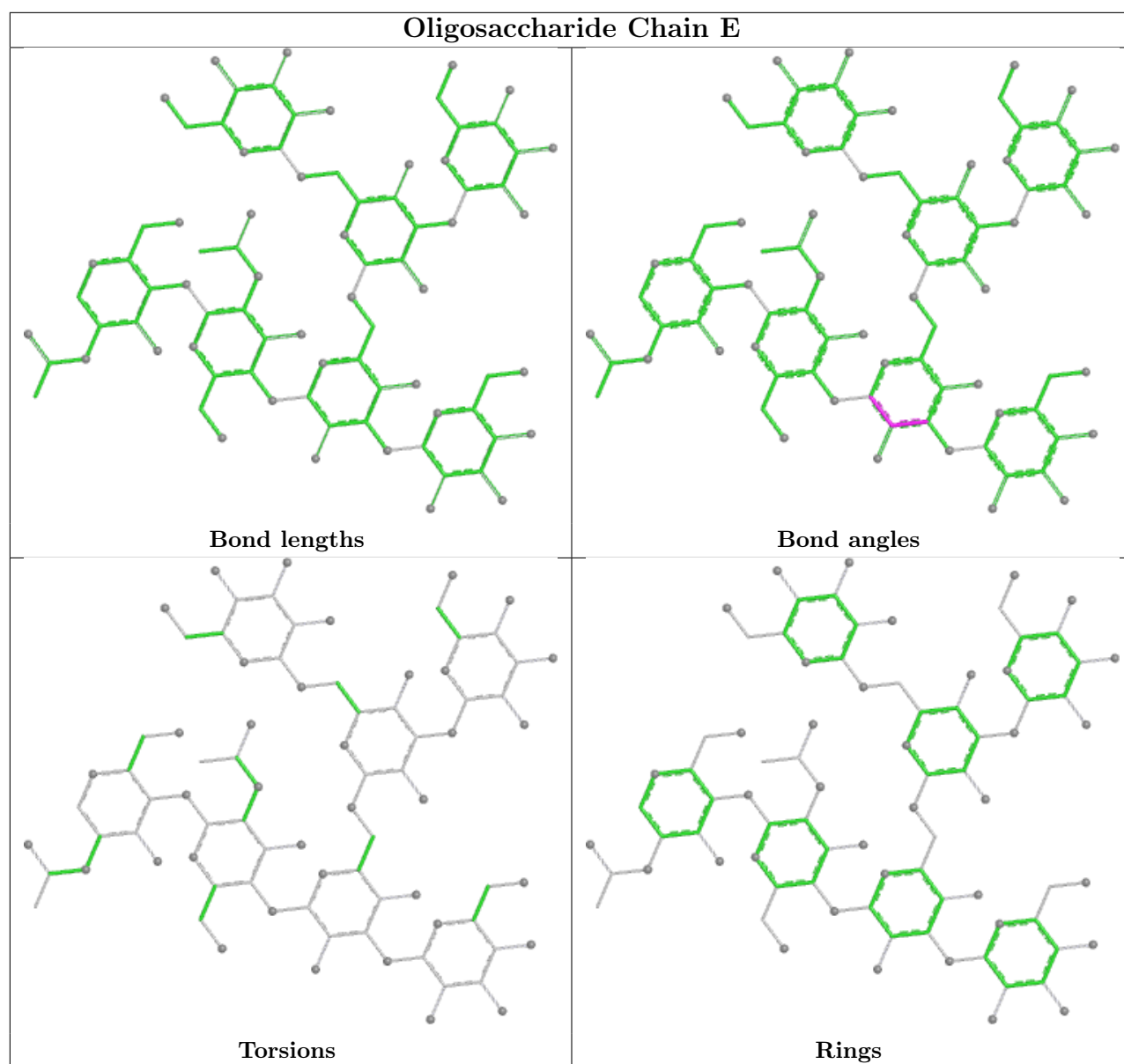
Mol	Chain	Res	Type	Atoms
4	D	7	MAN	C4-C5-C6-O6
4	D	7	MAN	O5-C5-C6-O6

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](#)

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SO4	B	214	-	4,4,4	0.24	0	6,6,6	0.12	0
6	BU3	B	213	-	4,5,5	1.27	1 (25%)	6,6,6	0.70	0
5	SO4	L	214	-	4,4,4	0.25	0	6,6,6	0.11	0
5	SO4	H	226	-	4,4,4	0.27	0	6,6,6	0.10	0
7	NAG	G	573	3	14,14,15	0.54	0	17,19,21	0.73	0
7	NAG	C	573	3	14,14,15	0.56	0	17,19,21	0.97	1 (5%)
5	SO4	B	215	-	4,4,4	0.24	0	6,6,6	0.14	0
6	BU3	L	213	-	4,5,5	1.27	1 (25%)	6,6,6	0.70	0
5	SO4	B	216	-	4,4,4	0.24	0	6,6,6	0.18	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
6	BU3	B	213	-	-	0/4/4/4	-
7	NAG	G	573	3	-	0/6/23/26	0/1/1/1
7	NAG	C	573	3	-	0/6/23/26	0/1/1/1
6	BU3	L	213	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L	213	BU3	O6-C3	-2.14	1.37	1.43
6	B	213	BU3	O6-C3	-2.14	1.37	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	573	NAG	O5-C1-C2	-2.43	107.53	111.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	L	214	SO4	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	224/248 (90%)	0.63	35 (15%) 5 4	18, 40, 109, 148	0
1	H	227/248 (91%)	0.06	7 (3%) 51 51	18, 33, 69, 96	0
2	B	211/216 (97%)	0.37	5 (2%) 59 60	20, 44, 76, 101	0
2	L	211/216 (97%)	0.17	11 (5%) 33 31	17, 38, 77, 101	0
3	C	90/124 (72%)	1.72	36 (40%) 1 0	34, 76, 117, 138	0
3	G	90/124 (72%)	1.98	43 (47%) 0 0	41, 72, 112, 140	0
All	All	1053/1176 (89%)	0.57	137 (13%) 7 6	17, 44, 98, 148	0

The worst 5 of 137 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	181	VAL	5.9
1	A	136	ALA	5.6
1	A	213	PRO	5.6
3	G	235	PHE	5.5
3	G	210	ALA	5.3

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	TYS	H	100(H)	16/17	0.91	0.11	30,36,91,100	0
1	PCA	H	2	8/9	0.92	0.08	30,39,48,54	0
1	TYS	H	100(G)	16/17	0.93	0.10	30,38,88,90	0
1	TYS	A	100(G)	16/17	0.94	0.10	27,32,75,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	TYS	A	100(H)	16/17	0.95	0.10	24,34,64,67	0
1	PCA	A	2	8/9	0.96	0.06	19,24,33,33	0

6.3 Carbohydrates

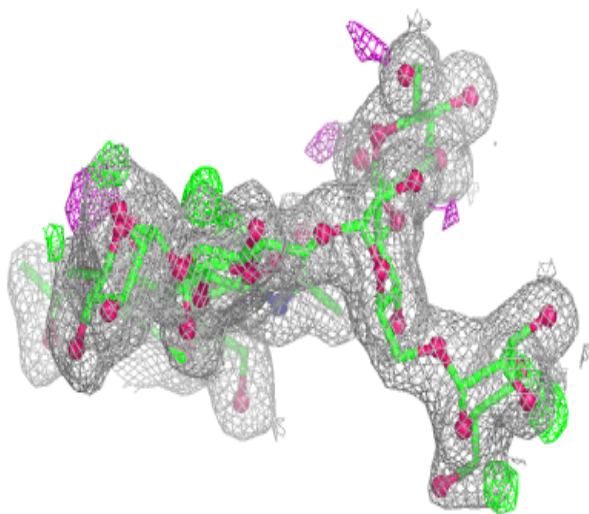
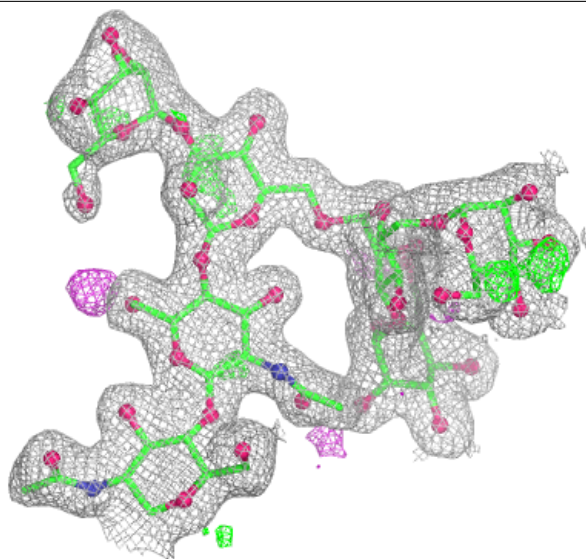
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

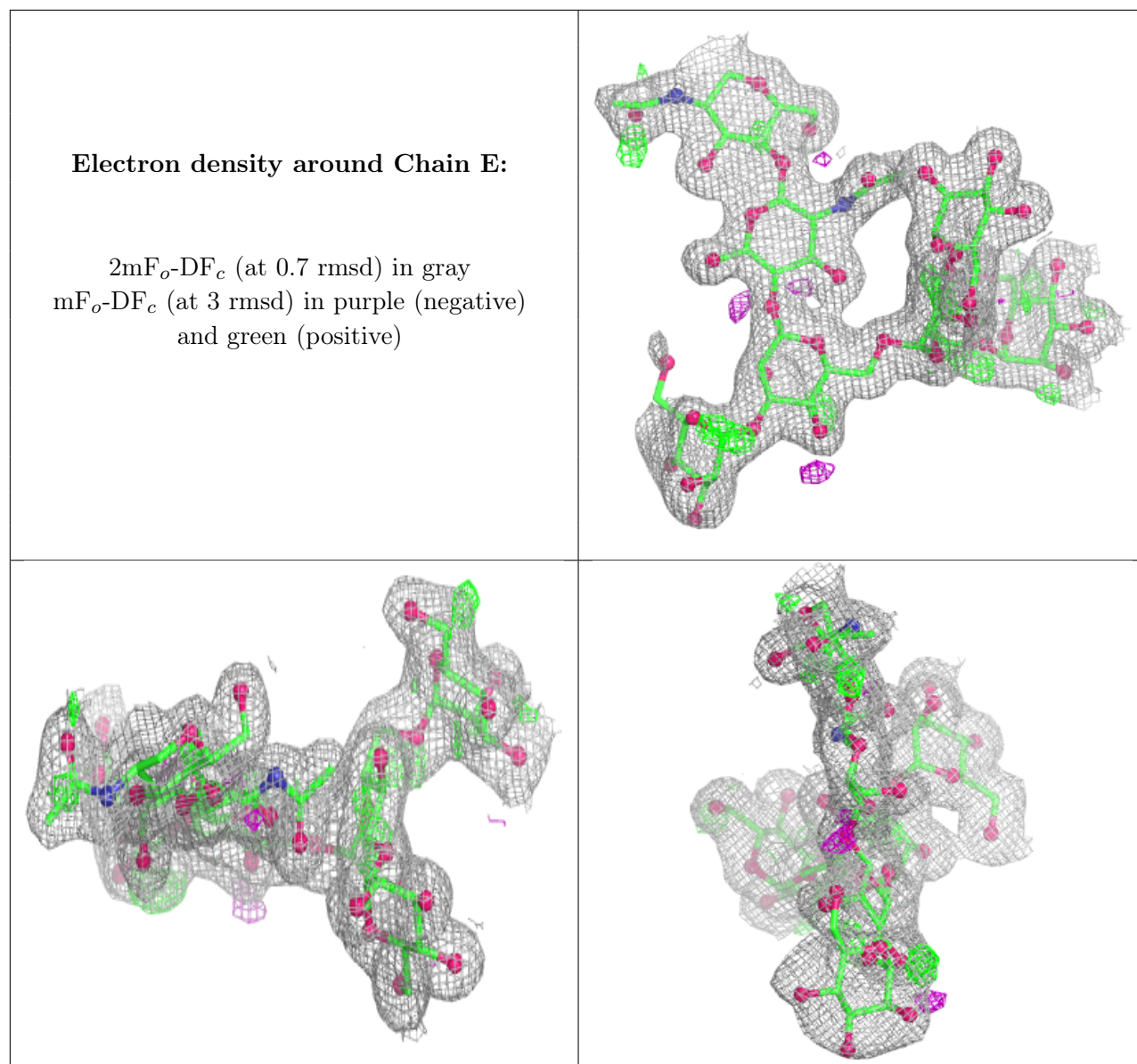
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	D	1	14/15	-	-	26,32,41,43	0
4	NAG	D	2	14/15	-	-	24,31,43,44	0
4	BMA	D	3	11/12	-	-	25,34,42,52	0
4	MAN	D	4	11/12	-	-	21,25,28,28	0
4	MAN	D	5	11/12	-	-	20,23,26,27	0
4	MAN	D	6	11/12	-	-	21,25,38,54	0
4	MAN	D	7	11/12	-	-	52,68,75,80	0
4	NAG	E	1	14/15	-	-	28,34,43,50	0
4	NAG	E	2	14/15	-	-	28,36,45,53	0
4	BMA	E	3	11/12	-	-	31,43,53,60	0
4	MAN	E	4	11/12	-	-	25,28,33,37	0
4	MAN	E	5	11/12	-	-	22,26,32,36	0
4	MAN	E	6	11/12	-	-	23,27,40,43	0
4	MAN	E	7	11/12	-	-	74,80,87,88	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 D:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	BU3	L	213	6/6	0.77	0.16	34,44,58,60	0
5	SO4	H	226	5/5	0.79	0.11	76,80,89,94	0
5	SO4	B	216	5/5	0.82	0.11	59,62,77,88	0
5	SO4	B	215	5/5	0.82	0.12	40,57,71,83	0
7	NAG	G	573	14/15	0.86	0.13	39,46,53,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SO4	L	214	5/5	0.87	0.10	78,83,90,96	0
5	SO4	B	214	5/5	0.87	0.09	67,79,95,103	0
6	BU3	B	213	6/6	0.91	0.10	34,44,58,60	0
7	NAG	C	573	14/15	0.92	0.10	31,38,47,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.