



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 13, 2024 – 11:06 pm GMT

PDB ID : 6QZK  
Title : Structure of Clostridium butyricum Argonaute bound to a guide DNA (5' deoxycytidine) and a 19-mer target DNA  
Authors : Swarts, D.C.; Jinek, M.; Hegge, J.W.; Van der Oost, J.  
Deposited on : 2019-03-11  
Resolution : 3.55 Å (reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

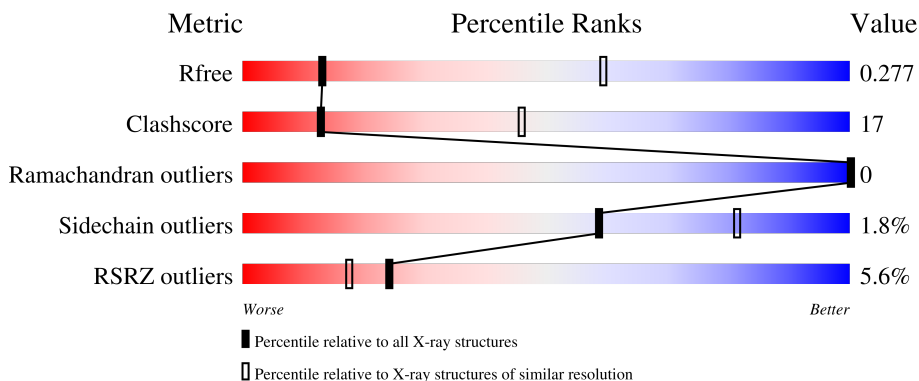
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


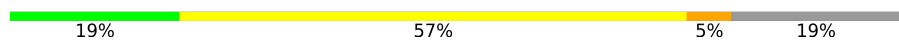

The reported resolution of this entry is 3.55 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1028 (3.60-3.48)
Clashscore	141614	1109 (3.60-3.48)
Ramachandran outliers	138981	1073 (3.60-3.48)
Sidechain outliers	138945	1074 (3.60-3.48)
RSRZ outliers	127900	1079 (3.62-3.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	748	
2	B	21	
3	C	19	

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
5	FMT	B	101	-	-	-	X

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6749 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 Clostridium butyricum Argonaute.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	746	6033	3916	985	1112	20	0	0	0

- Molecule 2 is a DNA chain called siDNA guide (5'-D(P\*CP\*GP\*AP\*GP\*GP\*TP\*AP\*GP\*TP\*AP\*GP\*GP\*TP\*TP\*GP\*TP\*AP\*TP\*AP\*GP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	17	341	159	63	102	17	0	0	0

- Molecule 3 is a DNA chain called DNA target (5'-D(T\*AP\*TP\*AP\*CP\*AP\*AP\*CP\*CP\*TP\*AP\*CP\*TP\*AP\*CP\*CP\*TP\*CP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	C	18	356	173	61	105	17	0	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is FORMIC ACID (three-letter code: FMT) (formula: CH<sub>2</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 3 1 2	0	0
5	A	1	Total C O 3 1 2	0	0
5	B	1	Total C O 3 1 2	0	0
5	B	1	Total C O 3 1 2	0	0
5	B	1	Total C O 3 1 2	0	0

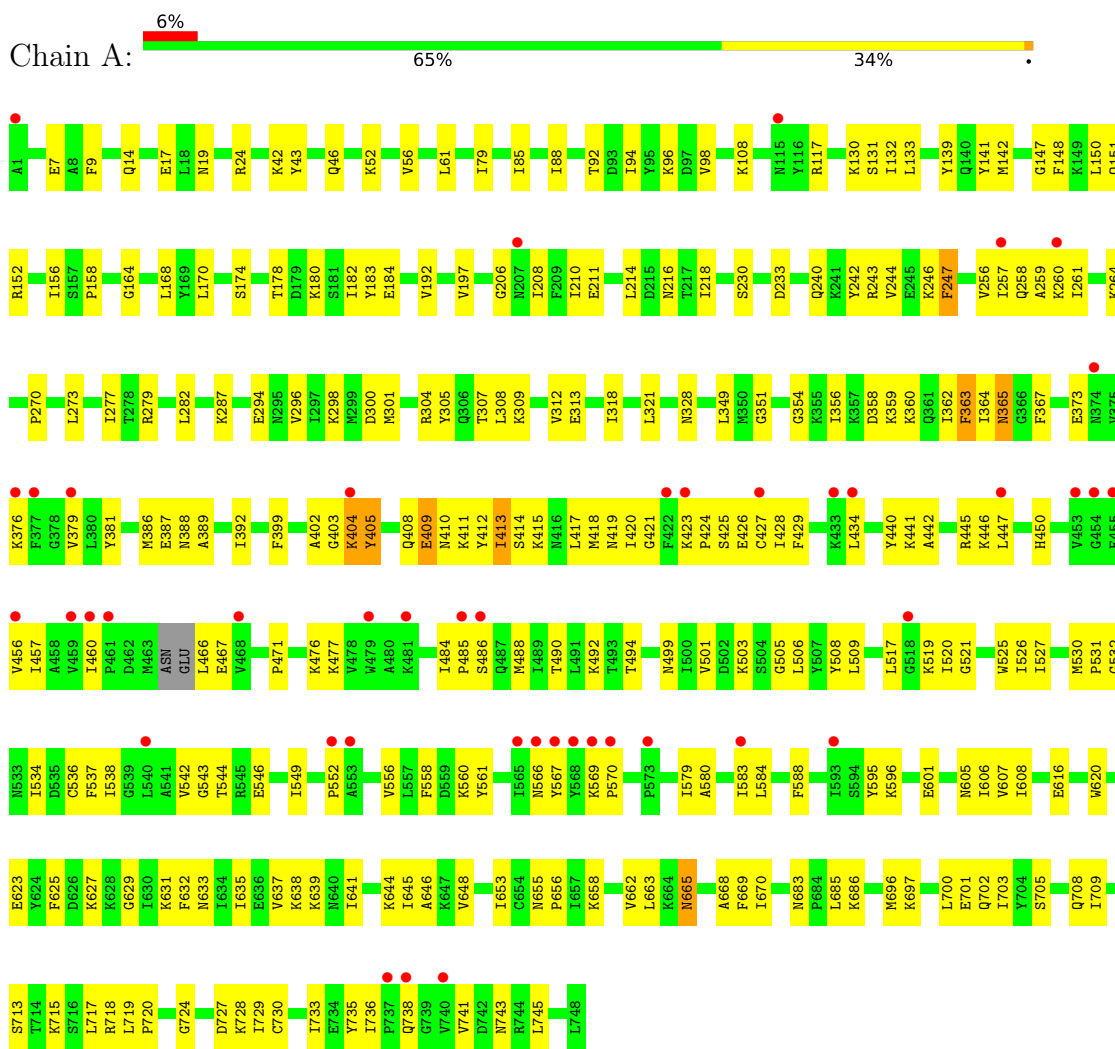
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	2	Total O 2 2	0	0
6	B	1	Total O 1 1	0	0

### 3 Residue-property plots

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

- Molecule 1: Clostridium butyricum Argonaute



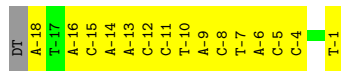
- Molecule 2: siDNA guide (5'-D(P\*CP\*GP\*AP\*GP\*GP\*TP\*AP\*GP\*TP\*AP\*GP\*GP\*TP\*TP\*GP\*TP\*AP\*TP\*AP\*GP\*T)-3')





- Molecule 3: DNA target (5'-D(T\*AP\*TP\*AP\*CP\*AP\*AP\*CP\*CP\*TP\*AP\*CP\*TP\*AP\*CP\*CP\*TP\*CP\*T)-3')

Chain C: 16% 79% 5%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	181.45Å 181.45Å 142.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	76.61 – 3.55 90.73 – 3.55	Depositor EDS
% Data completeness (in resolution range)	99.9 (76.61-3.55) 100.0 (90.73-3.55)	Depositor EDS
$R_{merge}$	0.28	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.48 (at 3.58Å)	Xtrriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.247 , 0.276 0.247 , 0.277	Depositor DCC
$R_{free}$ test set	1732 reflections (9.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	98.6	Xtrriage
Anisotropy	0.723	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 60.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.84	EDS
Total number of atoms	6749	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	110.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.66% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality i

### 5.1 Standard geometry i

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

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	0.28	0/6157	0.46	0/8298
2	B	0.87	1/382 (0.3%)	1.08	0/589
3	C	0.73	0/397	1.05	0/608
All	All	0.37	1/6936 (0.0%)	0.58	0/9495

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	DC	OP3-P	-10.75	1.48	1.61

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts i

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6033	0	6162	197	0
2	B	341	0	181	17	0
3	C	356	0	205	18	0
4	A	1	0	0	0	0
5	A	6	0	2	0	0
5	B	9	0	3	0	0
6	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	1	0	0	0	0
All	All	6749	0	6553	220	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 (220) 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:532:GLY:HA2	1:A:560:LYS:HD3	1.57	0.85
1:A:156:ILE:HG22	1:A:158:PRO:HD3	1.63	0.80
1:A:373:GLU:HA	1:A:419:ASN:HB3	1.64	0.78
1:A:132:ILE:HD12	1:A:296:VAL:HG11	1.68	0.76
1:A:367:PHE:HE2	1:A:417:LEU:HB3	1.50	0.75
1:A:424:PRO:HG2	1:A:426:GLU:HG2	1.69	0.74
1:A:257:ILE:HD11	1:A:273:LEU:HD12	1.69	0.73
1:A:178:THR:HG22	1:A:180:LYS:H	1.56	0.71
1:A:506:LEU:HD12	3:C:-1:DT:H71	1.71	0.70
2:B:14:DT:H2'	2:B:15:DG:C8	2.25	0.70
1:A:147:GLY:HA3	1:A:178:THR:HG23	1.76	0.68
1:A:279:ARG:HG2	3:C:-6:DA:H1'	1.76	0.68
1:A:410:ASN:ND2	1:A:413:ILE:O	2.25	0.67
1:A:132:ILE:HD11	1:A:142:MET:SD	2.35	0.67
1:A:379:VAL:HA	1:A:457:ILE:O	1.94	0.67
1:A:206:GLY:N	2:B:11:DG:OP1	2.28	0.67
3:C:-15:DC:H2'	3:C:-14:DA:C8	2.30	0.67
1:A:607:VAL:HG22	1:A:633:ASN:HB2	1.76	0.66
2:B:13:DT:H2'	2:B:14:DT:H71	1.78	0.65
3:C:-7:DT:H2''	3:C:-6:DA:O5'	1.96	0.65
1:A:362:ILE:HG12	1:A:525:TRP:HE3	1.62	0.65
1:A:42:LYS:HE3	1:A:43:TYR:HE1	1.61	0.64
2:B:2:DG:H2'	2:B:3:DA:C8	2.32	0.64
1:A:527:ILE:HD11	1:A:705:SER:HB3	1.79	0.64
1:A:392:ILE:HB	1:A:494:THR:HG23	1.78	0.64
1:A:639:LYS:HD3	1:A:719:LEU:HD21	1.80	0.63
1:A:409:GLU:HG2	1:A:411:LYS:H	1.64	0.63
1:A:530:MET:HG2	1:A:560:LYS:HB2	1.81	0.62
1:A:705:SER:O	1:A:709:ILE:HG13	1.99	0.62
1:A:309:LYS:HE3	1:A:328:ASN:HB2	1.82	0.62
1:A:638:LYS:HA	3:C:-11:DC:OP1	1.99	0.62
1:A:596:LYS:HG3	1:A:601:GLU:HA	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:LYS:HA	1:A:421:GLY:HA3	1.82	0.61
2:B:3:DA:H2'	2:B:4:DG:C8	2.35	0.61
1:A:381:TYR:HE2	1:A:429:PHE:HD1	1.49	0.60
1:A:526:ILE:HG23	1:A:561:TYR:HB3	1.84	0.60
1:A:240:GLN:HG2	1:A:243:ARG:HG2	1.84	0.60
3:C:-10:DT:H2'	3:C:-9:DA:C8	2.37	0.60
1:A:399:PHE:HA	1:A:404:LYS:HA	1.84	0.60
1:A:558:PHE:HE2	1:A:566:ASN:HB2	1.65	0.60
1:A:543:GLY:HA3	1:A:730:CYS:HB3	1.84	0.59
1:A:414:SER:HB2	1:A:417:LEU:HD21	1.84	0.59
3:C:-10:DT:H2'	3:C:-9:DA:H8	1.65	0.59
1:A:499:ASN:HB3	1:A:501:VAL:HG23	1.85	0.59
1:A:376:LYS:H	1:A:376:LYS:HD2	1.68	0.57
1:A:427:CYS:HA	1:A:450:HIS:CD2	2.39	0.57
1:A:670:ILE:HD11	1:A:700:LEU:HB3	1.86	0.57
1:A:354:GLY:N	1:A:365:ASN:OD1	2.37	0.57
1:A:442:ALA:O	1:A:446:LYS:HG2	2.05	0.57
1:A:362:ILE:HG21	1:A:525:TRP:HZ3	1.69	0.57
1:A:381:TYR:CE2	1:A:429:PHE:HD1	2.22	0.57
1:A:440:TYR:CD1	1:A:471:PRO:HB3	2.39	0.57
1:A:663:LEU:HD12	1:A:668:ALA:HB2	1.86	0.56
2:B:4:DG:H2'	2:B:5:DG:C8	2.39	0.56
1:A:648:VAL:HG22	1:A:653:ILE:HG12	1.86	0.56
2:B:7:DA:H2'	2:B:8:DG:C8	2.42	0.55
3:C:-14:DA:H2'	3:C:-13:DA:C8	2.42	0.55
1:A:412:TYR:O	1:A:413:ILE:HD13	2.06	0.55
1:A:703:ILE:HG23	1:A:720:PRO:HG3	1.88	0.55
1:A:663:LEU:HD11	1:A:696:MET:HB3	1.89	0.55
1:A:318:ILE:HG22	1:A:321:LEU:H	1.71	0.54
1:A:558:PHE:HD1	1:A:595:TYR:HD2	1.54	0.54
1:A:427:CYS:HA	1:A:450:HIS:HD2	1.72	0.54
3:C:-7:DT:H2'	3:C:-6:DA:C8	2.43	0.54
1:A:14:GLN:HG3	1:A:164:GLY:HA3	1.89	0.54
1:A:240:GLN:HG3	1:A:242:TYR:CE2	2.43	0.53
1:A:645:ILE:HG23	1:A:656:PRO:HG3	1.91	0.53
1:A:17:GLU:HG2	1:A:88:ILE:HG13	1.90	0.53
1:A:530:MET:SD	1:A:531:PRO:HD2	2.49	0.53
1:A:569:LYS:HG3	1:A:570:PRO:HD2	1.90	0.53
1:A:182:ILE:HG23	1:A:192:VAL:HG21	1.91	0.52
1:A:298:LYS:HE2	2:B:6:DT:O2	2.10	0.52
1:A:305:TYR:CD2	1:A:653:ILE:HD12	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:LYS:HE3	1:A:43:TYR:CE1	2.43	0.52
1:A:490:THR:HG22	1:A:492:LYS:H	1.75	0.52
1:A:197:VAL:HG11	1:A:273:LEU:HB3	1.91	0.52
1:A:261:ILE:N	1:A:264:LYS:O	2.36	0.52
1:A:180:LYS:HG3	1:A:184:GLU:HB2	1.91	0.52
1:A:364:ILE:HA	1:A:414:SER:OG	2.08	0.52
1:A:536:CYS:O	1:A:606:ILE:HG13	2.09	0.52
1:A:387:GLU:H	1:A:387:GLU:CD	2.14	0.51
1:A:300:ASP:OD1	1:A:301:MET:N	2.44	0.51
1:A:405:TYR:CD1	1:A:506:LEU:HD21	2.46	0.51
1:A:170:LEU:HD23	1:A:308:LEU:HB3	1.92	0.51
1:A:623:GLU:HG3	1:A:627:LYS:HE3	1.93	0.51
1:A:724:GLY:O	1:A:728:LYS:HG2	2.11	0.51
1:A:183:TYR:HB2	1:A:270:PRO:HB2	1.93	0.51
1:A:379:VAL:HG23	1:A:423:LYS:HE2	1.94	0.50
1:A:367:PHE:CD1	1:A:517:LEU:HD22	2.46	0.50
1:A:520:ILE:HG22	1:A:521:GLY:H	1.75	0.50
1:A:441:LYS:O	1:A:445:ARG:HG2	2.11	0.50
1:A:505:GLY:O	1:A:508:TYR:HB2	2.12	0.50
1:A:133:LEU:HB3	1:A:307:THR:HG22	1.93	0.50
1:A:131:SER:HB3	1:A:141:TYR:CE2	2.47	0.50
1:A:404:LYS:HE2	1:A:506:LEU:HD22	1.93	0.50
1:A:538:ILE:HG12	1:A:556:VAL:HG22	1.92	0.50
1:A:208:ILE:HA	1:A:261:ILE:HG22	1.95	0.49
1:A:152:ARG:NH2	1:A:174:SER:OG	2.46	0.49
1:A:362:ILE:HG21	1:A:525:TRP:CZ3	2.47	0.49
1:A:363:PHE:N	1:A:363:PHE:HD1	2.10	0.49
1:A:637:VAL:HG13	1:A:685:LEU:HD13	1.94	0.49
1:A:358:ASP:OD1	1:A:359:LYS:N	2.46	0.49
1:A:381:TYR:CD1	1:A:386:MET:HB2	2.47	0.49
1:A:569:LYS:NZ	1:A:733:ILE:O	2.28	0.49
1:A:519:LYS:NZ	2:B:1:DC:OP3	2.41	0.49
1:A:566:ASN:OD1	1:A:743:ASN:HA	2.13	0.49
1:A:538:ILE:O	1:A:608:ILE:HA	2.13	0.49
1:A:519:LYS:HA	1:A:745:LEU:O	2.13	0.48
1:A:46:GLN:HB2	1:A:52:LYS:O	2.14	0.48
1:A:363:PHE:N	1:A:363:PHE:CD1	2.80	0.48
1:A:434:LEU:HA	1:A:440:TYR:HE2	1.78	0.48
1:A:108:LYS:O	1:A:117:ARG:HD2	2.13	0.48
1:A:349:LEU:HB2	1:A:356:ILE:HD11	1.94	0.48
1:A:418:MET:HG3	1:A:420:ILE:HG23	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:658:LYS:HD3	1:A:708:GLN:NE2	2.29	0.48
1:A:718:ARG:HB2	2:B:5:DG:H5''	1.95	0.48
1:A:131:SER:HB3	1:A:141:TYR:CD2	2.48	0.48
1:A:218:ILE:HG22	1:A:256:VAL:HG22	1.96	0.48
1:A:405:TYR:HE1	1:A:503:LYS:HA	1.79	0.47
3:C:-7:DT:H2'	3:C:-6:DA:H8	1.78	0.47
2:B:5:DG:H2'	2:B:6:DT:C6	2.49	0.47
1:A:460:ILE:HD11	1:A:488:MET:HG2	1.95	0.47
1:A:7:GLU:OE2	1:A:304:ARG:NH2	2.44	0.47
1:A:713:SER:HB2	2:B:3:DA:H4'	1.97	0.47
1:A:425:SER:O	1:A:428:ILE:HG13	2.15	0.47
1:A:665:ASN:O	1:A:665:ASN:ND2	2.44	0.47
1:A:294:GLU:OE2	1:A:298:LYS:NZ	2.32	0.46
1:A:178:THR:HG22	1:A:180:LYS:N	2.26	0.46
1:A:373:GLU:HA	1:A:419:ASN:CB	2.38	0.46
1:A:132:ILE:HG21	1:A:150:LEU:HD22	1.98	0.46
1:A:309:LYS:HE2	1:A:313:GLU:OE2	2.16	0.46
1:A:542:VAL:HG12	1:A:552:PRO:HA	1.97	0.46
1:A:403:GLY:HA3	1:A:408:GLN:HG2	1.97	0.46
1:A:697:LYS:O	1:A:701:GLU:HG3	2.16	0.46
2:B:15:DG:H2''	2:B:16:DT:H5''	1.98	0.46
1:A:388:ASN:HB3	1:A:494:THR:CG2	2.46	0.46
1:A:717:LEU:HD21	1:A:724:GLY:HA3	1.98	0.46
1:A:404:LYS:HB2	1:A:413:ILE:HG21	1.97	0.45
1:A:605:ASN:HB3	1:A:631:LYS:HB3	1.99	0.45
1:A:24:ARG:HB2	1:A:79:ILE:HG13	1.98	0.45
1:A:527:ILE:CD1	1:A:705:SER:HB3	2.45	0.45
1:A:211:GLU:HB2	1:A:260:LYS:H	1.81	0.45
1:A:240:GLN:HB2	3:C:-18:DA:O5'	2.16	0.45
1:A:404:LYS:CE	1:A:413:ILE:HG12	2.46	0.45
1:A:309:LYS:CE	1:A:328:ASN:HB2	2.45	0.45
1:A:216:ASN:O	1:A:256:VAL:HG23	2.16	0.44
1:A:56:VAL:HG22	1:A:61:LEU:HD22	1.99	0.44
1:A:584:LEU:HB3	1:A:620:TRP:CH2	2.51	0.44
1:A:641:ILE:HG21	1:A:686:LYS:HG3	1.98	0.44
1:A:644:LYS:NZ	1:A:683:ASN:OD1	2.46	0.44
1:A:9:PHE:CD2	1:A:312:VAL:HG21	2.52	0.44
1:A:304:ARG:CZ	1:A:644:LYS:HD2	2.48	0.44
1:A:645:ILE:HD11	1:A:662:VAL:HG21	1.99	0.44
1:A:301:MET:CE	1:A:646:ALA:HB2	2.48	0.44
1:A:531:PRO:HG2	1:A:702:GLN:HG3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:580:ALA:HB3	1:A:583:ILE:HG12	2.00	0.44
1:A:389:ALA:O	1:A:392:ILE:HG22	2.18	0.44
1:A:409:GLU:HG3	1:A:411:LYS:NZ	2.33	0.44
1:A:402:ALA:HB1	1:A:415:LYS:HE2	1.99	0.43
1:A:466:LEU:HG	1:A:467:GLU:H	1.84	0.43
1:A:460:ILE:HG23	1:A:471:PRO:HD2	1.99	0.43
1:A:538:ILE:HG13	1:A:606:ILE:HD11	2.00	0.43
1:A:607:VAL:HG13	1:A:635:ILE:HD12	2.00	0.43
1:A:447:LEU:HD11	1:A:456:VAL:HG21	2.00	0.43
1:A:606:ILE:O	1:A:632:PHE:HA	2.18	0.43
2:B:4:DG:H2'	2:B:5:DG:H8	1.81	0.43
1:A:301:MET:HG2	1:A:655:ASN:OD1	2.18	0.43
1:A:404:LYS:HE2	1:A:506:LEU:CD2	2.49	0.43
1:A:639:LYS:NZ	1:A:727:ASP:OD1	2.50	0.43
1:A:244:VAL:HG12	1:A:247:PHE:CZ	2.53	0.43
1:A:486:SER:OG	1:A:488:MET:SD	2.63	0.43
1:A:729:ILE:HA	1:A:736:ILE:CD1	2.49	0.43
1:A:546:GLU:HB3	1:A:549:ILE:HD12	2.01	0.43
1:A:94:ILE:O	1:A:98:VAL:HG23	2.19	0.42
1:A:608:ILE:HG22	1:A:633:ASN:O	2.19	0.42
2:B:7:DA:H2'	2:B:8:DG:H8	1.82	0.42
1:A:130:LYS:HA	1:A:151:GLN:HG2	2.02	0.42
1:A:569:LYS:HE2	1:A:738:GLN:HB3	2.02	0.42
1:A:214:LEU:HD11	1:A:258:GLN:HB2	2.02	0.42
1:A:625:PHE:O	1:A:629:GLY:N	2.53	0.42
1:A:569:LYS:HB3	1:A:738:GLN:HA	2.00	0.42
1:A:476:LYS:HB3	1:A:735:TYR:HB3	2.01	0.42
1:A:641:ILE:HD12	1:A:669:PHE:HE1	1.83	0.42
1:A:279:ARG:HA	1:A:282:LEU:HD12	2.01	0.42
1:A:399:PHE:HD1	1:A:509:LEU:HD13	1.84	0.42
1:A:715:LYS:HZ2	2:B:3:DA:H2	1.66	0.42
1:A:19:ASN:HA	1:A:85:ILE:HG22	2.03	0.41
1:A:230:SER:HB3	1:A:233:ASP:OD2	2.21	0.41
1:A:304:ARG:O	1:A:308:LEU:HG	2.20	0.41
1:A:381:TYR:HE2	1:A:429:PHE:CD1	2.33	0.41
1:A:605:ASN:HA	1:A:631:LYS:O	2.20	0.41
1:A:686:LYS:HE3	3:C:-12:DC:OP2	2.20	0.41
3:C:-16:DA:H2''	3:C:-15:DC:H6	1.85	0.41
1:A:156:ILE:HG23	1:A:168:LEU:HD11	2.02	0.41
1:A:360:LYS:NZ	3:C:-1:DT:P	2.93	0.41
1:A:534:ILE:HD13	1:A:537:PHE:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:709:ILE:HG22	1:A:709:ILE:O	2.21	0.41
1:A:211:GLU:H	1:A:259:ALA:HA	1.85	0.41
1:A:526:ILE:HG22	1:A:527:ILE:N	2.36	0.41
3:C:-5:DC:H2''	3:C:-4:DC:H5'	2.02	0.41
1:A:558:PHE:CD1	1:A:595:TYR:HD2	2.35	0.41
3:C:-12:DC:H2'	3:C:-11:DC:C6	2.56	0.41
1:A:246:LYS:HD2	1:A:246:LYS:HA	1.87	0.41
1:A:364:ILE:HA	1:A:414:SER:HG	1.83	0.41
1:A:588:PHE:CE1	1:A:625:PHE:HZ	2.39	0.41
1:A:92:THR:O	1:A:96:LYS:HG3	2.21	0.41
1:A:287:LYS:HB2	1:A:287:LYS:HE3	1.88	0.41
1:A:404:LYS:HG2	1:A:405:TYR:N	2.33	0.41
1:A:526:ILE:CG2	1:A:527:ILE:N	2.83	0.41
3:C:-9:DA:H2'	3:C:-8:DC:C6	2.56	0.41
3:C:-6:DA:H2'	3:C:-5:DC:C6	2.56	0.41
1:A:569:LYS:HE3	1:A:733:ILE:HG23	2.03	0.40
1:A:351:GLY:HA3	1:A:365:ASN:HB3	2.04	0.40
1:A:484:ILE:HA	1:A:485:PRO:HD3	1.95	0.40
1:A:192:VAL:HG22	1:A:210:ILE:HD12	2.03	0.40
1:A:544:THR:O	1:A:544:THR:OG1	2.38	0.40
1:A:638:LYS:HB2	1:A:686:LYS:HB2	2.02	0.40
1:A:579:ILE:HB	1:A:616:GLU:OE2	2.21	0.40
1:A:148:PHE:HE1	1:A:277:ILE:HD12	1.86	0.40
2:B:5:DG:H2'	2:B:6:DT:H6	1.86	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	742/748 (99%)	710 (96%)	32 (4%)	0	<b>100</b> <b>100</b>

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	670/672 (100%)	658 (98%)	12 (2%)	59 81

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	139	TYR
1	A	247	PHE
1	A	363	PHE
1	A	365	ASN
1	A	404	LYS
1	A	405	TYR
1	A	409	GLU
1	A	413	ILE
1	A	477	LYS
1	A	567	TYR
1	A	665	ASN
1	A	741	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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



## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	$\# Z  > 2$
5	FMT	B	102	-	2,2,2	0.71	0	1,1,1	0.23	0
5	FMT	B	103	-	2,2,2	0.70	0	1,1,1	0.24	0
5	FMT	A	803	-	2,2,2	0.71	0	1,1,1	0.25	0
5	FMT	B	101	-	2,2,2	0.71	0	1,1,1	0.24	0
5	FMT	A	802	-	2,2,2	0.70	0	1,1,1	0.24	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	746/748 (99%)	0.49	44 (5%) 22 16	70, 110, 148, 196	0
2	B	17/21 (80%)	0.07	0 100 100	81, 100, 136, 160	0
3	C	18/19 (94%)	0.19	0 100 100	88, 103, 140, 144	0
All	All	781/788 (99%)	0.48	44 (5%) 24 17	70, 109, 148, 196	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	374	ASN	5.1
1	A	453	VAL	3.7
1	A	427	CYS	3.7
1	A	567	TYR	3.6
1	A	455	PHE	3.6
1	A	553	ALA	3.4
1	A	1	ALA	3.4
1	A	454	GLY	3.3
1	A	486	SER	3.2
1	A	485	PRO	3.2
1	A	422	PHE	3.2
1	A	434	LEU	3.0
1	A	423	LYS	2.9
1	A	460	ILE	2.8
1	A	552	PRO	2.8
1	A	456	VAL	2.7
1	A	115	ASN	2.7
1	A	518	GLY	2.5
1	A	566	ASN	2.5
1	A	404	LYS	2.4
1	A	377	PHE	2.4
1	A	738	GLN	2.4
1	A	479	TRP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	593	ILE	2.3
1	A	481	LYS	2.3
1	A	459	VAL	2.3
1	A	740	VAL	2.2
1	A	737	PRO	2.2
1	A	468	VAL	2.2
1	A	568	TYR	2.2
1	A	379	VAL	2.2
1	A	207	ASN	2.2
1	A	447	LEU	2.1
1	A	570	PRO	2.1
1	A	461	PRO	2.1
1	A	565	ILE	2.1
1	A	540	LEU	2.1
1	A	257	ILE	2.1
1	A	569	LYS	2.0
1	A	583	ILE	2.0
1	A	260	LYS	2.0
1	A	376	LYS	2.0
1	A	433	LYS	2.0
1	A	573	PRO	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 monosaccharides 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. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	FMT	B	101	3/3	0.48	0.50	113,113,113,113	0
5	FMT	A	803	3/3	0.51	0.31	125,125,126,126	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	FMT	B	103	3/3	0.54	0.31	119,119,119,120	0
5	FMT	B	102	3/3	0.74	0.36	99,99,99,99	0
5	FMT	A	802	3/3	0.74	0.39	100,100,101,101	0
4	MG	A	801	1/1	0.95	0.07	97,97,97,97	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.