



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 12:02 PM UTC

PDB ID : 6QZK / pdb_00006qzk
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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

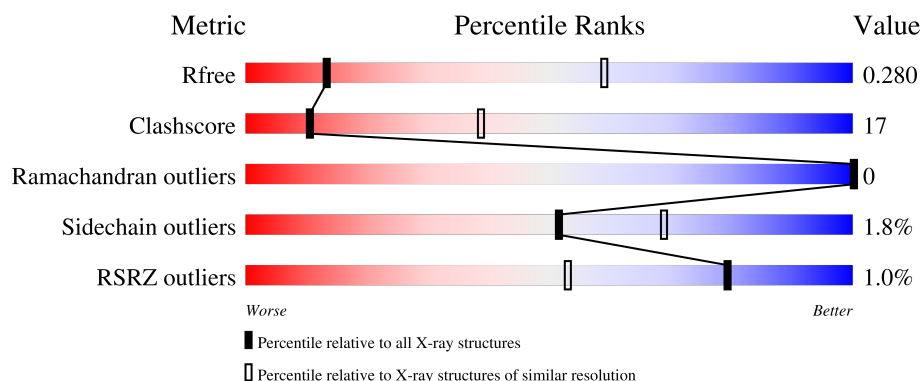
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

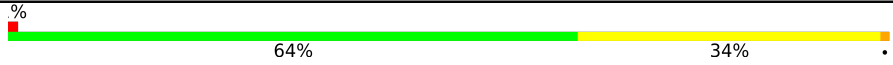
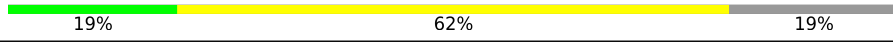
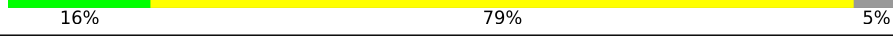
The reported resolution of this entry is 3.55 Å.

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}	180053	1008 (3.58-3.50)
Clashscore	190562	1062 (3.58-3.50)
Ramachandran outliers	187476	1033 (3.58-3.50)
Sidechain outliers	187428	1034 (3.58-3.50)
RSRZ outliers	180081	1007 (3.58-3.50)

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

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

2 Entry composition

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 Protein argonaute.

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

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	ALA	MET	conflict	UNP A0ABF7PLX0
A	188	ASP	ASN	conflict	UNP A0ABF7PLX0
A	191	GLY	ASP	conflict	UNP A0ABF7PLX0
A	200	LYS	ARG	conflict	UNP A0ABF7PLX0
A	204	ALA	SER	conflict	UNP A0ABF7PLX0
A	212	LYS	GLU	conflict	UNP A0ABF7PLX0
A	216	ASN	LYS	conflict	UNP A0ABF7PLX0
A	217	THR	SER	conflict	UNP A0ABF7PLX0
A	220	ASP	GLU	conflict	UNP A0ABF7PLX0
A	253	ASN	LYS	conflict	UNP A0ABF7PLX0
A	258	GLN	LYS	conflict	UNP A0ABF7PLX0
A	343	VAL	ILE	conflict	UNP A0ABF7PLX0
A	466	LEU	SER	conflict	UNP A0ABF7PLX0
A	541	ALA	ASP	engineered mutation	UNP A0ABF7PLX0
A	611	ALA	ASP	engineered mutation	UNP A0ABF7PLX0

- 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
2	B	17	Total	C	N	O	P	0	0	0
			341	159	63	102	17			

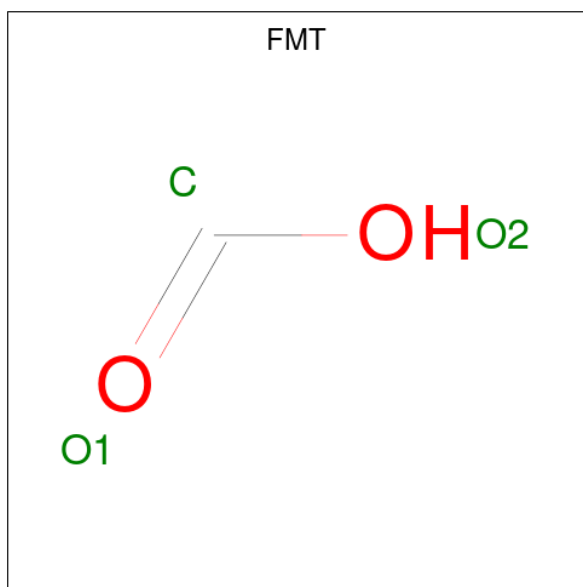
- Molecule 3 is a DNA chain called DNA target (5'-D(T*AP*TP*AP*CP*AP*AP*CP*CP*TP*AP*CP*TP*AP*CP*CP*TP*CP*T)-3').

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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



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

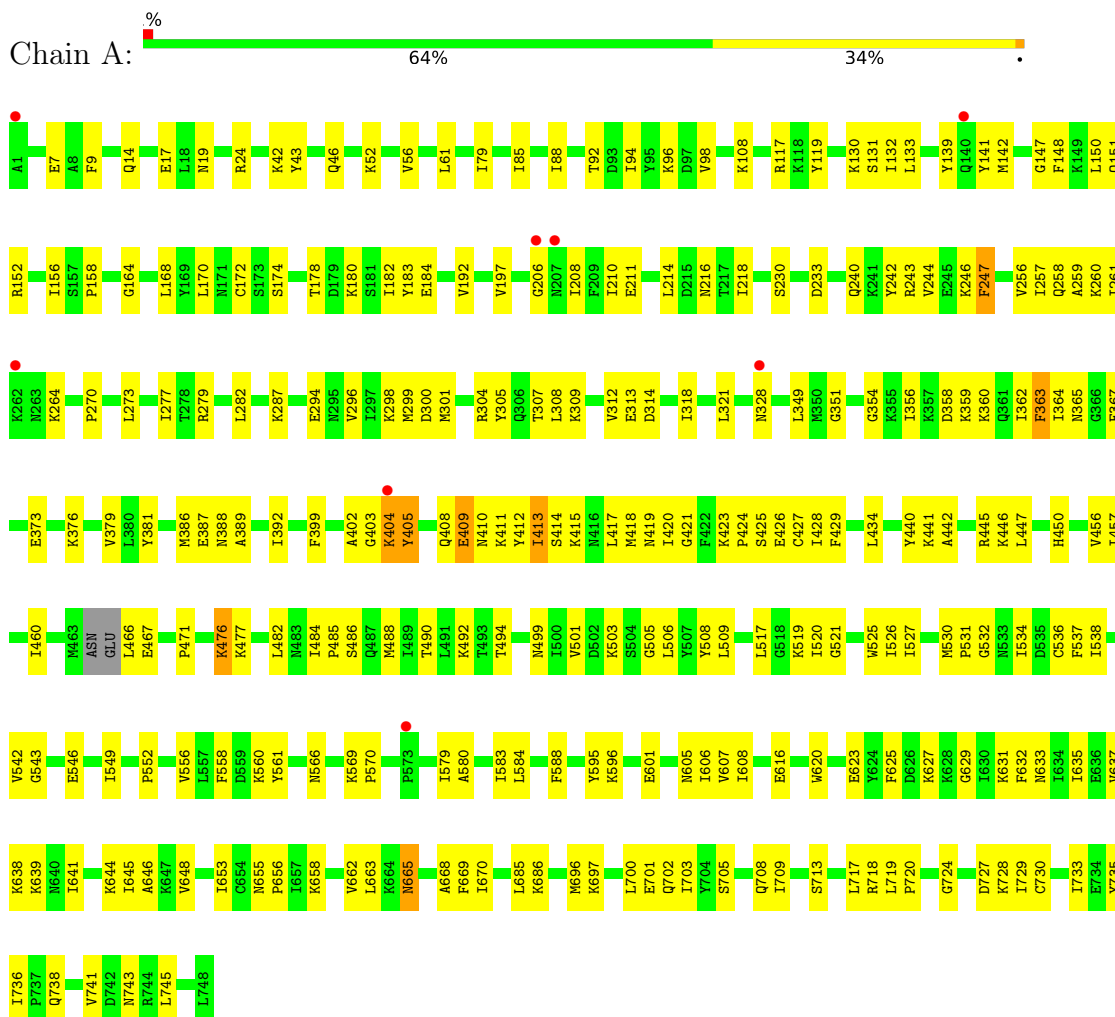
- Molecule 6 is water.

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

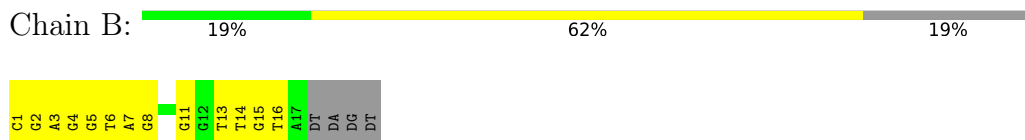
3 Residue-property plots

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.

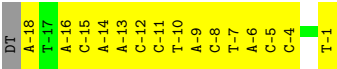
• Molecule 1: Protein argonaute



• 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')



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	181.45Å 181.45Å 142.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	76.61 – 3.55 76.61 – 3.55	Depositor EDS
% Data completeness (in resolution range)	99.9 (76.61-3.55) 99.9 (76.61-3.55)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 3.58Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.247 , 0.276 0.251 , 0.280	Depositor DCC
R_{free} test set	1732 reflections (6.79%)	wwPDB-VP
Wilson B-factor (Å ²)	98.6	Xtriage
Anisotropy	0.723	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 106.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6749	wwPDB-VP
Average B, all atoms (Å ²)	110.0	wwPDB-VP

Xtriage'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.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.19	0/6157	0.42	0/8298
2	B	0.29	0/382	0.56	0/589
3	C	0.32	0/397	0.62	0/608
All	All	0.20	0/6936	0.44	0/9495

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 [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	200	0
2	B	341	0	181	16	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
6	B	1	0	0	0	0
All	All	6749	0	6553	223	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 (223) 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
1:A:299:MET:HE3	1:A:304:ARG:HG2	1.80	0.61
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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:569:LYS:NZ	1:A:733:ILE:O	2.28	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
1:A:387:GLU:CD	1:A:387:GLU:H	2.14	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
1:A:301:MET:HE2	1:A:646:ALA:HB2	1.87	0.55
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:486:SER:OG	1:A:488:MET:SD	2.63	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
1:A:42:LYS:HE3	1:A:43:TYR:CE1	2.43	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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:172:CYS:SG	1:A:299:MET:HE1	2.53	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
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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:527:ILE:HD12	1:A:530:MET:HE2	1.95	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:142:MET:HE2	1:A:142:MET:HB3	1.88	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:9:PHE:CD2	1:A:312:VAL:HG21	2.52	0.44
1:A:119:TYR:OH	1:A:314:ASP:OD2	2.29	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

Continued on next page...

Continued from previous page...

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: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: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:364:ILE:HA	1:A:414:SER:HG	1.84	0.41
1:A:534:ILE:HD13	1:A:537:PHE:CE1	2.55	0.41
1:A:709:ILE:HG22	1:A:709:ILE:O	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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 ⓘ

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	742/748 (99%)	710 (96%)	32 (4%)	0	100	100

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%)	51 69

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	404	LYS
1	A	405	TYR
1	A	409	GLU
1	A	413	ILE
1	A	476	LYS
1	A	477	LYS
1	A	482	LEU
1	A	665	ASN
1	A	741	VAL

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

Mol	Chain	Res	Type
1	A	48	ASN
1	A	267	ASN
1	A	470	ASN
1	A	550	HIS

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 oligosaccharides 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	0.73	0	1,1,1	0.22	0
5	FMT	B	103	-	2,2,2	0.73	0	1,1,1	0.24	0
5	FMT	B	101	-	2,2,2	0.74	0	1,1,1	0.24	0
5	FMT	A	803	-	2,2,2	0.73	0	1,1,1	0.24	0
5	FMT	A	802	-	2,2,2	0.73	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	746/748 (99%)	0.36	8 (1%) 78 51	70, 110, 148, 196	0
2	B	17/21 (80%)	0.32	0 100 100	81, 100, 136, 160	0
3	C	18/19 (94%)	0.46	0 100 100	88, 103, 140, 144	0
All	All	781/788 (99%)	0.37	8 (1%) 79 53	70, 109, 148, 196	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	207	ASN	2.6
1	A	1	ALA	2.6
1	A	404	LYS	2.4
1	A	262	LYS	2.4
1	A	140	GLN	2.3
1	A	206	GLY	2.2
1	A	328	ASN	2.2
1	A	573	PRO	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	FMT	B	103	3/3	0.46	0.15	119,119,119,120	0
5	FMT	A	803	3/3	0.62	0.10	125,125,126,126	0
5	FMT	A	802	3/3	0.70	0.16	100,100,101,101	0
5	FMT	B	102	3/3	0.83	0.12	99,99,99,99	0
5	FMT	B	101	3/3	0.85	0.08	113,113,113,113	0
4	MG	A	801	1/1	0.97	0.05	97,97,97,97	0

6.5 Other polymers [i](#)

There are no such residues in this entry.