



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2020 – 06:25 PM BST

PDB ID : 4RF7
Title : Crystal structure of double-domain arginine kinase from *Anthopleura japonica* in complex with substrate L-arginine
Authors : Wang, Z.; Qiao, Z.; Ye, S.; Zhang, R.
Deposited on : 2014-09-25
Resolution : 2.10 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.4.dev1
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) : 2.14.4.dev1

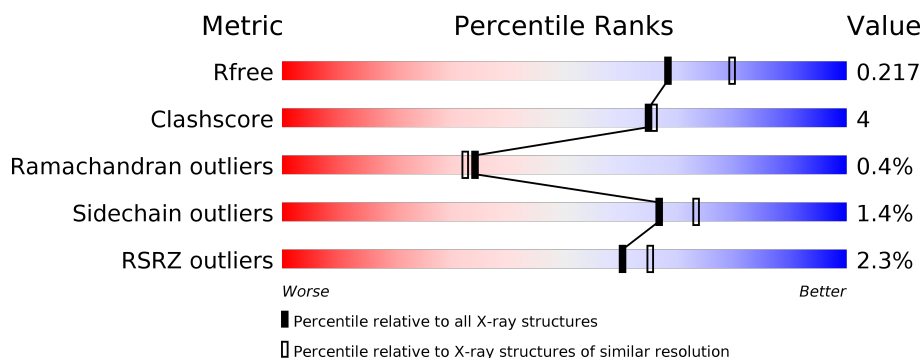
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 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	718	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>7%</div> <div>.</div> </div> </div>
1	B	718	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div>..</div> </div> </div>

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
2	ACT	B	1003	-	-	X	-
3	ARG	A	807	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 12184 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.

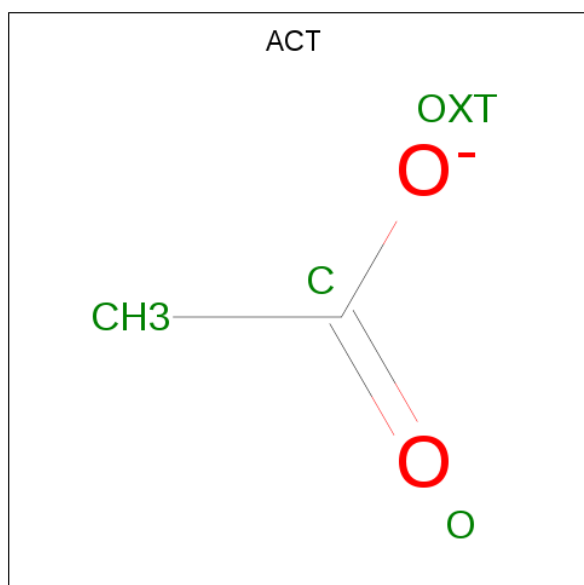
- Molecule 1 is a protein called Arginine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	692	Total	C	N	O	S	0	1	0
			5470	3455	964	1026	25			
1	B	700	Total	C	N	O	S	0	1	0
			5538	3493	979	1042	24			

There are 6 discrepancies between the modelled and reference sequences:

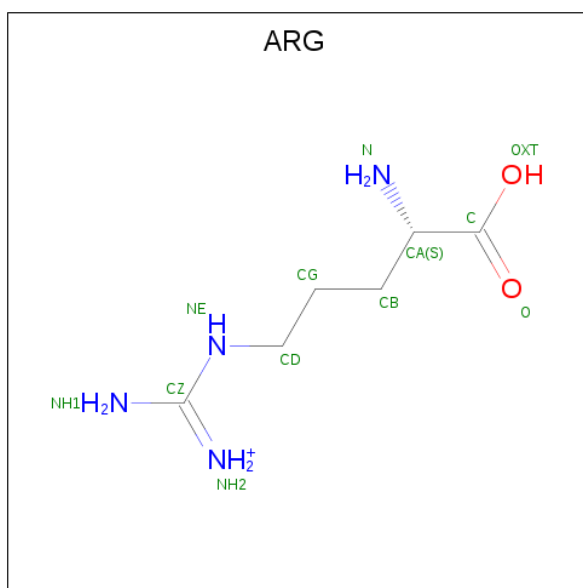
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP O15992
A	-1	PRO	-	EXPRESSION TAG	UNP O15992
A	0	HIS	-	EXPRESSION TAG	UNP O15992
B	-2	GLY	-	EXPRESSION TAG	UNP O15992
B	-1	PRO	-	EXPRESSION TAG	UNP O15992
B	0	HIS	-	EXPRESSION TAG	UNP O15992

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0

- Molecule 3 is ARGinine (three-letter code: ARG) (formula: $C_6H_{15}N_4O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 12 6 4 2	0	0
3	A	1	Total C N O 12 6 4 2	0	0
3	A	1	Total C N O 12 6 4 2	0	0
3	B	1	Total C N O 12 6 4 2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			12	6	4	2		
3	B	1	Total	C	N	O	0	0
			12	6	4	2		

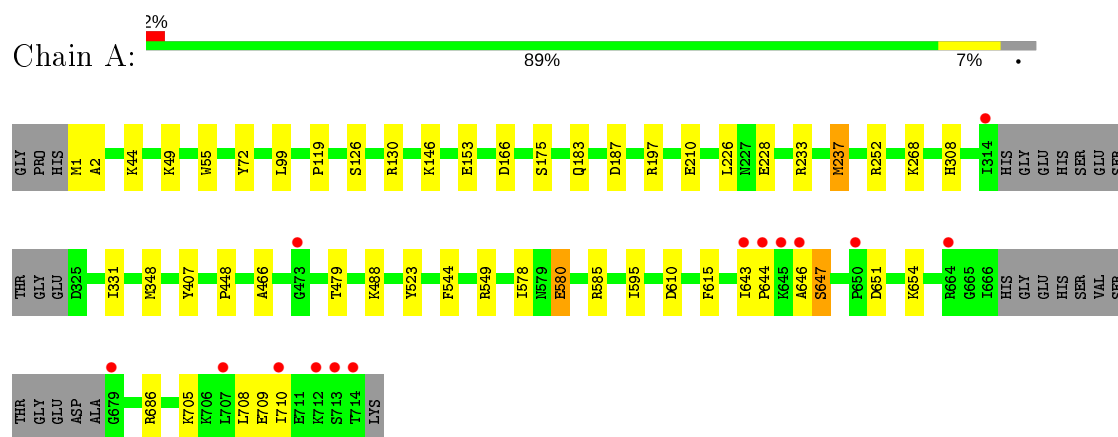
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	545	Total	O	0	0
			545	545		
4	B	527	Total	O	0	0
			527	527		

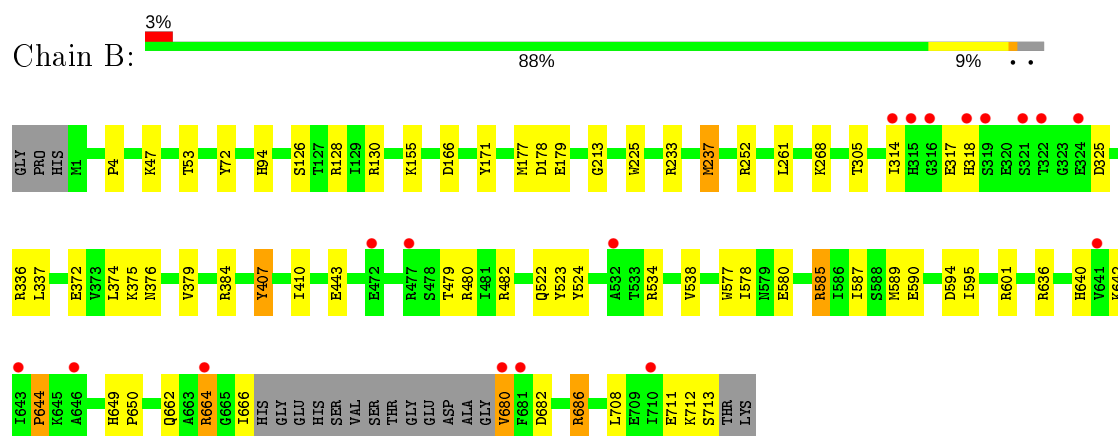
3 Residue-property plots

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: Arginine kinase



• Molecule 1: Arginine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.50Å 59.27Å 162.26Å 90.00° 91.31° 90.00°	Depositor
Resolution (Å)	47.85 – 2.10 47.85 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.85-2.10) 90.9 (47.85-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.10Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.180 , 0.217 0.180 , 0.217	Depositor DCC
R_{free} test set	4401 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12184	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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: ACT

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	0.25	0/5582	0.42	0/7527
1	B	0.26	0/5653	0.43	0/7622
All	All	0.25	0/11235	0.42	0/15149

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

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	5470	0	5469	40	0
1	B	5538	0	5522	48	0
2	A	16	0	12	1	0
2	B	16	0	12	5	0
3	A	36	0	36	9	0
3	B	36	0	36	7	0
4	A	545	0	0	13	1
4	B	527	0	0	12	1
All	All	12184	0	11087	91	1

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 4.

All (91) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:ARG:HE	2:B:1003:ACT:H1	1.22	1.01
1:B:177:MET:O	4:B:1580:HOH:O	1.98	0.81
1:B:590:GLU:HG3	1:B:601:ARG:HH22	1.47	0.78
1:B:128:ARG:HH22	2:B:1003:ACT:H3	1.52	0.73
1:B:72:TYR:HH	3:B:1005:ARG:N	1.87	0.71
1:B:374:LEU:O	1:B:384:ARG:NH1	2.23	0.69
1:A:1:MET:O	4:A:1192:HOH:O	2.12	0.68
1:B:594:ASP:OD1	4:B:1337:HOH:O	2.11	0.68
1:B:664:ARG:NH1	4:B:1516:HOH:O	2.25	0.68
1:A:175:SER:OG	4:A:1269:HOH:O	2.11	0.67
1:B:662:GLN:HB3	1:B:686:ARG:HD3	1.76	0.66
1:B:126:SER:HB2	3:B:1006:ARG:HD2	1.79	0.65
1:A:166:ASP:OD1	1:A:252:ARG:NH2	2.31	0.64
1:A:44:LYS:NZ	4:A:1285:HOH:O	2.23	0.64
3:B:1007:ARG:N	4:B:1353:HOH:O	2.31	0.62
1:B:636:ARG:NH2	1:B:682:ASP:OD1	2.33	0.61
1:A:705:LYS:NZ	1:A:709:GLU:OE2	2.33	0.61
1:A:549:ARG:NH1	4:A:1334:HOH:O	2.34	0.61
1:B:642:LYS:HA	1:B:680:VAL:HB	1.84	0.60
1:A:615:PHE:H	3:A:807:ARG:HG3	1.67	0.59
1:B:577:TRP:HB2	1:B:585:ARG:HB3	1.85	0.59
3:B:1006:ARG:NH1	4:B:1151:HOH:O	2.36	0.58
1:B:166:ASP:OD1	1:B:252:ARG:NH2	2.37	0.58
1:A:448:PRO:O	4:A:1044:HOH:O	2.17	0.58
1:B:268:LYS:NZ	4:B:1268:HOH:O	2.37	0.57
1:A:651:ASP:OD1	1:A:654:LYS:NZ	2.38	0.57
1:B:590:GLU:HG3	1:B:601:ARG:NH2	2.19	0.57
1:B:376:ASN:HB3	1:B:379:VAL:HG23	1.87	0.56
1:A:72:TYR:HH	3:A:805:ARG:N	2.03	0.55
1:B:336:ARG:NH1	4:B:1380:HOH:O	2.39	0.54
1:B:128:ARG:HH12	2:B:1003:ACT:H2	1.72	0.53
1:B:305:THR:OG1	4:B:1236:HOH:O	2.18	0.53
1:B:649:HIS:CG	1:B:650:PRO:HD2	2.45	0.52
1:B:155:LYS:HD3	1:B:261:LEU:HD22	1.90	0.52
1:A:308:HIS:ND1	4:A:1260:HOH:O	2.34	0.51
1:B:372:GLU:OE2	1:B:375:LYS:NZ	2.36	0.51
1:A:130:ARG:HD2	1:A:233:ARG:CZ	2.41	0.51
1:B:225:TRP:HB2	1:B:233[A]:ARG:HB3	1.93	0.51
1:A:153:GLU:HA	1:A:226:LEU:HD21	1.93	0.50
1:A:466:ALA:O	4:A:970:HOH:O	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:443:GLU:OE1	4:B:1442:HOH:O	2.20	0.50
1:B:4:PRO:HD2	4:B:1545:HOH:O	2.11	0.49
1:A:331:ILE:HB	1:A:348:MET:HE1	1.94	0.49
1:B:179:GLU:OE1	4:B:1584:HOH:O	2.19	0.49
1:B:534:ARG:O	1:B:538:VAL:HG23	2.13	0.48
2:A:803:ACT:O	3:A:807:ARG:NH1	2.42	0.48
1:B:522:GLN:OE1	1:B:524:TYR:OH	2.13	0.48
1:B:94:HIS:CG	1:B:337:LEU:HD12	2.49	0.48
1:B:479:THR:HG21	1:B:595:ILE:HA	1.95	0.48
1:A:99:LEU:HD22	1:A:268:LYS:HG2	1.97	0.47
1:A:126:SER:OG	3:A:806:ARG:NH1	2.39	0.47
1:A:233:ARG:CZ	4:A:1307:HOH:O	2.63	0.47
1:B:237:MET:N	1:B:237:MET:SD	2.87	0.47
1:B:47:LYS:HG2	1:B:53:THR:HG22	1.97	0.47
1:A:610:ASP:OD1	3:A:807:ARG:HD3	2.15	0.47
1:A:610:ASP:OD2	3:A:807:ARG:NH2	2.48	0.47
1:B:126:SER:CB	3:B:1006:ARG:HD2	2.45	0.46
1:B:480:ARG:HH21	1:B:587:ILE:HD13	1.81	0.46
1:B:644:PRO:HD2	1:B:711:GLU:OE2	2.16	0.46
1:A:643:ILE:O	1:A:647:SER:HB3	2.16	0.46
1:B:523:TYR:CE1	1:B:578:ILE:HD12	2.51	0.46
1:B:128:ARG:HH22	2:B:1003:ACT:CH3	2.26	0.45
1:B:317:GLU:OE1	1:B:318:HIS:NE2	2.50	0.45
1:A:233:ARG:NH2	4:A:1366:HOH:O	2.50	0.45
1:A:615:PHE:H	3:A:807:ARG:CD	2.30	0.45
1:A:119:PRO:HD2	4:A:945:HOH:O	2.17	0.44
1:A:646:ALA:HB2	1:A:710:ILE:HB	1.98	0.44
1:A:146:LYS:NZ	1:A:210:GLU:OE1	2.46	0.44
1:A:523:TYR:CE1	1:A:578:ILE:HD12	2.52	0.44
1:B:708:LEU:O	1:B:712:LYS:HG2	2.19	0.43
1:A:49:LYS:HE2	1:A:49:LYS:HB3	1.72	0.43
1:A:544:PHE:CD1	1:A:580:GLU:HG3	2.54	0.43
1:A:615:PHE:H	3:A:807:ARG:CG	2.31	0.42
1:B:171:TYR:OH	1:B:213:GLY:HA3	2.19	0.42
1:A:166:ASP:CG	1:A:252:ARG:HH22	2.22	0.42
1:A:2:ALA:HB2	1:A:197:ARG:HD2	2.02	0.42
1:A:488:LYS:NZ	3:A:807:ARG:O	2.30	0.42
1:B:233[B]:ARG:HH21	2:B:1003:ACT:H2	1.84	0.42
1:A:183:GLN:NE2	1:A:187:ASP:OD1	2.53	0.42
1:B:482:ARG:HD3	1:B:585:ARG:HD2	2.01	0.42
1:A:479:THR:HG21	1:A:595:ILE:HA	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:ARG:HD2	4:A:1307:HOH:O	2.18	0.41
1:B:237:MET:HB2	3:B:1006:ARG:HG2	2.01	0.41
1:B:237:MET:HB2	3:B:1006:ARG:CG	2.51	0.41
1:A:585:ARG:NH2	4:A:994:HOH:O	2.30	0.41
1:B:47:LYS:HE2	4:B:1405:HOH:O	2.19	0.41
1:A:233:ARG:NH2	4:A:1307:HOH:O	2.54	0.41
1:A:708:LEU:HA	1:A:708:LEU:HD23	1.92	0.41
1:B:407:TYR:HA	1:B:410:ILE:HG12	2.02	0.41
1:B:314:ILE:HD11	1:B:325:ASP:CG	2.42	0.40
1:A:237:MET:N	1:A:237:MET:SD	2.94	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1407:HOH:O	4:B:1617:HOH:O[2_546]	1.97	0.23

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	687/718 (96%)	666 (97%)	18 (3%)	3 (0%)	34 32
1	B	697/718 (97%)	677 (97%)	18 (3%)	2 (0%)	41 41
All	All	1384/1436 (96%)	1343 (97%)	36 (3%)	5 (0%)	34 32

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	644	PRO
1	A	580	GLU
1	A	644	PRO

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Mol	Chain	Res	Type
1	A	228	GLU
1	B	580	GLU

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	597/616 (97%)	592 (99%)	5 (1%)	81	86
1	B	604/616 (98%)	593 (98%)	11 (2%)	59	65
All	All	1201/1232 (98%)	1185 (99%)	16 (1%)	67	75

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	TRP
1	A	237	MET
1	A	407	TYR
1	A	647	SER
1	A	686	ARG
1	B	178	ASP
1	B	237	MET
1	B	407	TYR
1	B	585	ARG
1	B	589	MET
1	B	640	HIS
1	B	664	ARG
1	B	666	ILE
1	B	680	VAL
1	B	686	ARG
1	B	713	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	183	GLN

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](#)

14 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	ACT	B	1002	-	1,3,3	1.26	0	0,3,3	0.00	-
2	ACT	A	802	-	1,3,3	1.35	0	0,3,3	0.00	-
3	ARG	B	1006	-	7,11,11	0.23	0	6,13,13	0.15	0
2	ACT	B	1001	-	1,3,3	1.30	0	0,3,3	0.00	-
2	ACT	B	1003	-	1,3,3	0.40	0	0,3,3	0.00	-
3	ARG	A	805	-	7,11,11	0.28	0	6,13,13	0.30	0
3	ARG	B	1007	-	7,11,11	0.30	0	6,13,13	0.24	0
2	ACT	A	801	-	1,3,3	1.25	0	0,3,3	0.00	-
3	ARG	A	806	-	7,11,11	0.29	0	6,13,13	0.19	0
2	ACT	A	804	-	1,3,3	1.23	0	0,3,3	0.00	-
2	ACT	B	1004	-	1,3,3	1.32	0	0,3,3	0.00	-
3	ARG	A	807	-	7,11,11	0.25	0	6,13,13	0.28	0
3	ARG	B	1005	-	7,11,11	0.29	0	6,13,13	0.29	0
2	ACT	A	803	-	1,3,3	1.28	0	0,3,3	0.00	-

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
3	ARG	B	1006	-	-	1/7/11/11	-
3	ARG	A	805	-	-	0/7/11/11	-
3	ARG	B	1007	-	-	2/7/11/11	-
3	ARG	A	806	-	-	1/7/11/11	-
3	ARG	A	807	-	-	1/7/11/11	-
3	ARG	B	1005	-	-	1/7/11/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1007	ARG	N-CA-CB-CG
3	B	1007	ARG	C-CA-CB-CG
3	A	806	ARG	NE-CD-CG-CB
3	B	1006	ARG	CA-CB-CG-CD
3	A	807	ARG	CA-CB-CG-CD
3	B	1005	ARG	NE-CD-CG-CB

There are no ring outliers.

8 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1006	ARG	5	0
2	B	1003	ACT	5	0
3	A	805	ARG	1	0
3	B	1007	ARG	1	0
3	A	806	ARG	1	0
3	A	807	ARG	7	0
3	B	1005	ARG	1	0
2	A	803	ACT	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	692/718 (96%)	-0.22	14 (2%) 65 69	22, 37, 78, 127	0
1	B	700/718 (97%)	-0.22	18 (2%) 56 61	24, 37, 79, 118	0
All	All	1392/1436 (96%)	-0.22	32 (2%) 60 65	22, 37, 79, 127	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	646	ALA	8.1
1	A	650	PRO	6.3
1	A	643	ILE	5.3
1	B	680	VAL	4.9
1	B	710	ILE	4.8
1	B	315	HIS	4.5
1	B	318	HIS	4.2
1	B	681	PHE	3.8
1	A	710	ILE	3.8
1	B	314	ILE	3.7
1	B	322	THR	3.5
1	B	319	SER	3.2
1	B	643	ILE	3.2
1	B	532	ALA	3.0
1	A	679	GLY	2.9
1	B	321	SER	2.9
1	B	664	ARG	2.7
1	A	712	LYS	2.7
1	A	644	PRO	2.7
1	A	645	LYS	2.4
1	B	324	GLU	2.3
1	B	472	GLU	2.3
1	A	707	LEU	2.3
1	B	477	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	714	THR	2.2
1	A	713	SER	2.2
1	A	664	ARG	2.1
1	B	641	VAL	2.1
1	A	473	GLY	2.0
1	A	314	ILE	2.0
1	B	646	ALA	2.0
1	B	316	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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
3	ARG	B	1007	12/12	0.62	0.26	78,88,94,97	0
2	ACT	B	1004	4/4	0.69	0.15	83,85,85,85	0
2	ACT	A	802	4/4	0.70	0.21	62,63,64,65	0
2	ACT	A	801	4/4	0.71	0.26	54,54,56,59	0
2	ACT	A	803	4/4	0.71	0.19	60,63,64,67	0
3	ARG	B	1006	12/12	0.73	0.30	61,63,69,69	0
3	ARG	A	806	12/12	0.75	0.30	60,81,103,107	0
3	ARG	A	807	12/12	0.76	0.31	39,62,78,80	0
2	ACT	A	804	4/4	0.84	0.16	71,75,75,77	0
3	ARG	B	1005	12/12	0.85	0.15	50,54,59,59	0
2	ACT	B	1003	4/4	0.88	0.12	28,33,35,43	0
3	ARG	A	805	12/12	0.88	0.15	53,56,58,59	0
2	ACT	B	1001	4/4	0.92	0.24	48,53,53,56	0
2	ACT	B	1002	4/4	0.95	0.07	72,72,72,73	0

6.5 Other polymers

There are no such residues in this entry.