



Full wwPDB X-ray Structure Validation Report ⓘ

May 3, 2025 – 03:42 PM EDT

PDB ID : 3A1J / pdb_00003a1j
Title : Crystal structure of the human Rad9-Hus1-Rad1 complex
Authors : Sohn, S.Y.; Cho, Y.
Deposited on : 2009-04-08
Resolution : 2.50 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 Phenix2.0rc1
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

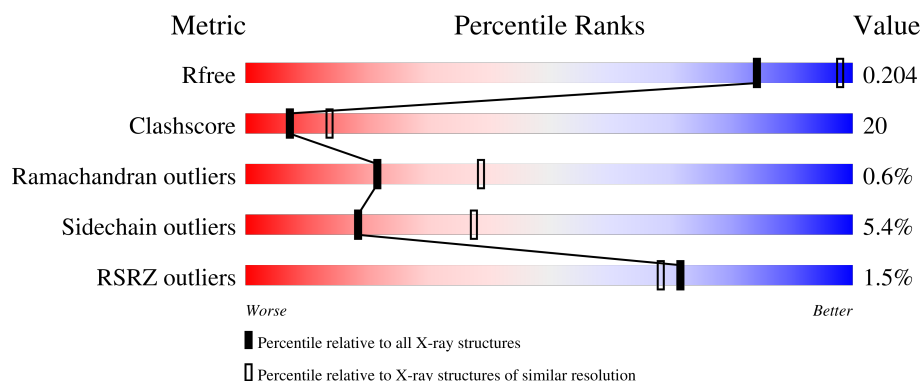
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	164625	5504 (2.50-2.50)
Clashscore	180529	6282 (2.50-2.50)
Ramachandran outliers	177936	6191 (2.50-2.50)
Sidechain outliers	177891	6193 (2.50-2.50)
RSRZ outliers	164620	5504 (2.50-2.50)

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	266	<div> <div>0%</div> <div> <div>64%</div> <div>33%</div> <div>.</div> </div> </div>
2	B	281	<div> <div>2%</div> <div> <div>54%</div> <div>37%</div> <div>...</div> </div> </div>
3	C	263	<div> <div>2%</div> <div> <div>62%</div> <div>34%</div> <div>.</div> </div> </div>
4	D	2	<div> <div>100%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6467 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 Cell cycle checkpoint control protein RAD9A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	Se	0	0	0
			2044	1299	356	373	10	6			

- Molecule 2 is a protein called Checkpoint protein HUS1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	269	Total	C	N	O	S	Se	0	0	0
			2138	1366	364	392	7	9			

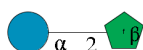
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	HIS	-	expression tag	UNP O60921

- Molecule 3 is a protein called Cell cycle checkpoint protein RAD1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	263	Total	C	N	O	S	Se	0	0	0
			2076	1315	341	402	11	7			

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



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

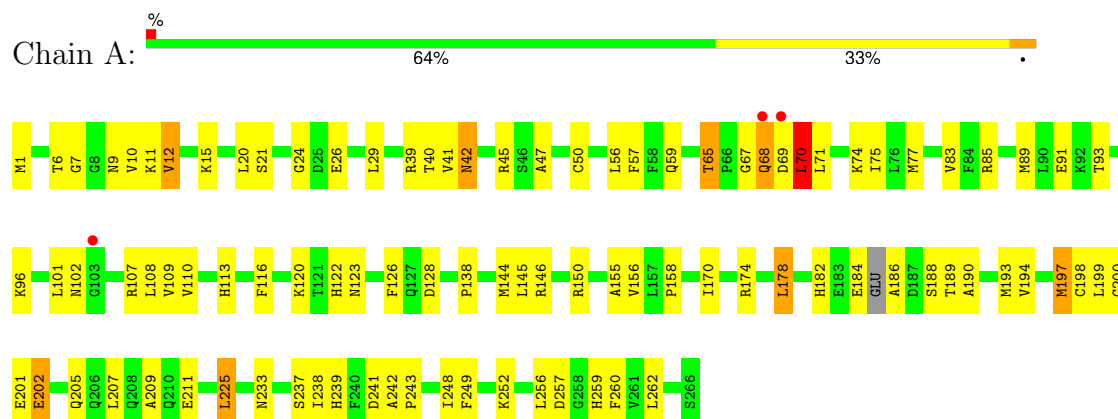
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	64	Total 64	O 64	0	0
5	B	54	Total 54	O 54	0	0
5	C	68	Total 68	O 68	0	0

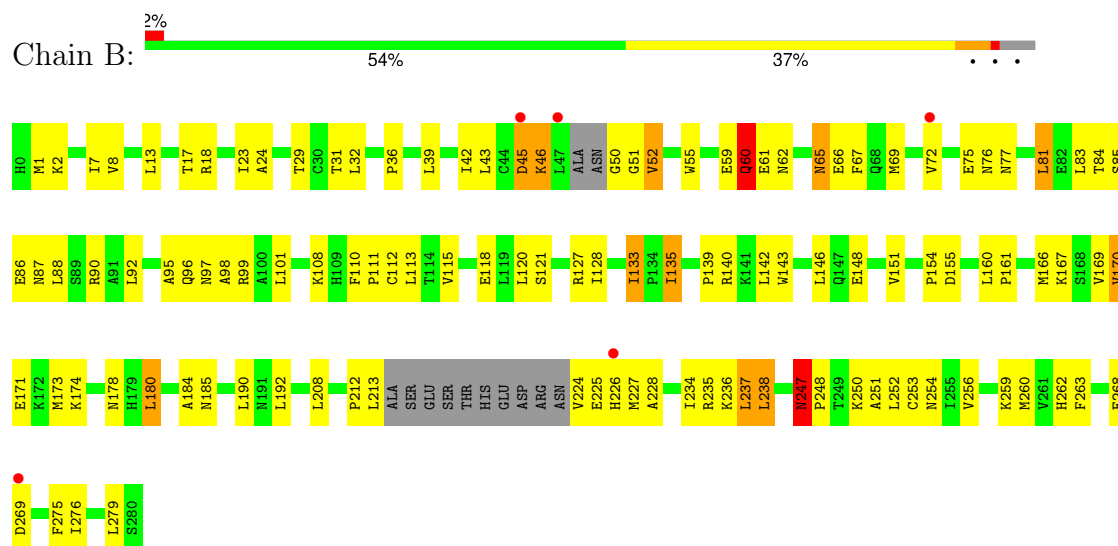
3 Residue-property plots [i](#)

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

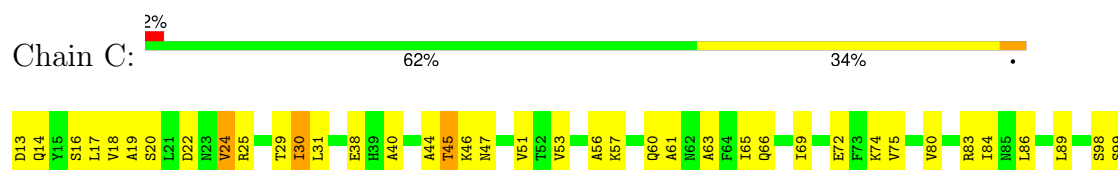
- Molecule 1: Cell cycle checkpoint control protein RAD9A

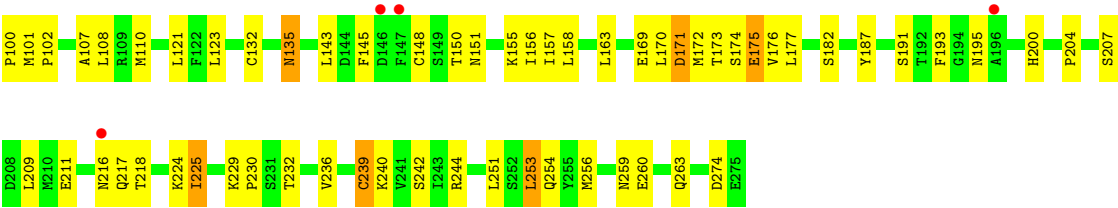


- Molecule 2: Checkpoint protein HUS1

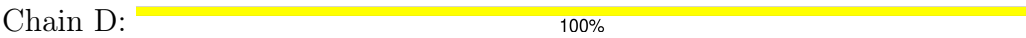


- Molecule 3: Cell cycle checkpoint protein RAD1





● Molecule 4: beta-D-fructofuranose-(2-1)-alpha-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	77.15Å 70.78Å 86.66Å 90.00° 99.20° 90.00°	Depositor
Resolution (Å)	19.90 – 2.50 19.90 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.2 (19.90-2.50) 95.0 (19.90-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.16 (at 2.50Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.209 , 0.282 0.202 , 0.204	Depositor DCC
R_{free} test set	1555 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	43.3	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6467	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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: GLC, FRU

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	0.44	0/2075	0.94	6/2788 (0.2%)
2	B	0.43	0/2166	0.94	9/2912 (0.3%)
3	C	0.42	0/2107	0.90	4/2837 (0.1%)
All	All	0.43	0/6348	0.93	19/8537 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	THR	N-CA-C	7.91	120.99	111.82
2	B	113	LEU	N-CA-C	-7.05	95.51	107.99
1	A	109	VAL	N-CA-C	6.37	117.29	107.99
3	C	56	ALA	N-CA-C	6.24	119.97	111.17
1	A	71	LEU	N-CA-C	6.18	123.96	110.80
1	A	57	PHE	N-CA-C	-5.94	104.75	112.23
3	C	22	ASP	N-CA-C	-5.88	106.06	113.18
3	C	225	ILE	N-CA-C	5.64	117.05	110.62
2	B	247	ASN	CA-C-N	5.64	125.59	119.78
2	B	247	ASN	C-N-CA	5.64	125.59	119.78
2	B	42	ILE	N-CA-C	5.58	116.15	108.12
2	B	228	ALA	N-CA-C	-5.51	100.25	109.24
2	B	226	HIS	N-CA-C	-5.41	102.87	110.50
1	A	59	GLN	N-CA-C	-5.38	105.45	112.23
3	C	239	CYS	N-CA-C	-5.35	106.64	114.39
2	B	155	ASP	N-CA-C	-5.30	105.50	111.28
1	A	128	ASP	N-CA-C	-5.24	102.64	110.23
2	B	133	ILE	N-CA-C	5.19	113.00	107.76
2	B	65	ASN	N-CA-C	-5.18	107.01	113.38

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2044	0	2079	82	0
2	B	2138	0	2205	103	0
3	C	2076	0	2050	82	0
4	D	23	0	21	0	0
5	A	64	0	0	7	0
5	B	54	0	0	1	0
5	C	68	0	0	4	0
All	All	6467	0	6355	257	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 20.

All (257) 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)
2:B:169:VAL:HG12	2:B:173:MSE:HE2	1.39	1.05
1:A:197:MSE:HE2	1:A:198:CYS:N	1.82	0.95
2:B:13:LEU:HD21	2:B:95:ALA:HB1	1.56	0.88
3:C:177:LEU:HB2	3:C:225:ILE:HD13	1.52	0.88
3:C:173:THR:HG21	3:C:195:ASN:H	1.39	0.87
1:A:6:THR:HG22	1:A:96:LYS:HG2	1.55	0.87
2:B:81:LEU:HD23	2:B:135:ILE:HD13	1.54	0.86
3:C:239:CYS:SG	3:C:256:MSE:HE2	2.18	0.83
1:A:107:ARG:HD3	1:A:123:ASN:HB3	1.59	0.82
2:B:43:LEU:HD13	2:B:45:ASP:H	1.43	0.82
1:A:197:MSE:HE3	2:B:128:ILE:O	1.79	0.82
1:A:197:MSE:HE2	1:A:198:CYS:H	1.46	0.81
2:B:43:LEU:HD11	2:B:45:ASP:HB3	1.65	0.79
3:C:172:MSE:HE2	3:C:229:LYS:HD2	1.62	0.79
1:A:11:LYS:HE2	1:A:15:LYS:HE2	1.64	0.79
2:B:76:ASN:HD22	2:B:140:ARG:NH1	1.80	0.79
2:B:13:LEU:CD2	2:B:95:ALA:HB1	2.15	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:GLU:HG2	1:A:74:LYS:HE3	1.68	0.75
1:A:7:GLY:O	1:A:10:VAL:HG12	1.85	0.75
3:C:171:ASP:OD1	3:C:173:THR:HB	1.88	0.74
2:B:69:MSE:HE1	2:B:77:ASN:ND2	2.02	0.74
3:C:45:THR:HG22	3:C:47:ASN:H	1.51	0.74
2:B:247:ASN:HD22	2:B:248:PRO:HD2	1.52	0.73
3:C:31:LEU:HD21	3:C:51:VAL:HG21	1.69	0.73
3:C:66:GLN:HG2	3:C:69:ILE:HG12	1.72	0.72
3:C:148:CYS:HB2	3:C:151:ASN:ND2	2.04	0.71
2:B:55:TRP:CD2	2:B:148:GLU:HB3	2.26	0.71
2:B:174:LYS:HD2	2:B:238:LEU:HD12	1.73	0.69
1:A:170:ILE:HG12	1:A:209:ALA:HB3	1.75	0.68
2:B:46:LYS:HA	2:B:51:GLY:HA2	1.75	0.68
1:A:193:MSE:CE	2:B:90:ARG:CZ	2.71	0.68
2:B:166:MSE:O	2:B:170:VAL:HG13	1.93	0.68
2:B:256:VAL:HG12	2:B:259:LYS:HB2	1.76	0.67
2:B:60:GLN:HG3	2:B:67:PHE:CG	2.31	0.65
1:A:146:ARG:HE	1:A:205:GLN:NE2	1.94	0.65
3:C:173:THR:CG2	3:C:195:ASN:H	2.09	0.65
2:B:84:THR:O	2:B:87:ASN:HB2	1.97	0.65
2:B:60:GLN:HG3	2:B:67:PHE:CD2	2.33	0.64
2:B:75:GLU:HG3	2:B:76:ASN:N	2.12	0.64
2:B:170:VAL:HA	2:B:173:MSE:HE3	1.80	0.64
2:B:248:PRO:HG2	2:B:251:ALA:HB2	1.81	0.63
3:C:224:LYS:HE3	3:C:274:ASP:OD1	1.99	0.63
1:A:89:MSE:HE3	3:C:169:GLU:HG3	1.82	0.62
1:A:197:MSE:HE3	2:B:128:ILE:C	2.23	0.62
1:A:182:HIS:HD2	5:A:275:HOH:O	1.82	0.62
1:A:138:PRO:HG3	1:A:248:ILE:HD13	1.83	0.61
3:C:155:LYS:NZ	3:C:157:ILE:HD11	2.16	0.61
2:B:254:ASN:HD22	2:B:262:HIS:CD2	2.18	0.61
1:A:150:ARG:HG2	5:A:287:HOH:O	2.00	0.61
3:C:44:ALA:HB3	3:C:80:VAL:CG1	2.31	0.61
2:B:32:LEU:HD11	2:B:39:LEU:HG	1.82	0.60
1:A:68:GLN:C	1:A:68:GLN:HE21	2.08	0.60
1:A:83:VAL:HG11	1:A:110:VAL:HG11	1.82	0.60
1:A:259:HIS:HB2	5:A:278:HOH:O	2.00	0.60
3:C:29:THR:HB	5:C:294:HOH:O	2.02	0.60
1:A:184:GLU:HG2	1:A:186:ALA:HB3	1.84	0.59
3:C:24:VAL:HG11	3:C:123:LEU:HD23	1.84	0.59
1:A:197:MSE:CE	2:B:128:ILE:O	2.51	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:31:THR:HG21	2:B:146:LEU:HD13	1.85	0.59
2:B:254:ASN:HD22	2:B:262:HIS:HD2	1.51	0.58
3:C:65:ILE:HD12	3:C:65:ILE:N	2.19	0.58
3:C:155:LYS:HZ1	3:C:157:ILE:HD11	1.68	0.58
3:C:45:THR:CG2	3:C:47:ASN:H	2.14	0.57
2:B:154:PRO:HB3	2:B:254:ASN:HB3	1.85	0.57
2:B:69:MSE:HE1	2:B:77:ASN:HD21	1.68	0.57
2:B:224:VAL:HG12	2:B:225:GLU:N	2.20	0.57
2:B:174:LYS:HE3	2:B:235:ARG:HG2	1.85	0.57
1:A:116:PHE:CD2	3:C:209:LEU:HD11	2.40	0.56
3:C:158:LEU:HB3	3:C:163:LEU:HD11	1.86	0.56
2:B:75:GLU:HG3	2:B:76:ASN:H	1.70	0.56
1:A:144:MSE:HA	1:A:238:ILE:O	2.05	0.56
3:C:211:GLU:OE1	3:C:240:LYS:HE2	2.06	0.56
1:A:174:ARG:HD3	1:A:201:GLU:CD	2.31	0.56
3:C:45:THR:HG23	3:C:46:LYS:N	2.21	0.56
1:A:256:LEU:C	1:A:256:LEU:HD12	2.30	0.56
1:A:42:ASN:HB2	1:A:47:ALA:H	1.71	0.55
3:C:99:SER:N	3:C:100:PRO:HD3	2.21	0.55
2:B:24:ALA:HA	2:B:85:SER:HB3	1.87	0.55
2:B:185:ASN:ND2	2:B:227:MSE:HE3	2.22	0.55
3:C:40:ALA:HB2	3:C:53:VAL:HG12	1.89	0.55
2:B:160:LEU:HD21	2:B:253:CYS:HB2	1.89	0.54
3:C:98:SER:C	3:C:100:PRO:HD3	2.32	0.54
1:A:145:LEU:CD2	1:A:207:LEU:HG	2.37	0.54
1:A:174:ARG:NH2	1:A:207:LEU:O	2.41	0.54
1:A:101:LEU:C	1:A:102:ASN:HD22	2.16	0.54
1:A:26:GLU:HG2	1:A:74:LYS:CE	2.36	0.53
2:B:81:LEU:CD2	2:B:135:ILE:HD13	2.35	0.53
1:A:1:MSE:N	1:A:65:THR:HB	2.24	0.53
2:B:268:GLU:O	2:B:269:ASP:CG	2.51	0.53
2:B:43:LEU:HD13	2:B:43:LEU:C	2.34	0.53
2:B:260:MSE:HG3	2:B:276:ILE:O	2.09	0.53
3:C:19:ALA:HA	3:C:72:GLU:O	2.09	0.52
2:B:52:VAL:HG23	2:B:275:PHE:O	2.09	0.52
3:C:135:ASN:N	3:C:135:ASN:HD22	2.07	0.52
3:C:216:ASN:HB2	5:C:322:HOH:O	2.08	0.52
2:B:8:VAL:O	2:B:8:VAL:HG22	2.10	0.52
2:B:160:LEU:CD2	2:B:253:CYS:HB2	2.40	0.52
3:C:110:MSE:HG2	3:C:121:LEU:CD2	2.40	0.52
2:B:160:LEU:HD11	2:B:263:PHE:HD1	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:250:LYS:HE3	2:B:252:LEU:HD12	1.92	0.52
2:B:59:GLU:HB2	2:B:62:ASN:ND2	2.25	0.52
1:A:193:MSE:HE1	2:B:90:ARG:CZ	2.38	0.51
2:B:161:PRO:HB3	2:B:208:LEU:HD21	1.91	0.51
1:A:1:MSE:N	1:A:65:THR:HG22	2.25	0.51
2:B:99:ARG:HD3	2:B:118:GLU:OE2	2.11	0.51
1:A:146:ARG:HE	1:A:205:GLN:HE21	1.58	0.51
2:B:65:ASN:OD1	2:B:66:GLU:HG3	2.11	0.51
3:C:24:VAL:HG11	3:C:123:LEU:CD2	2.40	0.51
1:A:39:ARG:HG2	1:A:50:CYS:SG	2.51	0.51
1:A:197:MSE:CE	1:A:198:CYS:H	2.21	0.50
2:B:1:MSE:HE2	2:B:72:VAL:HG12	1.92	0.50
3:C:158:LEU:CB	3:C:163:LEU:HD11	2.41	0.50
2:B:2:LYS:O	2:B:69:MSE:HA	2.11	0.50
2:B:7:ILE:HD12	2:B:101:LEU:HD23	1.94	0.50
1:A:68:GLN:C	1:A:68:GLN:NE2	2.69	0.50
2:B:151:VAL:HG23	2:B:151:VAL:O	2.12	0.50
2:B:178:ASN:HA	2:B:234:ILE:CG2	2.41	0.49
1:A:170:ILE:HG22	1:A:211:GLU:HA	1.94	0.49
3:C:83:ARG:HG3	3:C:83:ARG:HH11	1.76	0.49
3:C:17:LEU:HD12	3:C:18:VAL:N	2.28	0.49
3:C:155:LYS:HG3	3:C:244:ARG:HG2	1.95	0.49
2:B:160:LEU:HD11	2:B:263:PHE:CD1	2.47	0.49
3:C:24:VAL:CG1	3:C:108:LEU:HB2	2.42	0.48
3:C:13:ASP:O	3:C:13:ASP:OD1	2.31	0.48
1:A:150:ARG:NH2	1:A:233:ASN:HB2	2.29	0.48
2:B:184:ALA:HA	2:B:227:MSE:HE2	1.96	0.48
3:C:110:MSE:HG2	3:C:121:LEU:HD23	1.95	0.48
1:A:225:LEU:HG	1:A:260:PHE:CD1	2.49	0.48
2:B:121:SER:HB2	5:B:318:HOH:O	2.12	0.48
1:A:120:LYS:HE2	1:A:122:HIS:CE1	2.49	0.48
3:C:45:THR:CG2	3:C:46:LYS:N	2.76	0.47
1:A:249:PHE:CE2	1:A:262:LEU:HD12	2.49	0.47
3:C:57:LYS:HA	3:C:143:LEU:HD11	1.97	0.47
3:C:187:TYR:HD2	3:C:204:PRO:HA	1.79	0.47
3:C:175:GLU:HG3	3:C:176:VAL:HG13	1.97	0.47
2:B:173:MSE:HE1	2:B:192:LEU:HB3	1.97	0.47
2:B:268:GLU:C	2:B:269:ASP:CG	2.81	0.47
3:C:148:CYS:C	3:C:150:THR:H	2.20	0.47
2:B:36:PRO:CA	2:B:69:MSE:HE2	2.45	0.47
3:C:174:SER:CB	3:C:193:PHE:H	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MSE:H1	1:A:65:THR:CB	2.28	0.46
1:A:170:ILE:N	1:A:170:ILE:HD12	2.30	0.46
2:B:139:PRO:HG2	2:B:142:LEU:HG	1.97	0.46
1:A:10:VAL:HG13	1:A:11:LYS:N	2.30	0.46
1:A:42:ASN:CG	1:A:47:ALA:HB3	2.39	0.46
2:B:76:ASN:HB2	2:B:140:ARG:NH1	2.30	0.46
2:B:247:ASN:ND2	2:B:248:PRO:HD2	2.26	0.46
3:C:18:VAL:HB	3:C:74:LYS:HG2	1.97	0.46
3:C:121:LEU:HB2	3:C:132:CYS:HB2	1.97	0.46
2:B:190:LEU:C	2:B:190:LEU:HD23	2.41	0.46
2:B:76:ASN:HD22	2:B:140:ARG:CZ	2.27	0.46
1:A:197:MSE:HE1	2:B:127:ARG:HB3	1.98	0.46
1:A:207:LEU:C	1:A:207:LEU:HD23	2.40	0.46
2:B:259:LYS:O	2:B:260:MSE:HB2	2.14	0.46
3:C:89:LEU:HD12	3:C:89:LEU:O	2.15	0.46
3:C:232:THR:O	3:C:236:VAL:HG23	2.15	0.46
1:A:252:LYS:HG3	1:A:257:ASP:OD1	2.16	0.46
3:C:148:CYS:C	3:C:150:THR:N	2.74	0.46
2:B:268:GLU:C	2:B:269:ASP:OD1	2.58	0.46
1:A:41:VAL:HG21	1:A:45:ARG:HD3	1.98	0.46
1:A:242:ALA:O	1:A:243:PRO:C	2.57	0.45
1:A:24:GLY:HA3	1:A:40:THR:OG1	2.16	0.45
1:A:102:ASN:HD22	1:A:102:ASN:N	2.12	0.45
3:C:17:LEU:HD12	3:C:18:VAL:H	1.80	0.45
1:A:70:LEU:HD23	5:A:303:HOH:O	2.16	0.45
2:B:112:CYS:HA	2:B:133:ILE:O	2.16	0.45
3:C:259:ASN:ND2	3:C:263:GLN:HB2	2.31	0.45
3:C:191:SER:OG	3:C:200:HIS:HD2	1.99	0.45
1:A:200:GLY:HA3	1:A:202:GLU:OE2	2.17	0.45
2:B:81:LEU:N	2:B:81:LEU:HD12	2.30	0.45
2:B:101:LEU:HD11	2:B:115:VAL:HG13	1.98	0.45
3:C:254:GLN:HG3	5:C:307:HOH:O	2.14	0.45
1:A:9:ASN:HA	1:A:12:VAL:HG13	1.99	0.45
3:C:24:VAL:HG13	3:C:108:LEU:HB2	1.99	0.45
1:A:158:PRO:HG2	1:A:197:MSE:HG3	1.99	0.45
2:B:76:ASN:ND2	2:B:140:ARG:HD3	2.31	0.45
3:C:74:LYS:HG3	5:C:292:HOH:O	2.16	0.45
3:C:83:ARG:HG3	3:C:83:ARG:NH1	2.32	0.45
1:A:85:ARG:HD3	3:C:169:GLU:OE1	2.17	0.45
2:B:212:PRO:O	2:B:213:LEU:HB2	2.15	0.45
1:A:146:ARG:HG3	1:A:237:SER:OG	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:178:ASN:HA	2:B:234:ILE:HG22	1.99	0.45
2:B:43:LEU:CD1	2:B:45:ASP:H	2.23	0.44
3:C:25:ARG:HE	3:C:260:GLU:CD	2.25	0.44
1:A:10:VAL:CG1	5:A:310:HOH:O	2.66	0.44
2:B:43:LEU:CD1	2:B:45:ASP:HB3	2.41	0.44
3:C:229:LYS:N	3:C:230:PRO:HD2	2.32	0.44
3:C:156:ILE:O	3:C:242:SER:HA	2.18	0.44
1:A:102:ASN:N	1:A:102:ASN:ND2	2.65	0.44
1:A:155:ALA:CB	1:A:199:LEU:HD11	2.47	0.44
2:B:60:GLN:CG	2:B:67:PHE:CG	2.99	0.44
1:A:150:ARG:HG3	1:A:150:ARG:HH11	1.82	0.44
2:B:110:PHE:HB2	2:B:111:PRO:HD2	2.00	0.44
1:A:193:MSE:HE3	2:B:90:ARG:NH1	2.33	0.44
2:B:96:GLN:HG3	2:B:97:ASN:N	2.32	0.44
2:B:167:LYS:O	2:B:171:GLU:HB2	2.17	0.44
3:C:84:ILE:HD12	3:C:84:ILE:C	2.42	0.44
3:C:172:MSE:CE	3:C:229:LYS:HD2	2.41	0.44
1:A:1:MSE:N	1:A:65:THR:CG2	2.81	0.44
3:C:60:GLN:HB2	3:C:145:PHE:HE2	1.83	0.44
3:C:107:ALA:O	3:C:123:LEU:HA	2.18	0.44
3:C:135:ASN:HD22	3:C:135:ASN:H	1.66	0.44
2:B:224:VAL:CG1	2:B:225:GLU:N	2.81	0.43
3:C:259:ASN:HD21	3:C:263:GLN:HB2	1.82	0.43
2:B:101:LEU:HD11	2:B:115:VAL:CG1	2.48	0.43
3:C:156:ILE:HG22	3:C:158:LEU:HD22	2.00	0.43
3:C:204:PRO:HG2	3:C:207:SER:HB2	1.99	0.43
2:B:250:LYS:CE	2:B:252:LEU:HD12	2.49	0.43
1:A:1:MSE:N	1:A:65:THR:CB	2.81	0.43
3:C:57:LYS:HA	3:C:143:LEU:CD1	2.49	0.43
1:A:29:LEU:O	1:A:29:LEU:HD23	2.18	0.43
2:B:36:PRO:CB	2:B:69:MSE:HE2	2.49	0.43
2:B:60:GLN:HG3	2:B:67:PHE:CD1	2.54	0.43
1:A:21:SER:HA	1:A:77:MSE:HE3	2.01	0.42
1:A:156:VAL:O	1:A:156:VAL:HG12	2.19	0.42
2:B:140:ARG:HA	2:B:143:TRP:CE2	2.54	0.42
1:A:239:HIS:HD2	5:A:313:HOH:O	2.02	0.42
3:C:16:SER:O	3:C:75:VAL:HA	2.19	0.42
2:B:29:THR:O	2:B:43:LEU:HD23	2.19	0.42
2:B:46:LYS:HD2	2:B:46:LYS:C	2.44	0.42
3:C:171:ASP:OD1	3:C:173:THR:CB	2.64	0.42
2:B:36:PRO:HB3	2:B:69:MSE:HE2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:20:SER:HA	3:C:108:LEU:O	2.19	0.42
3:C:30:ILE:CG2	3:C:63:ALA:HB1	2.49	0.42
3:C:157:ILE:HG12	3:C:242:SER:CB	2.49	0.42
3:C:101:MSE:HE3	3:C:102:PRO:HD2	2.02	0.42
3:C:182:SER:HB2	3:C:218:THR:HG23	2.02	0.42
2:B:23:ILE:C	2:B:23:ILE:HD12	2.45	0.42
3:C:53:VAL:HG22	3:C:61:ALA:HB3	2.01	0.42
2:B:169:VAL:O	2:B:173:MSE:HG3	2.20	0.41
1:A:189:THR:O	1:A:190:ALA:HB3	2.20	0.41
2:B:36:PRO:HD3	2:B:69:MSE:CE	2.51	0.41
2:B:88:LEU:O	2:B:92:LEU:HG	2.19	0.41
1:A:96:LYS:HB2	1:A:113:HIS:HB2	2.01	0.41
1:A:186:ALA:C	1:A:188:SER:H	2.28	0.41
3:C:155:LYS:HG3	3:C:244:ARG:CG	2.51	0.41
1:A:241:ASP:HB3	1:A:242:ALA:H	1.59	0.41
2:B:248:PRO:CG	2:B:251:ALA:HB2	2.49	0.41
3:C:174:SER:HB2	3:C:193:PHE:H	1.85	0.41
1:A:42:ASN:HD22	1:A:42:ASN:HA	1.52	0.41
1:A:75:ILE:CD1	1:A:108:LEU:HD22	2.51	0.41
2:B:95:ALA:HA	2:B:98:ALA:HB2	2.02	0.41
3:C:251:LEU:HG	3:C:253:LEU:HD13	2.03	0.41
1:A:67:GLY:O	1:A:69:ASP:OD1	2.38	0.41
1:A:145:LEU:HD22	1:A:207:LEU:HG	2.01	0.41
2:B:17:THR:HA	2:B:92:LEU:HD13	2.03	0.41
3:C:157:ILE:HG12	3:C:242:SER:HB2	2.02	0.41
2:B:50:GLY:HA3	2:B:236:LYS:NZ	2.35	0.41
2:B:76:ASN:HD22	2:B:140:ARG:HH11	1.63	0.41
1:A:75:ILE:HG12	1:A:126:PHE:HB3	2.03	0.40
1:A:178:LEU:HD12	1:A:178:LEU:HA	1.87	0.40
1:A:197:MSE:HE2	1:A:197:MSE:C	2.45	0.40
2:B:18:ARG:NH2	2:B:268:GLU:OE2	2.45	0.40
2:B:180:LEU:HD22	2:B:237:LEU:HD12	2.04	0.40
3:C:216:ASN:O	3:C:217:GLN:HB2	2.20	0.40
1:A:174:ARG:HA	5:A:307:HOH:O	2.22	0.40
2:B:192:LEU:HD23	2:B:192:LEU:HA	1.94	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	261/266 (98%)	244 (94%)	15 (6%)	2 (1%)	16	31
2	B	263/281 (94%)	243 (92%)	18 (7%)	2 (1%)	16	31
3	C	261/263 (99%)	245 (94%)	15 (6%)	1 (0%)	30	49
All	All	785/810 (97%)	732 (93%)	48 (6%)	5 (1%)	22	39

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	171	ASP
1	A	70	LEU
1	A	91	GLU
2	B	60	GLN
2	B	108	LYS

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	224/219 (102%)	212 (95%)	12 (5%)	18	37
2	B	248/249 (100%)	232 (94%)	16 (6%)	14	29
3	C	237/230 (103%)	227 (96%)	10 (4%)	25	49
All	All	709/698 (102%)	671 (95%)	38 (5%)	18	37

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	20	LEU
1	A	42	ASN
1	A	56	LEU
1	A	65	THR
1	A	68	GLN
1	A	70	LEU
1	A	178	LEU
1	A	194	VAL
1	A	197	MSE
1	A	202	GLU
1	A	225	LEU
2	B	45	ASP
2	B	46	LYS
2	B	52	VAL
2	B	60	GLN
2	B	61	GLU
2	B	81	LEU
2	B	83	LEU
2	B	86	GLU
2	B	120	LEU
2	B	135	ILE
2	B	170	VAL
2	B	180	LEU
2	B	237	LEU
2	B	238	LEU
2	B	247	ASN
2	B	279	LEU
3	C	14	GLN
3	C	24	VAL
3	C	30	ILE
3	C	38	GLU
3	C	45	THR
3	C	86	LEU
3	C	135	ASN
3	C	170	LEU
3	C	175	GLU
3	C	253	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN

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Mol	Chain	Res	Type
1	A	59	GLN
1	A	62	GLN
1	A	68	GLN
1	A	102	ASN
1	A	127	GLN
1	A	143	HIS
1	A	205	GLN
1	A	208	GLN
1	A	210	GLN
2	B	0	HIS
2	B	15	HIS
2	B	62	ASN
2	B	68	GLN
2	B	76	ASN
2	B	226	HIS
2	B	231	HIS
2	B	247	ASN
2	B	258	ASN
2	B	262	HIS
2	B	267	HIS
3	C	35	HIS
3	C	39	HIS
3	C	76	GLN
3	C	135	ASN
3	C	151	ASN
3	C	200	HIS
3	C	221	ASN
3	C	247	ASN

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	$\# Z > 2$
4	GLC	D	1	4	11,11,12	1.24	1 (9%)	15,15,17	0.76	0
4	FRU	D	2	4	11,12,12	1.79	1 (9%)	10,18,18	0.76	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	GLC	D	1	4	-	0/2/19/22	0/1/1/1
4	FRU	D	2	4	-	1/5/24/24	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	2	FRU	O2-C2	5.23	1.49	1.40
4	D	1	GLC	C1-C2	2.02	1.57	1.52

There are no bond angle outliers.

There are no chirality outliers.

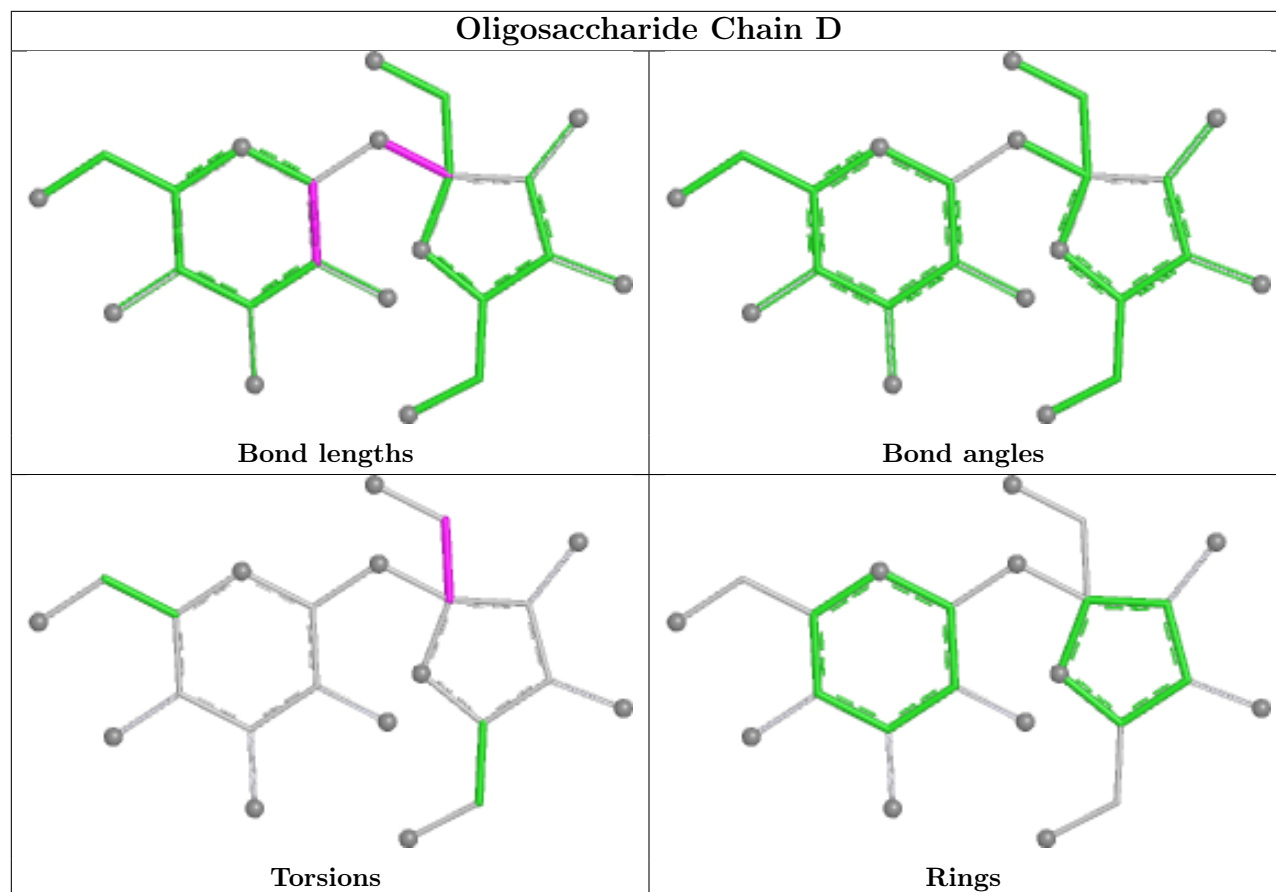
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	2	FRU	O1-C1-C2-C3

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](#)

There are no ligands in this entry.

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	259/266 (97%)	-0.34	3 (1%) 76 73	24, 40, 68, 90	0
2	B	260/281 (92%)	-0.18	5 (1%) 66 63	23, 42, 73, 85	0
3	C	256/263 (97%)	-0.17	4 (1%) 70 67	24, 44, 76, 87	0
All	All	775/810 (95%)	-0.23	12 (1%) 71 68	23, 42, 73, 90	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	103	GLY	4.9
2	B	45	ASP	4.3
3	C	146	ASP	4.2
2	B	47	LEU	3.5
2	B	269	ASP	3.1
2	B	72	VAL	2.9
3	C	216	ASN	2.9
3	C	196	ALA	2.7
1	A	68	GLN	2.7
3	C	147	PHE	2.4
2	B	226	HIS	2.2
1	A	69	ASP	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.

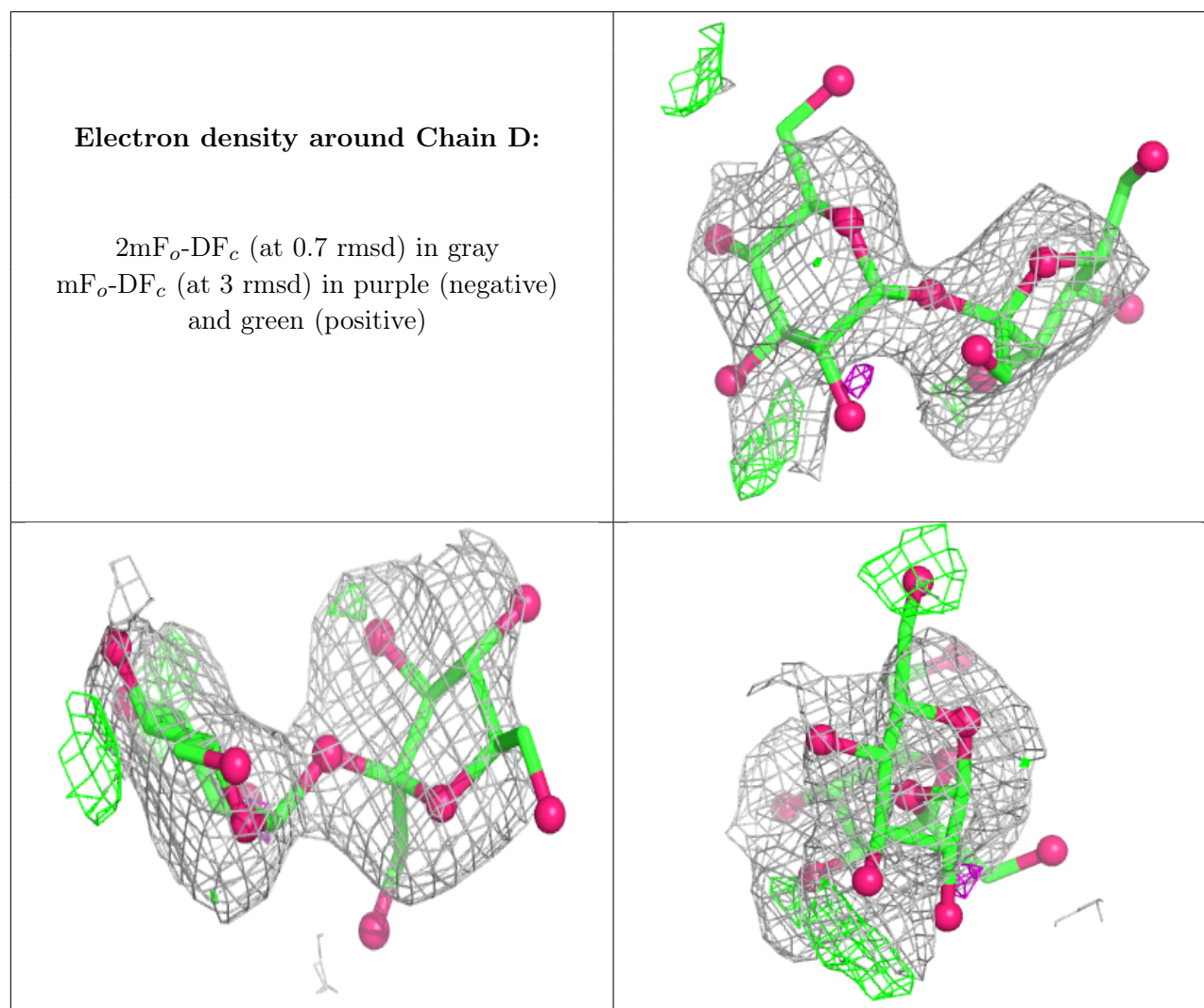
6.3 Carbohydrates [i](#)

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GLC	D	1	11/12	0.61	0.17	89,91,93,93	0
4	FRU	D	2	12/12	0.68	0.15	89,92,94,94	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.