



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 23, 2024 – 04:09 PM JST

PDB ID : 8Y06
Title : Crystal structure of LbCas12a in complex with crRNA and 12nt target DNA
Authors : Lin, X.; Chen, J.; Liu, L.
Deposited on : 2024-01-22
Resolution : 3.99 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	:	4.02b-467
Xtriage (Phenix)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

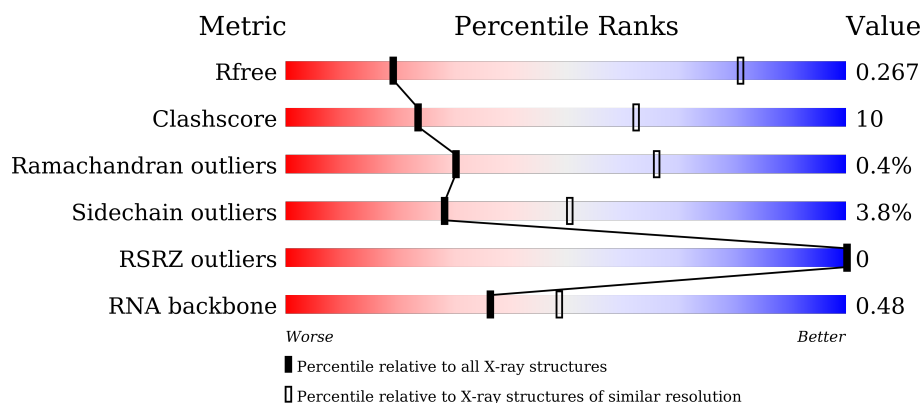
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.99 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}	164625	1028 (4.22-3.78)
Clashscore	180529	1055 (4.20-3.80)
Ramachandran outliers	177936	1004 (4.20-3.80)
Sidechain outliers	177891	1027 (4.22-3.78)
RSRZ outliers	164620	1029 (4.22-3.78)
RNA backbone	3690	1144 (5.00-3.00)

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	1228	<div> <div></div> <div>74%</div> <div>23%</div> <div>..</div> </div>
2	B	40	<div> <div>38%</div> <div>28%</div> <div>15%</div> <div>5%</div> <div>15%</div> </div>
3	C	21	<div> <div>43%</div> <div>52%</div> <div>5%</div> </div>
4	D	11	<div> <div>64%</div> <div>36%</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 11215 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 LbCas12a.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1202	Total	C	N	O	S	0	0	0
			9841	6333	1608	1871	29			

- Molecule 2 is a RNA chain called RNA (34-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	34	Total	C	N	O	P	0	0	0
			719	324	128	234	33			

- Molecule 3 is a DNA chain called DNA (5'-D(P*TP*TP*AP*CP*TP*GP*GP*AP*TP*GP*CP*GP*TP*AP*AP*AP*GP*GP*AP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	21	Total	C	N	O	P	0	0	0
			437	207	84	125	21			

- Molecule 4 is a DNA chain called DNA (5'-D(*CP*GP*TP*CP*CP*TP*TP*TP*AP*TP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	11	Total	C	N	O	P	0	0	0
			217	107	31	69	10			

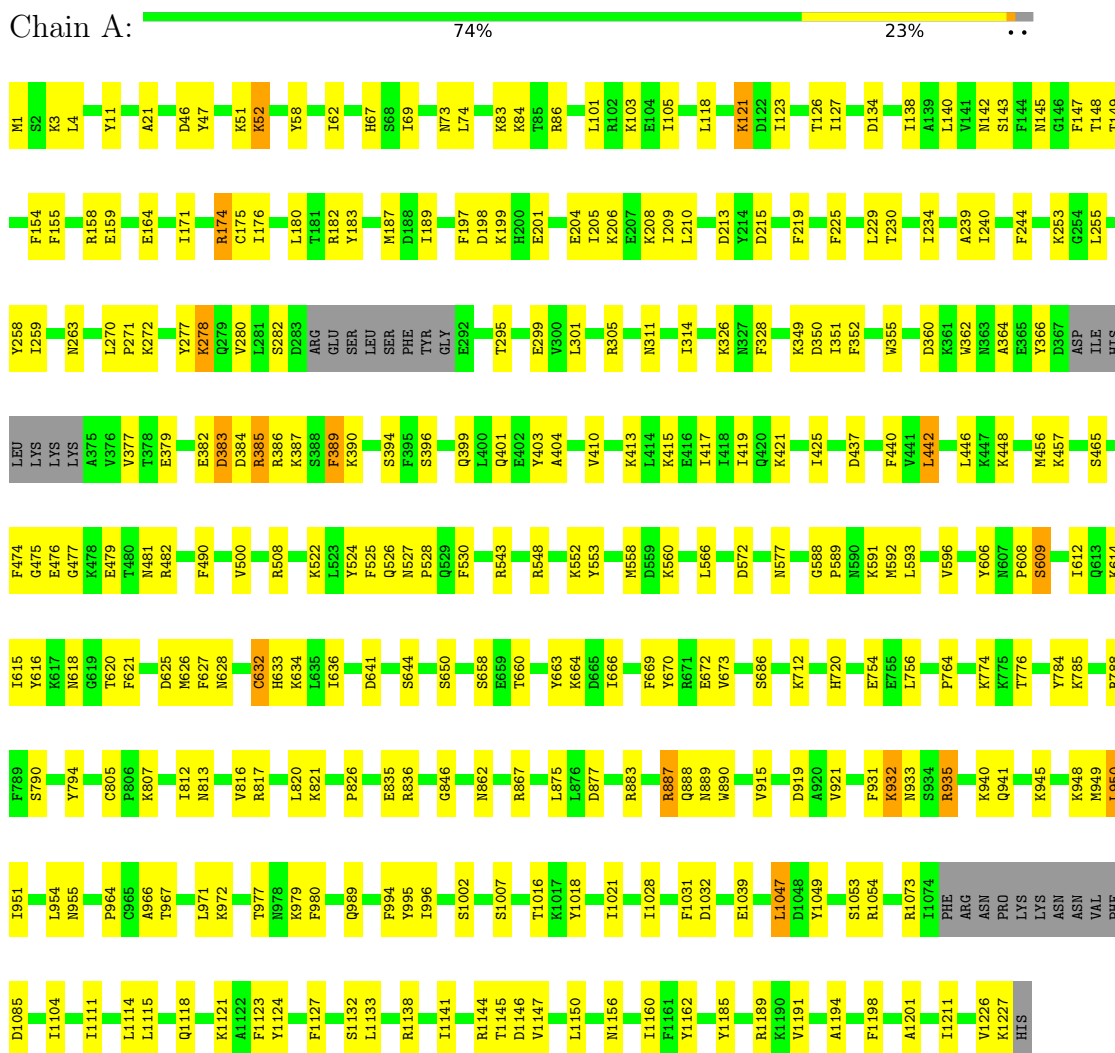
- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

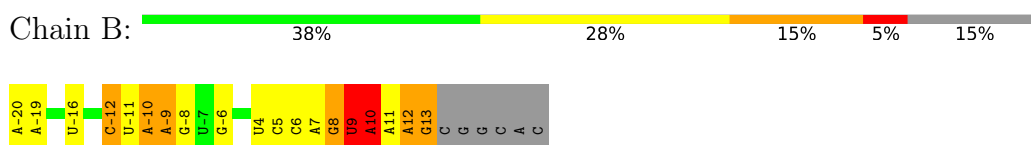
3 Residue-property plots

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: LbCas12a

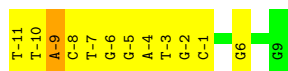


• Molecule 2: RNA (34-MER)



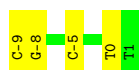
- Molecule 3: DNA (5'-D(P*TP*TP*AP*CP*TP*GP*GP*AP*TP*GP*CP*GP*TP*AP*AP*AP*GP*GP*AP*CP*G)-3')

Chain C:  43% 52% 5%



- Molecule 4: DNA (5'-D(*CP*GP*TP*CP*CP*TP*TP*TP*AP*TP*T)-3')

Chain D:  64% 36%



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	155.86Å 155.86Å 212.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.60 – 3.99 49.60 – 3.99	Depositor EDS
% Data completeness (in resolution range)	94.9 (49.60-3.99) 95.1 (49.60-3.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.02 (at 4.00Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.246 , 0.272 0.246 , 0.267	Depositor DCC
R_{free} test set	1241 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	2.3	Xtriage
Anisotropy	1.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.22 , 0.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	0.148 for -h,-k,l	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	11215	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.28	0/10046	0.45	0/13514
2	B	0.32	0/804	0.96	2/1250 (0.2%)
3	C	0.65	0/491	0.98	1/757 (0.1%)
4	D	0.61	0/240	1.13	0/368
All	All	0.32	0/11581	0.56	3/15889 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	-9	DA	O4'-C1'-N9	6.27	112.39	108.00
2	B	9	U	C2-N1-C1'	6.26	125.21	117.70
2	B	10	A	O5'-P-OP1	5.62	117.44	110.70

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	9841	0	9681	190	0
2	B	719	0	366	19	0
3	C	437	0	237	10	0
4	D	217	0	129	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	1	0	0	0	0
All	All	11215	0	10413	212	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 10.

All (212) 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)
3:C:-11:DT:H2''	3:C:-10:DT:H5''	1.43	1.01
1:A:305:ARG:NH2	1:A:437:ASP:O	2.07	0.88
1:A:835:GLU:HG2	1:A:935:ARG:HB2	1.53	0.86
1:A:774:LYS:NZ	2:B:-8:G:O6	2.10	0.85
1:A:263:ASN:HD21	1:A:270:LEU:HB2	1.42	0.84
1:A:377:VAL:HG13	1:A:385:ARG:HH12	1.45	0.80
3:C:-4:DA:H2''	3:C:-3:DT:H5''	1.63	0.79
1:A:351:ILE:HD12	1:A:352:PHE:H	1.48	0.78
1:A:633:HIS:HD2	1:A:660:THR:HG22	1.49	0.78
1:A:390:LYS:O	1:A:390:LYS:NZ	2.16	0.77
1:A:616:TYR:OH	4:D:0:DT:OP1	2.03	0.77
1:A:1016:THR:HG1	1:A:1132:SER:HG	1.26	0.76
1:A:887:ARG:NH1	1:A:889:ASN:OD1	2.19	0.75
1:A:74:LEU:HD21	1:A:225:PHE:HB3	1.70	0.74
1:A:349:LYS:HB2	1:A:355:TRP:HD1	1.55	0.72
1:A:996:ILE:HD11	1:A:1191:VAL:HG23	1.71	0.72
1:A:941:GLN:NE2	3:C:-8:DC:OP1	2.23	0.71
1:A:121:LYS:HA	1:A:148:THR:HG21	1.71	0.71
1:A:932:LYS:HE2	1:A:940:LYS:HD2	1.73	0.71
1:A:821:LYS:NZ	1:A:1201:ALA:O	2.24	0.70
1:A:142:ASN:O	1:A:145:ASN:ND2	2.24	0.70
1:A:210:LEU:HD21	1:A:240:ILE:HD11	1.74	0.70
2:B:10:A:H4'	2:B:10:A:OP1	1.91	0.69
3:C:-6:DG:H1'	3:C:-5:DG:H5''	1.72	0.69
1:A:349:LYS:HB2	1:A:355:TRP:CD1	2.29	0.68
1:A:83:LYS:HG2	1:A:86:ARG:HG2	1.76	0.68
1:A:1115:LEU:HD12	1:A:1123:PHE:HZ	1.60	0.67
1:A:389:PHE:H	1:A:389:PHE:HD1	1.43	0.67
1:A:525:PHE:O	1:A:543:ARG:NH2	2.28	0.66
1:A:835:GLU:OE2	1:A:1138:ARG:NH1	2.29	0.65
1:A:305:ARG:HH12	1:A:440:PHE:HB3	1.62	0.65
1:A:591:LYS:HG2	4:D:0:DT:H1'	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1185:TYR:CZ	1:A:1189:ARG:HD2	2.34	0.62
3:C:-5:DG:H2''	3:C:-4:DA:H5''	1.81	0.61
1:A:350:ASP:HB3	1:A:417:ILE:HD13	1.83	0.61
1:A:401:GLN:HG2	1:A:410:VAL:HB	1.82	0.60
1:A:785:LYS:HB2	2:B:-20:A:H5''	1.83	0.60
1:A:836:ARG:NH2	1:A:1147:VAL:O	2.35	0.60
1:A:836:ARG:NH2	1:A:1144:ARG:O	2.35	0.59
1:A:377:VAL:O	1:A:1053:SER:HB2	2.03	0.59
1:A:477:GLY:HA3	1:A:479:GLU:HB2	1.84	0.58
1:A:206:LYS:HA	1:A:210:LEU:HB2	1.86	0.57
1:A:383:ASP:O	1:A:387:LYS:N	2.36	0.57
1:A:171:ILE:HG12	1:A:280:VAL:HG21	1.87	0.57
1:A:263:ASN:OD1	1:A:270:LEU:N	2.30	0.57
1:A:205:ILE:HG23	1:A:209:ILE:HD12	1.85	0.56
2:B:10:A:H3'	2:B:11:A:C8	2.40	0.56
1:A:508:ARG:HG3	1:A:890:TRP:CD1	2.41	0.56
1:A:1007:SER:O	1:A:1226:VAL:HG21	2.06	0.56
1:A:588:GLY:HA3	1:A:591:LYS:HD2	1.87	0.56
1:A:278:LYS:NZ	1:A:282:SER:O	2.35	0.55
1:A:140:LEU:O	1:A:143:SER:OG	2.20	0.55
2:B:4:U:H2'	2:B:5:C:C6	2.41	0.55
1:A:572:ASP:O	1:A:577:ASN:ND2	2.33	0.54
1:A:1194:ALA:HB1	1:A:1211:ILE:HD12	1.88	0.54
1:A:625:ASP:OD1	1:A:625:ASP:N	2.39	0.54
1:A:351:ILE:HD12	1:A:352:PHE:N	2.21	0.54
1:A:189:ILE:HG23	1:A:271:PRO:HG2	1.89	0.54
1:A:47:TYR:CZ	1:A:51:LYS:HE3	2.43	0.53
1:A:362:TRP:HB2	1:A:403:TYR:OH	2.09	0.53
1:A:11:TYR:OH	1:A:977:THR:O	2.21	0.53
1:A:149:THR:HB	1:A:528:PRO:HB2	1.91	0.53
1:A:474:PHE:HE1	1:A:490:PHE:HE2	1.58	0.52
2:B:-20:A:O2'	2:B:-11:U:O4	2.22	0.52
2:B:10:A:H5'	2:B:11:A:OP2	2.10	0.52
1:A:351:ILE:HG22	1:A:413:LYS:HG3	1.91	0.52
1:A:1031:PHE:CE2	1:A:1047:LEU:HD23	2.45	0.52
1:A:46:ASP:HB3	1:A:140:LEU:HD11	1.91	0.52
1:A:548:ARG:HD3	1:A:553:TYR:CE2	2.44	0.52
1:A:3:LYS:NZ	1:A:919:ASP:O	2.33	0.51
1:A:641:ASP:O	1:A:644:SER:OG	2.22	0.51
1:A:263:ASN:CG	1:A:270:LEU:H	2.13	0.51
1:A:1198:PHE:CZ	1:A:1211:ILE:HD11	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:887:ARG:CG	1:A:888:GLN:H	2.24	0.51
1:A:210:LEU:HD11	1:A:219:PHE:CZ	2.46	0.51
1:A:360:ASP:O	1:A:364:ALA:N	2.41	0.51
1:A:812:ILE:O	1:A:816:VAL:HG23	2.11	0.51
1:A:158:ARG:HH12	2:B:5:C:H5'	1.76	0.51
1:A:826:PRO:HD2	1:A:846:GLY:HA3	1.93	0.51
1:A:967:THR:HA	1:A:972:LYS:HB3	1.92	0.51
1:A:415:LYS:HG2	1:A:419:ILE:HD12	1.93	0.50
2:B:9:U:H2'	2:B:10:A:O4'	2.11	0.50
1:A:255:LEU:O	1:A:259:ILE:HG13	2.12	0.50
1:A:836:ARG:HE	1:A:1145:THR:HA	1.77	0.50
2:B:-20:A:O2'	2:B:-19:A:OP2	2.30	0.50
1:A:663:TYR:OH	1:A:672:GLU:OE2	2.21	0.50
1:A:240:ILE:HG22	1:A:255:LEU:HG	1.93	0.50
1:A:1141:ILE:HB	1:A:1144:ARG:HB2	1.93	0.49
1:A:385:ARG:O	1:A:387:LYS:N	2.45	0.49
1:A:351:ILE:HG22	1:A:413:LYS:CG	2.43	0.49
1:A:955:ASN:HD21	1:A:979:LYS:HA	1.76	0.49
4:D:-9:DC:H2'	4:D:-8:DG:C8	2.47	0.49
1:A:1039:GLU:OE2	1:A:1039:GLU:N	2.37	0.49
1:A:174:ARG:HH21	1:A:277:TYR:HB3	1.77	0.49
1:A:932:LYS:HD2	1:A:932:LYS:N	2.27	0.49
1:A:326:LYS:HD3	1:A:419:ILE:HD11	1.95	0.49
2:B:12:A:O2'	2:B:13:G:OP2	2.29	0.49
3:C:-7:DT:H4'	3:C:-6:DG:OP1	2.11	0.49
1:A:862:ASN:ND2	1:A:867:ARG:HG2	2.28	0.48
1:A:658:SER:OG	1:A:672:GLU:OE1	2.31	0.48
1:A:835:GLU:HG2	1:A:935:ARG:CB	2.36	0.48
1:A:4:LEU:HD11	1:A:915:VAL:HG12	1.96	0.48
1:A:394:SER:HB3	1:A:475:GLY:HA2	1.96	0.48
1:A:620:THR:O	1:A:627:PHE:HA	2.14	0.48
1:A:1115:LEU:HD12	1:A:1123:PHE:CZ	2.44	0.48
1:A:69:ILE:HD11	1:A:105:ILE:HA	1.96	0.48
1:A:314:ILE:HD13	1:A:500:VAL:HG12	1.96	0.48
2:B:9:U:H3'	2:B:10:A:C5'	2.43	0.48
1:A:1104:ILE:HD11	1:A:1118:GLN:HG3	1.96	0.48
1:A:592:MET:O	1:A:596:VAL:HG23	2.14	0.47
1:A:1111:ILE:O	1:A:1114:LEU:N	2.45	0.47
1:A:51:LYS:HE2	1:A:154:PHE:CD1	2.49	0.47
1:A:615:ILE:HD11	1:A:634:LYS:NZ	2.30	0.47
1:A:21:ALA:HB3	1:A:794:TYR:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:LEU:O	1:A:305:ARG:HG3	2.13	0.47
1:A:366:TYR:CE2	1:A:385:ARG:HG3	2.48	0.47
1:A:508:ARG:HG3	1:A:890:TRP:CG	2.50	0.47
1:A:1018:TYR:HE1	1:A:1124:TYR:HB2	1.79	0.47
1:A:589:PRO:O	1:A:593:LEU:HB2	2.14	0.47
1:A:615:ILE:HD11	1:A:634:LYS:HZ2	1.80	0.47
1:A:950:LEU:O	1:A:954:LEU:HD13	2.15	0.47
1:A:457:LYS:HE3	1:A:888:GLN:O	2.15	0.47
1:A:474:PHE:CE1	1:A:490:PHE:HE2	2.33	0.46
1:A:875:LEU:HD21	2:B:-12:C:OP1	2.15	0.46
1:A:362:TRP:CH2	1:A:389:PHE:HA	2.50	0.46
1:A:1021:ILE:HD11	1:A:1121:LYS:HB3	1.98	0.46
1:A:176:ILE:O	1:A:180:LEU:HB3	2.15	0.46
1:A:527:ASN:HB3	1:A:530:PHE:HB2	1.98	0.46
3:C:6:DG:N2	4:D:-5:DC:O2	2.49	0.46
1:A:836:ARG:NE	1:A:1145:THR:HA	2.31	0.46
1:A:377:VAL:HG11	1:A:1054:ARG:HH21	1.80	0.46
1:A:820:LEU:HD21	1:A:921:VAL:HG21	1.98	0.46
1:A:175:CYS:HB3	1:A:229:LEU:HD22	1.97	0.45
1:A:379:GLU:HG2	1:A:1054:ARG:HG2	1.98	0.45
1:A:305:ARG:HH22	1:A:440:PHE:HB3	1.81	0.45
1:A:103:LYS:HG2	1:A:164:GLU:OE1	2.17	0.45
1:A:183:TYR:O	1:A:187:MET:HG3	2.17	0.45
1:A:204:GLU:HG2	1:A:208:LYS:HG3	1.97	0.45
1:A:239:ALA:HB1	1:A:244:PHE:HD2	1.82	0.45
1:A:966:ALA:O	1:A:972:LYS:HB2	2.17	0.45
1:A:666:ILE:HG13	1:A:670:TYR:CE2	2.52	0.44
1:A:989:GLN:HB2	1:A:994:PHE:CD2	2.53	0.44
1:A:199:LYS:H	1:A:199:LYS:HG2	1.53	0.44
1:A:225:PHE:CE2	1:A:229:LEU:HD21	2.51	0.44
1:A:764:PRO:HB2	1:A:776:THR:CG2	2.47	0.44
1:A:67:HIS:CE1	1:A:230:THR:HG21	2.53	0.44
1:A:933:ASN:ND2	1:A:933:ASN:H	2.16	0.44
1:A:1028:ILE:HG12	1:A:1127:PHE:CE2	2.52	0.44
1:A:669:PHE:O	1:A:673:VAL:HG23	2.17	0.44
2:B:8:G:O2'	2:B:9:U:O4'	2.27	0.44
1:A:118:LEU:HA	1:A:123:ILE:HD13	2.00	0.44
1:A:813:ASN:O	1:A:817:ARG:HG3	2.18	0.44
1:A:58:TYR:HE1	1:A:118:LEU:HB3	1.82	0.43
1:A:479:GLU:HB3	1:A:482:ARG:HD2	2.00	0.43
1:A:756:LEU:HD11	1:A:790:SER:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:PHE:O	1:A:159:GLU:HG3	2.18	0.43
1:A:788:ARG:HD2	2:B:-16:U:OP2	2.18	0.43
1:A:945:LYS:O	1:A:949:MET:HG3	2.19	0.43
3:C:-2:DG:H2'	3:C:-1:DC:C6	2.54	0.43
1:A:389:PHE:CD1	1:A:389:PHE:N	2.85	0.43
1:A:566:LEU:HD23	1:A:566:LEU:HA	1.86	0.43
1:A:134:ASP:O	1:A:138:ILE:HG13	2.18	0.43
1:A:396:SER:OG	1:A:399:GLN:HG3	2.19	0.43
1:A:543:ARG:HG2	1:A:558:MET:HB2	2.01	0.43
1:A:278:LYS:HG2	2:B:7:A:OP1	2.19	0.42
1:A:1146:ASP:OD1	1:A:1146:ASP:N	2.44	0.42
2:B:-10:A:H5'	2:B:-9:A:C2	2.54	0.42
1:A:382:GLU:O	1:A:385:ARG:HB3	2.20	0.42
1:A:413:LYS:HZ3	1:A:417:ILE:HD11	1.84	0.42
1:A:614:LYS:O	1:A:618:ASN:N	2.33	0.42
1:A:636:ILE:HG23	1:A:669:PHE:CG	2.54	0.42
2:B:6:C:H2'	2:B:7:A:O4'	2.20	0.42
1:A:805:CYS:O	1:A:807:LYS:HE2	2.20	0.42
1:A:197:PHE:CZ	1:A:255:LEU:HB3	2.54	0.42
1:A:784:TYR:O	2:B:-19:A:H5'	2.19	0.42
1:A:377:VAL:CG1	1:A:385:ARG:HH12	2.26	0.42
1:A:526:GLN:HB2	1:A:560:LYS:HD2	2.02	0.42
1:A:1150:LEU:HD22	1:A:1162:TYR:HE2	1.83	0.42
1:A:1211:ILE:HD13	1:A:1211:ILE:HA	1.74	0.42
3:C:-9:DA:H1'	3:C:-8:DC:O5'	2.20	0.42
1:A:609:SER:OG	1:A:612:ILE:HG13	2.19	0.42
1:A:1039:GLU:H	1:A:1039:GLU:CD	2.18	0.41
1:A:52:LYS:N	1:A:52:LYS:HD2	2.34	0.41
1:A:201:GLU:HB3	1:A:258:TYR:CE2	2.55	0.41
1:A:572:ASP:HB2	1:A:686:SER:HB2	2.01	0.41
1:A:754:GLU:H	1:A:754:GLU:CD	2.23	0.41
3:C:-9:DA:OP2	3:C:-9:DA:H8	2.03	0.41
1:A:404:ALA:HB2	1:A:410:VAL:HG23	2.02	0.41
1:A:875:LEU:HD23	1:A:875:LEU:HA	1.84	0.41
1:A:126:THR:OG1	1:A:127:ILE:HD12	2.20	0.41
1:A:142:ASN:HA	1:A:145:ASN:HD21	1.85	0.41
1:A:421:LYS:O	1:A:425:ILE:HG13	2.21	0.41
1:A:948:LYS:HG3	1:A:980:PHE:CD1	2.56	0.41
1:A:522:LYS:HD3	1:A:524:TYR:CZ	2.56	0.41
1:A:606:TYR:O	1:A:608:PRO:HD3	2.21	0.41
1:A:384:ASP:N	1:A:384:ASP:OD1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:TYR:O	1:A:62:ILE:HG13	2.21	0.41
1:A:101:LEU:HB3	1:A:176:ILE:HD12	2.03	0.41
1:A:198:ASP:OD1	1:A:199:LYS:N	2.54	0.41
1:A:1156:ASN:CG	1:A:1160:ILE:HG22	2.41	0.40
1:A:1160:ILE:HA	1:A:1160:ILE:HD12	1.83	0.40
1:A:263:ASN:ND2	1:A:270:LEU:HB2	2.21	0.40
1:A:442:LEU:HD11	1:A:446:LEU:CD1	2.51	0.40
1:A:820:LEU:HD23	1:A:820:LEU:HA	1.89	0.40
1:A:230:THR:O	1:A:234:ILE:HG13	2.21	0.40
1:A:712:LYS:HA	1:A:712:LYS:HD3	1.87	0.40
1:A:964:PRO:O	1:A:971:LEU:HB2	2.21	0.40
1:A:1073:ARG:HG3	1:A:1133:LEU:HD21	2.02	0.40
1:A:127:ILE:HD12	1:A:127:ILE:H	1.86	0.40
1:A:951:ILE:HD11	1:A:995:TYR:OH	2.21	0.40
1:A:295:THR:HG22	1:A:299:GLU:OE1	2.22	0.40
1:A:621:PHE:HA	1:A:632:CYS:HB2	2.03	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	1194/1228 (97%)	1153 (97%)	36 (3%)	5 (0%)	30 66

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	385	ARG
1	A	386	ARG
1	A	481	ASN
1	A	476	GLU
1	A	389	PHE

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	1068/1115 (96%)	1027 (96%)	41 (4%)	28 51

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	52	LYS
1	A	73	ASN
1	A	84	LYS
1	A	121	LYS
1	A	147	PHE
1	A	174	ARG
1	A	182	ARG
1	A	213	ASP
1	A	215	ASP
1	A	253	LYS
1	A	272	LYS
1	A	278	LYS
1	A	311	ASN
1	A	328	PHE
1	A	383	ASP
1	A	442	LEU
1	A	448	LYS
1	A	456	MET
1	A	465	SER
1	A	552	LYS
1	A	609	SER
1	A	626	MET
1	A	628	ASN
1	A	632	CYS
1	A	650	SER
1	A	664	LYS
1	A	720	HIS
1	A	877	ASP
1	A	883	ARG

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Mol	Chain	Res	Type
1	A	887	ARG
1	A	931	PHE
1	A	932	LYS
1	A	935	ARG
1	A	950	LEU
1	A	1002	SER
1	A	1032	ASP
1	A	1047	LEU
1	A	1049	TYR
1	A	1085	ASP
1	A	1227	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	468	ASN
1	A	862	ASN
1	A	888	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	33/40 (82%)	9 (27%)	1 (3%)

All (9) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	-12	C
2	B	-10	A
2	B	-9	A
2	B	-6	G
2	B	8	G
2	B	9	U
2	B	10	A
2	B	12	A
2	B	13	G

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	8	G

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 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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	1202/1228 (97%)	-1.21	0 100 100	3, 38, 157, 195	0
2	B	34/40 (85%)	-1.31	0 100 100	11, 27, 187, 206	0
3	C	21/21 (100%)	-1.16	0 100 100	11, 35, 176, 186	0
4	D	11/11 (100%)	-1.51	0 100 100	15, 23, 41, 42	0
All	All	1268/1300 (97%)	-1.21	0 100 100	3, 38, 158, 206	0

There are no RSRZ outliers to report.

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, 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	MG	A	1301	1/1	0.99	0.12	1,1,1,1	0

6.5 Other polymers [i](#)

There are no such residues in this entry.