



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 1, 2023 – 11:01 PM EDT

PDB ID : 3OGX  
Title : Crystal structure of the complex of Peptidoglycan Recognition protein (PGRP-s) with Heparin-Dissacharide at 2.8 Å resolution  
Authors : Sharma, P.; Dube, D.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.  
Deposited on : 2010-08-17  
Resolution : 2.80 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](mailto: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>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

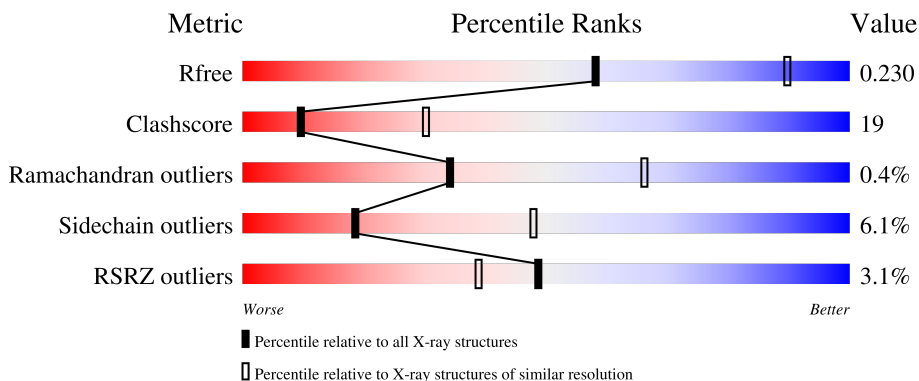
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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	171	
1	B	171	
1	C	171	
1	D	171	
2	E	2	

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:

<b>Mol</b>	<b>Type</b>	<b>Chain</b>	<b>Res</b>	<b>Chirality</b>	<b>Geometry</b>	<b>Clashes</b>	<b>Electron density</b>
2	UAP	E	2	-	-	-	X

## 2 Entry composition [i](#)

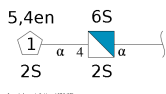
There are 5 unique types of molecules in this entry. The entry contains 5693 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Peptidoglycan recognition protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	171	1337	834	254	241	8	0	0	0
1	B	171	1336	834	254	240	8	0	0	0
1	C	171	1337	834	254	241	8	0	0	0
1	D	171	1336	834	254	240	8	0	0	0

- Molecule 2 is an oligosaccharide called 4-deoxy-2-O-sulfo-alpha-L-threo-hex-4-enopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose.



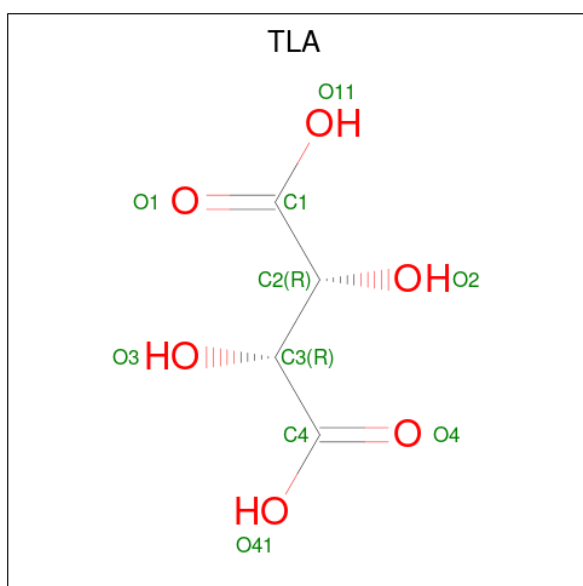
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	E	2	35	12	1	19	3	0	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula:  $C_4H_6O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			10	4	6		

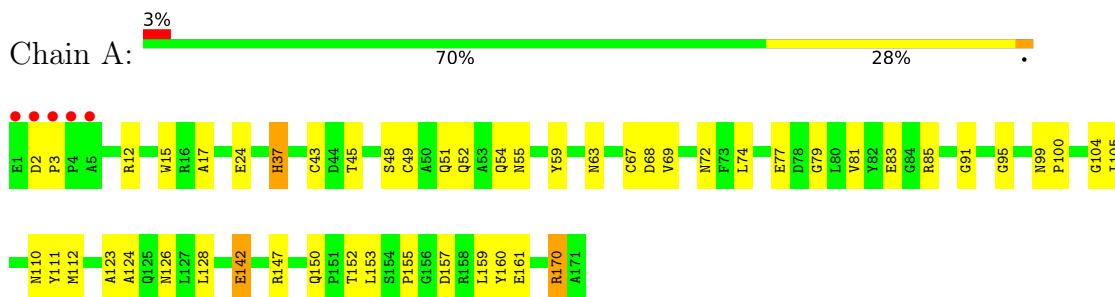
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	75	Total O 75 75	0	0
5	B	72	Total O 72 72	0	0
5	C	72	Total O 72 72	0	0
5	D	71	Total O 71 71	0	0

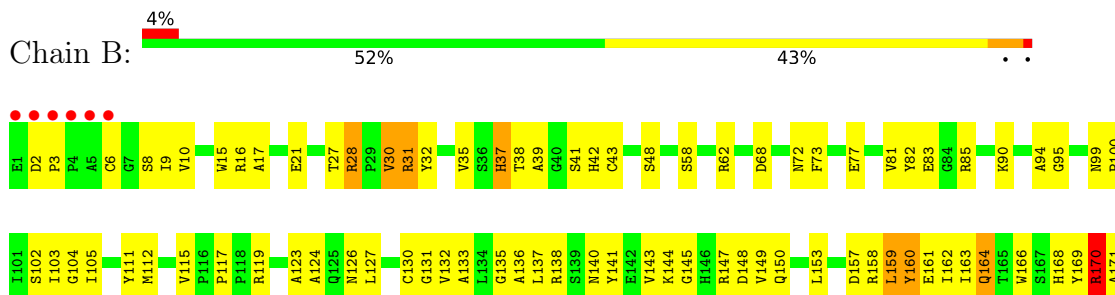
### 3 Residue-property plots

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

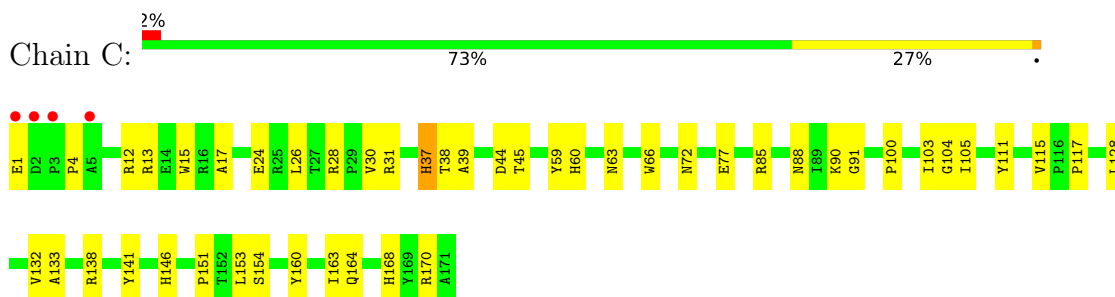
- Molecule 1: Peptidoglycan recognition protein 1



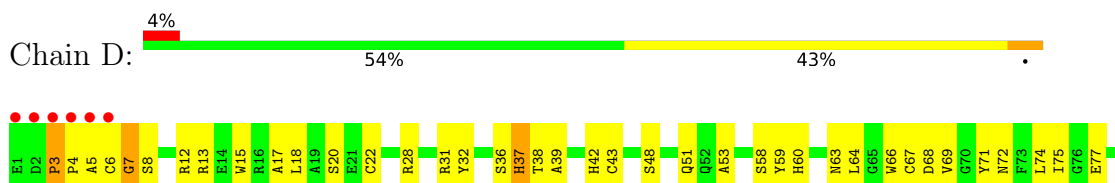
- Molecule 1: Peptidoglycan recognition protein 1

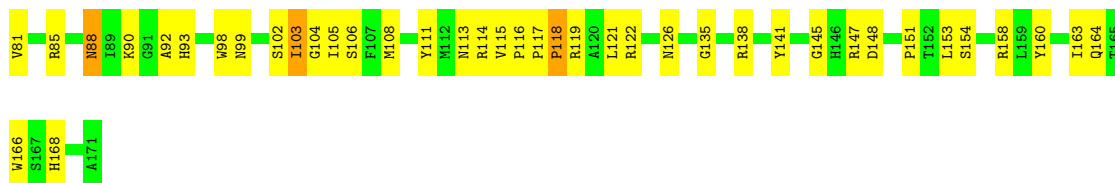


- Molecule 1: Peptidoglycan recognition protein 1



- Molecule 1: Peptidoglycan recognition protein 1





- Molecule 2: 4-deoxy-2-O-sulfo- $\alpha$ -L-threo-hex-4-enopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)- $\alpha$ -D-glucopyranose

Chain E:

100%

S0H1  
UAP2



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.42Å 101.76Å 162.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.88 – 2.80 47.89 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.2 (47.88-2.80) 97.0 (47.89-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.34 (at 2.81Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.205 , 0.234 0.197 , 0.230	Depositor DCC
$R_{free}$ test set	911 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.4	Xtrriage
Anisotropy	1.059	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 63.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5693	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.27% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: UAP, SGN, GOL, TLA

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	0.57	0/1374	0.81	0/1871
1	B	0.62	2/1373 (0.1%)	0.81	0/1871
1	C	0.48	0/1374	0.78	1/1871 (0.1%)
1	D	0.52	0/1373	0.82	1/1871 (0.1%)
All	All	0.55	2/5494 (0.0%)	0.81	2/7484 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	160	TYR	CD1-CE1	-5.50	1.31	1.39
1	B	160	TYR	CD2-CE2	-5.16	1.31	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	7	GLY	N-CA-C	5.64	127.21	113.10
1	C	115	VAL	N-CA-C	-5.01	97.46	111.00

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	1337	0	1288	38	0
1	B	1336	0	1288	68	0
1	C	1337	0	1288	44	0
1	D	1336	0	1288	59	0
2	E	35	0	14	3	0
3	A	6	0	8	0	0
3	D	6	0	8	0	0
4	C	10	0	4	2	0
5	A	75	0	0	0	0
5	B	72	0	0	0	0
5	C	72	0	0	0	0
5	D	71	0	0	0	0
All	All	5693	0	5186	206	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 19.

All (206) 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:D:72:ASN:ND2	1:D:104:GLY:H	1.36	1.21
1:D:72:ASN:HD22	1:D:104:GLY:N	1.55	1.03
1:B:72:ASN:HD22	1:B:104:GLY:N	1.57	1.01
1:B:72:ASN:ND2	1:B:104:GLY:H	1.59	0.99
1:D:28:ARG:HH21	1:D:88:ASN:ND2	1.63	0.96
1:D:28:ARG:HH21	1:D:88:ASN:HD21	1.13	0.94
1:C:72:ASN:HD22	1:C:104:GLY:H	1.12	0.93
1:B:159:LEU:HD12	1:B:159:LEU:O	1.72	0.89
1:B:37:HIS:HD2	1:B:111:TYR:H	1.23	0.87
1:B:73:PHE:HB2	1:B:105:ILE:HG22	1.56	0.86
1:C:59:TYR:HA	1:C:63:ASN:HD22	1.38	0.86
1:D:28:ARG:NH2	1:D:88:ASN:HD21	1.71	0.86
1:A:72:ASN:HD22	1:A:104:GLY:H	1.20	0.85
1:C:77:GLU:HG2	1:C:117:PRO:HD2	1.59	0.85
1:C:72:ASN:ND2	1:C:104:GLY:H	1.75	0.84
1:C:31:ARG:HD2	1:C:138:ARG:HH12	1.41	0.83
1:D:115:VAL:HG12	1:D:158:ARG:HB3	1.61	0.83
1:C:28:ARG:HH21	1:C:88:ASN:HD21	1.25	0.82
1:C:37:HIS:HD2	1:C:111:TYR:H	1.26	0.82
1:A:72:ASN:ND2	1:A:104:GLY:H	1.81	0.79
1:B:43:CYS:HB2	1:B:48:SER:OG	1.84	0.77
2:E:2:UAP:H3	2:E:2:UAP:O1S	1.84	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:31:ARG:HD2	1:C:138:ARG:NH1	2.01	0.75
1:D:72:ASN:ND2	1:D:104:GLY:N	2.22	0.75
1:B:169:TYR:CG	1:B:170:ARG:N	2.56	0.73
1:D:72:ASN:HD22	1:D:104:GLY:H	0.78	0.73
1:D:37:HIS:HD2	1:D:111:TYR:H	1.37	0.73
1:B:169:TYR:CD2	1:B:170:ARG:N	2.57	0.72
1:C:170:ARG:HG2	1:C:170:ARG:NH1	2.05	0.71
1:A:52:GLN:HA	1:A:55:ASN:HD22	1.55	0.71
1:A:126:ASN:HD21	1:B:8:SER:H	1.39	0.71
1:D:3:PRO:HB2	1:D:4:PRO:HD3	1.72	0.71
1:C:170:ARG:HG2	1:C:170:ARG:HH11	1.56	0.70
1:B:37:HIS:CD2	1:B:111:TYR:H	2.07	0.70
1:A:142:GLU:HG2	1:A:170:ARG:HG3	1.73	0.69
1:B:169:TYR:O	1:B:170:ARG:HB2	1.89	0.69
1:D:38:THR:O	1:D:39:ALA:HB3	1.93	0.67
1:C:38:THR:O	1:C:39:ALA:HB3	1.93	0.67
1:A:95:GLY:HA3	1:A:150:GLN:NE2	2.11	0.66
1:C:85:ARG:HD2	1:C:91:GLY:HA2	1.77	0.66
1:D:36:SER:O	1:D:106:SER:HA	1.96	0.65
1:D:48:SER:O	1:D:51:GLN:HB3	1.97	0.65
1:C:60:HIS:O	1:C:66:TRP:HB2	1.97	0.64
1:D:72:ASN:HA	1:D:85:ARG:H	1.61	0.64
1:B:143:VAL:HB	1:B:169:TYR:HA	1.79	0.64
1:A:157:ASP:O	1:A:161:GLU:HG3	1.99	0.63
1:B:124:ALA:O	1:B:127:LEU:HB3	1.98	0.63
1:B:159:LEU:HD12	1:B:159:LEU:C	2.18	0.62
1:D:53:ALA:N	1:D:108:MET:HE1	2.15	0.62
1:B:10:VAL:HG23	1:B:81:VAL:O	2.00	0.62
1:B:115:VAL:CG1	1:B:162:ILE:HD11	2.30	0.62
1:A:147:ARG:HD2	1:A:160:TYR:CE1	2.35	0.61
1:B:81:VAL:HG23	1:B:123:ALA:HB1	1.82	0.61
1:A:142:GLU:CG	1:A:170:ARG:HG3	2.30	0.61
1:B:2:ASP:HB2	1:B:3:PRO:HD2	1.83	0.60
1:A:142:GLU:HG2	1:A:170:ARG:CG	2.32	0.60
4:C:172:TLA:H3	1:D:153:LEU:HD12	1.84	0.60
1:C:59:TYR:HA	1:C:63:ASN:ND2	2.14	0.59
1:B:72:ASN:HD22	1:B:104:GLY:H	0.76	0.59
1:D:15:TRP:CZ2	1:D:17:ALA:HB2	2.38	0.58
1:B:112:MET:SD	1:B:157:ASP:OD1	2.62	0.58
1:D:77:GLU:HG2	1:D:117:PRO:HD2	1.86	0.58
1:B:6:CYS:O	1:B:130:CYS:SG	2.61	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:ARG:HG3	1:B:32:TYR:CD2	2.39	0.57
1:B:131:GLY:HA3	1:B:137:LEU:HD12	1.87	0.57
1:C:4:PRO:HG3	1:C:133:ALA:HB1	1.86	0.57
1:B:90:LYS:HG3	1:B:100:PRO:HA	1.86	0.57
1:C:37:HIS:CD2	1:C:111:TYR:H	2.15	0.56
1:A:85:ARG:HD2	1:A:91:GLY:HA2	1.87	0.56
1:C:72:ASN:HD22	1:C:104:GLY:N	1.94	0.55
1:D:37:HIS:CD2	1:D:111:TYR:H	2.22	0.55
1:A:74:LEU:HD12	1:A:74:LEU:N	2.21	0.55
1:D:28:ARG:HD3	1:D:135:GLY:O	2.06	0.55
1:D:59:TYR:HD1	1:D:63:ASN:HB2	1.71	0.55
1:D:71:TYR:CD1	1:D:106:SER:HB2	2.42	0.55
1:C:24:GLU:OE2	1:C:90:LYS:HE2	2.07	0.54
1:D:147:ARG:NH2	1:D:153:LEU:O	2.40	0.54
1:C:26:LEU:HD13	1:C:30:VAL:HG21	1.88	0.54
1:A:37:HIS:HD2	1:A:111:TYR:H	1.55	0.54
1:A:12:ARG:HG3	1:A:83:GLU:OE1	2.08	0.54
1:D:74:LEU:HD12	1:D:74:LEU:N	2.23	0.54
2:E:1:SGN:O1S	2:E:1:SGN:H3	2.06	0.54
1:A:68:ASP:OD1	1:A:69:VAL:N	2.38	0.54
1:B:99:ASN:N	1:B:100:PRO:HD3	2.23	0.54
1:D:105:ILE:O	1:D:105:ILE:HG13	2.07	0.54
1:C:170:ARG:HH11	1:C:170:ARG:CG	2.21	0.53
1:D:85:ARG:NH1	1:D:90:LYS:O	2.39	0.53
1:B:30:VAL:HG23	1:B:136:ALA:O	2.09	0.53
1:B:99:ASN:N	1:B:100:PRO:CD	2.71	0.53
1:B:132:VAL:O	1:B:135:GLY:HA2	2.09	0.53
1:C:24:GLU:OE2	1:C:90:LYS:CE	2.57	0.53
1:C:28:ARG:HH21	1:C:88:ASN:ND2	1.99	0.53
1:D:43:CYS:O	1:D:77:GLU:HB2	2.09	0.53
1:C:128:LEU:O	1:C:132:VAL:HG23	2.08	0.53
1:D:53:ALA:HA	1:D:108:MET:CE	2.39	0.53
1:B:149:VAL:O	1:B:150:GLN:HG2	2.08	0.52
1:D:22:CYS:SG	1:D:68:ASP:HB3	2.49	0.52
1:A:43:CYS:O	1:A:77:GLU:HB2	2.09	0.52
1:B:43:CYS:O	1:B:77:GLU:HB2	2.09	0.52
1:D:38:THR:O	1:D:39:ALA:CB	2.58	0.52
1:C:59:TYR:O	1:C:63:ASN:HB2	2.09	0.51
1:A:95:GLY:CA	1:A:150:GLN:NE2	2.74	0.50
1:D:141:TYR:O	1:D:168:HIS:HD2	1.95	0.50
1:A:99:ASN:N	1:A:100:PRO:HD2	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:VAL:O	1:B:135:GLY:N	2.44	0.50
1:C:141:TYR:O	1:C:168:HIS:HD2	1.95	0.50
1:D:59:TYR:CE1	1:D:64:LEU:HD21	2.47	0.49
1:D:92:ALA:HA	1:D:99:ASN:ND2	2.27	0.49
1:B:28:ARG:HB2	1:B:136:ALA:HA	1.93	0.49
1:D:32:TYR:O	1:D:102:SER:HA	2.13	0.49
1:C:1:GLU:O	1:C:1:GLU:HG2	2.13	0.48
1:C:38:THR:O	1:C:39:ALA:CB	2.58	0.48
1:D:122:ARG:O	1:D:126:ASN:HB2	2.13	0.48
1:D:93:HIS:CE1	1:D:104:GLY:HA3	2.48	0.48
1:B:144:LYS:NZ	1:B:171:ALA:CB	2.77	0.48
1:B:72:ASN:ND2	1:B:103:ILE:HA	2.29	0.48
1:A:95:GLY:CA	1:A:150:GLN:HE22	2.26	0.48
1:B:144:LYS:NZ	1:B:171:ALA:HB3	2.28	0.47
1:C:44:ASP:OD1	1:C:45:THR:HG23	2.14	0.47
1:C:151:PRO:HD3	1:D:66:TRP:CH2	2.49	0.47
1:B:95:GLY:HA3	1:B:150:GLN:NE2	2.30	0.47
1:B:35:VAL:HG22	1:B:105:ILE:HD11	1.97	0.47
1:B:81:VAL:HG11	1:B:127:LEU:HD13	1.96	0.47
1:D:145:GLY:O	1:D:148:ASP:HB2	2.15	0.47
1:C:153:LEU:HD12	4:C:172:TLA:H2	1.97	0.47
1:D:5:ALA:O	1:D:6:CYS:HB3	2.15	0.46
1:D:116:PRO:HG2	1:D:121:LEU:HD21	1.96	0.46
1:B:144:LYS:HZ3	1:B:171:ALA:HB3	1.81	0.46
1:B:160:TYR:CZ	1:B:164:GLN:NE2	2.84	0.46
1:C:12:ARG:HB3	1:C:17:ALA:HB3	1.96	0.46
1:D:160:TYR:O	1:D:164:GLN:HG3	2.16	0.46
1:B:147:ARG:NH2	1:B:153:LEU:O	2.45	0.46
1:C:160:TYR:O	1:C:164:GLN:HG3	2.15	0.46
1:D:60:HIS:O	1:D:66:TRP:HB2	2.16	0.46
1:A:126:ASN:ND2	1:B:8:SER:H	2.11	0.46
1:B:123:ALA:O	1:B:127:LEU:HB2	2.16	0.46
1:C:90:LYS:HD2	1:C:100:PRO:HB3	1.98	0.45
1:B:73:PHE:O	1:B:105:ILE:HA	2.16	0.45
1:B:94:ALA:O	1:B:150:GLN:NE2	2.49	0.45
1:B:28:ARG:HD3	1:B:136:ALA:HA	1.98	0.45
1:C:15:TRP:CZ3	1:C:17:ALA:HB2	2.51	0.45
1:D:67:CYS:O	1:D:68:ASP:HB2	2.16	0.45
1:D:31:ARG:NH2	1:D:138:ARG:NE	2.64	0.45
1:D:42:HIS:CD2	1:D:114:ARG:CZ	3.00	0.44
1:D:98:TRP:O	1:D:102:SER:OG	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:103:ILE:HG22	1:D:103:ILE:O	2.18	0.44
1:D:68:ASP:O	1:D:69:VAL:C	2.55	0.44
1:A:112:MET:CE	1:A:157:ASP:OD2	2.65	0.44
1:B:95:GLY:HA3	1:B:150:GLN:HE22	1.83	0.44
1:D:75:ILE:HD12	1:D:81:VAL:HG22	2.00	0.44
1:D:163:ILE:O	1:D:166:TRP:HB2	2.18	0.44
1:A:126:ASN:ND2	1:B:8:SER:N	2.65	0.44
1:B:32:TYR:HB2	1:B:102:SER:HB3	2.00	0.44
1:B:38:THR:O	1:B:39:ALA:HB3	2.18	0.43
1:C:153:LEU:HD23	1:C:153:LEU:HA	1.86	0.43
1:B:131:GLY:O	1:B:136:ALA:N	2.43	0.43
1:D:160:TYR:CE2	1:D:164:GLN:NE2	2.87	0.43
1:A:155:PRO:HB2	1:A:159:LEU:HD23	2.00	0.43
1:A:126:ASN:HD21	1:B:8:SER:N	2.09	0.43
1:C:24:GLU:HG2	1:C:90:LYS:HE3	2.01	0.43
1:A:37:HIS:CD2	1:A:110:ASN:HA	2.54	0.43
1:C:59:TYR:OH	1:D:151:PRO:HA	2.19	0.43
1:B:145:GLY:O	1:B:148:ASP:HB2	2.17	0.43
1:C:103:ILE:HD13	1:C:103:ILE:HA	1.77	0.43
1:B:77:GLU:HG2	1:B:117:PRO:HD2	2.00	0.42
1:A:67:CYS:O	1:A:68:ASP:HB2	2.19	0.42
1:B:138:ARG:O	1:B:141:TYR:HB3	2.19	0.42
1:A:124:ALA:O	1:A:128:LEU:HG	2.20	0.42
1:D:158:ARG:HA	1:D:158:ARG:HD2	1.74	0.42
1:B:133:ALA:C	1:B:135:GLY:N	2.72	0.42
1:A:99:ASN:N	1:A:100:PRO:CD	2.83	0.42
1:A:2:ASP:HA	1:A:3:PRO:HD3	1.88	0.42
1:A:43:CYS:HB2	1:A:48:SER:OG	2.20	0.41
1:C:15:TRP:CE3	1:C:17:ALA:HB2	2.55	0.41
1:D:12:ARG:NH2	1:D:20:SER:HB2	2.35	0.41
1:A:105:ILE:O	1:A:105:ILE:HG13	2.20	0.41
1:B:82:TYR:N	1:B:82:TYR:CD2	2.88	0.41
1:C:37:HIS:CE1	1:C:154:SER:HA	2.55	0.41
1:D:3:PRO:CB	1:D:4:PRO:HD3	2.48	0.41
1:A:81:VAL:HG23	1:A:123:ALA:HB1	2.02	0.41
1:B:159:LEU:HD12	1:B:163:ILE:HG23	2.01	0.41
1:B:2:ASP:HB2	1:B:3:PRO:CD	2.48	0.41
1:B:132:VAL:O	1:B:135:GLY:CA	2.69	0.41
1:B:144:LYS:HZ3	1:B:171:ALA:CB	2.33	0.41
1:B:166:TRP:HB2	1:B:169:TYR:HB2	2.01	0.41
1:D:53:ALA:HA	1:D:108:MET:HE3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:THR:OG1	1:A:153:LEU:N	2.53	0.41
1:D:3:PRO:HB2	1:D:4:PRO:CD	2.39	0.41
1:D:15:TRP:CH2	1:D:17:ALA:HB2	2.56	0.41
1:A:51:GLN:O	1:A:54:GLN:HB3	2.21	0.41
1:A:59:TYR:O	1:A:63:ASN:HB2	2.21	0.41
1:B:58:SER:O	1:B:62:ARG:HB2	2.20	0.41
1:C:28:ARG:NH2	1:C:88:ASN:HD21	2.04	0.41
1:A:77:GLU:C	1:A:79:GLY:H	2.21	0.41
1:B:9:ILE:CG2	1:B:83:GLU:HB2	2.51	0.40
1:B:68:ASP:OD2	1:B:85:ARG:HG2	2.21	0.40
1:B:140:ASN:OD1	1:B:140:ASN:N	2.52	0.40
1:B:158:ARG:O	1:B:161:GLU:HG2	2.20	0.40
1:C:146:HIS:HB3	1:C:154:SER:HB3	2.02	0.40
1:D:117:PRO:O	1:D:118:PRO:C	2.59	0.40
1:A:45:THR:O	1:A:49:CYS:HB2	2.21	0.40
1:B:15:TRP:CZ2	1:B:17:ALA:HB2	2.56	0.40
1:A:15:TRP:CZ2	1:A:17:ALA:HB2	2.56	0.40
1:C:77:GLU:HG2	1:C:117:PRO:CD	2.41	0.40
1:C:146:HIS:ND1	1:C:154:SER:HB3	2.36	0.40
2:E:1:SGN:O1S	2:E:1:SGN:C3	2.70	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	169/171 (99%)	160 (95%)	9 (5%)	0	100 100
1	B	169/171 (99%)	150 (89%)	18 (11%)	1 (1%)	25 56
1	C	169/171 (99%)	154 (91%)	15 (9%)	0	100 100
1	D	169/171 (99%)	148 (88%)	19 (11%)	2 (1%)	13 39
All	All	676/684 (99%)	612 (90%)	61 (9%)	3 (0%)	34 66



All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	7	GLY
1	D	3	PRO
1	B	170	ARG

### 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	139/139 (100%)	135 (97%)	4 (3%)	42 76
1	B	139/139 (100%)	124 (89%)	15 (11%)	6 19
1	C	139/139 (100%)	135 (97%)	4 (3%)	42 76
1	D	139/139 (100%)	128 (92%)	11 (8%)	12 34
All	All	556/556 (100%)	522 (94%)	34 (6%)	18 48

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	GLU
1	A	37	HIS
1	A	142	GLU
1	A	170	ARG
1	B	16	ARG
1	B	21	GLU
1	B	27	THR
1	B	28	ARG
1	B	30	VAL
1	B	31	ARG
1	B	37	HIS
1	B	41	SER
1	B	42	HIS
1	B	119	ARG
1	B	126	ASN
1	B	159	LEU
1	B	164	GLN

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Mol	Chain	Res	Type
1	B	168	HIS
1	B	170	ARG
1	C	13	ARG
1	C	37	HIS
1	C	105	ILE
1	C	163	ILE
1	D	8	SER
1	D	13	ARG
1	D	18	LEU
1	D	37	HIS
1	D	58	SER
1	D	88	ASN
1	D	103	ILE
1	D	113	ASN
1	D	118	PRO
1	D	119	ARG
1	D	154	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	HIS
1	A	42	HIS
1	A	55	ASN
1	A	72	ASN
1	A	126	ASN
1	A	150	GLN
1	A	168	HIS
1	B	37	HIS
1	B	72	ASN
1	B	150	GLN
1	C	37	HIS
1	C	63	ASN
1	C	72	ASN
1	C	88	ASN
1	C	168	HIS
1	D	37	HIS
1	D	63	ASN
1	D	72	ASN
1	D	88	ASN
1	D	164	GLN
1	D	168	HIS

### 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](#)

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
2	SGN	E	1	2	19,20,20	2.16	2 (10%)	24,31,31	1.75	2 (8%)
2	UAP	E	2	2	15,15,16	2.06	2 (13%)	18,22,24	3.67	7 (38%)

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	SGN	E	1	2	-	4/11/31/31	0/1/1/1
2	UAP	E	2	2	-	1/9/22/25	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	2	UAP	O5-C5	7.09	1.47	1.37
2	E	1	SGN	O2S-S1	6.47	1.49	1.42
2	E	1	SGN	O1S-S1	5.78	1.48	1.42
2	E	2	UAP	O2-C2	-2.02	1.44	1.47

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	2	UAP	O5-C5-C4	-11.45	115.14	124.81
2	E	2	UAP	O5-C5-C6	8.01	123.54	111.52
2	E	1	SGN	O1S-S1-O2S	-6.05	105.85	120.16
2	E	2	UAP	C1-C2-C3	4.01	112.69	108.98
2	E	1	SGN	C1-C2-C3	3.81	115.74	110.54
2	E	2	UAP	O6B-C6-C5	2.73	121.02	114.20
2	E	2	UAP	C3-C4-C5	-2.73	116.98	121.60
2	E	2	UAP	C2-O2-S	2.22	120.81	117.91
2	E	2	UAP	O6B-C6-O6A	-2.01	119.02	123.61

There are no chirality outliers.

All (5) torsion outliers are listed below:

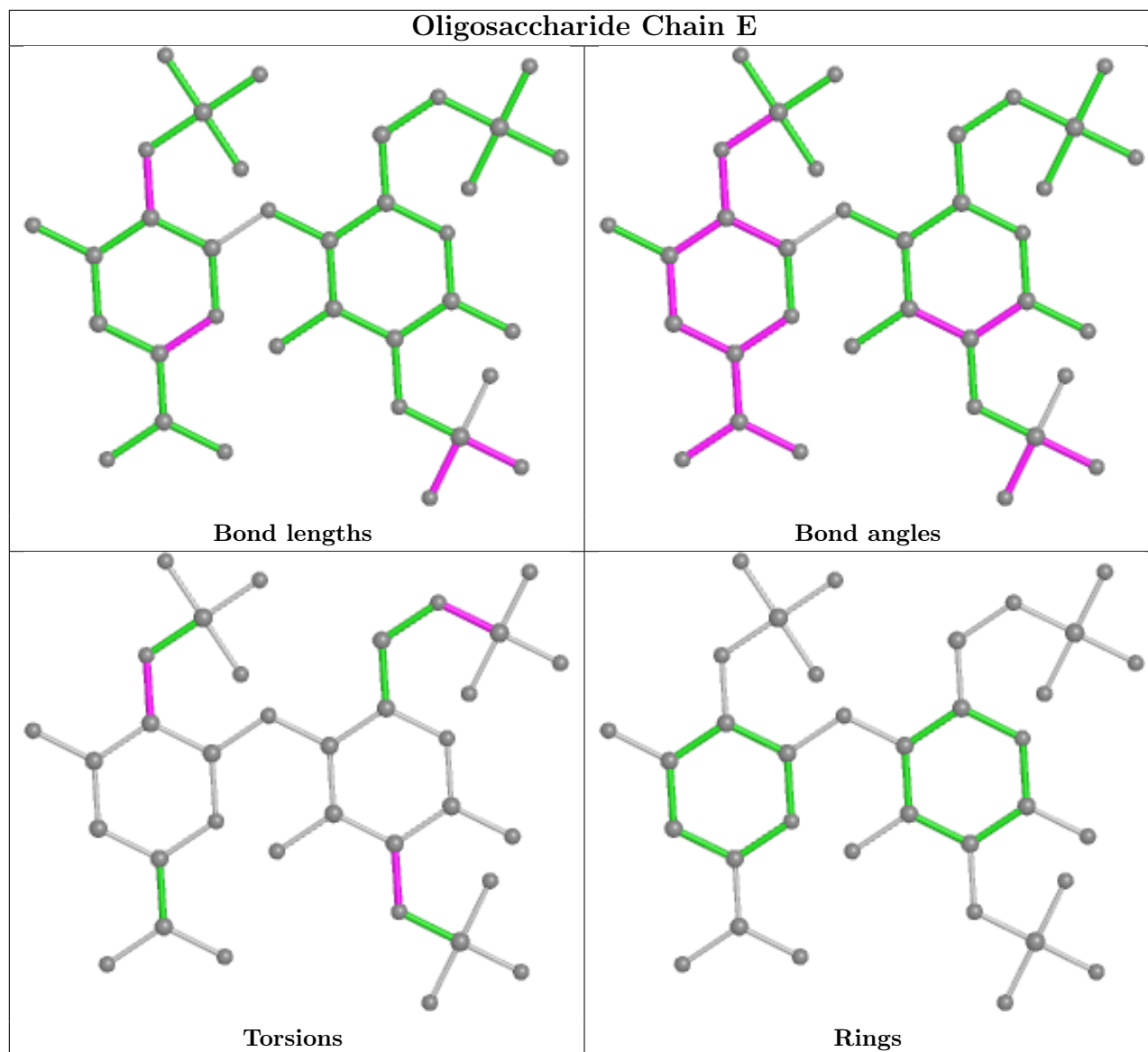
Mol	Chain	Res	Type	Atoms
2	E	1	SGN	C3-C2-N2-S1
2	E	2	UAP	C3-C2-O2-S
2	E	1	SGN	C6-O6-S2-O4S
2	E	1	SGN	C6-O6-S2-O6S
2	E	1	SGN	C6-O6-S2-O5S

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	2	UAP	1	0
2	E	1	SGN	2	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](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	GOL	D	172	-	5,5,5	0.35	0	5,5,5	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	TLA	C	172	-	9,9,9	2.21	4 (44%)	12,12,12	1.98	4 (33%)
3	GOL	A	172	-	5,5,5	0.36	0	5,5,5	0.26	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	GOL	D	172	-	-	1/4/4/4	-
4	TLA	C	172	-	-	2/12/12/12	-
3	GOL	A	172	-	-	0/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	172	TLA	C3-C4	-3.74	1.47	1.52
4	C	172	TLA	C2-C1	-3.62	1.47	1.52
4	C	172	TLA	O11-C1	-2.85	1.21	1.30
4	C	172	TLA	O41-C4	-2.37	1.22	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	172	TLA	O3-C3-C4	-3.50	103.32	110.66
4	C	172	TLA	O41-C4-C3	3.23	122.00	113.27
4	C	172	TLA	O11-C1-C2	2.56	120.20	113.27
4	C	172	TLA	O4-C4-C3	-2.42	115.28	121.63

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	172	TLA	C2-C3-C4-O41
4	C	172	TLA	C2-C3-C4-O4
3	D	172	GOL	C1-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	172	TLA	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	171/171 (100%)	-0.18	5 (2%) 51 41	15, 33, 57, 99	0
1	B	171/171 (100%)	-0.16	6 (3%) 44 34	21, 43, 72, 99	0
1	C	171/171 (100%)	-0.46	4 (2%) 60 51	12, 26, 52, 99	0
1	D	171/171 (100%)	-0.31	6 (3%) 44 34	18, 32, 56, 99	0
All	All	684/684 (100%)	-0.28	21 (3%) 49 39	12, 33, 69, 99	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	PRO	9.2
1	D	2	ASP	8.1
1	B	2	ASP	7.4
1	D	1	GLU	7.1
1	A	2	ASP	6.7
1	B	4	PRO	6.5
1	B	3	PRO	6.2
1	C	5	ALA	5.5
1	A	1	GLU	5.4
1	C	2	ASP	3.7
1	B	5	ALA	3.2
1	D	6	CYS	3.1
1	D	5	ALA	3.0
1	A	5	ALA	2.8
1	B	1	GLU	2.6
1	C	1	GLU	2.5
1	A	4	PRO	2.4
1	C	3	PRO	2.3
1	D	3	PRO	2.1
1	B	6	CYS	2.1
1	D	4	PRO	2.1



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

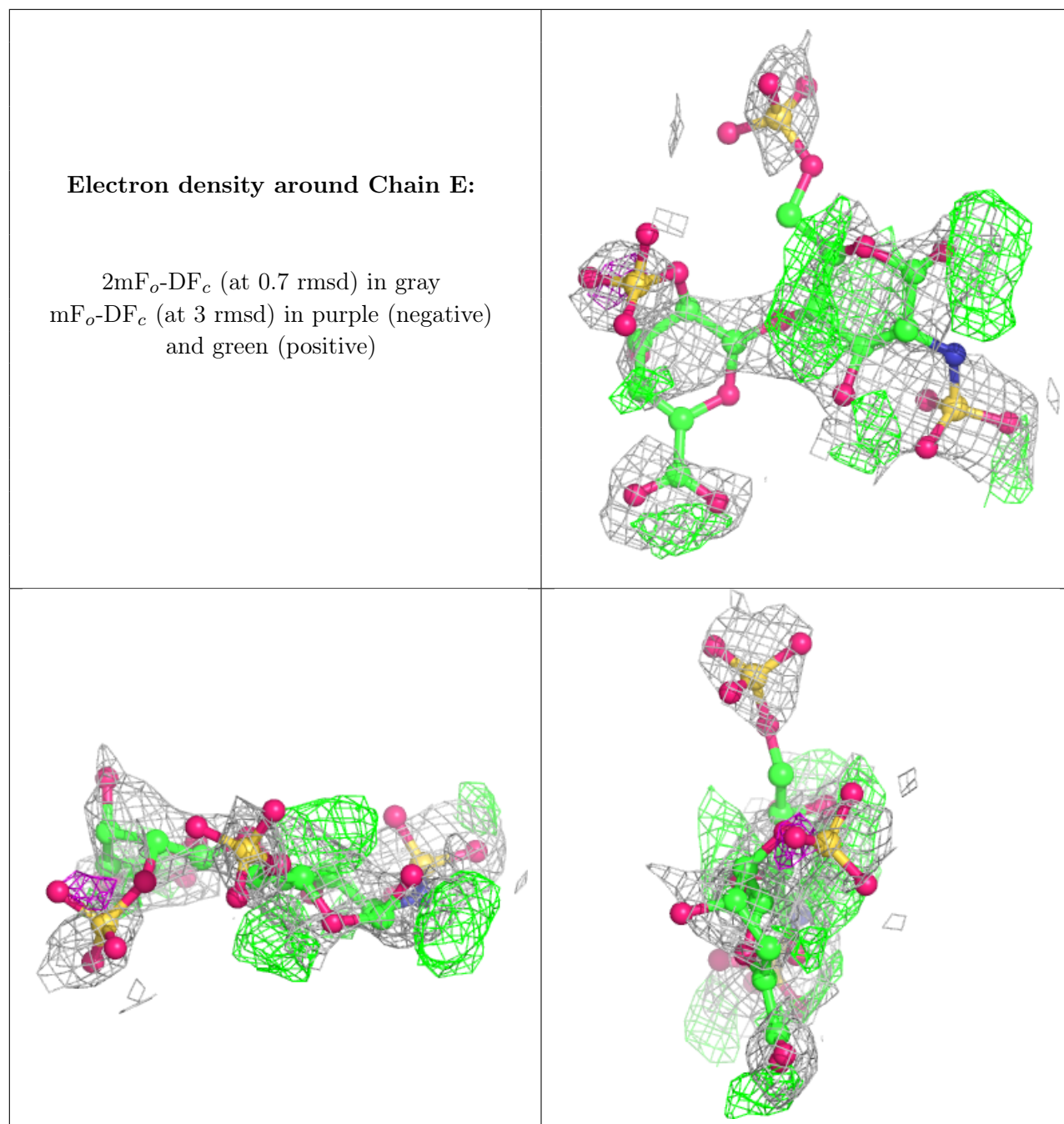
There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	UAP	E	2	15/16	0.69	0.43	35,40,46,46	15
2	SGN	E	1	20/20	0.76	0.38	38,42,49,49	20

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	GOL	A	172	6/6	0.72	0.27	66,68,69,72	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	TLA	C	172	10/10	0.89	0.19	26,38,44,48	2
3	GOL	D	172	6/6	0.92	0.22	36,40,43,50	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.