



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:34 AM UTC

PDB ID : 8AG0 / pdb_00008ag0
Title : Crystal structure of mutant PRELID3a-TRIAP1 complex - R53E
Authors : Milara, X.; Perez-Dorado, J.I.; Matthews, S.J.
Deposited on : 2022-07-18
Resolution : 2.70 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

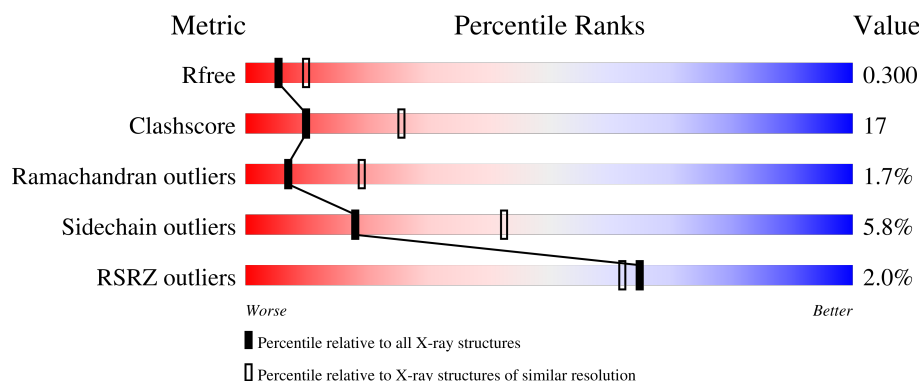
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	186	
2	B	446	
3	D	2	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4594 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 PRELI domain containing protein 3A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	165	Total	C	N	O	S	0	0	0
			1277	813	222	235	7			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP Q96N28
A	-12	ALA	-	expression tag	UNP Q96N28
A	-11	HIS	-	expression tag	UNP Q96N28
A	-10	HIS	-	expression tag	UNP Q96N28
A	-9	HIS	-	expression tag	UNP Q96N28
A	-8	HIS	-	expression tag	UNP Q96N28
A	-7	HIS	-	expression tag	UNP Q96N28
A	-6	HIS	-	expression tag	UNP Q96N28
A	-5	VAL	-	expression tag	UNP Q96N28
A	-4	ASP	-	expression tag	UNP Q96N28
A	-3	ASP	-	expression tag	UNP Q96N28
A	-2	ASP	-	expression tag	UNP Q96N28
A	-1	ASP	-	expression tag	UNP Q96N28
A	0	LYS	-	expression tag	UNP Q96N28
A	53	GLU	ARG	engineered mutation	UNP Q96N28

- Molecule 2 is a protein called Maltose/maltodextrin-binding periplasmic protein,TP53-regulated inhibitor of apoptosis 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	426	Total	C	N	O	S	0	0	0
			3245	2080	529	624	12			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P0AEX9
B	82	ALA	ASP	conflict	UNP P0AEX9
B	83	ALA	LYS	conflict	UNP P0AEX9
B	172	ALA	GLU	conflict	UNP P0AEX9
B	173	ALA	ASN	conflict	UNP P0AEX9
B	239	ALA	LYS	conflict	UNP P0AEX9
B	359	ALA	GLU	conflict	UNP P0AEX9
B	362	ALA	-	linker	UNP P0AEX9
B	363	ALA	-	linker	UNP P0AEX9
B	364	ALA	-	linker	UNP P0AEX9
B	365	GLN	-	linker	UNP P0AEX9
B	366	THR	-	linker	UNP P0AEX9
B	367	ASN	-	linker	UNP P0AEX9
B	368	ALA	-	linker	UNP P0AEX9
B	369	ALA	-	linker	UNP P0AEX9
B	370	ALA	-	linker	UNP P0AEX9

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose.



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

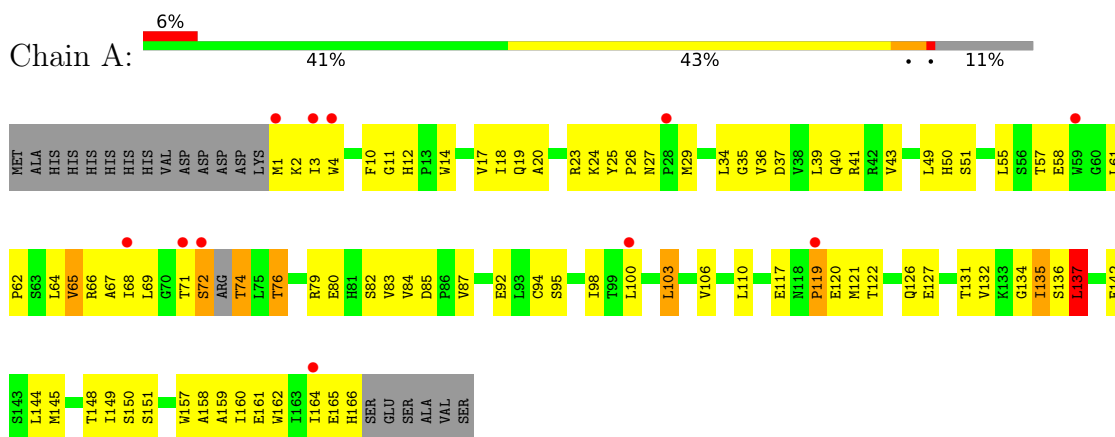
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	O	0	0
			12	12		
4	B	37	Total	O	0	0
			37	37		

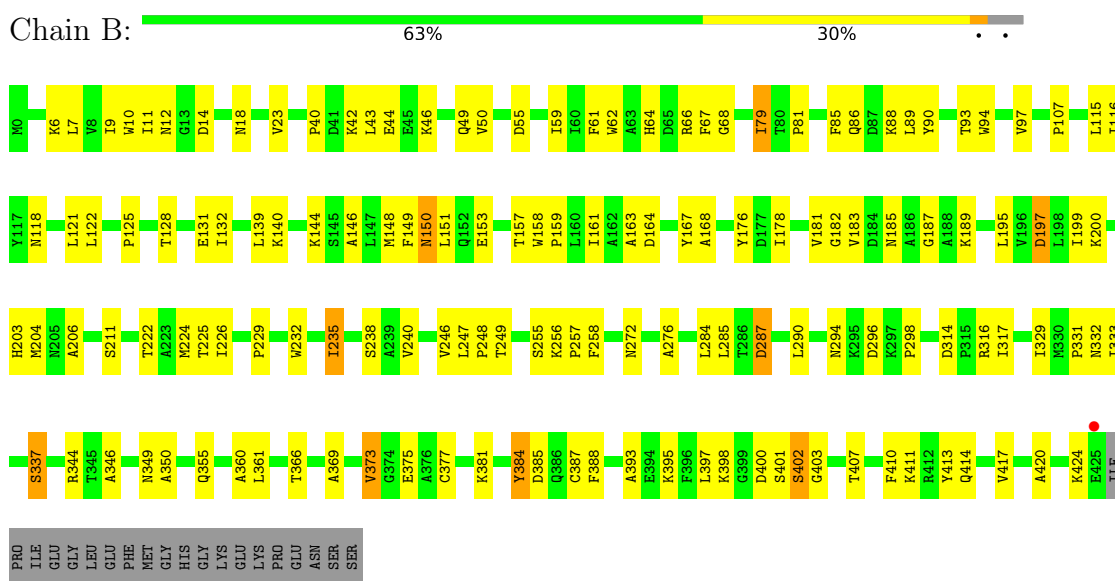
3 Residue-property plots

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: PRELI domain containing protein 3A



- Molecule 2: Maltose/maltodextrin-binding periplasmic protein, TP53-regulated inhibitor of apoptosis 1



- Molecule 3: alpha-D-glucopyranose-(1-4)-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	62.30Å 66.81Å 82.14Å 90.00° 96.46° 90.00°	Depositor
Resolution (Å)	45.41 – 2.70 45.41 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.9 (45.41-2.70) 99.1 (45.41-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.213 , 0.304 0.215 , 0.300	Depositor DCC
R_{free} test set	970 reflections (5.21%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.823	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4594	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.17% 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, BGC

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	0.36	0/1306	0.58	0/1783
2	B	0.39	0/3322	0.58	0/4517
All	All	0.38	0/4628	0.58	0/6300

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1277	0	1284	54	0
2	B	3245	0	3129	104	0
3	D	23	0	21	1	0
4	A	12	0	0	0	0
4	B	37	0	0	3	0
All	All	4594	0	4434	156	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 17.

All (156) 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:119:PRO:HB2	1:A:121:MET:HE3	1.56	0.88
1:A:62:PRO:HB2	1:A:76:THR:HG21	1.58	0.83
1:A:66:ARG:HE	1:A:71:THR:HA	1.45	0.82
2:B:122:LEU:HD21	2:B:125:PRO:HA	1.59	0.82
2:B:79:ILE:HD12	2:B:81:PRO:HD3	1.60	0.82
2:B:89:LEU:HD23	2:B:107:PRO:HG2	1.62	0.80
1:A:26:PRO:HG3	1:A:35:GLY:HA2	1.69	0.74
1:A:142:GLU:HA	1:A:145:MET:HE2	1.71	0.72
1:A:117:GLU:O	1:A:119:PRO:HD3	1.91	0.71
2:B:115:LEU:HD11	2:B:224:MET:HE2	1.76	0.68
2:B:181:VAL:HG12	2:B:183:VAL:HG13	1.79	0.65
1:A:17:VAL:HG11	1:A:122:THR:HG21	1.78	0.65
1:A:79:ARG:HB2	1:A:98:ILE:HD11	1.79	0.65
2:B:395:LYS:HA	2:B:400:ASP:HB3	1.79	0.64
2:B:68:GLY:HA3	2:B:332:ASN:O	1.98	0.63
2:B:116:ILE:HB	2:B:225:THR:HG23	1.83	0.61
2:B:40:PRO:HD2	2:B:43:LEU:HD12	1.82	0.60
1:A:27:ASN:HD21	1:A:159:ALA:HA	1.64	0.60
2:B:384:TYR:CE1	2:B:388:PHE:HB2	2.38	0.58
1:A:67:ALA:O	2:B:49:GLN:HG2	2.04	0.57
1:A:120:GLU:O	1:A:120:GLU:HG3	2.04	0.57
2:B:85:PHE:HA	2:B:88:LYS:HE3	1.86	0.57
1:A:110:LEU:HG	1:A:126:GLN:HG3	1.87	0.56
1:A:165:GLU:O	1:A:166:HIS:ND1	2.38	0.56
2:B:400:ASP:OD1	2:B:401:SER:N	2.38	0.56
2:B:50:VAL:HB	2:B:55:ASP:HB3	1.88	0.55
2:B:90:TYR:O	2:B:93:THR:OG1	2.25	0.54
2:B:185:ASN:O	2:B:189:LYS:HG3	2.08	0.54
1:A:144:LEU:O	1:A:148:THR:HG23	2.08	0.53
1:A:36:VAL:HG22	1:A:55:LEU:HD22	1.90	0.53
2:B:384:TYR:O	2:B:387:CYS:N	2.41	0.53
2:B:395:LYS:CA	2:B:400:ASP:HB3	2.38	0.53
1:A:24:LYS:HE3	1:A:25:TYR:CZ	2.43	0.53
1:A:39:LEU:C	1:A:40:GLN:HG2	2.34	0.53
2:B:43:LEU:HD22	2:B:62:TRP:HE3	1.74	0.52
2:B:384:TYR:HE1	2:B:388:PHE:HB2	1.74	0.52
2:B:287:ASP:N	2:B:287:ASP:OD1	2.43	0.52
2:B:148:MET:HB2	2:B:222:THR:HG21	1.93	0.51
1:A:57:THR:OG1	1:A:76:THR:O	2.25	0.51
2:B:195:LEU:O	2:B:199:ILE:HG12	2.10	0.51
2:B:349:ASN:HB3	2:B:355:GLN:HB2	1.93	0.51
2:B:153:GLU:OE2	2:B:344:ARG:NH1	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:118:ASN:ND2	2:B:121:LEU:HB2	2.26	0.51
2:B:159:PRO:HG3	2:B:257:PRO:HA	1.92	0.51
2:B:118:ASN:HD22	2:B:121:LEU:HB2	1.76	0.50
1:A:161:GLU:HB3	1:A:165:GLU:OE2	2.11	0.50
2:B:232:TRP:CH2	2:B:316:ARG:HB3	2.47	0.49
1:A:83:VAL:HG12	1:A:92:GLU:HB3	1.94	0.49
2:B:115:LEU:HD21	2:B:224:MET:HE1	1.95	0.49
2:B:167:TYR:CE1	2:B:182:GLY:HA3	2.48	0.49
2:B:395:LYS:C	2:B:400:ASP:HB3	2.37	0.49
1:A:14:TRP:CZ2	1:A:18:ILE:HG13	2.47	0.49
2:B:64:HIS:HE2	2:B:329:ILE:HG22	1.77	0.49
2:B:235:ILE:HA	2:B:238:SER:HB3	1.95	0.49
1:A:85:ASP:OD1	1:A:87:VAL:HG22	2.13	0.49
2:B:12:ASN:HD22	2:B:14:ASP:CG	2.21	0.49
2:B:121:LEU:HD22	2:B:139:LEU:HD22	1.94	0.49
2:B:149:PHE:CD2	2:B:226:ILE:HD12	2.48	0.49
2:B:10:TRP:O	2:B:61:PHE:HB2	2.13	0.48
2:B:199:ILE:HD13	2:B:204:MET:HB2	1.95	0.48
2:B:146:ALA:O	2:B:224:MET:HG3	2.14	0.48
1:A:14:TRP:HA	1:A:122:THR:OG1	2.13	0.48
2:B:43:LEU:HD22	2:B:62:TRP:CE3	2.49	0.48
2:B:115:LEU:HD22	2:B:248:PRO:HD3	1.96	0.48
2:B:149:PHE:HE1	2:B:157:THR:HG22	1.77	0.48
1:A:157:TRP:N	1:A:157:TRP:CD1	2.80	0.48
2:B:246:VAL:HG22	2:B:247:LEU:O	2.14	0.48
2:B:384:TYR:CE2	2:B:410:PHE:CD1	3.02	0.48
1:A:157:TRP:O	1:A:160:ILE:N	2.47	0.48
2:B:7:LEU:HD11	2:B:276:ALA:HA	1.96	0.47
1:A:76:THR:HG23	1:A:100:LEU:HD12	1.95	0.47
2:B:66:ARG:NH1	2:B:337:SER:HB3	2.29	0.47
1:A:49:LEU:O	1:A:83:VAL:HA	2.14	0.47
2:B:397:LEU:HA	2:B:397:LEU:HD23	1.71	0.47
1:A:162:TRP:O	1:A:166:HIS:ND1	2.47	0.47
2:B:18:ASN:HB2	2:B:296:ASP:OD2	2.14	0.47
2:B:176:TYR:CE2	2:B:331:PRO:HB3	2.49	0.47
2:B:255:SER:O	2:B:257:PRO:HD3	2.15	0.47
2:B:413:TYR:CZ	2:B:417:VAL:HG11	2.50	0.47
1:A:19:GLN:O	1:A:23:ARG:HG2	2.15	0.47
2:B:183:VAL:O	2:B:361:LEU:HD22	2.15	0.47
2:B:411:LYS:HA	2:B:414:GLN:HB2	1.97	0.46
1:A:72:SER:O	1:A:74:THR:OG1	2.24	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:178:ILE:HG22	2:B:333:ILE:HD12	1.97	0.46
2:B:314:ASP:HB3	2:B:317:ILE:HG12	1.97	0.46
1:A:119:PRO:HG2	1:A:121:MET:HB2	1.97	0.46
2:B:203:HIS:O	4:B:501:HOH:O	2.21	0.46
2:B:384:TYR:HE2	2:B:410:PHE:CD1	2.34	0.46
1:A:27:ASN:HD21	1:A:159:ALA:CA	2.27	0.46
2:B:151:LEU:HD12	2:B:206:ALA:HA	1.96	0.46
2:B:375:GLU:N	4:B:502:HOH:O	2.49	0.46
1:A:43:VAL:HG22	1:A:49:LEU:HD12	1.98	0.46
1:A:80:GLU:HG3	1:A:95:SER:HB3	1.97	0.46
2:B:46:LYS:N	2:B:46:LYS:HD2	2.31	0.46
2:B:140:LYS:HA	2:B:144:LYS:O	2.16	0.45
1:A:36:VAL:HG13	1:A:55:LEU:HD23	1.98	0.45
2:B:149:PHE:O	2:B:150:ASN:HB3	2.17	0.45
2:B:153:GLU:HG3	3:D:1:BGC:O6	2.16	0.45
2:B:178:ILE:O	2:B:369:ALA:O	2.35	0.45
1:A:65:VAL:O	1:A:68:ILE:HG13	2.16	0.45
2:B:384:TYR:C	2:B:384:TYR:CD1	2.95	0.45
2:B:6:LYS:O	2:B:272:ASN:ND2	2.49	0.45
2:B:128:THR:O	2:B:131:GLU:HG2	2.16	0.45
2:B:373:VAL:HG12	2:B:424:LYS:HE3	1.97	0.45
2:B:44:GLU:N	2:B:44:GLU:OE1	2.46	0.44
1:A:41:ARG:HA	1:A:50:HIS:O	2.18	0.44
1:A:12:HIS:CE1	1:A:164:ILE:HD11	2.53	0.44
2:B:232:TRP:HB2	2:B:298:PRO:HG2	2.00	0.44
2:B:413:TYR:O	2:B:417:VAL:HG13	2.17	0.44
1:A:136:SER:OG	1:A:137:LEU:N	2.50	0.43
2:B:381:LYS:HB2	2:B:413:TYR:CE2	2.53	0.43
1:A:34:LEU:HD11	1:A:58:GLU:HG2	2.01	0.43
2:B:314:ASP:OD1	2:B:316:ARG:HB2	2.19	0.43
1:A:134:GLY:O	1:A:136:SER:N	2.51	0.43
1:A:1:MET:HA	1:A:131:THR:HG22	2.01	0.43
2:B:285:LEU:HD23	2:B:290:LEU:HD21	2.01	0.42
2:B:62:TRP:HB3	2:B:67:PHE:HE1	1.84	0.42
2:B:116:ILE:HB	2:B:225:THR:CG2	2.48	0.42
2:B:346:ALA:HA	2:B:360:ALA:HB1	2.01	0.42
2:B:284:LEU:HD12	2:B:284:LEU:HA	1.85	0.42
2:B:294:ASN:OD1	2:B:298:PRO:HA	2.19	0.42
2:B:94:TRP:HA	2:B:97:VAL:HG22	2.02	0.42
2:B:9:ILE:HG12	2:B:59:ILE:HB	2.01	0.42
2:B:420:ALA:O	2:B:424:LYS:HE2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:118:ASN:OD1	2:B:240:VAL:HG13	2.19	0.42
2:B:398:LYS:HE3	2:B:398:LYS:HB3	1.88	0.42
2:B:149:PHE:O	2:B:150:ASN:CB	2.67	0.41
1:A:2:LYS:O	1:A:3:ILE:HD12	2.21	0.41
1:A:10:PHE:N	1:A:122:THR:O	2.47	0.41
1:A:11:GLY:HA2	1:A:121:MET:HG2	2.01	0.41
2:B:393:ALA:O	2:B:398:LYS:HD2	2.20	0.41
1:A:69:LEU:HD13	1:A:103:LEU:HD21	2.03	0.41
2:B:350:ALA:HA	2:B:355:GLN:O	2.21	0.41
2:B:158:TRP:HA	2:B:161:ILE:HD12	2.03	0.41
2:B:229:PRO:HA	2:B:232:TRP:CE2	2.56	0.41
2:B:64:HIS:NE2	2:B:329:ILE:HG22	2.35	0.41
2:B:122:LEU:HD23	2:B:122:LEU:O	2.20	0.41
2:B:164:ASP:O	2:B:187:GLY:HA3	2.21	0.41
2:B:168:ALA:HB3	4:B:520:HOH:O	2.20	0.41
2:B:181:VAL:HG21	2:B:369:ALA:HA	2.02	0.41
2:B:204:MET:HE2	2:B:204:MET:HB3	1.97	0.41
1:A:14:TRP:CE2	1:A:18:ILE:HG13	2.56	0.41
1:A:37:ASP:HA	2:B:413:TYR:OH	2.21	0.41
1:A:61:LEU:HD23	1:A:149:ILE:HG12	2.02	0.41
2:B:11:ILE:HG13	2:B:61:PHE:HB3	2.03	0.41
2:B:163:ALA:O	2:B:256:LYS:NZ	2.45	0.41
2:B:384:TYR:O	2:B:385:ASP:C	2.64	0.41
2:B:121:LEU:HD21	2:B:144:LYS:HG3	2.03	0.40
1:A:10:PHE:CE2	1:A:160:ILE:HD12	2.57	0.40
1:A:20:ALA:HB2	1:A:164:ILE:CG2	2.51	0.40
2:B:42:LYS:HD3	2:B:42:LYS:HA	1.82	0.40
1:A:4:TRP:O	1:A:127:GLU:HA	2.22	0.40
1:A:132:VAL:HG13	1:A:135:ILE:HA	2.02	0.40
2:B:168:ALA:O	2:B:181:VAL:HA	2.21	0.40
1:A:29:MET:HE2	1:A:158:ALA:HB1	2.04	0.40
2:B:197:ASP:HA	2:B:200:LYS:HB2	2.04	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	163/186 (88%)	136 (83%)	21 (13%)	6 (4%)	2	6
2	B	424/446 (95%)	382 (90%)	38 (9%)	4 (1%)	14	35
All	All	587/632 (93%)	518 (88%)	59 (10%)	10 (2%)	7	19

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	PRO
2	B	150	ASN
1	A	74	THR
1	A	135	ILE
2	B	384	TYR
2	B	402	SER
1	A	103	LEU
1	A	72	SER
1	A	137	LEU
2	B	403	GLY

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	142/167 (85%)	131 (92%)	11 (8%)	12	30
2	B	323/353 (92%)	307 (95%)	16 (5%)	22	48
All	All	465/520 (89%)	438 (94%)	27 (6%)	18	42

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	SER
1	A	64	LEU

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Mol	Chain	Res	Type
1	A	65	VAL
1	A	76	THR
1	A	82	SER
1	A	84	VAL
1	A	94	CYS
1	A	106	VAL
1	A	137	LEU
1	A	150	SER
1	A	151	SER
2	B	23	VAL
2	B	79	ILE
2	B	86	GLN
2	B	132	ILE
2	B	197	ASP
2	B	211	SER
2	B	235	ILE
2	B	249	THR
2	B	258	PHE
2	B	287	ASP
2	B	337	SER
2	B	366	THR
2	B	373	VAL
2	B	377	CYS
2	B	402	SER
2	B	407	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
2	B	39	HIS
2	B	355	GLN
2	B	389	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	$\# Z > 2$
3	BGC	D	1	3	12,12,12	0.63	0	17,17,17	2.06	5 (29%)
3	GLC	D	2	3	11,11,12	0.76	0	15,15,17	1.10	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	BGC	D	1	3	-	0/2/22/22	0/1/1/1
3	GLC	D	2	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	BGC	C1-C2-C3	-5.02	100.11	110.36
3	D	1	BGC	O5-C1-C2	-3.52	104.11	110.30
3	D	1	BGC	O3-C3-C2	2.81	117.01	110.38
3	D	1	BGC	O5-C5-C4	2.38	113.98	109.70
3	D	1	BGC	O2-C2-C3	2.22	115.62	110.38

There are no chirality outliers.

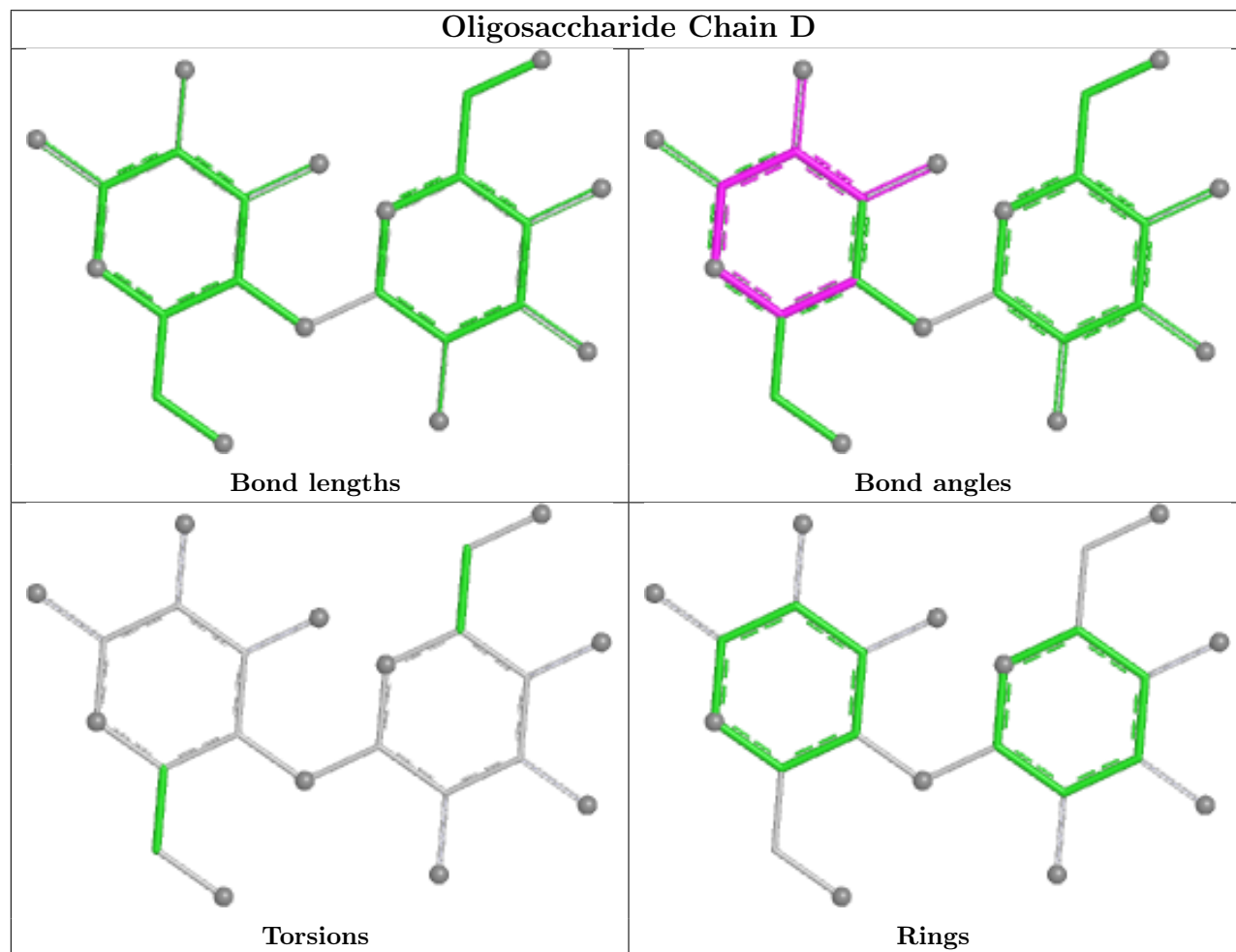
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1	BGC	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](#)

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 ⓘ

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	165/186 (88%)	0.57	11 (6%) 24 21	53, 72, 111, 130	0
2	B	426/446 (95%)	-0.01	1 (0%) 91 90	37, 54, 85, 119	0
All	All	591/632 (93%)	0.15	12 (2%) 65 62	37, 59, 97, 130	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	LEU	4.4
2	B	425	GLU	2.5
1	A	4	TRP	2.4
1	A	119	PRO	2.3
1	A	1	MET	2.3
1	A	59	TRP	2.3
1	A	3	ILE	2.3
1	A	68	ILE	2.2
1	A	164	ILE	2.2
1	A	72	SER	2.2
1	A	28	PRO	2.1
1	A	71	THR	2.0

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

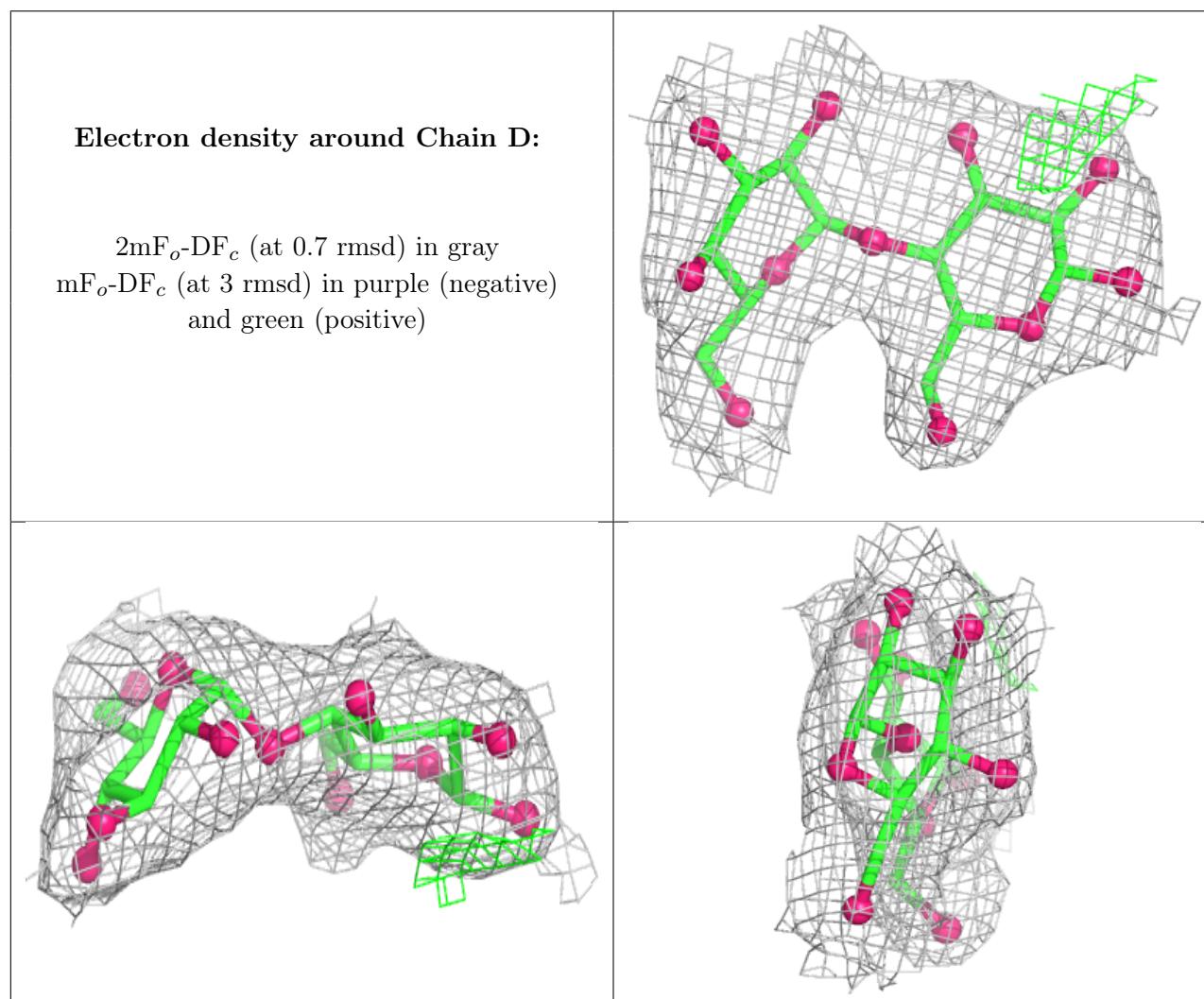
There are no non-standard protein/DNA/RNA residues in this entry.

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
3	BGC	D	1	12/12	-	-	32,37,40,42	0
3	GLC	D	2	11/12	-	-	35,38,39,43	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.