



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2016 – 05:23 AM EDT

PDB ID : 5EJZ
Title : Bacterial Cellulose Synthase Product-Bound State
Authors : Morgan, J.L.W.; Zimmer, J.
Deposited on : 2015-11-02
Resolution : 2.94 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0122
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107

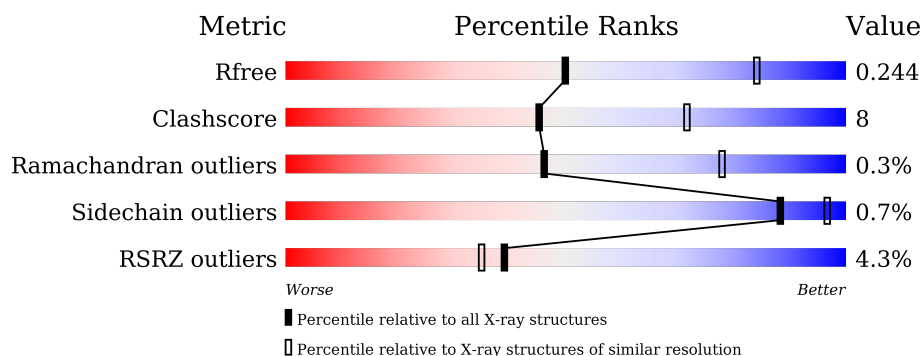
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.94 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}	91344	2057 (2.98-2.90)
Clashscore	102246	2308 (2.98-2.90)
Ramachandran outliers	100387	2245 (2.98-2.90)
Sidechain outliers	100360	2247 (2.98-2.90)
RSRZ outliers	91569	2065 (2.98-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	803	<div> <div>5%</div> <div> <div></div> <div>76%</div> <div>14%</div> <div>9%</div> </div> </div>
2	B	724	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>14%</div> <div>10%</div> </div> </div>
3	D	9	<div> <div></div> <div>100%</div> </div>

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:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
11	UND	B	802	-	-	-	X
4	BGC	A	910	-	-	-	X
4	BGC	A	917	-	-	-	X
5	SHG	A	918	-	-	-	X
8	MG	A	922	-	-	-	X
9	PLC	A	923	-	-	-	X
9	PLC	B	804	-	-	-	X

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 11078 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 Putative cellulose synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	728	Total	C	N	O	S	0	1	0
			5739	3725	1000	982	32			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q3J125
A	1	GLY	-	expression tag	UNP Q3J125
A	789	HIS	-	expression tag	UNP Q3J125
A	790	HIS	-	expression tag	UNP Q3J125
A	791	HIS	-	expression tag	UNP Q3J125
A	792	HIS	-	expression tag	UNP Q3J125
A	793	HIS	-	expression tag	UNP Q3J125
A	794	HIS	-	expression tag	UNP Q3J125
A	795	LYS	-	expression tag	UNP Q3J125
A	796	LEU	-	expression tag	UNP Q3J125
A	797	HIS	-	expression tag	UNP Q3J125
A	798	HIS	-	expression tag	UNP Q3J125
A	799	HIS	-	expression tag	UNP Q3J125
A	800	HIS	-	expression tag	UNP Q3J125
A	801	HIS	-	expression tag	UNP Q3J125
A	802	HIS	-	expression tag	UNP Q3J125

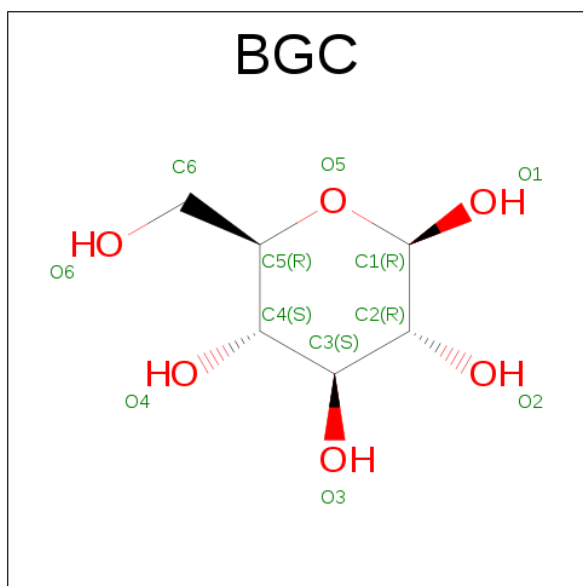
- Molecule 2 is a protein called Putative cellulose synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	655	Total	C	N	O	S	0	0	0
			4887	3100	864	907	16			

- Molecule 3 is a protein called poly(unk).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	9	Total	C	N	O	0	0	0
			45	27	9	9			

- Molecule 4 is BETA-D-GLUCOSE (three-letter code: BGC) (formula: $C_6H_{12}O_6$).



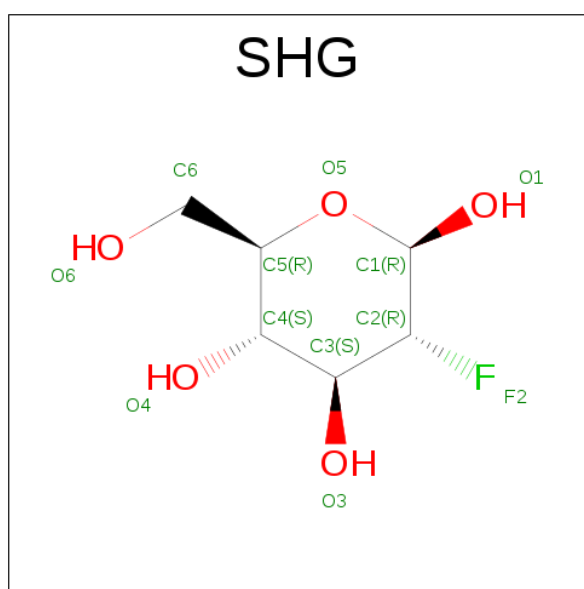
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			12	6	6		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		

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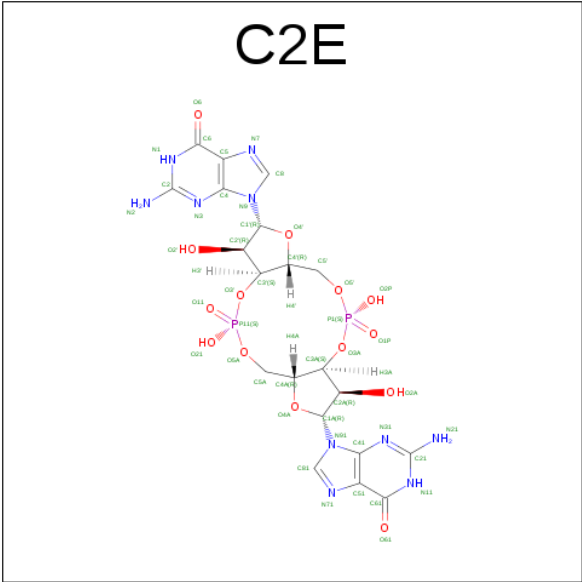
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is 2-deoxy-2-fluoro-beta-D-glucopyranose (three-letter code: SHG) (formula: $C_6H_{11}FO_5$).



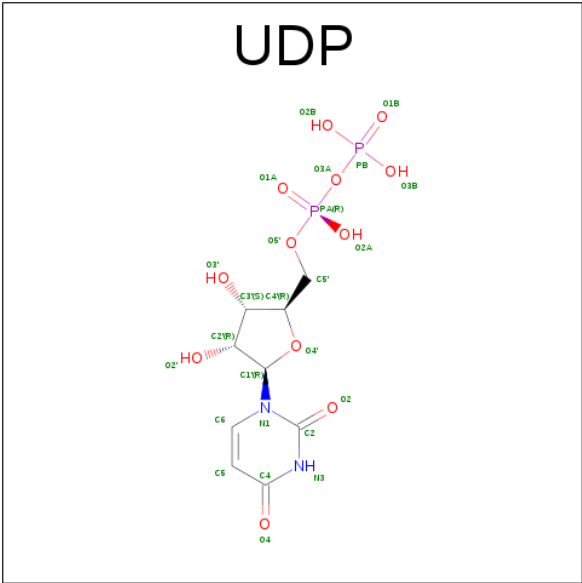
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	F	O	0	0
			11	6	1	4		

- Molecule 6 is 9,9'-[(2R,3R,3aS,5S,7aR,9R,10R,10aS,12S,14aR)-3,5,10,12-tetrahydroxy-5,12-dioxidoctahydro-2H,7H-difuro[3,2-d:3',2'-j][1,3,7,9,2,8]tetraoxadiphosphacyclododecine-2,9-diyl]bis(2-amino-1,9-dihydro-6H-purin-6-one) (three-letter code: C2E) (formula: $C_{20}H_{24}N_{10}O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total 46	C 20	N 10	O 14	P 2	0	0
6	A	1	Total 46	C 20	N 10	O 14	P 2	0	0

- Molecule 7 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C₉H₁₄N₂O₁₂P₂).

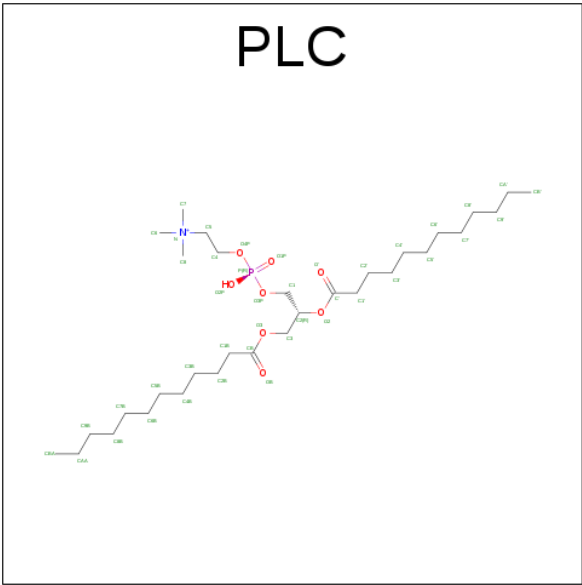


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

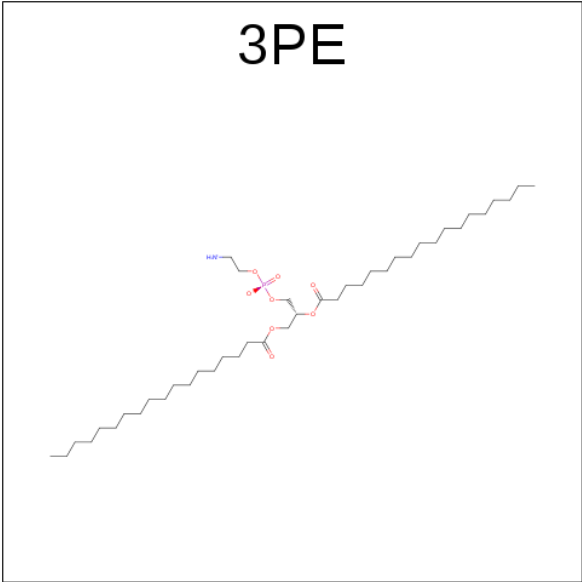
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	1	Total Mg 1 1	0	0
8	A	1	Total Mg 1 1	0	0

- Molecule 9 is DIUNDECYL PHOSPHATIDYL CHOLINE (three-letter code: PLC) (formula: C₃₂H₆₅NO₈P).



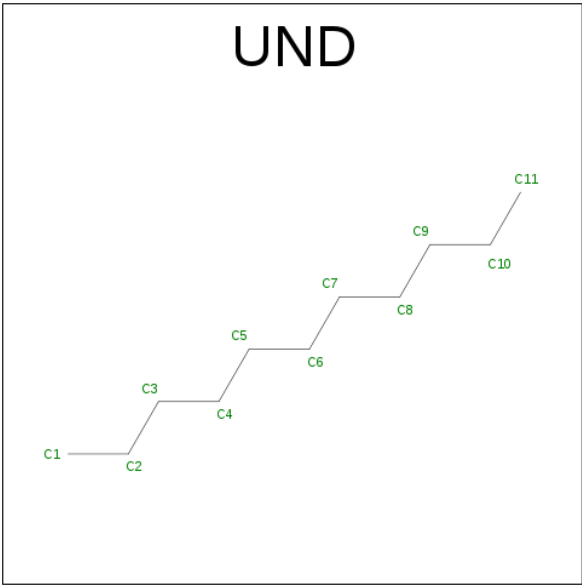
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total C N O P 38 28 1 8 1	0	0
9	B	1	Total C 9 9	0	0
9	B	1	Total C 11 11	0	0

- Molecule 10 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOETHANOLAMINE (three-letter code: 3PE) (formula: C₄₁H₈₂NO₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	A	1	Total	C	N	O	P	0	0
			20	10	1	8	1		

- Molecule 11 is UNDECANE (three-letter code: UND) (formula: C₁₁H₂₄).

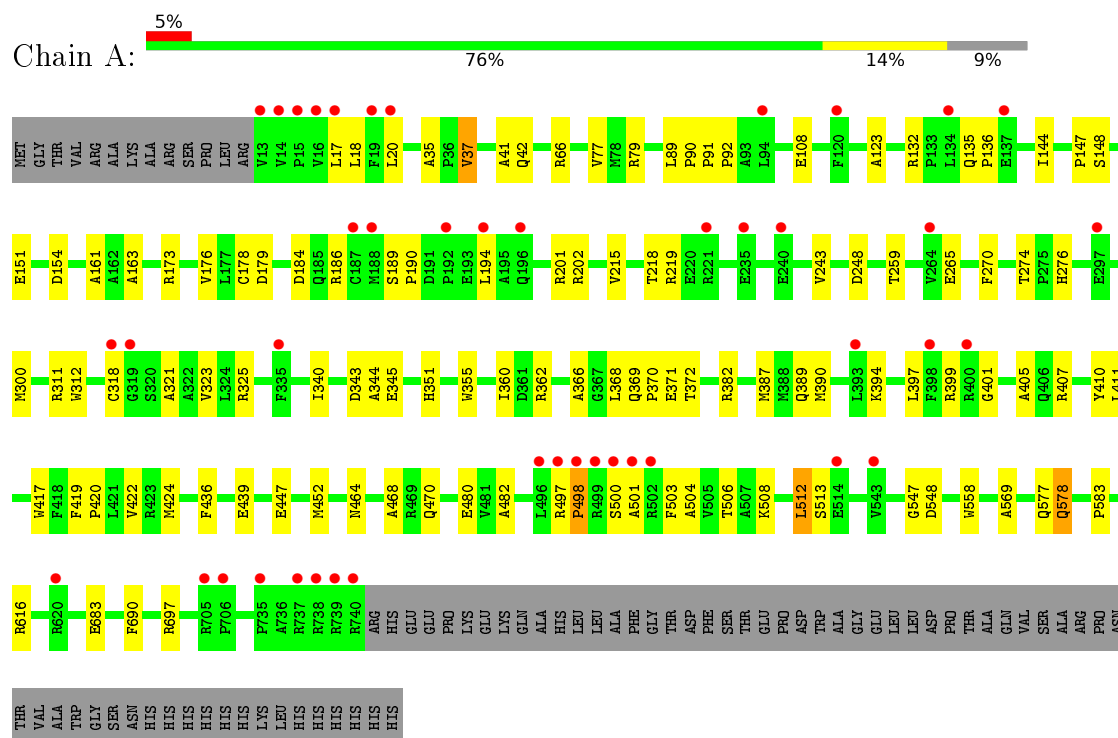


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	1	Total	C	0	0
			11	11		

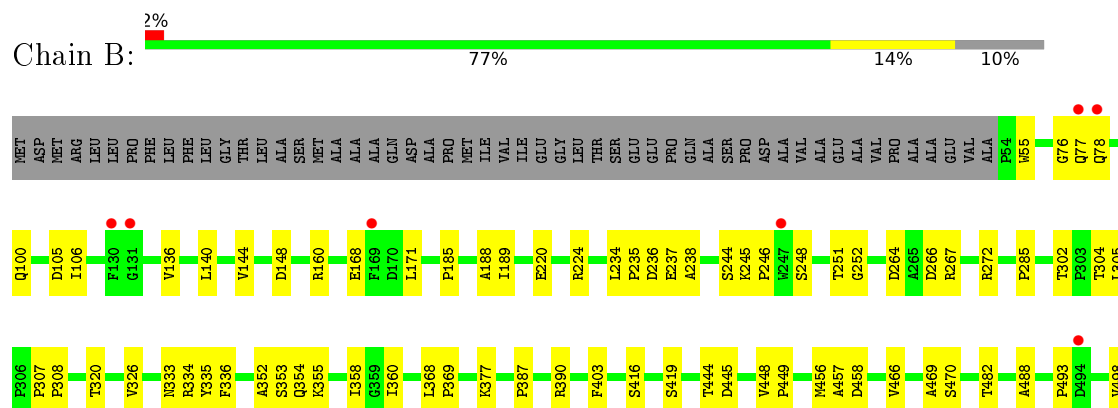
3 Residue-property plots

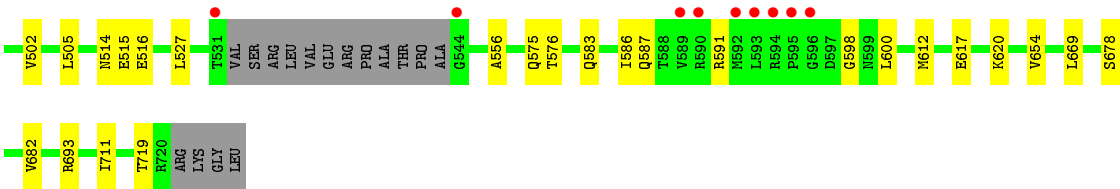
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Putative cellulose synthase



• Molecule 2: Putative cellulose synthase





● Molecule 3: poly(unk)

Chain D:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.27Å 216.84Å 221.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.35 – 2.94 38.70 – 2.94	Depositor EDS
% Data completeness (in resolution range)	98.5 (34.35-2.94) 98.5 (38.70-2.94)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.95Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.206 , 0.233 0.223 , 0.244	Depositor DCC
R_{free} test set	3431 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	73.0	Xtriage
Anisotropy	0.549	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.5	EDS
Estimated twinning fraction	0.017 for -h,l,k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	2 of 68819 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11078	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, BGC, C2E, UDP, PLC, SHG, UND, 3PE

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	0.23	0/5888	0.40	0/8007
2	B	0.24	0/5006	0.43	0/6865
All	All	0.23	0/10894	0.41	0/14872

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	5739	0	5855	89	0
2	B	4887	0	4966	85	0
3	D	45	0	13	0	0
4	A	188	0	154	14	0
5	A	11	0	9	0	0
6	A	92	0	44	1	0
7	A	25	0	11	4	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
9	A	38	0	53	2	0
9	B	20	0	38	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	A	20	0	14	0	0
11	B	11	0	24	2	0
All	All	11078	0	11181	174	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 8.

All (174) 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:77:GLN:NE2	2:B:168:GLU:CB	2.16	1.09
1:A:173:ARG:O	1:A:173:ARG:HD2	1.57	1.05
2:B:77:GLN:HG3	2:B:168:GLU:OE1	1.58	1.03
2:B:77:GLN:NE2	2:B:168:GLU:HB3	1.71	1.02
2:B:77:GLN:CD	2:B:168:GLU:HB3	1.82	1.00
1:A:340:ILE:CD1	1:A:501:ALA:HB3	1.92	0.99
1:A:340:ILE:HD13	1:A:501:ALA:HB1	1.46	0.97
1:A:340:ILE:CD1	1:A:501:ALA:CB	2.45	0.94
1:A:480:GLU:OE2	4:A:914:BGC:O6	1.86	0.92
2:B:352:ALA:HB3	2:B:354:GLN:HE22	1.36	0.90
2:B:77:GLN:NE2	2:B:168:GLU:HB2	1.85	0.89
1:A:340:ILE:HD13	1:A:501:ALA:CB	2.04	0.87
2:B:352:ALA:HB3	2:B:354:GLN:NE2	1.90	0.86
2:B:77:GLN:HE22	2:B:335:TYR:HB3	1.41	0.85
1:A:340:ILE:HD11	1:A:501:ALA:HB3	1.59	0.82
2:B:78:GLN:HG3	2:B:336:PHE:CD1	2.15	0.80
2:B:516:GLU:OE2	2:B:598:GLY:HA3	1.81	0.80
1:A:616:ARG:NH2	6:A:919:C2E:O61	2.15	0.80
2:B:516:GLU:CG	2:B:600:LEU:HB3	2.12	0.80
2:B:77:GLN:NE2	2:B:335:TYR:HB3	1.98	0.79
1:A:548:ASP:OD1	4:A:907:BGC:O6	2.01	0.79
2:B:77:GLN:NE2	2:B:168:GLU:OE1	2.18	0.76
2:B:77:GLN:CG	2:B:168:GLU:OE1	2.33	0.76
2:B:354:GLN:OE1	2:B:354:GLN:N	2.18	0.75
1:A:371:GLU:OE2	1:A:578:GLN:NE2	2.20	0.74
1:A:243:VAL:HG22	1:A:323:VAL:HG22	1.70	0.73
1:A:66:ARG:NH1	1:A:123:ALA:O	2.25	0.70
2:B:336:PHE:HB3	2:B:419:SER:OG	1.92	0.70
1:A:340:ILE:HG22	1:A:340:ILE:O	1.90	0.69
1:A:382:ARG:NH1	1:A:504:ALA:O	2.24	0.69
2:B:516:GLU:HG3	2:B:600:LEU:HB3	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:508:LYS:NZ	7:A:921:UDP:O2A	2.26	0.68
2:B:516:GLU:HG2	2:B:600:LEU:CB	2.25	0.67
2:B:77:GLN:O	2:B:78:GLN:HB3	1.95	0.67
1:A:399:ARG:O	1:A:407:ARG:NH1	2.27	0.67
2:B:352:ALA:HB3	2:B:354:GLN:OE1	1.95	0.67
2:B:77:GLN:HE22	2:B:168:GLU:HB2	1.56	0.67
2:B:251:THR:HG22	2:B:252:GLY:N	2.11	0.66
1:A:343:ASP:OD1	1:A:344:ALA:N	2.28	0.66
1:A:173:ARG:C	1:A:173:ARG:HD2	2.15	0.65
2:B:583:GLN:HA	2:B:586:ILE:HG12	1.79	0.65
2:B:482:THR:HG22	2:B:502:VAL:HB	1.78	0.65
1:A:108:GLU:OE2	4:A:912:BGC:O2	2.12	0.64
2:B:352:ALA:HB3	2:B:354:GLN:CD	2.16	0.64
1:A:186:ARG:HG2	1:A:194:LEU:HD21	1.79	0.64
4:A:905:BGC:O6	2:B:387:PRO:HG3	1.98	0.64
2:B:516:GLU:HG3	2:B:516:GLU:O	1.98	0.64
1:A:382:ARG:NH2	7:A:921:UDP:O2B	2.31	0.63
2:B:516:GLU:HG2	2:B:600:LEU:HB3	1.78	0.63
1:A:372:THR:HG22	1:A:512:LEU:HD21	1.79	0.63
1:A:369:GLN:HG3	1:A:370:PRO:HD2	1.81	0.62
2:B:245:LYS:HB3	2:B:246:PRO:HD2	1.82	0.62
1:A:340:ILE:HD11	1:A:501:ALA:CB	2.20	0.62
4:A:905:BGC:C6	2:B:390:ARG:HH22	2.13	0.62
4:A:905:BGC:H6C1	2:B:390:ARG:HH22	1.64	0.62
1:A:345:GLU:HA	1:A:390:MET:HE2	1.83	0.60
2:B:235:PRO:O	2:B:236:ASP:HB2	2.00	0.60
1:A:351:HIS:HD1	1:A:410:TYR:HH	1.48	0.60
1:A:382:ARG:HH22	1:A:506:THR:HG1	1.48	0.59
1:A:259:THR:HG21	1:A:323:VAL:HG21	1.84	0.59
2:B:78:GLN:HG3	2:B:336:PHE:HD1	1.63	0.59
2:B:100:GLN:HE21	2:B:136:VAL:HG23	1.68	0.59
2:B:148:ASP:HB3	2:B:305:LEU:HG	1.85	0.58
2:B:244:SER:OG	2:B:248:SER:HB3	2.03	0.58
1:A:340:ILE:CG2	1:A:340:ILE:O	2.52	0.58
1:A:343:ASP:OD2	4:A:917:BGC:H4	2.03	0.58
2:B:185:PRO:HG2	2:B:188:ALA:HB2	1.84	0.58
2:B:360:ILE:HA	2:B:444:THR:HG23	1.86	0.57
1:A:439:GLU:HB2	4:A:908:BGC:H6C2	1.87	0.56
1:A:394:LYS:O	1:A:399:ARG:NH1	2.38	0.56
2:B:244:SER:OG	2:B:248:SER:CB	2.53	0.56
1:A:512:LEU:HD12	1:A:578:GLN:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:106:ILE:HG22	2:B:171:LEU:HD22	1.88	0.56
1:A:368:LEU:HD22	1:A:583:PRO:HG2	1.88	0.55
1:A:35:ALA:O	1:A:79:ARG:NH2	2.40	0.55
1:A:179:ASP:OD1	1:A:201:ARG:NH1	2.40	0.55
1:A:184:ASP:HB3	1:A:202:ARG:HH12	1.72	0.55
1:A:132:ARG:HG3	1:A:265:GLU:CD	2.27	0.55
1:A:311:ARG:HH22	2:B:719:THR:HG23	1.71	0.54
2:B:77:GLN:HG3	2:B:333:ASN:HB3	1.90	0.54
1:A:482:ALA:HB2	1:A:569:ALA:HB1	1.90	0.53
1:A:135:GLN:HG3	1:A:136:PRO:HD2	1.90	0.53
1:A:340:ILE:HG23	1:A:503:PHE:HD1	1.74	0.53
1:A:218:THR:OG1	1:A:219:ARG:N	2.42	0.53
2:B:352:ALA:CB	2:B:354:GLN:HE22	2.16	0.51
1:A:419:PHE:HA	1:A:422:VAL:HG22	1.92	0.51
2:B:245:LYS:HB3	2:B:246:PRO:CD	2.39	0.51
1:A:497:ARG:HD2	1:A:500:SER:HB3	1.91	0.51
1:A:154:ASP:N	1:A:154:ASP:OD1	2.43	0.51
1:A:300:MET:HA	1:A:470:GLN:HB3	1.92	0.51
1:A:345:GLU:HA	1:A:390:MET:CE	2.40	0.51
2:B:587:GLN:O	2:B:591:ARG:HB2	2.10	0.51
1:A:362:ARG:HH12	1:A:697:ARG:HD2	1.75	0.50
1:A:176:VAL:HG22	1:A:215:VAL:HB	1.93	0.50
2:B:320:THR:HA	2:B:445:ASP:HA	1.94	0.50
2:B:264:ASP:OD1	2:B:267:ARG:NH2	2.45	0.50
2:B:302:THR:O	2:B:304:THR:HG23	2.12	0.49
2:B:678:SER:O	2:B:682:VAL:HG23	2.12	0.49
2:B:77:GLN:N	2:B:77:GLN:OE1	2.45	0.49
2:B:77:GLN:OE1	2:B:168:GLU:O	2.30	0.49
1:A:436:PHE:HA	9:A:923:PLC:H1'2	1.95	0.49
1:A:161:ALA:HB1	1:A:683:GLU:HG2	1.93	0.49
2:B:189:ILE:O	2:B:272:ARG:NH1	2.45	0.49
1:A:173:ARG:C	1:A:173:ARG:CD	2.80	0.48
2:B:224:ARG:HD3	2:B:469:ALA:HA	1.95	0.48
2:B:575:GLN:HE21	11:B:802:UND:H21	1.78	0.48
2:B:514:ASN:O	2:B:515:GLU:HG3	2.14	0.48
2:B:238:ALA:HB2	2:B:488:ALA:HB2	1.94	0.48
2:B:516:GLU:HG2	2:B:600:LEU:HB2	1.94	0.48
1:A:512:LEU:O	1:A:578:GLN:O	2.31	0.48
2:B:140:LEU:HB3	2:B:144:VAL:HB	1.96	0.47
11:B:802:UND:H42	9:B:803:PLC:H4'2	1.95	0.47
1:A:419:PHE:CD2	1:A:420:PRO:HD3	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:ARG:NH1	1:A:697:ARG:HD2	2.29	0.47
2:B:220:GLU:OE2	2:B:224:ARG:NH2	2.48	0.47
2:B:251:THR:CG2	2:B:252:GLY:N	2.78	0.47
2:B:353:SER:O	2:B:355:LYS:HD2	2.13	0.47
1:A:270:PHE:HB3	1:A:355:TRP:HB3	1.96	0.47
1:A:274:THR:OG1	1:A:321:ALA:O	2.23	0.46
1:A:37:VAL:HG13	1:A:41:ALA:HB3	1.98	0.46
1:A:389:GLN:NE2	1:A:498:PRO:O	2.49	0.46
1:A:387:MET:HG3	1:A:417:TRP:CD1	2.50	0.46
1:A:464:ASN:O	1:A:468:ALA:HB2	2.16	0.46
2:B:377:LYS:HB2	2:B:416:SER:HB2	1.98	0.46
1:A:397:LEU:HD21	1:A:411:LEU:HD13	1.96	0.46
1:A:18:LEU:HD22	2:B:711:ILE:HG23	1.97	0.46
2:B:358:ILE:HG13	2:B:403:PHE:HE2	1.80	0.46
1:A:147:PRO:HA	1:A:178:CYS:HB2	1.98	0.46
1:A:452:MET:HE2	4:A:911:BGC:O6	2.15	0.45
2:B:466:VAL:HG13	2:B:470:SER:OG	2.17	0.45
2:B:617:GLU:HG3	2:B:620:LYS:HB2	1.98	0.45
1:A:390:MET:HE3	1:A:390:MET:HB2	1.61	0.45
4:A:910:BGC:O3	4:A:911:BGC:O5	2.25	0.45
2:B:105:ASP:HA	2:B:160:ARG:HE	1.81	0.45
2:B:251:THR:HG22	2:B:252:GLY:H	1.82	0.45
2:B:456:MET:O	2:B:458:ASP:N	2.49	0.45
1:A:42:GLN:OE1	1:A:79:ARG:NH1	2.51	0.44
2:B:76:GLY:N	2:B:171:LEU:O	2.40	0.44
2:B:505:LEU:HD21	2:B:527:LEU:HD12	2.00	0.44
1:A:452:MET:HE2	1:A:558:TRP:HE1	1.83	0.43
1:A:452:MET:CE	4:A:911:BGC:O6	2.67	0.43
2:B:612:MET:HG2	2:B:654:VAL:HG22	2.01	0.43
1:A:274:THR:HG22	1:A:360:ILE:HG23	1.99	0.43
1:A:345:GLU:HB2	1:A:390:MET:CE	2.49	0.43
2:B:55:TRP:CE3	2:B:189:ILE:HG13	2.54	0.43
2:B:334:ARG:HD3	2:B:334:ARG:HA	1.84	0.43
1:A:276:HIS:HE1	1:A:318:CYS:HB3	1.84	0.42
1:A:513:SER:O	1:A:577:GLN:HG3	2.18	0.42
4:A:905:BGC:H6	2:B:387:PRO:HB3	1.85	0.42
2:B:307:PRO:HA	2:B:308:PRO:HD3	1.85	0.42
1:A:189:SER:HA	1:A:190:PRO:HD3	1.92	0.42
2:B:266:ASP:OD1	2:B:285:PRO:HD3	2.20	0.42
1:A:91:PRO:HA	1:A:92:PRO:HD3	1.90	0.41
1:A:312:TRP:HB3	1:A:405:ALA:HB1	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:77:GLN:CD	2:B:168:GLU:OE1	2.58	0.41
2:B:556:ALA:HB2	2:B:576:THR:HG23	2.01	0.41
2:B:234:LEU:HB3	2:B:237:GLU:HB2	2.01	0.41
1:A:144:ILE:HG21	1:A:163:ALA:HB1	2.03	0.41
1:A:248:ASP:HB2	1:A:368:LEU:HG	2.02	0.41
1:A:148:SER:HA	7:A:921:UDP:O2'	2.20	0.41
1:A:447:GLU:OE2	2:B:355:LYS:NZ	2.53	0.41
2:B:368:LEU:HA	2:B:369:PRO:HD3	1.84	0.41
2:B:448:VAL:HA	2:B:449:PRO:HD3	1.89	0.41
1:A:186:ARG:HG2	1:A:194:LEU:CD2	2.50	0.41
1:A:547:GLY:O	4:A:907:BGC:O6	2.38	0.41
2:B:466:VAL:HG22	2:B:498:VAL:CG1	2.50	0.41
1:A:366:ALA:HB3	1:A:690:PHE:HD1	1.85	0.41
1:A:151:GLU:OE2	7:A:921:UDP:O2'	2.32	0.41
1:A:420:PRO:O	1:A:424:MET:HG2	2.21	0.41
1:A:497:ARG:HD2	1:A:500:SER:CB	2.50	0.41
4:A:906:BGC:O3	9:A:923:PLC:H71	2.21	0.41
1:A:17:LEU:HD23	1:A:20:LEU:HD12	2.02	0.40
1:A:89:LEU:HA	1:A:90:PRO:HD2	1.91	0.40
2:B:78:GLN:CG	2:B:78:GLN:O	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	727/803 (90%)	701 (96%)	24 (3%)	2 (0%)	46	78
2	B	651/724 (90%)	624 (96%)	25 (4%)	2 (0%)	46	78
All	All	1378/1527 (90%)	1325 (96%)	49 (4%)	4 (0%)	46	78

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	498	PRO
2	B	457	ALA
1	A	401	GLY
2	B	493	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	599/661 (91%)	594 (99%)	5 (1%)	86	96
2	B	520/572 (91%)	517 (99%)	3 (1%)	90	97
All	All	1119/1233 (91%)	1111 (99%)	8 (1%)	88	96

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	VAL
1	A	77	VAL
1	A	325	ARG
1	A	512	LEU
1	A	578	GLN
2	B	326	VAL
2	B	669	LEU
2	B	693	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	222	ASN
2	B	77	GLN
2	B	100	GLN
2	B	514	ASN
2	B	575	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 28 ligands modelled in this entry, 2 are monoatomic - leaving 26 for Mogul analysis.

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	BGC	A	901	4	12,12,12	1.39	1 (8%)	17,17,17	1.46	3 (17%)
4	BGC	A	902	4	11,11,12	1.77	3 (27%)	15,15,17	1.67	4 (26%)
4	BGC	A	903	4	11,11,12	1.68	2 (18%)	15,15,17	2.08	3 (20%)
4	BGC	A	904	4	11,11,12	1.69	2 (18%)	15,15,17	2.21	6 (40%)
4	BGC	A	905	4	11,11,12	1.57	2 (18%)	15,15,17	1.88	4 (26%)
4	BGC	A	906	4	11,11,12	2.11	3 (27%)	15,15,17	1.07	1 (6%)
4	BGC	A	907	4	11,11,12	1.72	3 (27%)	15,15,17	1.05	0
4	BGC	A	908	4	11,11,12	1.67	2 (18%)	15,15,17	1.74	4 (26%)
4	BGC	A	909	4	11,11,12	1.70	2 (18%)	15,15,17	1.29	1 (6%)
4	BGC	A	910	4	11,11,12	1.65	2 (18%)	15,15,17	1.69	1 (6%)
4	BGC	A	911	4	11,11,12	1.63	2 (18%)	15,15,17	1.15	1 (6%)
4	BGC	A	912	4	11,11,12	1.70	2 (18%)	15,15,17	0.73	0
4	BGC	A	913	4	11,11,12	1.57	2 (18%)	15,15,17	1.20	1 (6%)
4	BGC	A	914	4	11,11,12	1.66	2 (18%)	15,15,17	1.67	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	BGC	A	915	4	11,11,12	1.61	2 (18%)	15,15,17	1.55	4 (26%)
4	BGC	A	916	4	11,11,12	1.53	2 (18%)	15,15,17	2.19	3 (20%)
4	BGC	A	917	5,4	11,11,12	1.76	2 (18%)	15,15,17	1.00	1 (6%)
5	SHG	A	918	4	11,11,12	1.68	2 (18%)	12,15,17	0.80	0
6	C2E	A	919	-	44,52,52	1.18	4 (9%)	50,82,82	1.83	10 (20%)
6	C2E	A	920	-	44,52,52	1.19	4 (9%)	50,82,82	1.79	11 (22%)
7	UDP	A	921	8	20,26,26	1.23	2 (10%)	24,40,40	1.40	2 (8%)
9	PLC	A	923	-	37,37,41	1.10	4 (10%)	41,45,49	1.11	2 (4%)
10	3PE	A	924	-	19,19,50	1.36	4 (21%)	21,24,55	1.46	2 (9%)
11	UND	B	802	-	10,10,10	0.26	0	9,9,9	0.54	0
9	PLC	B	803	-	8,8,41	0.31	0	7,7,49	0.76	0
9	PLC	B	804	-	10,10,41	0.30	0	9,9,49	0.80	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	BGC	A	901	4	-	0/2/22/22	0/1/1/1
4	BGC	A	902	4	-	0/2/19/22	0/1/1/1
4	BGC	A	903	4	-	0/2/19/22	0/1/1/1
4	BGC	A	904	4	-	0/2/19/22	0/1/1/1
4	BGC	A	905	4	-	0/2/19/22	0/1/1/1
4	BGC	A	906	4	-	0/2/19/22	0/1/1/1
4	BGC	A	907	4	-	0/2/19/22	0/1/1/1
4	BGC	A	908	4	-	0/2/19/22	0/1/1/1
4	BGC	A	909	4	-	0/2/19/22	0/1/1/1
4	BGC	A	910	4	-	0/2/19/22	0/1/1/1
4	BGC	A	911	4	-	0/2/19/22	0/1/1/1
4	BGC	A	912	4	-	0/2/19/22	0/1/1/1
4	BGC	A	913	4	-	0/2/19/22	0/1/1/1
4	BGC	A	914	4	-	0/2/19/22	0/1/1/1
4	BGC	A	915	4	-	0/2/19/22	0/1/1/1
4	BGC	A	916	4	-	0/2/19/22	0/1/1/1
4	BGC	A	917	5,4	-	0/2/19/22	0/1/1/1
5	SHG	A	918	4	-	0/2/19/22	0/1/1/1
6	C2E	A	919	-	-	0/22/62/62	0/6/7/7
6	C2E	A	920	-	-	0/22/62/62	0/6/7/7
7	UDP	A	921	8	-	0/12/32/32	0/2/2/2
9	PLC	A	923	-	-	0/41/41/45	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	3PE	A	924	-	-	0/22/22/54	0/0/0/0
11	UND	B	802	-	-	0/8/8/8	0/0/0/0
9	PLC	B	803	-	-	0/6/6/45	0/0/0/0
9	PLC	B	804	-	-	0/8/8/45	0/0/0/0

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	906	BGC	C2-C3	-3.16	1.48	1.52
4	A	917	BGC	C2-C3	-3.14	1.48	1.52
4	A	902	BGC	C2-C3	-2.97	1.48	1.52
4	A	913	BGC	C2-C3	-2.77	1.48	1.52
4	A	910	BGC	C2-C3	-2.73	1.48	1.52
4	A	915	BGC	C2-C3	-2.71	1.48	1.52
4	A	907	BGC	C2-C3	-2.69	1.48	1.52
4	A	912	BGC	C2-C3	-2.63	1.49	1.52
4	A	911	BGC	C2-C3	-2.63	1.49	1.52
4	A	909	BGC	C2-C3	-2.61	1.49	1.52
4	A	903	BGC	C2-C3	-2.58	1.49	1.52
4	A	908	BGC	C2-C3	-2.47	1.49	1.52
4	A	914	BGC	C2-C3	-2.44	1.49	1.52
10	A	924	3PE	O21-C2	-2.35	1.40	1.46
4	A	904	BGC	C2-C3	-2.34	1.49	1.52
9	A	923	PLC	O2-C2	-2.28	1.40	1.46
5	A	918	SHG	C2-C3	-2.25	1.48	1.51
10	A	924	3PE	O31-C3	-2.19	1.40	1.45
9	A	923	PLC	O3-C3	-2.16	1.40	1.45
4	A	916	BGC	C2-C3	-2.15	1.49	1.52
7	A	921	UDP	PB-O2B	-2.11	1.47	1.54
4	A	902	BGC	O5-C5	2.07	1.48	1.43
9	A	923	PLC	O2-C'	2.07	1.40	1.34
4	A	907	BGC	O5-C5	2.09	1.48	1.43
4	A	905	BGC	O2-C2	2.16	1.48	1.43
10	A	924	3PE	O21-C21	2.33	1.40	1.35
4	A	906	BGC	O5-C5	2.39	1.48	1.43
9	A	923	PLC	O3-CB	2.39	1.40	1.33
10	A	924	3PE	O31-C31	2.41	1.40	1.33
6	A	919	C2E	C51-C41	2.93	1.47	1.40
6	A	920	C2E	C51-C41	3.02	1.47	1.40
6	A	920	C2E	C5-C4	3.05	1.47	1.40
6	A	919	C2E	C5-C4	3.05	1.47	1.40
7	A	921	UDP	C4-N3	3.14	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	920	C2E	C61-C51	3.56	1.48	1.41
6	A	920	C2E	C6-C5	3.60	1.48	1.41
6	A	919	C2E	C61-C51	3.61	1.48	1.41
4	A	913	BGC	O5-C1	3.68	1.49	1.43
6	A	919	C2E	C6-C5	3.71	1.48	1.41
4	A	905	BGC	O5-C1	3.85	1.50	1.43
4	A	915	BGC	O5-C1	3.91	1.50	1.43
4	A	901	BGC	O5-C1	3.92	1.50	1.43
4	A	916	BGC	O5-C1	3.94	1.50	1.43
4	A	910	BGC	O5-C1	3.98	1.50	1.43
4	A	911	BGC	O5-C1	4.03	1.50	1.43
4	A	908	BGC	O5-C1	4.15	1.50	1.43
4	A	902	BGC	O5-C1	4.19	1.50	1.43
4	A	917	BGC	O5-C1	4.19	1.50	1.43
4	A	907	BGC	O5-C1	4.21	1.50	1.43
4	A	903	BGC	O5-C1	4.25	1.50	1.43
4	A	912	BGC	O5-C1	4.26	1.50	1.43
5	A	918	SHG	O5-C1	4.28	1.50	1.43
4	A	904	BGC	O5-C1	4.30	1.50	1.43
4	A	909	BGC	O5-C1	4.31	1.50	1.43
4	A	914	BGC	O5-C1	4.32	1.50	1.43
4	A	906	BGC	O5-C1	5.43	1.52	1.43

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	920	C2E	C5-C6-N1	-4.23	117.99	123.52
6	A	920	C2E	C51-C61-N11	-4.17	118.07	123.52
6	A	919	C2E	C5-C6-N1	-4.12	118.14	123.52
6	A	919	C2E	C51-C61-N11	-4.06	118.21	123.52
6	A	919	C2E	N31-C21-N11	-3.35	123.00	127.56
6	A	919	C2E	C61-C51-C41	-3.31	117.07	120.86
6	A	919	C2E	N3-C2-N1	-3.17	123.24	127.56
6	A	919	C2E	C6-C5-C4	-3.11	117.30	120.86
6	A	920	C2E	N3-C2-N1	-3.11	123.32	127.56
6	A	920	C2E	N31-C21-N11	-3.08	123.36	127.56
6	A	920	C2E	C6-C5-C4	-3.05	117.38	120.86
6	A	920	C2E	C61-C51-C41	-3.01	117.42	120.86
6	A	920	C2E	C2A-C1A-N91	-2.43	106.97	113.47
4	A	905	BGC	O4-C4-C5	-2.36	103.01	109.23
4	A	905	BGC	C1-O5-C5	-2.34	108.71	112.14
4	A	904	BGC	C6-C5-C4	-2.26	107.33	112.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	919	C2E	C1A-N91-C41	-2.24	124.31	126.81
4	A	904	BGC	O3-C3-C4	-2.07	105.69	110.36
4	A	915	BGC	C1-O5-C5	-2.05	109.12	112.14
6	A	920	C2E	O4A-C1A-N91	2.03	111.94	108.11
4	A	916	BGC	C3-C4-C5	2.09	113.96	110.23
4	A	911	BGC	C3-C4-C5	2.13	114.02	110.23
4	A	902	BGC	C1-C2-C3	2.15	112.15	109.55
6	A	920	C2E	C3'-C2'-C1'	2.15	104.73	100.06
7	A	921	UDP	O4'-C1'-N1	2.17	112.22	108.10
4	A	901	BGC	C4-C3-C2	2.18	114.80	110.79
4	A	904	BGC	C3-C4-C5	2.23	114.21	110.23
4	A	914	BGC	O5-C1-C2	2.24	114.47	110.89
4	A	904	BGC	O5-C5-C4	2.27	113.89	110.13
4	A	915	BGC	C2-C3-C4	2.32	115.09	111.05
4	A	908	BGC	O5-C5-C4	2.32	113.98	110.13
4	A	917	BGC	O5-C5-C4	2.35	114.03	110.13
4	A	902	BGC	C2-C3-C4	2.42	115.27	111.05
4	A	913	BGC	O5-C5-C4	2.44	114.17	110.13
6	A	919	C2E	C3'-C2'-C1'	2.44	105.37	100.06
4	A	915	BGC	C1-C2-C3	2.47	112.54	109.55
4	A	906	BGC	C1-C2-C3	2.51	112.60	109.55
4	A	915	BGC	C3-C4-C5	2.65	114.95	110.23
9	A	923	PLC	O3-CB-C1B	2.66	120.05	111.85
4	A	908	BGC	C2-C3-C4	2.75	115.84	111.05
4	A	908	BGC	C3-C4-C5	2.88	115.37	110.23
4	A	909	BGC	C1-C2-C3	2.91	113.08	109.55
4	A	903	BGC	C3-C4-C5	2.94	115.47	110.23
10	A	924	3PE	O31-C31-C32	3.03	120.19	111.16
4	A	902	BGC	O5-C5-C4	3.03	115.15	110.13
4	A	902	BGC	C3-C4-C5	3.04	115.64	110.23
4	A	901	BGC	O5-C5-C4	3.05	115.49	109.67
4	A	908	BGC	C1-C2-C3	3.11	113.33	109.55
4	A	916	BGC	C2-C3-C4	3.30	116.80	111.05
4	A	901	BGC	C3-C4-C5	3.30	116.11	110.23
4	A	903	BGC	C2-C3-C4	3.59	117.31	111.05
4	A	904	BGC	C2-C3-C4	3.78	117.65	111.05
9	A	923	PLC	O2-C'-C1'	4.00	119.96	111.53
4	A	905	BGC	C2-C3-C4	4.03	118.07	111.05
4	A	905	BGC	C1-C2-C3	4.37	114.84	109.55
10	A	924	3PE	O21-C21-C22	4.67	119.98	111.09
7	A	921	UDP	C4-N3-C2	4.73	119.20	114.21
4	A	914	BGC	C1-C2-C3	4.79	115.35	109.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	910	BGC	C1-C2-C3	5.01	115.62	109.55
4	A	903	BGC	C1-C2-C3	5.01	115.63	109.55
6	A	920	C2E	C61-N11-C21	5.05	121.79	115.88
6	A	919	C2E	C6-N1-C2	5.09	121.84	115.88
6	A	919	C2E	C61-N11-C21	5.16	121.93	115.88
6	A	920	C2E	C6-N1-C2	5.19	121.96	115.88
4	A	904	BGC	C1-C2-C3	5.73	116.49	109.55
4	A	916	BGC	C1-C2-C3	6.14	116.99	109.55

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

14 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	905	BGC	4	0
4	A	906	BGC	1	0
4	A	907	BGC	2	0
4	A	908	BGC	1	0
4	A	910	BGC	1	0
4	A	911	BGC	3	0
4	A	912	BGC	1	0
4	A	914	BGC	1	0
4	A	917	BGC	1	0
6	A	919	C2E	1	0
7	A	921	UDP	4	0
9	A	923	PLC	2	0
11	B	802	UND	2	0
9	B	803	PLC	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	728/803 (90%)	0.20	44 (6%) 25 20	55, 79, 133, 176	0
2	B	655/724 (90%)	-0.13	16 (2%) 62 59	49, 72, 116, 161	0
3	D	0/9	-	-	-	-
All	All	1383/1536 (90%)	0.04	60 (4%) 39 35	49, 75, 126, 176	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	16	VAL	5.3
1	A	13	VAL	5.3
1	A	15	PRO	4.3
2	B	77	GLN	4.3
1	A	318	CYS	4.0
1	A	188	MET	3.7
1	A	499	ARG	3.7
2	B	593	LEU	3.6
2	B	595	PRO	3.5
2	B	78	GLN	3.5
2	B	594	ARG	3.5
1	A	192	PRO	3.4
1	A	398	PHE	3.4
1	A	739	ARG	3.3
1	A	496	LEU	3.3
1	A	221	ARG	3.2
1	A	502	ARG	3.2
1	A	393	LEU	3.2
1	A	17	LEU	3.1
1	A	134	LEU	3.0
1	A	94	LEU	2.9
1	A	740	ARG	2.9
2	B	590	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	187	CYS	2.8
1	A	514	GLU	2.8
2	B	169	PHE	2.8
1	A	497	ARG	2.7
1	A	543	VAL	2.7
1	A	264	VAL	2.6
1	A	14	VAL	2.6
2	B	531	THR	2.6
2	B	130	PHE	2.6
1	A	620	ARG	2.6
1	A	735	PRO	2.6
2	B	544	GLY	2.6
1	A	498	PRO	2.6
2	B	589	VAL	2.6
1	A	335	PHE	2.5
1	A	501	ALA	2.5
1	A	19	PHE	2.5
1	A	737	ARG	2.5
1	A	194	LEU	2.5
1	A	738	ARG	2.4
1	A	137	GLU	2.4
1	A	706	PRO	2.4
2	B	494	ASP	2.3
1	A	120	PHE	2.3
2	B	592	MET	2.3
1	A	705	ARG	2.3
1	A	319	GLY	2.3
1	A	235	GLU	2.2
1	A	297	GLU	2.2
2	B	596	GLY	2.1
1	A	20	LEU	2.1
2	B	131	GLY	2.1
1	A	240	GLU	2.1
2	B	247	TRP	2.0
1	A	500	SER	2.0
1	A	400	ARG	2.0
1	A	196	GLN	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
8	MG	A	922	1/1	0.73	0.94	14.11	98,98,98,98	0
9	PLC	B	804	11/42	0.86	0.42	10.83	72,83,94,97	0
9	PLC	A	923	38/42	0.83	0.46	4.88	87,124,155,157	0
11	UND	B	802	11/11	0.87	0.37	4.86	63,68,80,82	0
4	BGC	A	910	11/12	0.93	0.26	2.64	65,96,113,131	0
5	SHG	A	918	11/12	0.77	0.41	1.85	100,108,125,138	0
4	BGC	A	913	11/12	0.94	0.29	1.84	66,71,88,92	0
4	BGC	A	917	11/12	0.91	0.41	1.81	77,82,98,99	0
4	BGC	A	912	11/12	0.96	0.26	1.81	64,80,89,94	0
10	3PE	A	924	20/51	0.91	0.26	1.77	87,105,115,120	0
4	BGC	A	916	11/12	0.92	0.38	1.35	69,75,90,100	0
4	BGC	A	904	11/12	0.95	0.18	1.17	56,67,86,93	0
4	BGC	A	905	11/12	0.92	0.21	0.97	49,52,86,113	0
4	BGC	A	915	11/12	0.93	0.32	0.96	62,76,87,89	0
4	BGC	A	914	11/12	0.94	0.28	0.79	57,74,80,87	0
4	BGC	A	911	11/12	0.97	0.20	0.77	77,86,100,103	0
4	BGC	A	908	11/12	0.90	0.25	0.59	67,85,116,135	0
4	BGC	A	906	11/12	0.92	0.22	0.54	50,58,80,103	0
4	BGC	A	909	11/12	0.94	0.23	0.47	84,91,101,106	0
7	UDP	A	921	25/25	0.94	0.18	-0.04	77,102,116,116	0
6	C2E	A	919	46/46	0.96	0.14	-0.88	49,68,83,95	0
8	MG	B	801	1/1	0.81	0.13	-1.68	61,61,61,61	0
4	BGC	A	907	11/12	0.95	0.14	-1.79	49,57,76,89	0
6	C2E	A	920	46/46	0.95	0.13	-2.33	49,74,90,101	0
4	BGC	A	903	11/12	0.91	0.14	-	81,91,107,109	0
9	PLC	B	803	9/42	0.91	0.26	-	49,77,87,92	0
4	BGC	A	901	12/12	0.81	0.31	-	120,132,137,143	0
4	BGC	A	902	11/12	0.88	0.18	-	100,107,112,112	0

6.5 Other polymers [i](#)

There are no such residues in this entry.