



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:04 PM GMT

PDB ID : 4QTN
Title : Crystal structure of the Vitamin B3 transporter PnuC
Authors : Guskov, A.; Jaehme, M.; Slotboom, D.J.
Deposited on : 2014-07-08
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
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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

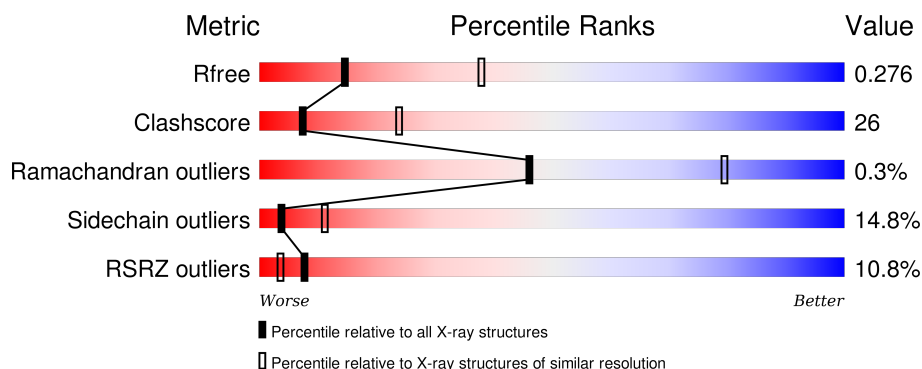
1 Overall quality at a glance ⓘ

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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (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 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	244	<div> <div>11%</div> <div>57%</div> <div>32%</div> <div>7%</div> <div>.</div> </div>
1	B	244	<div> <div>10%</div> <div>48%</div> <div>39%</div> <div>10%</div> <div>.</div> </div>
1	C	244	<div> <div>10%</div> <div>59%</div> <div>34%</div> <div>.</div> <div>.</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5808 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 Nicotinamide riboside transporter PnuC.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	Se	0	0	0
			1887	1269	298	310	2	8			
1	B	235	Total	C	N	O	S	Se	0	0	0
			1887	1269	298	310	2	8			
1	C	237	Total	C	N	O	S	Se	0	0	0
			1900	1276	301	313	2	8			

There are 24 discrepancies between the modelled and reference sequences:

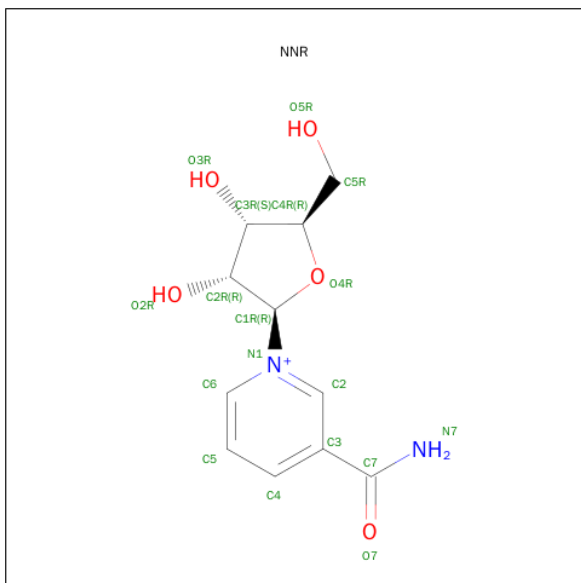
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	EXPRESSION TAG	UNP D2ZZC1
A	2	GLY	-	EXPRESSION TAG	UNP D2ZZC1
A	239	HIS	-	EXPRESSION TAG	UNP D2ZZC1
A	240	HIS	-	EXPRESSION TAG	UNP D2ZZC1
A	241	HIS	-	EXPRESSION TAG	UNP D2ZZC1
A	242	HIS	-	EXPRESSION TAG	UNP D2ZZC1
A	243	HIS	-	EXPRESSION TAG	UNP D2ZZC1
A	244	HIS	-	EXPRESSION TAG	UNP D2ZZC1
B	1	MSE	-	EXPRESSION TAG	UNP D2ZZC1
B	2	GLY	-	EXPRESSION TAG	UNP D2ZZC1
B	239	HIS	-	EXPRESSION TAG	UNP D2ZZC1
B	240	HIS	-	EXPRESSION TAG	UNP D2ZZC1
B	241	HIS	-	EXPRESSION TAG	UNP D2ZZC1
B	242	HIS	-	EXPRESSION TAG	UNP D2ZZC1
B	243	HIS	-	EXPRESSION TAG	UNP D2ZZC1
B	244	HIS	-	EXPRESSION TAG	UNP D2ZZC1
C	1	MSE	-	EXPRESSION TAG	UNP D2ZZC1
C	2	GLY	-	EXPRESSION TAG	UNP D2ZZC1
C	239	HIS	-	EXPRESSION TAG	UNP D2ZZC1
C	240	HIS	-	EXPRESSION TAG	UNP D2ZZC1
C	241	HIS	-	EXPRESSION TAG	UNP D2ZZC1
C	242	HIS	-	EXPRESSION TAG	UNP D2ZZC1
C	243	HIS	-	EXPRESSION TAG	UNP D2ZZC1

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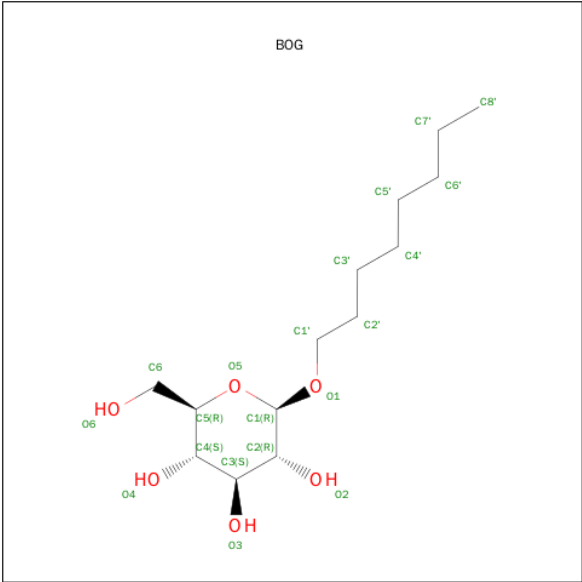
Chain	Residue	Modelled	Actual	Comment	Reference
C	244	HIS	-	EXPRESSION TAG	UNP D2ZZC1

- Molecule 2 is NICOTINAMIDE RIBOSIDE (three-letter code: NNR) (formula: $C_{11}H_{15}N_2O_5$).

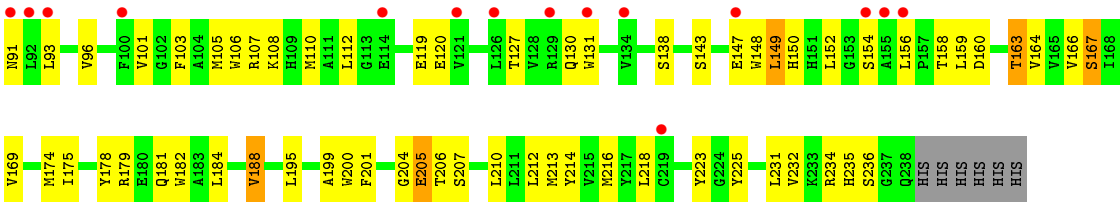


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			18	11	2	5		
2	B	1	Total	C	N	O	0	0
			18	11	2	5		
2	C	1	Total	C	N	O	0	0
			18	11	2	5		

- Molecule 3 is SUGAR (B-OCTYLGLUCOSIDE) (three-letter code: BOG) (formula: $C_{14}H_{28}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			20	14	6		
3	B	1	Total	C	O	0	0
			20	14	6		
3	B	1	Total	C	O	0	0
			20	14	6		
3	C	1	Total	C	O	0	0
			20	14	6		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.32Å 112.93Å 120.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.31 – 2.80 48.31 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.31-2.80) 99.8 (48.31-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.244 , 0.276 0.245 , 0.276	Depositor DCC
R_{free} test set	1597 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	102.7	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 79.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 31933 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5808	wwPDB-VP
Average B, all atoms (Å ²)	117.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NNR, BOG

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	0.61	0/1936	0.57	0/2633
1	B	0.63	0/1936	0.62	1/2633 (0.0%)
1	C	0.56	0/1949	0.58	1/2650 (0.0%)
All	All	0.60	0/5821	0.59	2/7916 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	110	MSE	CA-CB-CG	-5.94	103.20	113.30
1	C	231	LEU	CA-CB-CG	5.36	127.62	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	113	GLY	Peptide

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	1887	0	1931	109	0
1	B	1887	0	1931	120	0
1	C	1900	0	1942	107	0
2	A	18	0	14	1	0
2	B	18	0	14	0	0
2	C	18	0	14	1	0
3	A	20	0	28	1	0
3	B	40	0	56	2	0
3	C	20	0	28	0	0
All	All	5808	0	5958	302	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 26.

All (302) 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:B:25:MSE:HE1	1:C:184:LEU:CD1	1.67	1.22
1:A:154:SER:HB2	1:A:156:LEU:CD2	1.74	1.15
1:C:62:ILE:HD12	1:C:110:MSE:HE3	1.33	1.10
1:C:200:TRP:CD1	1:C:201:PHE:HE1	1.70	1.08
1:B:154:SER:HB2	1:B:156:LEU:HD22	1.31	1.08
1:A:154:SER:CB	1:A:156:LEU:HD22	1.86	1.04
1:C:200:TRP:CD1	1:C:201:PHE:CE1	2.45	1.04
1:A:25:MSE:HE1	1:B:183:ALA:HB1	1.39	1.03
1:C:200:TRP:HB2	1:C:206:THR:HG22	1.39	1.02
1:B:166:VAL:HG12	1:B:188:VAL:HG12	1.36	1.02
1:C:200:TRP:HD1	1:C:201:PHE:CE1	1.78	1.01
1:A:200:TRP:HD1	1:A:201:PHE:CD1	1.79	1.01
1:B:25:MSE:HE1	1:C:184:LEU:HD13	1.00	0.99
1:B:25:MSE:CE	1:C:184:LEU:HD13	1.93	0.99
1:A:159:LEU:HD23	1:A:195:LEU:CD2	1.96	0.95
1:C:234:ARG:O	1:C:235:HIS:HD2	1.50	0.94
1:A:154:SER:CB	1:A:156:LEU:CD2	2.46	0.93
1:A:184:LEU:HD13	1:C:25:MSE:HE2	1.48	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:TRP:HB2	1:C:206:THR:CG2	1.98	0.92
1:B:62:ILE:HD12	1:B:110:MSE:HE3	1.48	0.91
1:C:234:ARG:C	1:C:235:HIS:CD2	2.45	0.90
1:B:156:LEU:HD23	1:B:156:LEU:H	1.37	0.88
1:A:106:TRP:HB3	1:A:110:MSE:CE	2.04	0.88
1:A:159:LEU:HD23	1:A:195:LEU:HD23	1.56	0.87
1:C:166:VAL:HG12	1:C:188:VAL:HG12	1.56	0.87
1:A:154:SER:HB2	1:A:156:LEU:HD22	1.49	0.87
1:B:120:GLU:OE2	1:B:234:ARG:NH1	2.08	0.87
1:A:200:TRP:CD1	1:A:201:PHE:CD1	2.63	0.86
1:C:149:LEU:O	1:C:154:SER:OG	1.91	0.86
1:A:156:LEU:O	1:A:156:LEU:HG	1.75	0.84
1:A:154:SER:HB2	1:A:156:LEU:HD23	1.58	0.84
1:A:94:VAL:O	1:A:98:VAL:HG23	1.79	0.83
1:A:154:SER:HB3	1:A:156:LEU:HD22	1.58	0.82
1:A:25:MSE:CE	1:B:183:ALA:HB1	2.09	0.81
1:B:166:VAL:CG1	1:B:188:VAL:HG12	2.10	0.81
1:B:166:VAL:HG12	1:B:188:VAL:CG1	2.11	0.81
1:B:149:LEU:O	1:B:154:SER:OG	2.00	0.78
1:B:106:TRP:HB3	1:B:110:MSE:HE3	1.63	0.78
1:C:234:ARG:C	1:C:235:HIS:HD2	1.83	0.78
1:B:106:TRP:HB3	1:B:110:MSE:CE	2.13	0.78
1:C:131:TRP:HH2	1:C:184:LEU:CD2	1.95	0.78
1:C:159:LEU:CD2	1:C:195:LEU:HD21	2.12	0.78
1:A:25:MSE:HE1	1:B:183:ALA:CB	2.15	0.76
1:A:159:LEU:HD23	1:A:195:LEU:HD21	1.67	0.75
1:A:36:PHE:C	1:A:37:ASN:HD22	1.89	0.75
1:A:205:GLU:OE1	1:A:205:GLU:HA	1.84	0.75
1:A:159:LEU:CD2	1:A:195:LEU:HD23	2.17	0.74
1:B:155:ALA:O	1:B:157:PRO:HD3	1.87	0.73
1:C:205:GLU:OE1	1:C:206:THR:N	2.22	0.72
1:A:68:GLY:O	1:A:72:VAL:HG22	1.91	0.70
1:B:101:VAL:O	1:B:105:MSE:HE2	1.91	0.70
1:A:183:ALA:CB	1:C:25:MSE:HE1	2.21	0.70
1:C:10:GLU:HG2	1:C:63:SER:HB3	1.74	0.70
1:A:200:TRP:CD1	1:A:201:PHE:CE1	2.80	0.69
1:B:143:SER:O	1:B:147:GLU:HG3	1.93	0.69
1:B:4:LEU:HD22	1:B:4:LEU:O	1.93	0.68
1:A:149:LEU:HG	1:A:154:SER:OG	1.93	0.68
1:B:145:TYR:O	1:B:149:LEU:HD22	1.93	0.68
1:A:162:VAL:HG21	1:C:33:VAL:HG23	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:LEU:CD2	1:B:156:LEU:H	2.07	0.68
1:B:75:TYR:CD1	1:B:210:LEU:HD13	2.29	0.68
1:A:183:ALA:HB3	1:C:25:MSE:HE1	1.77	0.67
1:A:193:ILE:HG23	1:A:211:LEU:HD11	1.77	0.67
1:C:62:ILE:CD1	1:C:110:MSE:HE3	2.20	0.67
1:A:159:LEU:O	1:A:163:THR:HG22	1.95	0.66
1:C:166:VAL:HG13	1:C:184:LEU:HD12	1.77	0.66
1:C:234:ARG:O	1:C:235:HIS:CD2	2.40	0.66
1:B:10:GLU:HG2	1:B:63:SER:HB3	1.77	0.65
1:A:178:TYR:O	1:A:181:GLN:HG2	1.97	0.65
1:B:25:MSE:HE1	1:C:184:LEU:HD11	1.73	0.64
1:B:106:TRP:CZ2	1:B:174:MSE:HE2	2.32	0.64
1:A:93:LEU:O	1:A:97:PRO:HG2	1.99	0.63
1:C:57:VAL:HG11	2:C:301:NNR:H5	1.80	0.63
1:B:105:MSE:HE1	1:B:175:ILE:HD12	1.80	0.63
1:B:7:TRP:HH2	1:B:66:LEU:CD1	2.11	0.62
1:B:148:TRP:HE3	1:B:149:LEU:HB3	1.65	0.62
1:C:10:GLU:HG2	1:C:63:SER:CB	2.30	0.62
1:B:37:ASN:HD21	1:C:158:THR:CG2	2.11	0.62
1:A:201:PHE:HZ	1:B:201:PHE:CZ	2.17	0.62
1:B:36:PHE:C	1:B:37:ASN:HD22	2.03	0.62
1:B:212:LEU:HD22	1:B:216:MSE:HE3	1.81	0.62
1:B:3:SER:O	1:B:6:TRP:HB3	2.00	0.62
1:B:6:TRP:O	1:B:8:LYS:N	2.32	0.62
1:B:7:TRP:CH2	1:B:66:LEU:HD11	2.36	0.61
1:B:90:LEU:O	1:B:94:VAL:HB	2.00	0.61
1:C:159:LEU:CD2	1:C:195:LEU:CD2	2.79	0.61
1:A:77:TYR:O	1:A:80:TYR:HB3	2.01	0.61
1:B:52:LEU:HD13	1:B:67:PHE:CE1	2.36	0.61
1:C:159:LEU:C	1:C:159:LEU:HD23	2.22	0.60
1:B:85:TYR:HA	1:B:88:MSE:HG2	1.84	0.60
1:B:7:TRP:CH2	1:B:66:LEU:CD1	2.85	0.60
1:C:75:TYR:CD1	1:C:210:LEU:HD13	2.36	0.60
1:B:205:GLU:HA	1:B:205:GLU:OE1	2.02	0.59
1:C:68:GLY:O	1:C:72:VAL:HG22	2.03	0.59
1:C:105:MSE:HE1	1:C:175:ILE:HD12	1.85	0.59
1:B:156:LEU:O	1:B:156:LEU:HG	2.02	0.58
1:A:36:PHE:HE1	1:B:142:THR:CG2	2.16	0.58
1:A:166:VAL:HG12	1:A:188:VAL:HG12	1.84	0.58
1:C:131:TRP:HH2	1:C:184:LEU:HD21	1.68	0.58
1:A:36:PHE:HB3	1:B:158:THR:CG2	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:GLU:O	1:B:150:HIS:HB2	2.04	0.58
1:C:106:TRP:HB3	1:C:110:MSE:CE	2.34	0.58
1:B:37:ASN:HD22	1:B:37:ASN:N	2.02	0.57
1:C:105:MSE:HE1	1:C:175:ILE:HB	1.86	0.57
1:C:3:SER:O	1:C:6:TRP:HB3	2.04	0.57
1:B:159:LEU:O	1:B:163:THR:CG2	2.52	0.57
1:C:101:VAL:HG12	1:C:105:MSE:HE2	1.85	0.57
1:B:4:LEU:HD13	1:B:4:LEU:C	2.26	0.57
1:B:74:LEU:O	1:B:78:VAL:HG23	2.04	0.56
1:A:51:ILE:HD11	1:A:216:MSE:SE	2.55	0.56
1:C:232:VAL:O	1:C:236:SER:OG	2.20	0.56
1:A:159:LEU:CD2	1:A:195:LEU:CD2	2.78	0.56
1:C:31:ALA:O	1:C:35:VAL:HG23	2.05	0.56
1:C:159:LEU:O	1:C:163:THR:HG22	2.06	0.55
1:B:15:TRP:CZ3	1:B:59:LYS:HG3	2.42	0.55
1:C:159:LEU:HD23	1:C:195:LEU:HD21	1.87	0.55
1:A:158:THR:HG23	1:C:36:PHE:HB3	1.88	0.55
1:B:154:SER:CB	1:B:156:LEU:HD22	2.22	0.55
1:B:76:ALA:HB2	1:B:91:ASN:HB2	1.89	0.55
1:B:28:GLY:O	1:B:32:VAL:HG23	2.08	0.54
1:B:210:LEU:HD22	1:B:213:MSE:HE2	1.89	0.54
1:B:37:ASN:HD21	1:C:158:THR:HG21	1.70	0.54
1:A:184:LEU:N	1:C:25:MSE:HE1	2.21	0.54
1:C:75:TYR:HD1	1:C:210:LEU:HD13	1.70	0.54
1:B:172:VAL:O	1:B:176:LEU:HD12	2.08	0.54
1:A:184:LEU:HD13	1:C:25:MSE:CE	2.31	0.54
1:C:138:SER:OG	1:C:169:VAL:HG21	2.08	0.54
1:A:199:ALA:O	1:A:205:GLU:O	2.25	0.54
1:C:143:SER:O	1:C:147:GLU:HG2	2.08	0.54
1:C:103:PHE:O	1:C:107:ARG:HB2	2.08	0.53
1:C:29:ILE:HB	1:C:216:MSE:HE1	1.90	0.53
1:B:5:ALA:O	1:B:8:LYS:HB3	2.08	0.53
1:B:182:TRP:CZ3	1:B:221:SER:HA	2.43	0.53
1:B:202:LYS:HE3	1:B:203:ASN:HB2	1.90	0.53
1:A:75:TYR:CD1	1:A:210:LEU:HD13	2.44	0.53
1:C:105:MSE:CE	1:C:175:ILE:HB	2.38	0.53
1:B:106:TRP:CE3	1:B:110:MSE:HE1	2.44	0.53
1:C:178:TYR:O	1:C:181:GLN:HG2	2.09	0.53
1:C:4:LEU:HD13	1:C:4:LEU:O	2.08	0.53
1:B:85:TYR:O	1:B:89:MSE:HE3	2.09	0.53
1:A:66:LEU:HD23	1:A:67:PHE:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:131:TRP:CH2	1:C:184:LEU:HD21	2.44	0.53
1:A:106:TRP:HB3	1:A:110:MSE:HE1	1.87	0.52
1:B:181:GLN:HG3	1:B:182:TRP:N	2.24	0.52
1:B:189:ASN:O	1:B:193:ILE:HG13	2.09	0.52
1:A:63:SER:O	1:A:66:LEU:HD23	2.08	0.52
1:B:11:LEU:HD22	3:B:303:BOG:H8'1	1.92	0.52
1:B:202:LYS:HE3	1:B:202:LYS:HA	1.92	0.52
1:A:62:ILE:O	1:A:62:ILE:HG13	2.10	0.52
1:C:106:TRP:CE3	1:C:110:MSE:HE1	2.45	0.51
1:C:199:ALA:O	1:C:205:GLU:O	2.28	0.51
1:B:148:TRP:CE3	1:B:149:LEU:CB	2.93	0.51
1:B:7:TRP:CZ2	1:B:66:LEU:HD11	2.45	0.51
1:B:6:TRP:O	1:B:9:ARG:N	2.43	0.51
1:A:25:MSE:CE	1:B:183:ALA:CB	2.80	0.51
1:A:106:TRP:HB3	1:A:110:MSE:HE2	1.92	0.51
1:C:25:MSE:O	1:C:25:MSE:HG2	2.11	0.51
1:C:181:GLN:HG3	1:C:182:TRP:N	2.26	0.51
1:A:85:TYR:HA	1:A:88:MSE:HG2	1.92	0.51
1:B:25:MSE:CE	1:C:184:LEU:CD1	2.63	0.51
1:C:127:THR:OG1	1:C:130:GLN:HG3	2.10	0.51
1:B:178:TYR:O	1:B:181:GLN:HG2	2.11	0.50
1:B:159:LEU:O	1:B:163:THR:HG22	2.11	0.50
1:B:101:VAL:O	1:B:104:ALA:HB3	2.12	0.50
1:B:35:VAL:HG12	1:B:35:VAL:O	2.11	0.50
1:B:148:TRP:CE3	1:B:149:LEU:HB3	2.46	0.50
1:A:47:ALA:HA	1:A:213:MSE:SE	2.61	0.50
1:C:166:VAL:CG1	1:C:188:VAL:HG12	2.34	0.50
1:C:200:TRP:HB2	1:C:206:THR:HG21	1.92	0.50
1:C:148:TRP:O	1:C:149:LEU:HB3	2.11	0.50
1:A:90:LEU:HD11	1:A:167:SER:HB3	1.94	0.50
1:B:195:LEU:O	1:B:198:VAL:HG22	2.12	0.50
1:C:34:PHE:C	1:C:36:PHE:H	2.14	0.50
1:C:39:ASP:N	1:C:39:ASP:OD1	2.38	0.49
1:A:74:LEU:O	1:A:78:VAL:HG23	2.11	0.49
1:A:170:ALA:HB2	1:A:184:LEU:HB3	1.93	0.49
1:C:89:MSE:O	1:C:93:LEU:HG	2.12	0.49
1:B:149:LEU:HD21	1:B:156:LEU:HD13	1.95	0.49
1:B:149:LEU:HG	1:B:149:LEU:O	2.12	0.49
1:C:51:ILE:HD11	1:C:216:MSE:SE	2.63	0.49
1:C:200:TRP:O	1:C:204:GLY:HA2	2.13	0.48
1:A:101:VAL:HG12	1:A:105:MSE:HE1	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:91:ASN:O	1:C:96:VAL:HG23	2.13	0.48
1:A:106:TRP:C	1:A:110:MSE:HE2	2.34	0.48
1:A:37:ASN:N	1:A:37:ASN:HD22	2.10	0.48
1:C:235:HIS:CD2	1:C:235:HIS:N	2.79	0.48
1:A:29:ILE:HB	1:A:216:MSE:HE1	1.95	0.48
1:C:85:TYR:HA	1:C:88:MSE:HG2	1.95	0.48
1:C:200:TRP:CD1	1:C:201:PHE:CD1	3.00	0.48
1:B:85:TYR:O	1:B:89:MSE:HG2	2.14	0.48
1:B:27:LEU:HD22	1:B:48:VAL:HG13	1.94	0.48
1:A:15:TRP:CZ3	1:A:59:LYS:HG3	2.49	0.48
1:A:36:PHE:HB3	1:B:158:THR:HG22	1.94	0.48
1:B:159:LEU:HD23	1:B:195:LEU:HD23	1.96	0.48
1:C:88:MSE:HG3	1:C:89:MSE:N	2.29	0.48
1:B:111:ALA:O	1:B:119:GLU:HB3	2.14	0.47
1:B:170:ALA:HB2	1:B:184:LEU:HB3	1.95	0.47
1:A:37:ASN:C	1:A:39:ASP:H	2.17	0.47
1:B:87:GLU:O	1:B:91:ASN:ND2	2.41	0.47
1:B:191:LEU:O	1:B:194:SER:HB2	2.14	0.47
1:A:106:TRP:CE3	1:A:110:MSE:HE1	2.49	0.47
1:A:37:ASN:C	1:A:39:ASP:N	2.67	0.47
1:A:201:PHE:CZ	1:B:201:PHE:CZ	3.02	0.47
1:A:163:THR:O	1:A:167:SER:HB2	2.13	0.47
1:A:75:TYR:HD1	1:A:210:LEU:HD13	1.79	0.47
1:B:165:VAL:O	1:B:169:VAL:HG23	2.14	0.47
1:A:148:TRP:CE3	1:A:149:LEU:HA	2.50	0.47
1:C:106:TRP:HE3	1:C:110:MSE:HE1	1.80	0.47
1:C:25:MSE:O	1:C:29:ILE:HG12	2.15	0.47
1:A:106:TRP:O	1:A:110:MSE:HE2	2.14	0.46
1:C:179:ARG:O	1:C:179:ARG:HG3	2.15	0.46
1:C:200:TRP:HA	1:C:205:GLU:C	2.36	0.46
1:A:193:ILE:HG23	1:A:211:LEU:CD1	2.43	0.46
1:A:142:THR:CG2	1:C:36:PHE:CE1	2.98	0.46
1:B:6:TRP:C	1:B:8:LYS:N	2.68	0.46
1:C:49:THR:O	1:C:71:SER:HB2	2.16	0.46
1:B:151:HIS:O	1:B:152:LEU:HB2	2.15	0.46
1:A:180:GLU:OE1	1:A:180:GLU:N	2.48	0.46
1:A:57:VAL:HG11	2:A:301:NNR:H5	1.97	0.46
1:B:11:LEU:CD2	3:B:303:BOG:H8'1	2.45	0.46
1:C:101:VAL:O	1:C:105:MSE:HE2	2.16	0.46
1:A:158:THR:CG2	1:C:36:PHE:HB3	2.45	0.46
1:C:19:GLU:OE1	1:C:223:TYR:OH	2.32	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:TYR:HD1	1:B:210:LEU:HD13	1.76	0.46
1:A:106:TRP:HE3	1:A:110:MSE:HE1	1.80	0.46
1:C:201:PHE:C	1:C:204:GLY:H	2.19	0.45
1:B:11:LEU:HD21	1:B:67:PHE:HZ	1.80	0.45
1:C:90:LEU:HD11	1:C:167:SER:HB3	1.98	0.45
1:A:106:TRP:CZ2	1:A:174:MSE:HE2	2.52	0.45
1:B:39:ASP:OD1	1:B:39:ASP:N	2.46	0.45
1:A:176:LEU:O	1:A:177:ARG:HB2	2.16	0.45
1:B:75:TYR:CE1	1:B:210:LEU:HD13	2.51	0.45
1:C:212:LEU:HD23	1:C:212:LEU:HA	1.83	0.45
1:A:200:TRP:O	1:A:204:GLY:HA2	2.17	0.45
1:A:192:THR:HG21	1:A:214:TYR:CZ	2.52	0.45
1:A:183:ALA:C	1:C:25:MSE:HE1	2.36	0.44
1:B:83:LYS:HA	1:B:85:TYR:CZ	2.52	0.44
1:A:200:TRP:O	1:A:205:GLU:N	2.51	0.44
1:A:202:LYS:HA	1:A:203:ASN:HA	1.59	0.44
1:A:203:ASN:O	1:A:205:GLU:N	2.50	0.44
1:C:57:VAL:HG12	1:C:182:TRP:CZ2	2.53	0.44
1:A:166:VAL:CG1	1:A:188:VAL:HG12	2.46	0.44
1:B:77:TYR:O	1:B:80:TYR:HB3	2.17	0.44
1:B:105:MSE:CE	1:B:175:ILE:HB	2.46	0.44
1:A:181:GLN:HG3	1:A:182:TRP:N	2.32	0.44
1:A:76:ALA:HB2	1:A:91:ASN:HB3	1.99	0.44
1:B:4:LEU:HD13	1:B:5:ALA:N	2.32	0.44
1:B:202:LYS:CA	1:B:202:LYS:HE3	2.47	0.44
1:A:165:VAL:O	1:A:169:VAL:HG23	2.17	0.44
1:A:203:ASN:O	1:A:204:GLY:C	2.56	0.44
1:B:201:PHE:HZ	1:C:201:PHE:HE2	1.66	0.44
1:B:105:MSE:HE1	1:B:175:ILE:HB	1.98	0.44
1:A:37:ASN:ND2	1:B:158:THR:HG21	2.32	0.44
1:C:82:PHE:CZ	1:C:207:SER:HB3	2.53	0.43
1:B:105:MSE:HE3	1:B:105:MSE:HB2	1.78	0.43
1:C:201:PHE:CD1	1:C:201:PHE:N	2.85	0.43
1:A:63:SER:O	1:A:66:LEU:CD2	2.66	0.43
1:B:201:PHE:CZ	1:C:201:PHE:HE2	2.36	0.43
1:A:200:TRP:O	1:A:204:GLY:CA	2.67	0.43
1:B:194:SER:O	1:B:198:VAL:HG13	2.18	0.43
1:B:6:TRP:O	1:B:7:TRP:C	2.56	0.43
1:A:24:LEU:CD1	1:B:132:LEU:HB2	2.49	0.43
1:C:3:SER:HB2	1:C:6:TRP:H	1.84	0.43
1:C:131:TRP:CH2	1:C:184:LEU:CD2	2.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:THR:HG23	1:C:36:PHE:HE1	1.84	0.42
1:A:115:THR:HB	1:A:117:GLU:HG3	2.00	0.42
1:B:159:LEU:O	1:B:163:THR:HG23	2.19	0.42
1:A:59:LYS:HD2	1:A:59:LYS:HA	1.88	0.42
1:B:121:VAL:HG21	1:B:177:ARG:HG2	2.01	0.42
1:C:206:THR:OG1	1:C:207:SER:N	2.52	0.42
1:B:23:LEU:HD11	1:B:52:LEU:HD23	2.01	0.42
1:A:101:VAL:O	1:A:105:MSE:HE2	2.18	0.42
1:A:35:VAL:O	1:A:35:VAL:HG12	2.19	0.42
1:C:149:LEU:HG	1:C:149:LEU:O	2.19	0.42
1:C:156:LEU:HA	1:C:156:LEU:HD12	1.76	0.42
1:A:142:THR:CG2	1:C:36:PHE:HE1	2.33	0.42
1:A:105:MSE:CE	1:A:175:ILE:HB	2.50	0.42
1:A:131:TRP:HH2	1:A:184:LEU:HD22	1.84	0.42
1:A:101:VAL:HG12	1:A:105:MSE:CE	2.50	0.42
1:C:80:TYR:C	1:C:82:PHE:H	2.24	0.41
1:B:202:LYS:O	1:B:202:LYS:HG3	2.20	0.41
1:B:203:ASN:O	1:B:205:GLU:N	2.53	0.41
1:A:76:ALA:HB2	1:A:91:ASN:CB	2.51	0.41
1:C:90:LEU:HG	1:C:164:VAL:HG13	2.03	0.41
1:C:4:LEU:CD1	1:C:4:LEU:C	2.89	0.41
1:C:108:LYS:HB2	1:C:108:LYS:HE2	1.77	0.41
1:B:201:PHE:CZ	1:C:201:PHE:CE2	3.09	0.41
1:A:37:ASN:HD21	1:B:158:THR:CG2	2.33	0.41
1:B:214:TYR:HA	1:B:217:TYR:HB2	2.03	0.41
1:A:105:MSE:HB2	1:A:105:MSE:HE2	1.70	0.41
1:A:41:TRP:CE3	1:A:42:LEU:HA	2.55	0.41
1:B:168:ILE:O	1:B:172:VAL:HG23	2.22	0.40
1:A:105:MSE:HE3	1:A:175:ILE:HB	2.03	0.40
1:A:23:LEU:HD23	3:A:302:BOG:C5'	2.52	0.40
1:B:198:VAL:O	1:B:200:TRP:N	2.54	0.40
1:A:45:VAL:HG12	1:A:74:LEU:CD1	2.51	0.40
1:A:149:LEU:HD23	1:A:156:LEU:HD13	2.02	0.40
1:B:148:TRP:O	1:B:149:LEU:HB3	2.22	0.40
1:C:213:MSE:HE3	1:C:214:TYR:CE2	2.56	0.40
1:C:200:TRP:C	1:C:201:PHE:CD1	2.95	0.40
1:A:31:ALA:O	1:A:34:PHE:HB3	2.22	0.40
1:B:37:ASN:C	1:B:39:ASP:H	2.24	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	233/244 (96%)	220 (94%)	13 (6%)	0	100	100
1	B	233/244 (96%)	214 (92%)	18 (8%)	1 (0%)	39	74
1	C	235/244 (96%)	217 (92%)	17 (7%)	1 (0%)	39	74
All	All	701/732 (96%)	651 (93%)	48 (7%)	2 (0%)	46	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	7	TRP
1	C	35	VAL

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	200/199 (100%)	173 (86%)	27 (14%)	5	14
1	B	200/199 (100%)	159 (80%)	41 (20%)	1	4
1	C	201/199 (101%)	180 (90%)	21 (10%)	9	25
All	All	601/597 (101%)	512 (85%)	89 (15%)	4	11

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER

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Mol	Chain	Res	Type
1	A	4	LEU
1	A	42	LEU
1	A	44	SER
1	A	62	ILE
1	A	88	MSE
1	A	90	LEU
1	A	105	MSE
1	A	114	GLU
1	A	115	THR
1	A	143	SER
1	A	144	VAL
1	A	154	SER
1	A	156	LEU
1	A	158	THR
1	A	160	ASP
1	A	163	THR
1	A	167	SER
1	A	174	MSE
1	A	184	LEU
1	A	188	VAL
1	A	200	TRP
1	A	205	GLU
1	A	211	LEU
1	A	212	LEU
1	A	218	LEU
1	A	225	TYR
1	B	4	LEU
1	B	25	MSE
1	B	30	GLN
1	B	33	VAL
1	B	37	ASN
1	B	39	ASP
1	B	42	LEU
1	B	66	LEU
1	B	69	LEU
1	B	83	LYS
1	B	88	MSE
1	B	90	LEU
1	B	92	LEU
1	B	98	VAL
1	B	105	MSE
1	B	112	LEU

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Mol	Chain	Res	Type
1	B	115	THR
1	B	128	VAL
1	B	139	VAL
1	B	143	SER
1	B	149	LEU
1	B	152	LEU
1	B	154	SER
1	B	158	THR
1	B	163	THR
1	B	167	SER
1	B	174	MSE
1	B	176	LEU
1	B	184	LEU
1	B	188	VAL
1	B	200	TRP
1	B	202	LYS
1	B	205	GLU
1	B	207	SER
1	B	213	MSE
1	B	218	LEU
1	B	225	TYR
1	B	229	THR
1	B	231	LEU
1	B	233	LYS
1	B	236	SER
1	C	3	SER
1	C	4	LEU
1	C	37	ASN
1	C	42	LEU
1	C	52	LEU
1	C	88	MSE
1	C	90	LEU
1	C	112	LEU
1	C	119	GLU
1	C	120	GLU
1	C	149	LEU
1	C	150	HIS
1	C	152	LEU
1	C	160	ASP
1	C	163	THR
1	C	167	SER
1	C	174	MSE

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Mol	Chain	Res	Type
1	C	188	VAL
1	C	205	GLU
1	C	218	LEU
1	C	225	TYR

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	37	ASN
1	A	171	GLN
1	B	37	ASN
1	C	37	ASN
1	C	235	HIS

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 ⓘ

7 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
2	NNR	A	301	-	16,19,19	4.96	7 (43%)	20,27,27	1.27	4 (20%)
3	BOG	A	302	-	20,20,20	1.22	2 (10%)	25,25,25	1.18	3 (12%)
2	NNR	B	301	-	16,19,19	5.16	6 (37%)	20,27,27	1.58	3 (15%)
3	BOG	B	302	-	20,20,20	1.52	2 (10%)	25,25,25	1.35	3 (12%)
3	BOG	B	303	-	20,20,20	1.41	2 (10%)	25,25,25	0.88	0
2	NNR	C	301	-	16,19,19	4.79	5 (31%)	20,27,27	1.03	1 (5%)
3	BOG	C	302	-	20,20,20	1.33	3 (15%)	25,25,25	1.19	3 (12%)

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	NNR	A	301	-	-	0/6/26/26	0/2/2/2
3	BOG	A	302	-	-	0/11/31/31	0/1/1/1
2	NNR	B	301	-	-	0/6/26/26	0/2/2/2
3	BOG	B	302	-	-	0/11/31/31	0/1/1/1
3	BOG	B	303	-	-	0/11/31/31	0/1/1/1
2	NNR	C	301	-	-	0/6/26/26	0/2/2/2
3	BOG	C	302	-	-	0/11/31/31	0/1/1/1

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	NNR	O4R-C4R	-6.16	1.30	1.45
2	B	301	NNR	O4R-C4R	-6.07	1.31	1.45
2	A	301	NNR	O4R-C4R	-5.67	1.31	1.45
2	B	301	NNR	O3R-C3R	-2.99	1.35	1.43
2	A	301	NNR	O3R-C3R	-2.91	1.36	1.43
2	C	301	NNR	O3R-C3R	-2.84	1.36	1.43
2	A	301	NNR	O7-C7	-2.33	1.19	1.24
2	A	301	NNR	C5R-C4R	2.07	1.59	1.51
3	C	302	BOG	O2-C2	2.19	1.48	1.43
3	C	302	BOG	O5-C5	2.46	1.50	1.44
3	B	303	BOG	O5-C5	2.54	1.50	1.44
2	B	301	NNR	C2-C3	2.56	1.42	1.39
3	A	302	BOG	O2-C2	2.57	1.49	1.43
3	A	302	BOG	O5-C1	2.75	1.48	1.41
3	C	302	BOG	O5-C1	3.39	1.50	1.41
3	B	302	BOG	O5-C1	3.69	1.51	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	BOG	O2-C2	3.77	1.52	1.43
3	B	303	BOG	O5-C1	4.00	1.52	1.41
2	C	301	NNR	O2R-C2R	4.48	1.53	1.43
2	B	301	NNR	O2R-C2R	4.51	1.53	1.43
2	A	301	NNR	O2R-C2R	4.80	1.54	1.43
2	A	301	NNR	C7-N7	5.54	1.44	1.33
2	C	301	NNR	C7-N7	5.56	1.44	1.33
2	B	301	NNR	C7-N7	6.34	1.45	1.33
2	C	301	NNR	O4R-C1R	15.83	1.61	1.41
2	A	301	NNR	O4R-C1R	16.67	1.62	1.41
2	B	301	NNR	O4R-C1R	17.27	1.63	1.41

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	NNR	C4R-O4R-C1R	-2.79	106.65	109.72
2	A	301	NNR	C4R-O4R-C1R	-2.55	106.92	109.72
2	B	301	NNR	C5R-C4R-C3R	-2.45	109.29	115.08
2	A	301	NNR	C5R-C4R-C3R	-2.45	109.29	115.08
3	C	302	BOG	C6-C5-C4	-2.32	107.30	113.02
3	A	302	BOG	C6-C5-C4	-2.23	107.51	113.02
2	C	301	NNR	O7-C7-N7	-2.01	119.77	122.59
2	A	301	NNR	C2-C3-C4	2.06	120.58	118.29
3	B	302	BOG	C4-C3-C2	2.06	114.64	110.79
2	A	301	NNR	C3-C7-N7	2.12	120.14	117.82
3	B	302	BOG	C3-C4-C5	2.16	113.96	110.20
3	C	302	BOG	C3-C4-C5	2.55	114.64	110.20
3	A	302	BOG	O5-C5-C4	2.56	114.49	109.68
3	A	302	BOG	O1-C1-C2	3.23	112.11	108.04
3	C	302	BOG	O5-C5-C4	3.31	115.89	109.68
2	B	301	NNR	O4R-C1R-N1	4.00	112.53	108.13
3	B	302	BOG	O1-C1-C2	4.23	113.38	108.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	NNR	1	0
3	A	302	BOG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	303	BOG	2	0
2	C	301	NNR	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	227/244 (93%)	0.58	26 (11%)	6 3	70, 112, 168, 196	0
1	B	227/244 (93%)	0.37	24 (10%)	8 4	72, 108, 174, 211	0
1	C	229/244 (93%)	0.48	24 (10%)	8 4	72, 115, 179, 210	0
All	All	683/732 (93%)	0.48	74 (10%)	8 3	70, 112, 175, 211	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	85	TYR	5.0
1	C	155	ALA	5.0
1	B	112	LEU	4.9
1	A	219	CYS	4.8
1	C	147	GLU	4.8
1	B	152	LEU	4.7
1	A	85	TYR	4.6
1	A	4	LEU	4.5
1	B	219	CYS	4.5
1	C	7	TRP	4.2
1	B	160	ASP	4.0
1	C	129	ARG	3.9
1	C	154	SER	3.8
1	B	189	ASN	3.7
1	A	12	PHE	3.6
1	B	217	TYR	3.6
1	C	93	LEU	3.6
1	B	221	SER	3.4
1	A	221	SER	3.3
1	C	11	LEU	3.2
1	A	101	VAL	3.1
1	A	112	LEU	3.0
1	A	192	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	115	THR	3.0
1	B	220	ASN	2.9
1	B	214	TYR	2.9
1	A	220	ASN	2.9
1	A	154	SER	2.9
1	B	223	TYR	2.9
1	B	186	ILE	2.9
1	A	214	TYR	2.8
1	A	11	LEU	2.7
1	A	94	VAL	2.7
1	B	93	LEU	2.7
1	C	69	LEU	2.7
1	A	36	PHE	2.7
1	A	215	VAL	2.6
1	B	35	VAL	2.6
1	C	76	ALA	2.6
1	C	91	ASN	2.6
1	B	187	VAL	2.6
1	A	93	LEU	2.6
1	C	114	GLU	2.6
1	A	62	ILE	2.6
1	A	116	ALA	2.5
1	C	92	LEU	2.5
1	A	144	VAL	2.5
1	C	121	VAL	2.5
1	C	131	TRP	2.5
1	C	219	CYS	2.4
1	C	100	PHE	2.4
1	B	188	VAL	2.4
1	B	190	ILE	2.4
1	C	12	PHE	2.3
1	C	60	GLY	2.3
1	A	217	TYR	2.3
1	A	206	THR	2.3
1	C	126	LEU	2.3
1	B	66	LEU	2.3
1	B	36	PHE	2.3
1	C	80	TYR	2.2
1	B	218	LEU	2.2
1	C	156	LEU	2.2
1	B	148	TRP	2.2
1	A	186	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	118	THR	2.1
1	C	4	LEU	2.1
1	B	145	TYR	2.1
1	B	185	TRP	2.1
1	A	80	TYR	2.0
1	A	222	VAL	2.0
1	B	100	PHE	2.0
1	A	107	ARG	2.0
1	C	134	VAL	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](#)

There are no carbohydrates in this entry.

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. 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
3	BOG	B	303	20/20	0.94	0.34	1.19	92,99,116,126	0
3	BOG	A	302	20/20	0.95	0.31	0.34	72,102,127,142	0
2	NNR	A	301	18/18	0.95	0.28	0.25	97,105,122,129	0
3	BOG	B	302	20/20	0.88	0.26	0.15	76,100,122,135	0
3	BOG	C	302	20/20	0.93	0.25	-0.33	81,124,153,158	0
2	NNR	C	301	18/18	0.94	0.19	-0.35	77,99,117,125	0
2	NNR	B	301	18/18	0.93	0.24	-0.38	80,95,109,111	0

6.5 Other polymers [i](#)

There are no such residues in this entry.