



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:18 PM UTC

PDB ID : 2OK2 / pdb_00002ok2
Title : MutS C-terminal domain fused to Maltose Binding Protein
Authors : Putnam, C.D.; Mendillo, M.L.; Kolodner, R.D.
Deposited on : 2007-01-15
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

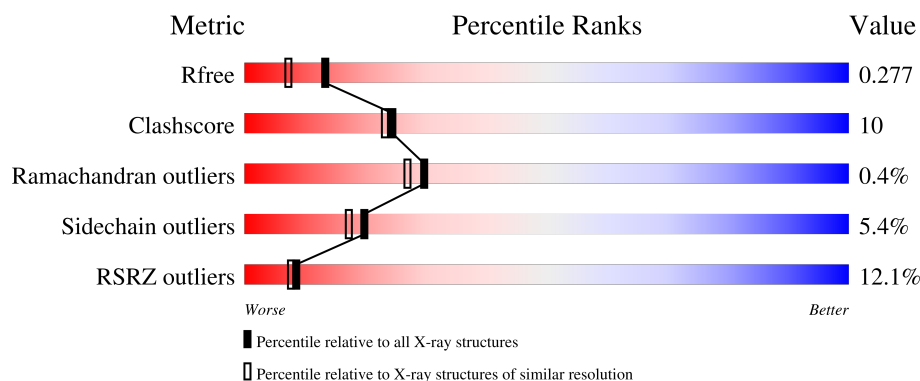
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	402	10% 77% 20% ..
1	B	402	14% 80% 17% .
2	C	2	50% 50%
2	D	2	50% 50%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6889 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Maltose-binding periplasmic protein, DNA mismatch repair protein mutS fusion protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	0	3	0
			3160	2032	512	608	8			
1	B	402	Total	C	N	O	S	0	4	0
			3167	2033	516	610	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	cloning artifact	UNP P0AEX9
A	367	HIS	-	cloning artifact	UNP P0AEX9
A	368	MET	-	cloning artifact	UNP P0AEX9
B	1	MET	-	cloning artifact	UNP P0AEX9
B	367	HIS	-	cloning artifact	UNP P0AEX9
B	368	MET	-	cloning artifact	UNP P0AEX9

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		

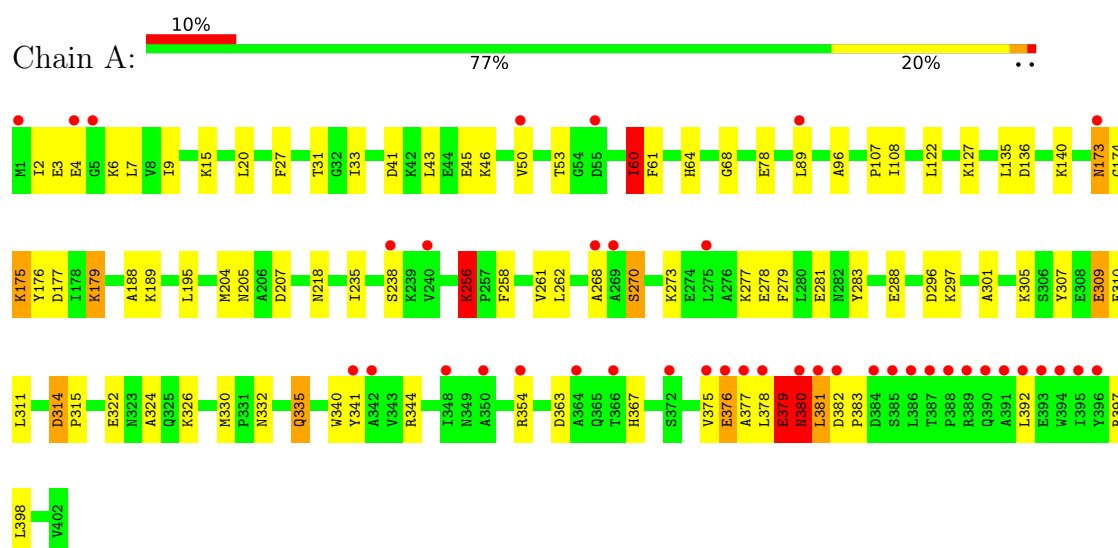
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	287	Total	O	0	0
			287	287		
4	B	224	Total	O	0	0
			224	224		

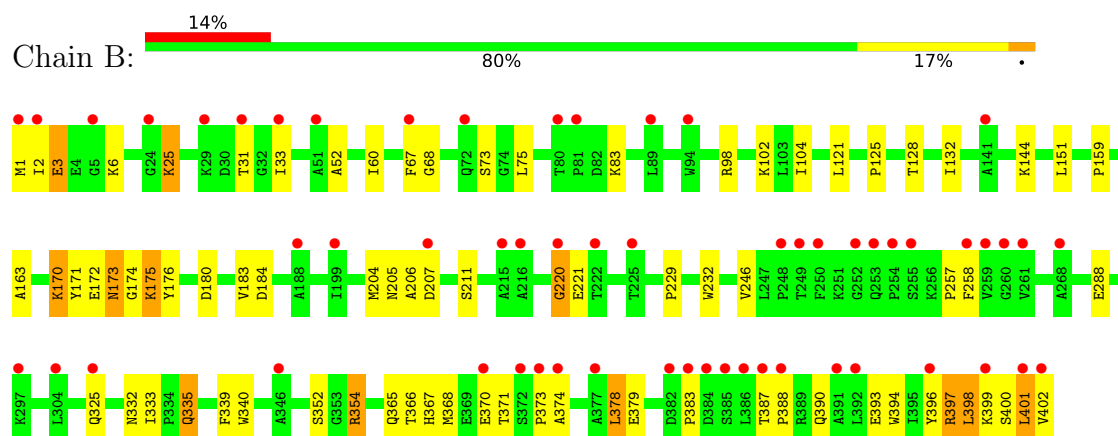
3 Residue-property plots [i](#)

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

- Molecule 1: Maltose-binding periplasmic protein, DNA mismatch repair protein mutS fusion protein



- Molecule 1: Maltose-binding periplasmic protein, DNA mismatch repair protein mutS fusion protein



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose





- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	174.29Å 88.36Å 61.35Å 90.00° 106.80° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.00) 97.9 (20.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.201 , 0.277 0.201 , 0.277	Depositor DCC
R_{free} test set	2998 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	34.4	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6889	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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: SO4, GLC

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	1.38	9/3237 (0.3%)	1.25	17/4398 (0.4%)
1	B	1.15	6/3243 (0.2%)	1.17	7/4404 (0.2%)
All	All	1.27	15/6480 (0.2%)	1.21	24/8802 (0.3%)

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
1	A	0	1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	376	GLU	CD-OE1	26.98	1.76	1.25
1	A	379	GLU	C-N	20.64	1.64	1.33
1	B	132	ILE	CA-CB	9.93	1.59	1.54
1	B	246	VAL	CA-CB	9.77	1.64	1.54
1	A	324	ALA	CA-CB	8.41	1.66	1.53
1	A	376	GLU	CD-OE2	7.59	1.39	1.25
1	A	380	ASN	CG-ND2	6.81	1.47	1.33
1	A	60	ILE	CA-CB	6.33	1.62	1.54
1	A	379	GLU	CD-OE1	6.02	1.36	1.25
1	B	163	ALA	CA-CB	5.98	1.63	1.53
1	B	220	GLY	C-O	5.54	1.31	1.24
1	B	374	ALA	N-CA	5.37	1.52	1.46
1	A	301	ALA	CA-CB	-5.27	1.46	1.53
1	B	128	THR	CA-C	-5.15	1.46	1.52
1	A	108	ILE	CA-CB	5.00	1.62	1.55

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	46	LYS	N-CA-C	-7.02	103.05	111.69
1	A	376	GLU	CG-CD-OE1	-6.74	102.89	118.40
1	B	125	PRO	CA-C-N	6.70	126.46	119.76
1	B	125	PRO	C-N-CA	6.70	126.46	119.76
1	A	256	LYS	CA-C-N	-6.66	113.43	120.03
1	A	256	LYS	C-N-CA	-6.66	113.43	120.03
1	B	132	ILE	N-CA-CB	6.33	115.39	110.45
1	A	270	SER	N-CA-C	6.11	118.28	109.48
1	B	374	ALA	N-CA-C	6.07	118.39	111.11
1	B	401	LEU	N-CA-C	5.83	120.67	113.50
1	A	179	LYS	N-CA-C	-5.77	106.37	113.41
1	A	380	ASN	N-CA-C	5.67	119.37	112.23
1	A	376	GLU	OE1-CD-OE2	5.66	136.49	122.90
1	A	41	ASP	CB-CA-C	5.61	119.87	109.54
1	A	78	GLU	N-CA-C	-5.57	101.60	109.96
1	A	397	ARG	N-CA-C	-5.56	105.22	111.28
1	A	392	LEU	N-CA-C	-5.51	105.36	111.36
1	A	262	LEU	N-CA-C	-5.39	101.66	109.59
1	A	314	ASP	CA-C-N	5.31	125.63	119.47
1	A	314	ASP	C-N-CA	5.31	125.63	119.47
1	B	393	GLU	N-CA-C	-5.22	105.48	111.07
1	A	188	ALA	N-CA-C	-5.17	105.65	111.28
1	B	221	GLU	CB-CG-CD	-5.15	103.85	112.60
1	A	296	ASP	N-CA-C	-5.00	103.91	111.56

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	379	GLU	Sidechain

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	3160	0	3118	76	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3167	0	3126	58	0
2	C	23	0	21	1	0
2	D	23	0	21	1	0
3	A	5	0	0	0	0
4	A	287	0	0	8	0
4	B	224	0	0	6	0
All	All	6889	0	6286	125	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (125) 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:376:GLU:OE1	1:A:376:GLU:CD	1.76	1.26
1:A:270:SER:O	1:A:273:LYS:HE2	1.65	0.97
1:A:175:LYS:HD3	1:A:176:TYR:H	1.29	0.96
1:A:31:THR:HG22	1:A:33:ILE:HD13	1.47	0.95
1:A:367:HIS:ND1	4:A:8336:HOH:O	2.04	0.91
1:B:31:THR:HG22	1:B:33:ILE:HG12	1.55	0.88
1:A:175:LYS:HA	1:A:175:LYS:HE2	1.56	0.86
1:A:378:LEU:HD11	1:B:398:LEU:HD21	1.57	0.86
1:B:399:LYS:O	1:B:402:VAL:HG23	1.76	0.86
1:B:172:GLU:HG3	1:B:173:ASN:H	1.43	0.84
1:A:268:ALA:O	1:A:273:LYS:NZ	2.15	0.79
1:A:288:GLU:HB2	4:A:8380:HOH:O	1.82	0.78
1:A:378:LEU:HD11	1:B:398:LEU:CD2	2.14	0.78
1:A:175:LYS:HD3	1:A:176:TYR:N	2.02	0.74
1:A:68:GLY:HA3	1:A:332:ASN:O	1.86	0.74
1:A:175:LYS:CD	1:A:176:TYR:H	2.02	0.72
1:B:387:THR:OG1	1:B:390:GLN:HG3	1.89	0.72
1:A:305:LYS:O	1:A:309:GLU:HG2	1.90	0.71
1:B:172:GLU:HG3	1:B:173:ASN:N	2.04	0.71
1:B:175:LYS:HE3	1:B:176:TYR:H	1.56	0.71
1:B:339:PHE:HA	1:B:368:MET:HE2	1.73	0.70
1:B:31:THR:CG2	1:B:33:ILE:HG12	2.21	0.70
1:B:205[B]:ASN:HD22	1:B:207:ASP:H	1.38	0.69
1:A:89:LEU:HD23	1:A:107:PRO:HG2	1.76	0.68
1:A:60:ILE:C	1:A:60:ILE:HD13	2.18	0.68
1:A:64:HIS:HB3	1:A:261[A]:VAL:HG13	1.74	0.68
1:A:31:THR:CG2	1:A:33:ILE:HD13	2.21	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:VAL:HG11	1:B:368:MET:HE1	1.76	0.67
1:A:363:ASP:OD2	4:A:8361:HOH:O	2.13	0.67
1:B:172:GLU:HG2	1:B:175:LYS:O	1.95	0.67
1:A:64:HIS:HB3	1:A:261[A]:VAL:CG1	2.26	0.65
1:B:366:THR:O	1:B:370:GLU:HG2	1.97	0.64
1:A:378:LEU:CD1	1:B:398:LEU:HD21	2.28	0.63
1:A:398:LEU:HB3	1:B:378:LEU:HD21	1.82	0.62
1:A:341[A]:TYR:CD1	1:A:367:HIS:NE2	2.68	0.62
1:B:354:ARG:HD2	4:B:8396:HOH:O	2.01	0.61
1:A:307:TYR:CZ	1:A:311:LEU:HD21	2.36	0.61
1:A:64:HIS:ND1	1:A:261[A]:VAL:HG12	2.15	0.60
1:A:341[B]:TYR:CD1	1:A:344:ARG:NH1	2.70	0.60
1:A:31:THR:HG22	1:A:33:ILE:CD1	2.29	0.58
1:B:25:LYS:HD3	1:B:25:LYS:C	2.27	0.58
1:B:170:LYS:HZ2	1:B:170:LYS:HB2	1.68	0.58
1:B:1:MET:C	1:B:3:GLU:H	2.12	0.57
1:B:333:ILE:HD12	1:B:335:GLN:HE21	1.69	0.57
1:B:67:PHE:HB3	1:B:104:ILE:HD12	1.87	0.57
1:A:9:ILE:HG21	1:A:20:LEU:HD21	1.85	0.57
1:A:27:PHE:O	1:A:31:THR:HB	2.05	0.56
1:B:398:LEU:C	1:B:398:LEU:HD23	2.31	0.56
1:A:381:LEU:HD12	1:A:381:LEU:H	1.69	0.56
1:B:383:PRO:HD2	4:B:8409:HOH:O	2.05	0.56
1:A:218:ASN:O	1:A:238:SER:OG	2.22	0.56
1:A:60:ILE:HD13	1:A:61:PHE:N	2.21	0.55
1:A:205:ASN:OD1	1:A:207:ASP:HB2	2.07	0.54
1:B:159:PRO:HG3	1:B:257:PRO:HA	1.90	0.54
1:A:195:LEU:CD1	1:A:204:MET:HE1	2.38	0.54
1:B:205[B]:ASN:ND2	1:B:207:ASP:H	2.03	0.54
1:B:333:ILE:HD12	1:B:335:GLN:NE2	2.24	0.53
1:B:172:GLU:CG	1:B:173:ASN:H	2.09	0.53
1:B:335:GLN:NE2	1:B:335:GLN:H	2.07	0.53
1:A:381:LEU:O	1:B:399:LYS:HE2	2.09	0.52
1:A:136:ASP:OD2	1:A:140:LYS:HE3	2.09	0.52
1:A:398:LEU:CB	1:B:378:LEU:HD11	2.40	0.52
1:A:195:LEU:HD11	1:A:204:MET:HE1	1.92	0.52
1:B:98:ARG:NH1	4:B:8366:HOH:O	2.42	0.52
1:A:341[B]:TYR:HD1	1:A:344:ARG:NH1	2.08	0.51
1:A:50:VAL:O	1:A:53:THR:HG22	2.10	0.50
1:B:52:ALA:HB1	1:B:397:ARG:NE	2.27	0.50
1:A:173[A]:ASN:ND2	1:A:173[A]:ASN:H	2.08	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205[B]:ASN:ND2	1:B:207:ASP:OD1	2.43	0.50
1:A:398:LEU:HB2	1:B:378:LEU:HD11	1.94	0.49
1:B:25:LYS:HD3	1:B:25:LYS:O	2.11	0.49
1:A:341[A]:TYR:CE1	1:A:367:HIS:CE1	3.01	0.48
1:B:102:LYS:NZ	4:B:8286:HOH:O	2.47	0.48
1:B:229:PRO:HA	1:B:232:TRP:CE2	2.49	0.48
1:B:352:SER:OG	1:B:354:ARG:HG2	2.13	0.48
1:A:6:LYS:HG2	1:A:7:LEU:N	2.29	0.47
1:A:43:LEU:C	1:A:43:LEU:HD12	2.39	0.47
1:A:136:ASP:O	1:A:140:LYS:HG3	2.14	0.47
1:A:45:GLU:HG2	1:A:380:ASN:HD22	1.80	0.46
1:A:175:LYS:HA	1:A:175:LYS:CE	2.36	0.46
1:B:365[B]:GLN:NE2	4:B:8210:HOH:O	2.48	0.46
1:A:64:HIS:CE1	1:A:330:MET:HB2	2.51	0.46
1:A:378:LEU:HA	1:A:381:LEU:CD1	2.46	0.46
1:B:367:HIS:CE1	1:B:379:GLU:OE1	2.68	0.46
1:A:64:HIS:CG	1:A:261[A]:VAL:HG12	2.51	0.46
1:A:189:LYS:NZ	4:A:8385:HOH:O	2.49	0.46
1:B:340:TRP:CE3	2:D:2:GLC:H61	2.51	0.45
1:A:15:LYS:HE2	4:A:8063:HOH:O	2.17	0.45
1:B:170:LYS:NZ	4:B:8393:HOH:O	2.48	0.45
1:B:205[B]:ASN:HD22	1:B:206:ALA:N	2.14	0.45
1:A:177:ASP:OD1	1:A:179:LYS:HB3	2.17	0.45
1:A:335:GLN:NE2	1:A:335:GLN:H	2.15	0.45
1:B:175:LYS:CE	1:B:176:TYR:H	2.26	0.45
1:A:314:ASP:HA	1:A:315:PRO:HD2	1.79	0.44
1:A:340:TRP:CD1	2:C:2:GLC:H4	2.53	0.44
1:B:68:GLY:HA3	1:B:332:ASN:O	2.17	0.44
1:B:144:LYS:HE2	1:B:220:GLY:O	2.16	0.44
1:B:151:LEU:HD11	1:B:204:MET:HE3	1.99	0.43
1:A:354:ARG:HA	1:A:354:ARG:HD2	1.73	0.43
1:A:122:LEU:HD11	1:A:135:LEU:HD11	2.00	0.43
1:A:382:ASP:HB3	4:A:8492:HOH:O	2.17	0.43
1:B:399:LYS:O	1:B:402:VAL:CG2	2.60	0.43
1:A:235:ILE:O	1:A:238:SER:HB3	2.18	0.43
1:A:341[B]:TYR:HD1	1:A:344:ARG:HH11	1.67	0.43
1:B:401:LEU:N	1:B:402:VAL:HA	2.33	0.43
1:B:171:TYR:HE1	1:B:174:GLY:C	2.27	0.42
1:B:394:TRP:O	1:B:398:LEU:HB3	2.19	0.42
1:A:173[A]:ASN:CG	1:A:174:GLY:N	2.77	0.42
1:A:96:ALA:HB3	1:A:261[A]:VAL:HG11	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:GLU:HG2	1:A:326:LYS:HE3	2.02	0.41
1:A:375:VAL:O	1:A:379:GLU:HG3	2.20	0.41
1:A:45:GLU:HG2	1:A:380:ASN:ND2	2.36	0.41
1:A:383:PRO:HB2	1:B:396:TYR:CE2	2.56	0.41
1:A:127:LYS:NZ	4:A:8288:HOH:O	2.51	0.41
1:B:184:ASP:HB2	1:B:365[B]:GLN:HB2	2.02	0.41
1:A:173[A]:ASN:H	1:A:173[A]:ASN:HD22	1.67	0.41
1:A:277:LYS:O	1:A:281:GLU:HB2	2.19	0.41
1:B:184:ASP:HB2	1:B:365[A]:GLN:HB2	2.02	0.41
1:A:279:PHE:O	1:A:283:TYR:HB2	2.21	0.41
1:B:73:SER:HB2	1:B:75:LEU:HG	2.02	0.41
1:B:170:LYS:NZ	1:B:180:ASP:OD2	2.53	0.41
1:A:377:ALA:O	1:A:380:ASN:HB2	2.21	0.40
1:A:398:LEU:HB3	1:B:378:LEU:HD11	2.03	0.40
1:A:256:LYS:HD3	4:A:8045:HOH: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	403/402 (100%)	391 (97%)	11 (3%)	1 (0%)	43	42
1	B	404/402 (100%)	393 (97%)	9 (2%)	2 (0%)	24	21
All	All	807/804 (100%)	784 (97%)	20 (2%)	3 (0%)	30	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2	ILE
1	A	2	ILE
1	B	373	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	329/326 (101%)	314 (95%)	15 (5%)	24	22
1	B	330/326 (101%)	309 (94%)	21 (6%)	16	12
All	All	659/652 (101%)	623 (94%)	36 (6%)	20	17

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLU
1	A	4	GLU
1	A	60	ILE
1	A	173[A]	ASN
1	A	173[B]	ASN
1	A	175	LYS
1	A	256	LYS
1	A	258	PHE
1	A	278	GLU
1	A	297	LYS
1	A	309	GLU
1	A	310	GLU
1	A	335	GLN
1	A	380	ASN
1	A	381	LEU
1	B	3	GLU
1	B	6	LYS
1	B	25	LYS
1	B	60	ILE
1	B	83	LYS
1	B	121	LEU
1	B	170	LYS
1	B	173	ASN
1	B	175	LYS
1	B	211	SER
1	B	258	PHE
1	B	288	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	325	GLN
1	B	335	GLN
1	B	354	ARG
1	B	371	THR
1	B	378	LEU
1	B	388	PRO
1	B	397	ARG
1	B	398	LEU
1	B	400	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	100	ASN
1	A	201	ASN
1	A	218	ASN
1	A	335	GLN
1	A	380	ASN
1	B	201	ASN
1	B	218	ASN
1	B	282	ASN
1	B	325	GLN
1	B	335	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

4 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	GLC	C	1	2	12,12,12	0.58	0	17,17,17	1.23	1 (5%)
2	GLC	C	2	2	11,11,12	0.82	0	15,15,17	0.98	1 (6%)
2	GLC	D	1	2	12,12,12	0.84	0	17,17,17	1.11	2 (11%)
2	GLC	D	2	2	11,11,12	0.69	0	15,15,17	2.07	3 (20%)

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	GLC	C	1	2	-	0/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	GLC	C1-O5-C5	6.20	120.49	112.19
2	D	2	GLC	C1-C2-C3	-3.33	104.79	109.64
2	C	1	GLC	O5-C1-C2	3.17	115.88	110.30
2	D	1	GLC	O1-C1-O5	-2.77	102.19	110.41
2	C	2	GLC	C1-O5-C5	2.24	115.19	112.19
2	D	2	GLC	O2-C2-C3	2.09	114.47	110.15
2	D	1	GLC	C4-C3-C2	-2.03	107.27	110.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

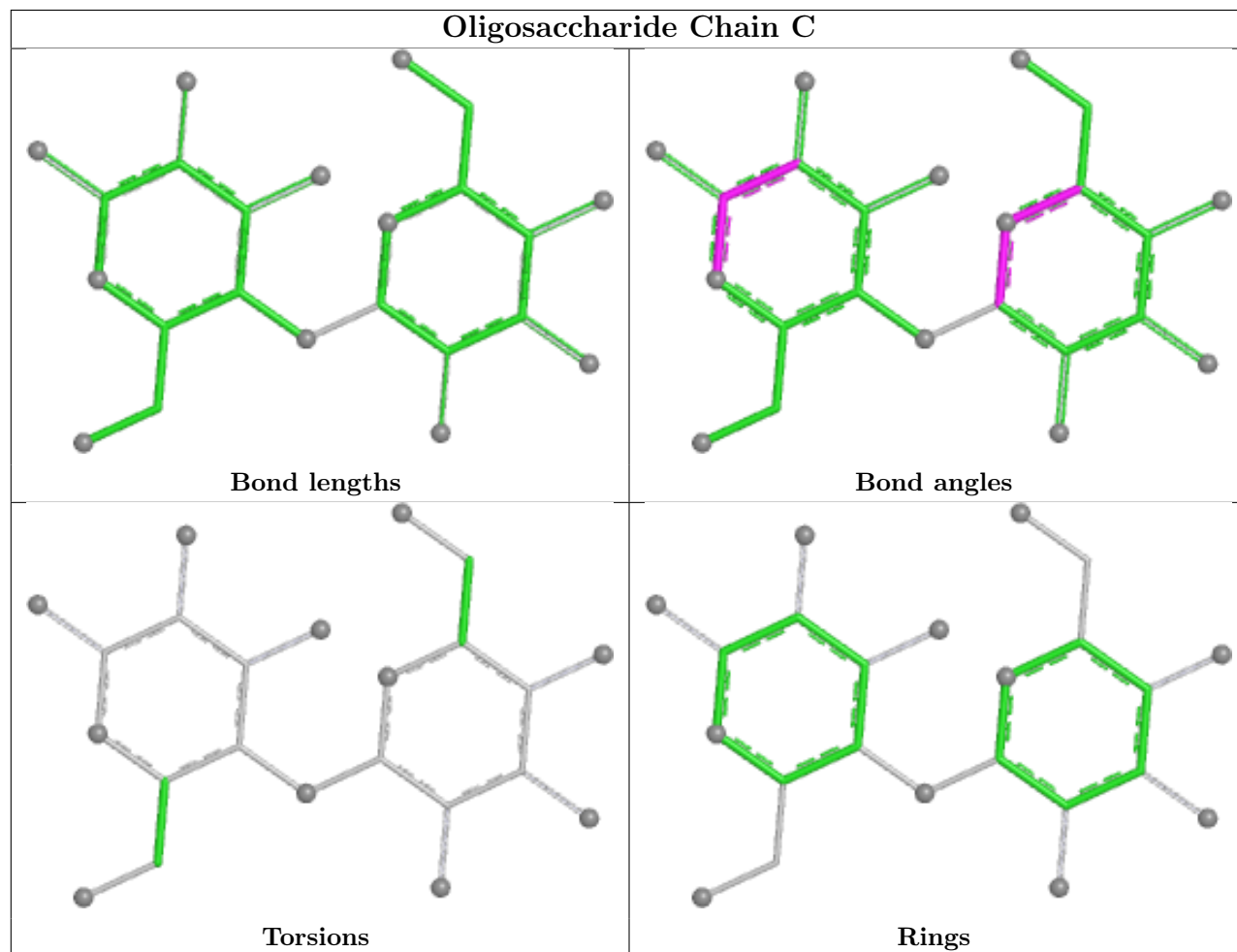
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	GLC	1	0

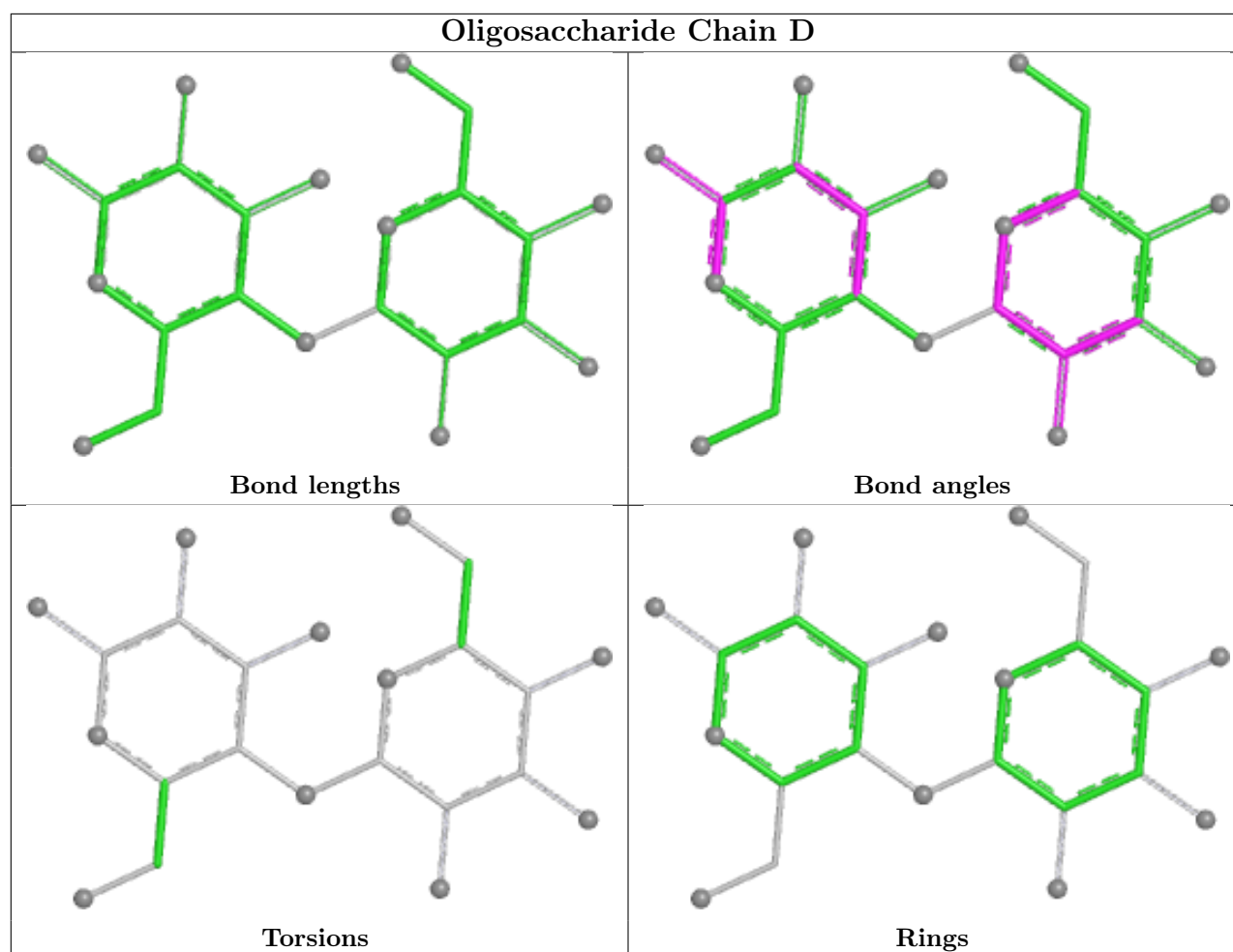
Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	GLC	1	0

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





5.6 Ligand geometry [i](#)

1 ligand is 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	SO4	A	5000	-	4,4,4	0.26	0	6,6,6	0.41	0

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 [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	379:GLU	C	380:ASN	N	1.64

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	402/402 (100%)	0.88	40 (9%)	12 11	19, 41, 47, 58	3 (0%)
1	B	402/402 (100%)	1.21	57 (14%)	6 5	19, 41, 46, 61	4 (0%)
All	All	804/804 (100%)	1.04	97 (12%)	8 8	19, 41, 46, 61	7 (0%)

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	388	PRO	6.1
1	A	391	ALA	5.0
1	A	387	THR	4.8
1	A	392	LEU	4.3
1	B	216	ALA	4.2
1	B	374	ALA	4.2
1	A	4	GLU	4.1
1	A	5	GLY	4.0
1	B	250	PHE	4.0
1	A	396	TYR	3.8
1	B	373	PRO	3.7
1	A	395	ILE	3.7
1	B	2	ILE	3.6
1	B	215	ALA	3.6
1	B	249	THR	3.5
1	A	389	ARG	3.5
1	B	51	ALA	3.3
1	A	394	TRP	3.1
1	A	378	LEU	3.1
1	B	392	LEU	3.1
1	B	1	MET	3.1
1	B	402	VAL	3.1
1	B	387	THR	3.0
1	B	384	ASP	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	67	PHE	3.0
1	B	33	ILE	3.0
1	B	31	THR	3.0
1	B	141	ALA	3.0
1	A	342	ALA	2.9
1	B	372	SER	2.8
1	A	268	ALA	2.8
1	B	89	LEU	2.8
1	A	366	THR	2.8
1	B	252	GLY	2.8
1	A	55	ASP	2.7
1	A	376	GLU	2.7
1	A	269	ALA	2.7
1	B	377	ALA	2.7
1	B	222	THR	2.7
1	B	401	LEU	2.7
1	A	382	ASP	2.7
1	A	393	GLU	2.7
1	A	341[A]	TYR	2.7
1	A	385	SER	2.7
1	A	377	ALA	2.6
1	B	207	ASP	2.6
1	B	268	ALA	2.6
1	B	386	LEU	2.6
1	A	386	LEU	2.6
1	B	383	PRO	2.6
1	A	348	ILE	2.6
1	B	24	GLY	2.6
1	B	248	PRO	2.5
1	A	1	MET	2.5
1	B	259	VAL	2.5
1	B	253	GLN	2.5
1	B	325	GLN	2.5
1	B	80	THR	2.5
1	B	255	SER	2.5
1	B	188	ALA	2.5
1	A	50	VAL	2.4
1	B	261	VAL	2.4
1	A	384	ASP	2.4
1	A	238	SER	2.4
1	A	390	GLN	2.4
1	B	297	LYS	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	391	ALA	2.3
1	A	275	LEU	2.3
1	A	375	VAL	2.3
1	B	258	PHE	2.3
1	B	385	SER	2.3
1	B	5	GLY	2.3
1	B	199	ILE	2.3
1	B	225	THR	2.3
1	B	254	PRO	2.3
1	B	29	LYS	2.3
1	B	382	ASP	2.3
1	B	94	TRP	2.3
1	B	388	PRO	2.2
1	A	173[A]	ASN	2.2
1	A	350	ALA	2.2
1	A	89	LEU	2.2
1	B	396	TYR	2.2
1	B	399	LYS	2.2
1	B	260	GLY	2.2
1	B	370	GLU	2.2
1	B	81	PRO	2.1
1	A	240	VAL	2.1
1	B	304	LEU	2.1
1	B	220	GLY	2.1
1	A	364	ALA	2.1
1	B	346	ALA	2.1
1	A	381	LEU	2.1
1	B	72	GLN	2.1
1	A	354	ARG	2.1
1	A	372	SER	2.0
1	A	380	ASN	2.0

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

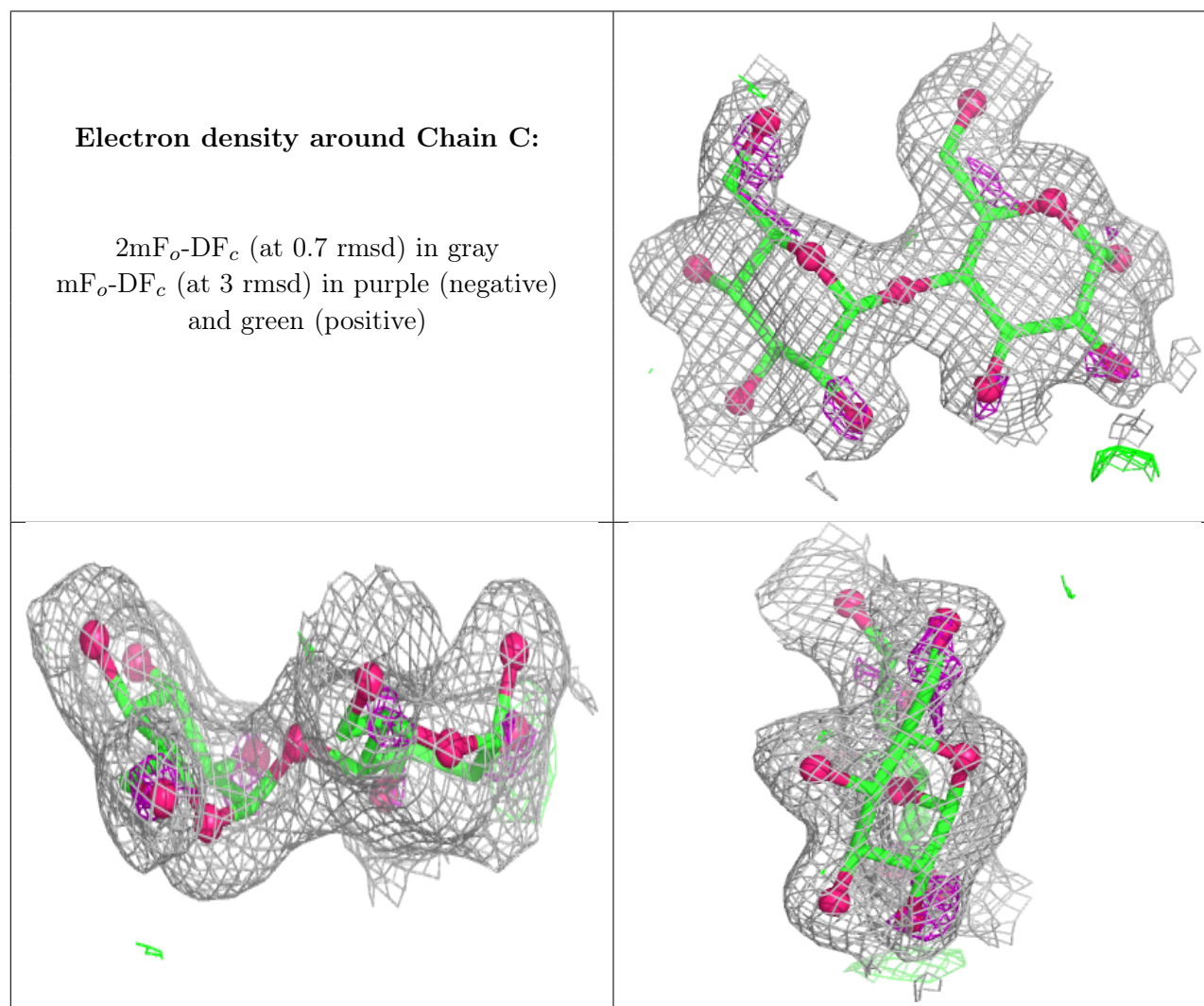
There are no non-standard protein/DNA/RNA residues in this entry.

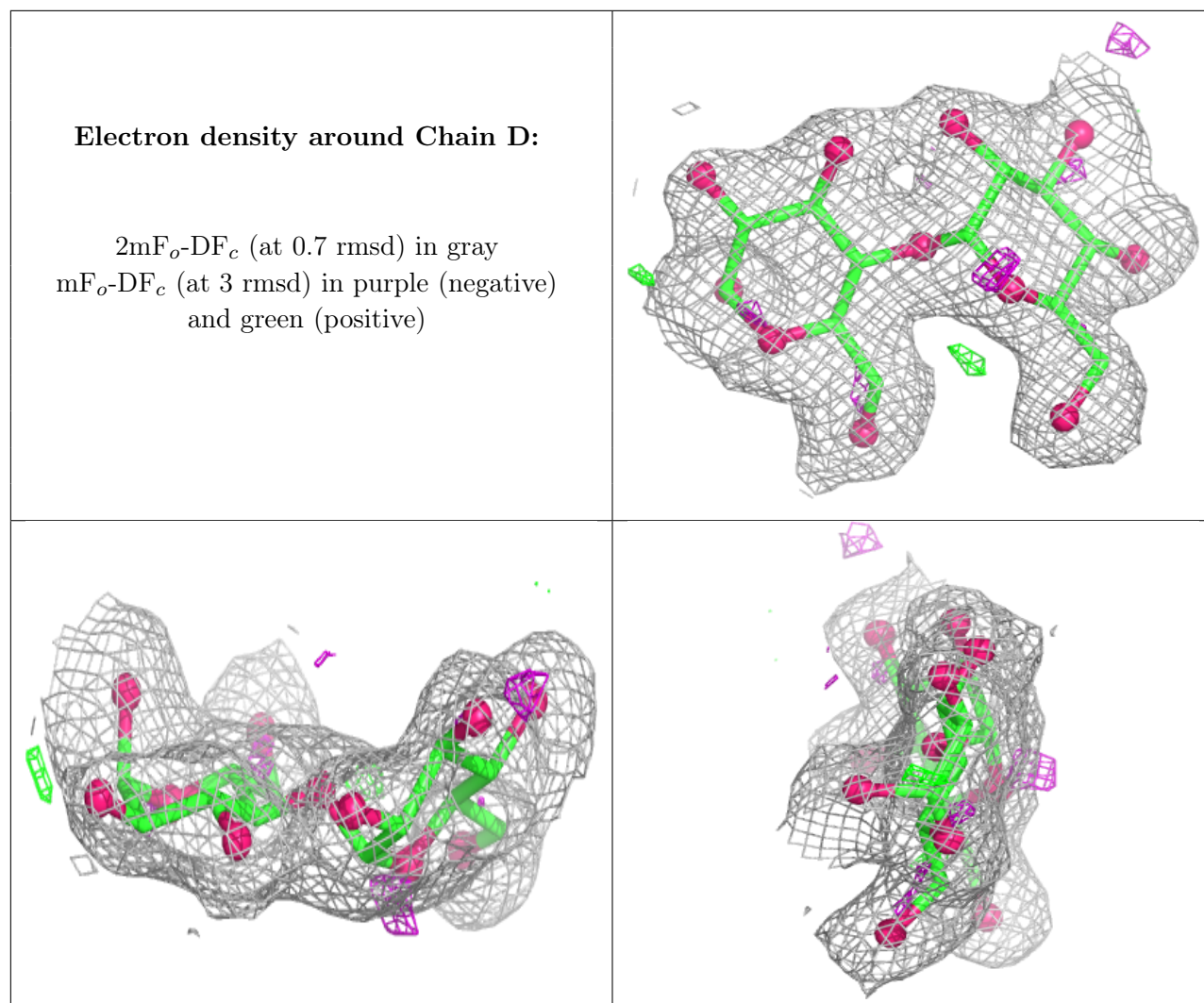
6.3 Carbohydrates [i](#)

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($\text{\AA}^2$)	Q<0.9
2	GLC	C	1	12/12	0.91	0.10	35,37,43,46	0
2	GLC	D	1	12/12	0.91	0.10	36,40,44,50	0
2	GLC	D	2	11/12	0.91	0.10	35,39,42,43	0
2	GLC	C	2	11/12	0.93	0.10	32,35,39,39	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($\text{\AA}^2$)	Q<0.9
3	SO4	A	5000	5/5	0.70	0.15	87,88,90,91	0

6.5 Other polymers [i](#)

There are no such residues in this entry.