



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:33 PM GMT

PDB ID : 4CIU
Title : Crystal structure of E. coli ClpB
Authors : Kopp, J.; Sinning, I.; Bukau, B.; Kummer, E.; Mogk, A.
Deposited on : 2013-12-16
Resolution : 3.50 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

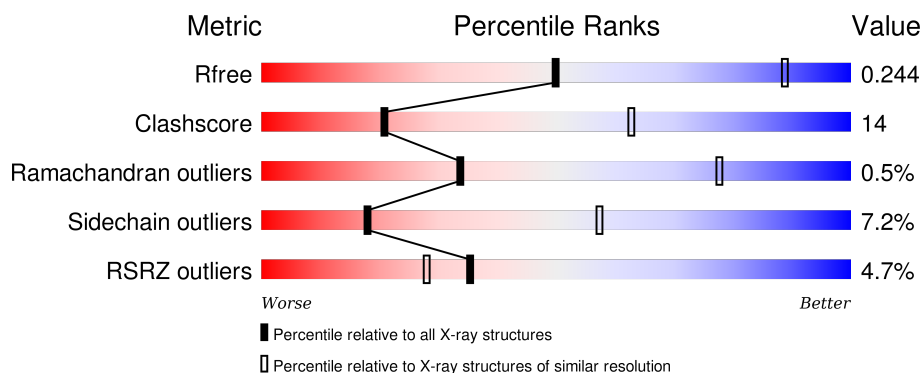
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	727	<div> <div>4%</div> <div>66%</div> <div>21%</div> <div>•</div> <div>10%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5257 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.

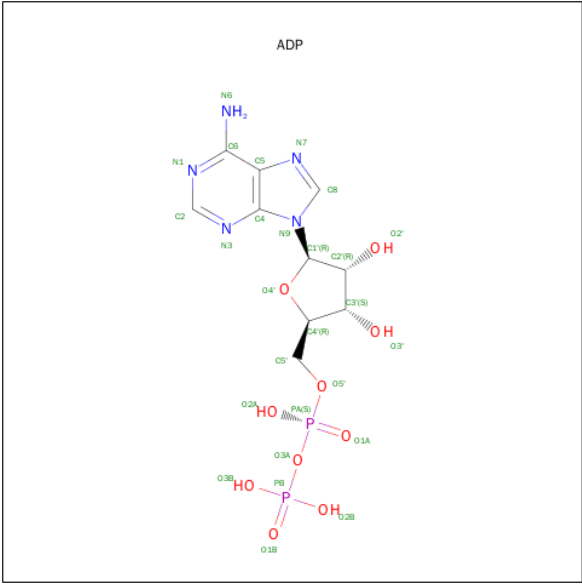
- Molecule 1 is a protein called CHAPERONE PROTEIN CLPB.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	657	5203	3260	939	985	2	17	0	0	0

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	858	HIS	-	EXPRESSION TAG	UNP P63284
A	859	SER	-	EXPRESSION TAG	UNP P63284
A	860	MSE	-	EXPRESSION TAG	UNP P63284
A	861	ALA	-	EXPRESSION TAG	UNP P63284
A	862	ARG	-	EXPRESSION TAG	UNP P63284
A	863	SER	-	EXPRESSION TAG	UNP P63284
A	864	HIS	-	EXPRESSION TAG	UNP P63284
A	865	HIS	-	EXPRESSION TAG	UNP P63284
A	866	HIS	-	EXPRESSION TAG	UNP P63284
A	867	HIS	-	EXPRESSION TAG	UNP P63284
A	868	HIS	-	EXPRESSION TAG	UNP P63284
A	869	HIS	-	EXPRESSION TAG	UNP P63284
A	279	ALA	GLU	ENGINEERED MUTATION	UNP P63284
A	432	ALA	GLU	ENGINEERED MUTATION	UNP P63284
A	678	ALA	GLU	ENGINEERED MUTATION	UNP P63284

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

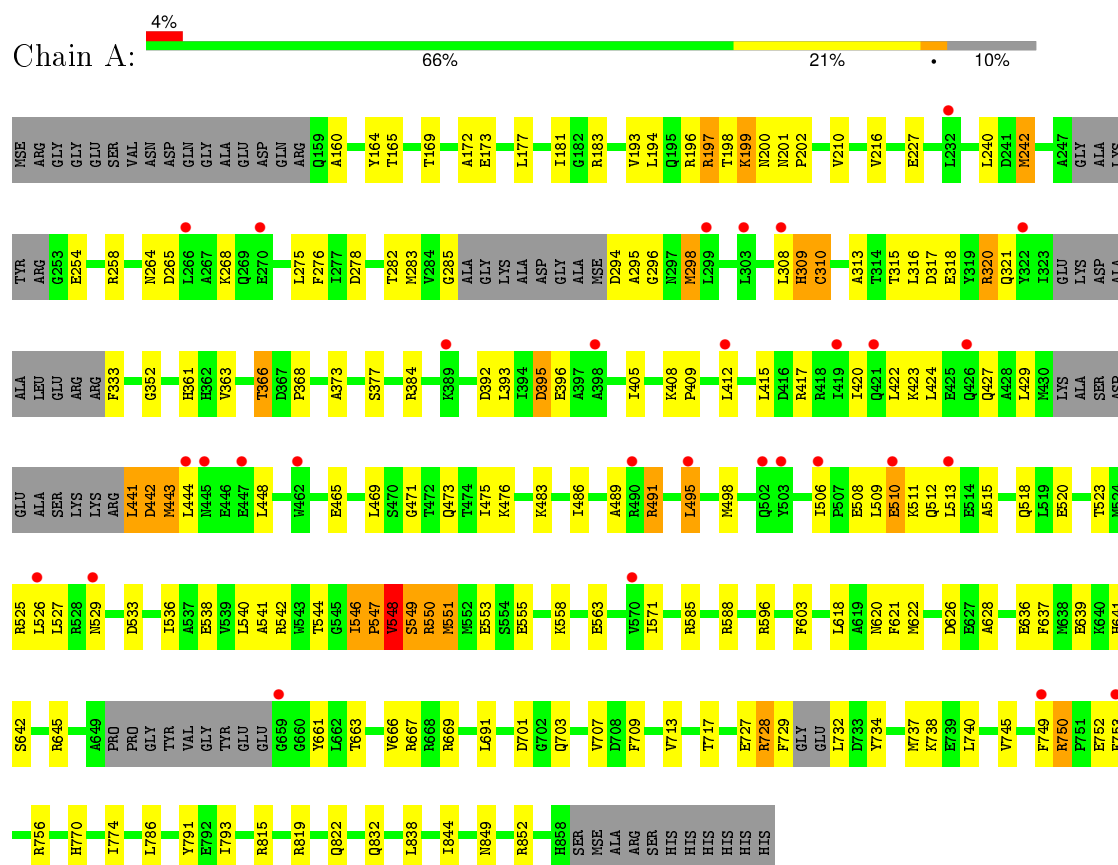


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). 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: CHAPERONE PROTEIN CLPB



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	127.33Å 127.33Å 119.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.80 – 3.50 81.15 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (36.80-3.50) 100.0 (81.15-3.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.28 (at 3.49Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.224 , 0.247 0.223 , 0.244	Depositor DCC
R_{free} test set	694 reflections (4.97%)	DCC
Wilson B-factor (Å ²)	121.0	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 69.7	EDS
Estimated twinning fraction	0.056 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 13992 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5257	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: ADP

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	0.26	1/5250 (0.0%)	0.47	1/7040 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	547	PRO	N-CD	5.15	1.55	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	546	ILE	C-N-CD	5.68	140.33	128.40

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	5203	0	5301	152	0
2	A	54	0	24	2	0
All	All	5257	0	5325	152	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 14.

All (152) 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:548:VAL:HG23	1:A:551:MSE:SE	1.96	1.14
1:A:198:THR:HG22	1:A:199:LYS:HD2	1.20	1.12
1:A:544:THR:HB	1:A:546:ILE:CD1	1.84	1.07
1:A:728:ARG:HD2	1:A:732:LEU:HD13	1.40	1.03
1:A:728:ARG:HD2	1:A:732:LEU:CD1	1.88	1.02
1:A:544:THR:OG1	1:A:546:ILE:HD13	1.65	0.97
1:A:198:THR:HG22	1:A:199:LYS:CD	1.94	0.95
1:A:196:ARG:HB3	1:A:200:ASN:HA	1.48	0.94
1:A:200:ASN:OD1	1:A:201:ASN:N	1.99	0.94
1:A:728:ARG:CD	1:A:732:LEU:HD13	1.98	0.93
1:A:544:THR:CB	1:A:546:ILE:HD13	1.98	0.92
1:A:728:ARG:HD2	1:A:732:LEU:HB2	1.50	0.92
1:A:728:ARG:HD2	1:A:732:LEU:CB	2.03	0.87
1:A:544:THR:CB	1:A:546:ILE:CD1	2.51	0.86
1:A:201:ASN:HB2	1:A:310:CYS:O	1.77	0.84
1:A:193:VAL:HG12	1:A:200:ASN:O	1.81	0.79
1:A:548:VAL:HA	1:A:551:MSE:HB2	1.68	0.76
1:A:199:LYS:N	1:A:199:LYS:HD2	2.01	0.75
1:A:366:THR:HG22	1:A:368:PRO:HD2	1.70	0.74
1:A:548:VAL:CG2	1:A:551:MSE:SE	2.82	0.73
1:A:729:PHE:CE2	1:A:734:TYR:HE1	2.07	0.72
1:A:197:ARG:HD3	1:A:197:ARG:H	1.54	0.72
1:A:637:PHE:HB3	1:A:642:SER:HB3	1.72	0.71
1:A:544:THR:HB	1:A:546:ILE:HD11	1.72	0.71
1:A:728:ARG:CD	1:A:732:LEU:HB2	2.20	0.70
1:A:396:GLU:OE2	1:A:542:ARG:NH2	2.25	0.69
1:A:546:ILE:N	1:A:546:ILE:HD12	2.08	0.69
1:A:196:ARG:CB	1:A:200:ASN:HA	2.22	0.69
1:A:199:LYS:H	1:A:199:LYS:HZ2	1.41	0.68
1:A:728:ARG:CB	1:A:737:MSE:SE	2.93	0.66
1:A:728:ARG:CD	1:A:732:LEU:CD1	2.66	0.66
1:A:728:ARG:HG3	1:A:737:MSE:SE	2.46	0.66
1:A:729:PHE:CZ	1:A:734:TYR:HE1	2.13	0.66
1:A:240:LEU:HB2	1:A:275:LEU:HD11	1.77	0.65
1:A:546:ILE:H	1:A:546:ILE:HD12	1.60	0.65
1:A:729:PHE:CE2	1:A:734:TYR:CE1	2.84	0.65
1:A:540:LEU:HG	1:A:551:MSE:HE1	1.78	0.65
1:A:405:ILE:HG23	1:A:527:LEU:HD23	1.81	0.63
1:A:315:THR:HB	1:A:318:GLU:OE1	1.98	0.62
1:A:628:ALA:HA	1:A:669:ARG:HD3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:596:ARG:NH2	1:A:756:ARG:O	2.33	0.61
1:A:201:ASN:OD1	1:A:333:PHE:CD1	2.54	0.61
1:A:728:ARG:HD2	1:A:732:LEU:CG	2.31	0.60
1:A:636:GLU:O	1:A:645:ARG:NH2	2.35	0.60
1:A:546:ILE:HG22	1:A:548:VAL:H	1.67	0.59
1:A:476:LYS:HG3	1:A:513:LEU:HD11	1.85	0.59
1:A:547:PRO:O	1:A:548:VAL:HG12	2.02	0.59
1:A:194:LEU:HD23	1:A:309:HIS:CE1	2.38	0.59
1:A:728:ARG:CG	1:A:732:LEU:HD13	2.33	0.58
1:A:745:VAL:HG13	1:A:749:PHE:HD2	1.70	0.57
1:A:547:PRO:HD2	1:A:550:ARG:HG3	1.87	0.57
1:A:393:LEU:HD11	1:A:540:LEU:HD22	1.86	0.57
1:A:193:VAL:O	1:A:196:ARG:HB2	2.06	0.56
1:A:384:ARG:NH2	1:A:392:ASP:OD2	2.40	0.55
1:A:198:THR:HG22	1:A:199:LYS:NZ	2.22	0.55
1:A:732:LEU:HD22	1:A:737:MSE:HG2	1.88	0.54
1:A:173:GLU:HG3	1:A:491:ARG:HH12	1.72	0.54
1:A:165:THR:HG22	1:A:240:LEU:HA	1.90	0.54
1:A:172:ALA:HB2	1:A:177:LEU:HD12	1.88	0.54
1:A:242:MSE:H	1:A:242:MSE:SE	2.40	0.54
1:A:663:THR:HG21	1:A:701:ASP:HB3	1.89	0.54
1:A:489:ALA:HB1	1:A:498:MSE:HG2	1.89	0.53
1:A:546:ILE:H	1:A:546:ILE:CD1	2.21	0.53
1:A:282:THR:O	1:A:294:ASP:N	2.41	0.53
1:A:555:GLU:OE2	1:A:588:ARG:NH1	2.43	0.52
1:A:728:ARG:CG	1:A:737:MSE:SE	3.07	0.52
1:A:540:LEU:CG	1:A:551:MSE:HE1	2.38	0.52
1:A:198:THR:CG2	1:A:199:LYS:NZ	2.73	0.52
1:A:508:GLU:HG3	1:A:509:LEU:N	2.25	0.52
1:A:786:LEU:HB3	1:A:791:TYR:HB2	1.91	0.51
1:A:199:LYS:N	1:A:199:LYS:CD	2.73	0.51
1:A:838:LEU:HD23	1:A:844:ILE:HD13	1.92	0.51
1:A:563:GLU:OE2	1:A:585:ARG:NH2	2.35	0.51
1:A:728:ARG:HG3	1:A:732:LEU:HD13	1.91	0.51
1:A:603:PHE:HB2	1:A:717:THR:HG22	1.93	0.50
1:A:815:ARG:NH2	2:A:902:ADP:O3B	2.45	0.50
1:A:315:THR:HG22	1:A:317:ASP:H	1.77	0.50
1:A:285:GLY:H	1:A:295:ALA:HA	1.76	0.49
1:A:193:VAL:CG1	1:A:200:ASN:O	2.57	0.49
1:A:520:GLU:HA	1:A:523:THR:HG22	1.94	0.49
1:A:819:ARG:HA	1:A:822:GLN:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:422:LEU:HB3	1:A:444:LEU:HD11	1.95	0.49
1:A:621:PHE:HD2	1:A:622:MSE:HG3	1.77	0.49
1:A:540:LEU:HB3	1:A:551:MSE:HE1	1.94	0.49
1:A:508:GLU:O	1:A:512:GLN:HG2	2.13	0.49
1:A:549:SER:O	1:A:553:GLU:HB3	2.13	0.49
1:A:515:ALA:HA	1:A:518:GLN:HG2	1.95	0.48
1:A:793:ILE:HG12	1:A:844:ILE:HB	1.94	0.48
1:A:622:MSE:HE1	1:A:713:VAL:HG21	1.95	0.48
1:A:546:ILE:N	1:A:546:ILE:CD1	2.76	0.48
1:A:544:THR:OG1	1:A:546:ILE:CD1	2.48	0.48
1:A:423:LYS:HG2	1:A:448:LEU:HD11	1.95	0.48
1:A:533:ASP:OD1	1:A:533:ASP:N	2.47	0.48
1:A:420:ILE:O	1:A:424:LEU:HG	2.13	0.48
1:A:750:ARG:HH11	1:A:750:ARG:HB2	1.78	0.48
1:A:639:GLU:N	1:A:642:SER:OG	2.40	0.47
1:A:198:THR:HG22	1:A:199:LYS:N	2.28	0.47
1:A:770:HIS:O	1:A:774:ILE:HG13	2.15	0.47
1:A:316:LEU:O	1:A:320:ARG:N	2.46	0.47
1:A:181:ILE:O	2:A:901:ADP:N6	2.46	0.47
1:A:471:GLY:O	1:A:475:ILE:HG22	2.15	0.47
1:A:160:ALA:HB1	1:A:164:TYR:HD2	1.80	0.46
1:A:728:ARG:HB3	1:A:737:MSE:SE	2.64	0.46
1:A:512:GLN:HA	1:A:515:ALA:HB3	1.97	0.46
1:A:283:MSE:O	1:A:296:GLY:N	2.38	0.46
1:A:483:LYS:O	1:A:486:ILE:HG22	2.15	0.46
1:A:506:ILE:O	1:A:510:GLU:HB2	2.16	0.46
1:A:752:GLU:O	1:A:756:ARG:HG2	2.16	0.45
1:A:526:LEU:HB2	1:A:527:LEU:HD12	1.99	0.45
1:A:377:SER:HB3	1:A:393:LEU:HD12	1.97	0.45
1:A:728:ARG:HA	1:A:728:ARG:HD3	1.59	0.45
1:A:832:GLN:HB3	1:A:838:LEU:HD13	1.99	0.45
1:A:417:ARG:O	1:A:420:ILE:HG22	2.17	0.45
1:A:637:PHE:HZ	1:A:661:TYR:HD2	1.65	0.44
1:A:361:HIS:NE2	1:A:395:ASP:OD1	2.51	0.44
1:A:278:ASP:HA	1:A:313:ALA:HB3	2.00	0.44
1:A:198:THR:CG2	1:A:199:LYS:HZ3	2.31	0.44
1:A:361:HIS:O	1:A:363:VAL:HG13	2.18	0.43
1:A:198:THR:O	1:A:199:LYS:HB2	2.19	0.43
1:A:183:ARG:NH2	1:A:210:VAL:O	2.51	0.43
1:A:558:LYS:HG2	1:A:621:PHE:CZ	2.53	0.43
1:A:738:LYS:HE2	1:A:738:LYS:HB3	1.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:ILE:HD11	1:A:529:ASN:HB3	2.01	0.43
1:A:728:ARG:HB2	1:A:737:MSE:SE	2.67	0.43
1:A:442:ASP:O	1:A:444:LEU:N	2.52	0.42
1:A:181:ILE:HG12	1:A:352:GLY:HA3	2.02	0.42
1:A:538:GLU:O	1:A:541:ALA:HB3	2.19	0.42
1:A:198:THR:CG2	1:A:199:LYS:HD2	2.14	0.42
1:A:422:LEU:HD22	1:A:444:LEU:HD11	2.02	0.42
1:A:429:LEU:HD13	1:A:441:LEU:HD12	2.02	0.42
1:A:465:GLU:OE2	1:A:525:ARG:N	2.51	0.42
1:A:508:GLU:O	1:A:511:LYS:HG2	2.20	0.41
1:A:442:ASP:HB3	1:A:443:MSE:H	1.63	0.41
1:A:196:ARG:HB3	1:A:200:ASN:CA	2.35	0.41
1:A:173:GLU:HG3	1:A:491:ARG:NH1	2.34	0.41
1:A:373:ALA:HA	1:A:536:ILE:HG21	2.02	0.41
1:A:495:LEU:HD22	1:A:495:LEU:HA	1.87	0.41
1:A:571:ILE:CG2	1:A:774:ILE:HG12	2.50	0.41
1:A:265:ASP:HA	1:A:268:LYS:HD2	2.03	0.41
1:A:409:PRO:HB2	1:A:412:LEU:HD13	2.02	0.41
1:A:618:LEU:HA	1:A:618:LEU:HD23	1.91	0.41
1:A:756:ARG:HA	1:A:756:ARG:HH11	1.85	0.41
1:A:666:VAL:HG21	1:A:709:PHE:CE2	2.55	0.41
1:A:540:LEU:CB	1:A:551:MSE:HE1	2.51	0.41
1:A:298:MSE:HE2	1:A:298:MSE:HB2	1.88	0.41
1:A:198:THR:HG22	1:A:199:LYS:CE	2.49	0.41
1:A:663:THR:HB	1:A:707:VAL:HG21	2.03	0.41
1:A:264:ASN:O	1:A:268:LYS:HG3	2.21	0.41
1:A:408:LYS:HA	1:A:409:PRO:HD3	1.94	0.41
1:A:791:TYR:CD1	1:A:844:ILE:HD11	2.56	0.41
1:A:216:VAL:HG21	1:A:276:PHE:CE1	2.56	0.40
1:A:164:TYR:OH	1:A:258:ARG:NE	2.55	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	643/727 (88%)	620 (96%)	20 (3%)	3 (0%)	34 78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	548	VAL
1	A	202	PRO
1	A	443	MSE

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	558/589 (95%)	518 (93%)	40 (7%)	18 57

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	THR
1	A	197	ARG
1	A	199	LYS
1	A	227	GLU
1	A	242	MSE
1	A	254	GLU
1	A	298	MSE
1	A	308	LEU
1	A	309	HIS
1	A	310	CYS
1	A	320	ARG
1	A	321	GLN
1	A	366	THR
1	A	395	ASP
1	A	415	LEU
1	A	427	GLN

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Mol	Chain	Res	Type
1	A	441	LEU
1	A	442	ASP
1	A	469	LEU
1	A	473	GLN
1	A	491	ARG
1	A	495	LEU
1	A	510	GLU
1	A	548	VAL
1	A	549	SER
1	A	550	ARG
1	A	551	MSE
1	A	620	ASN
1	A	626	ASP
1	A	641	HIS
1	A	667	ARG
1	A	691	LEU
1	A	703	GLN
1	A	727	GLU
1	A	728	ARG
1	A	740	LEU
1	A	750	ARG
1	A	753	PHE
1	A	849	ASN
1	A	852	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	272	ASN
1	A	736	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 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	ADP	A	901	-	22,29,29	0.99	1 (4%)	27,45,45	1.87	4 (14%)
2	ADP	A	902	-	22,29,29	1.00	1 (4%)	27,45,45	1.91	4 (14%)

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	ADP	A	901	-	-	0/12/32/32	0/3/3/3
2	ADP	A	902	-	-	0/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	ADP	C5-C4	3.10	1.47	1.40
2	A	902	ADP	C5-C4	3.14	1.47	1.40

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	ADP	N3-C2-N1	-6.93	123.59	128.89
2	A	902	ADP	N3-C2-N1	-6.81	123.68	128.89
2	A	902	ADP	PA-O3A-PB	-3.48	120.99	132.67
2	A	902	ADP	C2'-C1'-N9	-3.36	109.16	114.29
2	A	901	ADP	PA-O3A-PB	-3.30	121.59	132.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	ADP	C4-C5-N7	-3.02	106.70	109.48
2	A	902	ADP	C4-C5-N7	-2.95	106.77	109.48
2	A	901	ADP	C2'-C1'-N9	-2.08	111.11	114.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	ADP	1	0
2	A	902	ADP	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	640/727 (88%)	0.30	30 (4%) 35 28	37, 92, 164, 203	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	502	GLN	4.1
1	A	270	GLU	3.7
1	A	426	GLN	3.7
1	A	513	LEU	3.4
1	A	659	GLY	3.2
1	A	419	ILE	3.2
1	A	412	LEU	3.1
1	A	462	TRP	3.0
1	A	529	ASN	2.9
1	A	506	ILE	2.8
1	A	490	ARG	2.7
1	A	232	LEU	2.7
1	A	322	TYR	2.6
1	A	447	GLU	2.5
1	A	510	GLU	2.5
1	A	444	LEU	2.5
1	A	266	LEU	2.4
1	A	299	LEU	2.3
1	A	570	VAL	2.3
1	A	445	ASN	2.2
1	A	526	LEU	2.2
1	A	495	LEU	2.2
1	A	421	GLN	2.2
1	A	749	PHE	2.2
1	A	303	LEU	2.2
1	A	503	TYR	2.1
1	A	308	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	398	ALA	2.1
1	A	753	PHE	2.0
1	A	389	LYS	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(\AA^2)	Q<0.9
2	ADP	A	901	27/27	0.87	0.33	0.04	71,90,105,195	0
2	ADP	A	902	27/27	0.88	0.34	-0.33	57,79,93,95	0

6.5 Other polymers [i](#)

There are no such residues in this entry.