



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2017 – 01:25 PM EDT

PDB ID : 2OSB  
Title : Crystal structure of a thermostable mutant of Bacillus subtilis Adenylate Kinase (Q16L/Q199R/)  
Authors : Myers, J.; Counago, R.; Wilson, C.J.; Wu, G.; Shamoo, Y.  
Deposited on : unknown  
Resolution : 1.80 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : rb-20030345

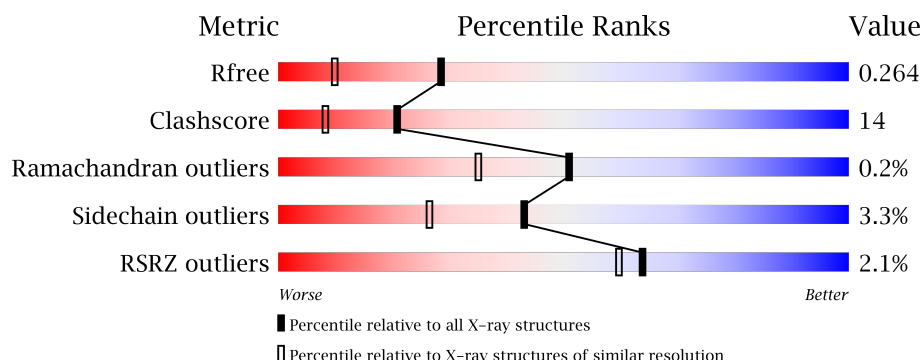
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	100719	4827 (1.80-1.80)
Clashscore	112137	5742 (1.80-1.80)
Ramachandran outliers	110173	5676 (1.80-1.80)
Sidechain outliers	110143	5675 (1.80-1.80)
RSRZ outliers	101464	4906 (1.80-1.80)

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  $\geq 3$ , 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	216	<div> <div>3%</div> <div>69%</div> <div>26%</div> <div>5%</div> </div>
1	B	216	<div> <div>%</div> <div>79%</div> <div>20%</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
3	MG	A	1223	-	-	-	X
3	MG	B	1222	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3759 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 Adenylate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1682	1056	281	335	10			
1	B	216	Total	C	N	O	S	0	0	0
			1682	1056	281	335	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	LEU	GLN	ENGINEERED	UNP P16304
A	199	ARG	GLN	ENGINEERED	UNP P16304
B	16	LEU	GLN	ENGINEERED	UNP P16304
B	199	ARG	GLN	ENGINEERED	UNP P16304

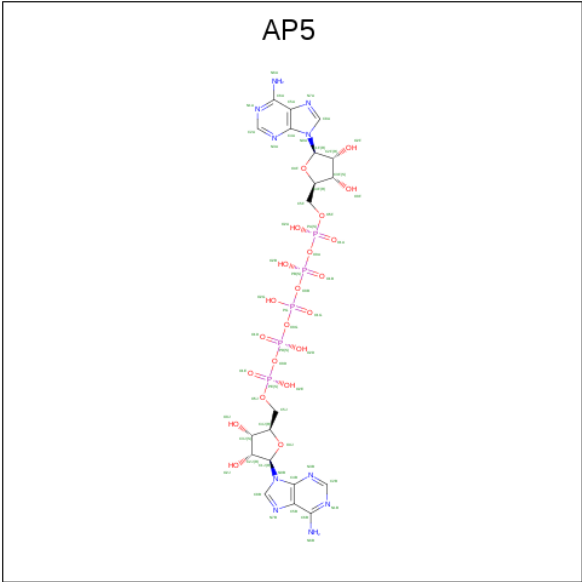
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is BIS(ADENOSINE)-5'-PENTAPHOSPHATE (three-letter code: AP5) (formula: C<sub>20</sub>H<sub>29</sub>N<sub>10</sub>O<sub>22</sub>P<sub>5</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			57	20	10	22	5		
4	B	1	Total	C	N	O	P	0	0
			57	20	10	22	5		

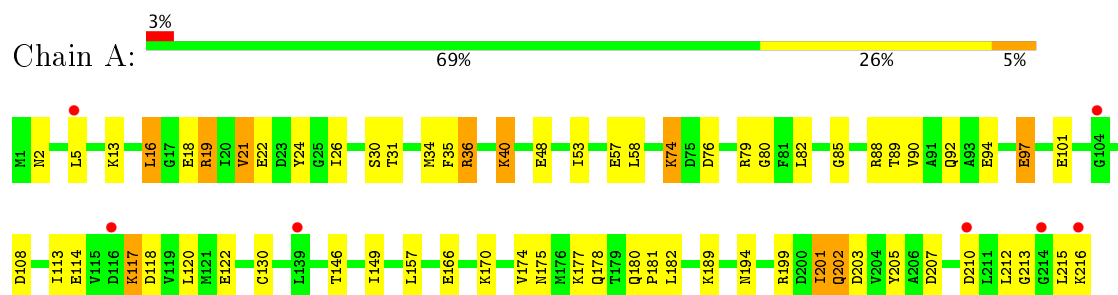
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	162	Total	O	0	0
			162	162		
5	B	115	Total	O	0	0
			115	115		

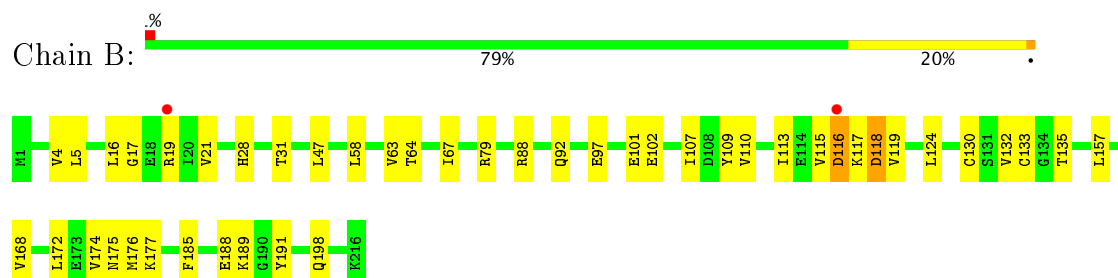
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Adenylate kinase



#### • Molecule 1: Adenylate kinase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	34.58 Å   74.32 Å   77.42 Å 90.00°   100.21°   90.00°	Depositor
Resolution (Å)	20.00 – 1.80 30.43 – 1.64	Depositor EDS
% Data completeness (in resolution range)	97.3 (20.00-1.80) 91.3 (30.43-1.64)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.50 (at 1.63 Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.245   ,   0.264 0.245   ,   0.264	Depositor DCC
$R_{free}$ test set	1781 reflections (5.12%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.2	Xtriage
Anisotropy	0.190	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 33.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.044 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3759	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.54 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.6319e-04.*

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

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MG, AP5

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	0.57	2/1707 (0.1%)	0.76	3/2301 (0.1%)
1	B	0.36	1/1707 (0.1%)	0.61	1/2301 (0.0%)
All	All	0.47	3/3414 (0.1%)	0.69	4/4602 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	202	GLN	C-N	15.10	1.68	1.34
1	B	116	ASP	C-N	-6.77	1.18	1.34
1	A	201	ILE	C-N	5.59	1.47	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	ILE	O-C-N	-10.44	106.00	122.70
1	A	201	ILE	CA-C-N	7.10	132.82	117.20
1	B	116	ASP	O-C-N	-7.00	111.50	122.70
1	A	201	ILE	C-N-CA	6.40	137.71	121.70

There are no chirality outliers.



All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	118	ASP	Mainchain
1	B	118	ASP	Mainchain

## 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	1682	0	1674	60	0
1	B	1682	0	1673	36	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	57	0	24	3	0
4	B	57	0	24	1	0
5	A	162	0	0	14	0
5	B	115	0	0	3	0
All	All	3759	0	3395	95	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 (95) 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:202:GLN:C	1:A:203:ASP:N	1.68	1.42
1:A:177:LYS:HE3	5:A:1270:HOH:O	1.43	1.19
1:A:114:GLU:OE2	1:A:194:ASN:HB3	1.48	1.11
1:A:21:VAL:HG23	1:A:26:ILE:O	1.50	1.11
1:B:21:VAL:HG11	1:B:28:HIS:HB2	1.36	1.03
1:A:57:GLU:OE2	5:A:1336:HOH:O	1.78	1.00
1:A:57:GLU:HG2	5:A:1336:HOH:O	1.62	0.99
1:A:166:GLU:OE2	5:A:1318:HOH:O	1.80	0.97
1:A:97:GLU:HA	1:A:97:GLU:OE1	1.63	0.94
1:A:57:GLU:CG	5:A:1336:HOH:O	2.18	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:ASP:OD1	5:B:1245:HOH:O	1.96	0.82
1:B:102:GLU:OE1	5:B:1236:HOH:O	1.97	0.80
1:A:213:GLY:O	5:A:1300:HOH:O	2.06	0.73
1:A:85:GLY:O	1:A:88:ARG:HD2	1.92	0.69
1:A:57:GLU:CD	5:A:1336:HOH:O	2.17	0.69
1:A:36:ARG:HD3	5:A:1259:HOH:O	1.93	0.69
1:B:117:LYS:HD2	1:B:176:MET:SD	2.34	0.68
1:A:16:LEU:HD12	1:A:113:ILE:CD1	2.24	0.67
1:A:202:GLN:C	1:A:203:ASP:CA	2.63	0.66
1:B:97:GLU:O	1:B:101:GLU:HG3	1.97	0.64
1:A:92:GLN:HE22	4:A:1218:AP5:H62B	1.46	0.63
1:A:216:LYS:HG3	1:A:216:LYS:OXT	1.98	0.62
1:B:16:LEU:HD12	1:B:113:ILE:HD13	1.82	0.62
1:A:5:LEU:HG	1:A:82:LEU:HD11	1.82	0.61
1:A:13:LYS:HG3	5:A:1292:HOH:O	2.00	0.61
1:A:178:GLN:O	1:A:181:PRO:HG2	2.01	0.61
1:A:212:LEU:HA	1:A:215:LEU:HD12	1.84	0.60
1:A:202:GLN:H	1:A:202:GLN:CD	2.05	0.59
1:B:63:VAL:O	1:B:67:ILE:HG23	2.04	0.57
1:A:88:ARG:HE	1:A:175:ASN:HD21	1.53	0.57
1:B:88:ARG:HH21	1:B:175:ASN:HD21	1.52	0.57
1:B:92:GLN:HE22	4:B:1219:AP5:H62B	1.51	0.57
1:A:202:GLN:CA	1:A:203:ASP:N	2.66	0.56
1:A:2:ASN:ND2	1:A:80:GLY:HA2	2.23	0.54
1:A:48:GLU:HG3	5:A:1262:HOH:O	2.08	0.53
1:A:16:LEU:HD12	1:A:113:ILE:HD11	1.88	0.53
1:A:16:LEU:HD12	1:A:113:ILE:HD13	1.89	0.53
1:B:188:GLU:OE2	1:B:189:LYS:HD3	2.09	0.53
1:A:89:THR:OG1	1:A:92:GLN:HG3	2.09	0.53
1:B:130:CYS:HA	1:B:157:LEU:HD23	1.89	0.53
1:A:18:GLU:O	1:A:22:GLU:HG3	2.08	0.52
1:A:97:GLU:OE1	5:A:1273:HOH:O	2.18	0.52
1:A:88:ARG:HA	1:A:182:LEU:HD11	1.92	0.52
1:B:172:LEU:O	1:B:176:MET:HB2	2.09	0.52
1:A:101:GLU:HG3	1:B:47:LEU:HD21	1.92	0.52
1:A:88:ARG:HB2	5:A:1272:HOH:O	2.10	0.52
1:B:88:ARG:HD3	1:B:175:ASN:ND2	2.26	0.51
1:B:58:LEU:HG	1:B:174:VAL:CG1	2.39	0.51
1:A:74:LYS:HB3	1:A:76:ASP:OD2	2.11	0.51
1:A:90:VAL:O	1:A:94:GLU:HG3	2.11	0.51
1:B:4:VAL:HG23	1:B:107:ILE:HD12	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:LEU:HG	1:B:174:VAL:HG11	1.92	0.49
1:A:88:ARG:HD3	4:A:1218:AP5:C5B	2.42	0.49
1:A:180:GLN:N	1:A:181:PRO:HD2	2.28	0.49
1:A:88:ARG:HD3	4:A:1218:AP5:C6B	2.43	0.48
1:B:64:THR:O	1:B:67:ILE:HG12	2.13	0.48
1:A:189:LYS:NZ	5:A:1323:HOH:O	2.46	0.48
1:B:88:ARG:HD3	1:B:175:ASN:HD21	1.79	0.48
1:A:88:ARG:HE	1:A:175:ASN:ND2	2.12	0.47
1:B:16:LEU:HD12	1:B:113:ILE:CD1	2.44	0.47
1:A:117:LYS:HA	1:A:120:LEU:HD12	1.97	0.47
1:A:215:LEU:O	1:A:216:LYS:HB3	2.15	0.47
1:A:21:VAL:CG2	1:A:26:ILE:O	2.41	0.46
1:A:170:LYS:O	1:A:174:VAL:HG23	2.15	0.46
1:A:199:ARG:HH22	1:A:207:ASP:CG	2.19	0.45
1:B:19:ARG:HG3	1:B:19:ARG:HH11	1.82	0.45
1:A:201:ILE:HG23	5:A:1285:HOH:O	2.15	0.45
1:B:119:VAL:HG21	1:B:198:GLN:HE21	1.82	0.45
1:B:5:LEU:HD11	1:B:113:ILE:HD11	1.99	0.45
1:A:146:THR:HB	1:A:149:ILE:HD11	2.00	0.44
1:B:177:LYS:HG3	5:B:1294:HOH:O	2.17	0.44
1:A:122:GLU:OE1	1:A:122:GLU:HA	2.18	0.44
1:B:58:LEU:HD22	1:B:88:ARG:HD2	1.98	0.44
1:A:18:GLU:O	1:A:21:VAL:HG12	2.18	0.43
1:B:133:CYS:SG	1:B:135:THR:HG22	2.59	0.43
1:B:185:PHE:O	1:B:189:LYS:HG2	2.19	0.43
1:B:17:GLY:O	1:B:21:VAL:HG13	2.18	0.43
1:A:35:PHE:HB3	1:A:53:ILE:HD11	2.00	0.43
1:A:5:LEU:HD23	1:A:5:LEU:HA	1.89	0.42
1:B:4:VAL:HB	1:B:110:VAL:HG22	2.01	0.42
1:A:19:ARG:HG3	1:A:205:TYR:CE2	2.54	0.42
1:B:177:LYS:HB3	1:B:177:LYS:HE2	1.77	0.42
1:B:117:LYS:C	1:B:119:VAL:H	2.24	0.42
1:B:124:LEU:HD12	1:B:168:VAL:HG13	2.02	0.42
1:B:79:ARG:NE	1:B:79:ARG:HA	2.35	0.42
1:A:130:CYS:HA	1:A:157:LEU:HD23	2.01	0.41
1:A:108:ASP:CG	1:A:215:LEU:HD22	2.40	0.41
1:A:24:TYR:CE1	1:A:212:LEU:HD23	2.55	0.41
1:B:109:TYR:HA	1:B:191:TYR:HB2	2.02	0.41
1:A:30:SER:O	1:A:34:MET:HG3	2.21	0.41
1:B:115:VAL:HG13	1:B:198:GLN:CG	2.50	0.41
1:A:58:LEU:HG	1:A:174:VAL:HG11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:LYS:HB3	1:A:40:LYS:HE3	1.50	0.40
1:A:19:ARG:HA	1:A:19:ARG:HD2	1.51	0.40
1:B:130:CYS:SG	1:B:132:VAL:HB	2.62	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	214/216 (99%)	209 (98%)	5 (2%)	0	100	100
1	B	214/216 (99%)	207 (97%)	6 (3%)	1 (0%)	32	17
All	All	428/432 (99%)	416 (97%)	11 (3%)	1 (0%)	51	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	118	ASP

### 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	183/183 (100%)	172 (94%)	11 (6%)	22	8
1	B	183/183 (100%)	182 (100%)	1 (0%)	91	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	366/366 (100%)	354 (97%)	12 (3%)	43 26

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

Mol	Chain	Res	Type
1	A	16	LEU
1	A	19	ARG
1	A	21	VAL
1	A	31	THR
1	A	36	ARG
1	A	40	LYS
1	A	74	LYS
1	A	79	ARG
1	A	97	GLU
1	A	117	LYS
1	A	210	ASP
1	B	31	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	92	GLN
1	A	164	ASN
1	A	175	ASN
1	B	92	GLN
1	B	112	ASN
1	B	142	ASN
1	B	175	ASN
1	B	180	GLN
1	B	198	GLN
1	B	202	GLN

### 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

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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
4	AP5	A	1218	3	48,62,62	1.16	3 (6%)	43,98,98	2.54	7 (16%)
4	AP5	B	1219	-	48,62,62	1.13	3 (6%)	43,98,98	2.60	8 (18%)

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
4	AP5	A	1218	3	-	0/36/76/76	0/6/6/6
4	AP5	B	1219	-	-	0/36/76/76	0/6/6/6

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1219	AP5	C2F-C1F	-2.31	1.50	1.53
4	A	1218	AP5	C2A-N3A	2.51	1.36	1.32
4	A	1218	AP5	O4F-C1F	2.56	1.44	1.41
4	B	1219	AP5	O4F-C1F	2.57	1.44	1.41
4	B	1219	AP5	O4J-C1J	2.71	1.45	1.41
4	A	1218	AP5	O4J-C1J	2.81	1.45	1.41

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1219	AP5	N3B-C2B-N1B	-10.74	119.50	128.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1218	AP5	N3B-C2B-N1B	-10.54	119.68	128.86
4	A	1218	AP5	N3A-C2A-N1A	-10.03	120.12	128.86
4	B	1219	AP5	N3A-C2A-N1A	-9.81	120.31	128.86
4	B	1219	AP5	C1F-N9A-C4A	-2.46	122.38	126.64
4	A	1218	AP5	C2B-N1B-C6B	2.34	122.86	118.77
4	A	1218	AP5	C4F-O4F-C1F	2.41	112.34	109.77
4	B	1219	AP5	C4F-O4F-C1F	2.53	112.46	109.77
4	A	1218	AP5	C2A-N1A-C6A	2.59	123.31	118.77
4	B	1219	AP5	C2B-N1B-C6B	2.67	123.44	118.77
4	A	1218	AP5	C5B-C6B-N6B	2.77	126.11	120.47
4	B	1219	AP5	C2A-N1A-C6A	2.84	123.75	118.77
4	B	1219	AP5	C5B-C6B-N6B	3.03	126.66	120.47
4	A	1218	AP5	C4J-O4J-C1J	3.48	113.47	109.77
4	B	1219	AP5	C4J-O4J-C1J	3.80	113.82	109.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1218	AP5	3	0
4	B	1219	AP5	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	202:GLN	C	203:ASP	N	1.68

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	116:ASP	C	117:LYS	N	1.18



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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	216/216 (100%)	0.16	7 (3%) 48 43	15, 24, 35, 44	0
1	B	216/216 (100%)	0.21	2 (0%) 84 82	17, 27, 36, 42	0
All	All	432/432 (100%)	0.19	9 (2%) 64 60	15, 25, 36, 44	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	116	ASP	3.0
1	A	5	LEU	2.8
1	A	216	LYS	2.7
1	A	214	GLY	2.4
1	B	116	ASP	2.4
1	A	104	GLY	2.1
1	B	19	ARG	2.1
1	A	210	ASP	2.1
1	A	139	LEU	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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	MG	B	1222	1/1	0.99	0.23	5.04	16,16,16,16	0
3	MG	A	1223	1/1	0.97	0.22	3.53	16,16,16,16	0
4	AP5	A	1218	57/57	0.94	0.12	-0.05	15,20,23,25	0
4	AP5	B	1219	57/57	0.93	0.10	-0.52	17,21,26,28	0
2	ZN	A	1220	1/1	0.99	0.07	-2.53	27,27,27,27	0
2	ZN	B	1221	1/1	0.97	0.03	-2.58	37,37,37,37	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.