



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:39 AM GMT

PDB ID : 2BAM  
Title : RESTRICTION ENDONUCLEASE BAMHI COMPLEX WITH DNA AND  
CALCIUM IONS (PRE-REACTIVE COMPLEX).  
Authors : Viadiu, H.; Aggarwal, A.K.  
Deposited on : 1998-08-19  
Resolution : 2.00 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

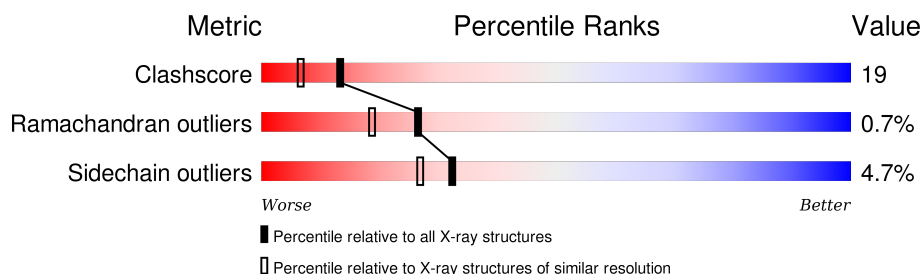
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	12	
1	D	12	
2	A	213	
2	B	213	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4164 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 DNA chain called DNA (5'-D(\*TP\*AP\*TP\*GP\*GP\*AP\*TP\*CP\*CP\*AP\*TP\*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	12	Total	C	N	O	P	0	0	0
			243	118	44	70	11			
1	D	11	Total	C	N	O	P	0	0	0
			222	108	39	65	10			

- Molecule 2 is a protein called PROTEIN (ENDONUCLEASE BAMHI).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	207	Total	C	N	O	S	0	0	0
			1678	1082	269	319	8			
2	B	210	Total	C	N	O	S	0	0	0
			1702	1097	273	324	8			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Ca	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	28	Total	O	0	0
			28	28		
4	D	25	Total	O	0	0
			25	25		
4	A	127	Total	O	0	0
			127	127		
4	B	137	Total	O	0	0
			137	137		

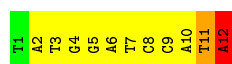
### 3 Residue-property plots

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 ( $RSRZ > 2$ ). 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.

Note EDS was not executed.

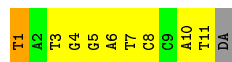
- Molecule 1: DNA (5'-D(\*TP\*AP\*TP\*GP\*GP\*AP\*TP\*CP\*CP\*AP\*TP\*A)-3')

Chain C: 



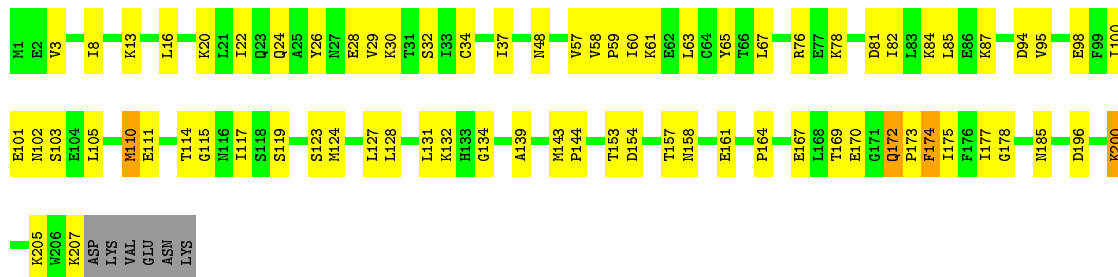
- Molecule 1: DNA (5'-D(\*TP\*AP\*TP\*GP\*GP\*AP\*TP\*CP\*CP\*AP\*TP\*A)-3')

Chain D: 



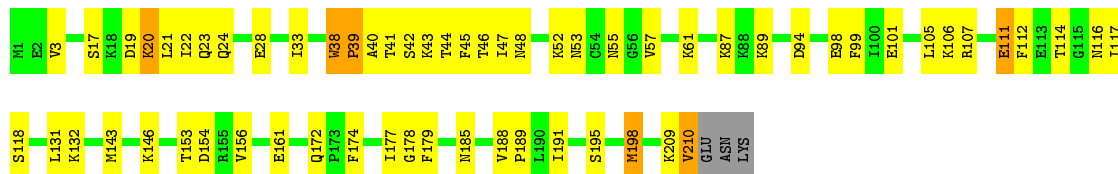
- Molecule 2: PROTEIN (ENDONUCLEASE BAMHI)

Chain A: 



- Molecule 2: PROTEIN (ENDONUCLEASE BAMHI)

Chain B: 



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	105.20 Å   79.40 Å   66.40 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	8.00 – 2.00	Depositor
% Data completeness (in resolution range)	81.6 (8.00-2.00)	Depositor
$R_{merge}$	0.07	Depositor
$R_{sym}$	7.40	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, $R_{free}$	0.233 , 0.251	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4164	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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

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	C	0.80	0/272	1.16	2/418 (0.5%)
1	D	0.79	0/248	1.08	1/381 (0.3%)
2	A	0.46	0/1713	0.67	0/2310
2	B	0.47	0/1737	0.69	0/2342
All	All	0.52	0/3970	0.76	3/5451 (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	C	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	C	11	DT	N1-C1'-C2'	6.13	124.25	112.60
1	D	1	DT	O4'-C1'-N1	5.36	111.75	108.00
1	C	12	DA	N9-C1'-C2'	5.13	122.35	112.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	12	DA	Sidechain

## 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	C	243	0	138	21	0
1	D	222	0	127	20	0
2	A	1678	0	1700	61	0
2	B	1702	0	1726	54	0
3	A	2	0	0	0	0
4	A	127	0	0	13	0
4	B	137	0	0	5	0
4	C	28	0	0	1	0
4	D	25	0	0	3	0
All	All	4164	0	3691	142	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 19.

All (142) 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:C:4:DG:H2''	1:C:5:DG:H5''	1.49	0.93
2:A:111:GLU:HG2	2:A:123:SER:HB3	1.55	0.89
1:D:6:DA:H2'	4:A:506:HOH:O	1.72	0.89
2:B:38:TRP:H	2:B:45:PHE:HA	1.41	0.85
1:C:6:DA:H2'	4:B:217:HOH:O	1.80	0.80
1:D:10:DA:H8	4:D:16:HOH:O	1.66	0.78
2:B:209:LYS:HG2	2:B:210:VAL:H	1.50	0.76
2:B:43:LYS:HG3	2:B:44:THR:HG23	1.68	0.74
1:D:4:DG:H1'	1:D:5:DG:H5''	1.69	0.74
1:D:7:DT:H5''	2:A:196:ASP:HB2	1.68	0.74
1:D:5:DG:H2''	1:D:6:DA:H5'	1.70	0.73
2:A:26:TYR:CZ	2:A:30:LYS:HD2	2.23	0.73
1:C:7:DT:H2'	2:B:153:THR:CG2	2.19	0.73
2:A:20:LYS:HB2	4:A:602:HOH:O	1.89	0.73
1:D:7:DT:H2'	2:A:153:THR:HG23	1.71	0.72
1:C:2:DA:OP2	1:C:2:DA:H2'	1.91	0.71
2:A:131:LEU:HD13	2:A:172:GLN:HG2	1.71	0.71
2:A:102:ASN:CG	2:A:103:SER:H	1.96	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:52:LYS:HE3	2:B:53:ASN:OD1	1.92	0.69
2:B:38:TRP:HB3	2:B:39:PRO:HD2	1.73	0.69
1:C:6:DA:H1'	1:C:7:DT:C5'	2.23	0.69
2:B:209:LYS:HG2	2:B:210:VAL:N	2.07	0.69
2:A:100:ILE:HG12	2:A:105:LEU:HD13	1.73	0.69
1:D:7:DT:H2'	2:A:153:THR:CG2	2.24	0.68
2:A:57:VAL:HG21	2:A:114:THR:HB	1.74	0.68
2:B:57:VAL:HG11	2:B:156:VAL:HG11	1.76	0.68
2:A:20:LYS:HE3	4:A:602:HOH:O	1.93	0.68
2:A:111:GLU:HG2	2:A:123:SER:CB	2.24	0.67
2:A:164:PRO:HA	2:B:87:LYS:HD3	1.76	0.67
2:A:169:THR:HG21	2:A:174:PHE:CB	2.26	0.66
1:D:3:DT:H6	4:D:33:HOH:O	1.79	0.65
1:C:7:DT:H2'	2:B:153:THR:HG23	1.78	0.65
2:B:153:THR:O	2:B:156:VAL:HG13	1.97	0.65
2:B:98:GLU:HB3	2:B:105:LEU:HD11	1.79	0.64
2:A:169:THR:HG21	2:A:174:PHE:HB2	1.79	0.64
2:A:3:VAL:HG13	2:A:177:ILE:HG23	1.80	0.63
1:C:6:DA:H1'	1:C:7:DT:H5'	1.79	0.63
2:A:169:THR:O	2:A:169:THR:HG22	1.97	0.62
1:C:6:DA:H2''	1:C:7:DT:H5'	1.81	0.62
2:A:32:SER:HA	2:A:67:LEU:HD23	1.82	0.61
2:B:38:TRP:O	2:B:39:PRO:C	2.38	0.60
2:A:34:CYS:SG	4:A:528:HOH:O	2.57	0.60
2:A:78:LYS:HE2	4:A:603:HOH:O	2.00	0.60
2:A:200:LYS:HE2	2:A:200:LYS:O	2.02	0.60
2:B:209:LYS:O	2:B:210:VAL:HG22	2.03	0.59
2:B:43:LYS:HG2	2:B:44:THR:N	2.18	0.59
2:B:38:TRP:CB	2:B:39:PRO:HD2	2.34	0.58
2:B:20:LYS:H	2:B:20:LYS:HZ2	1.49	0.58
2:B:39:PRO:O	2:B:42:SER:N	2.35	0.57
2:B:24:GLN:O	2:B:28:GLU:HG3	2.05	0.57
2:B:48:ASN:HA	2:B:185:ASN:O	2.04	0.57
1:C:9:DC:H2''	1:C:10:DA:C8	2.41	0.56
2:A:144:PRO:O	2:A:158:ASN:HA	2.06	0.56
2:B:55:ASN:HB3	2:B:191:ILE:HB	1.89	0.55
2:A:8:ILE:HG23	2:A:175:ILE:HG12	1.88	0.55
2:B:38:TRP:CE3	2:B:38:TRP:HA	2.42	0.55
2:B:94:ASP:HB3	2:B:111:GLU:HA	1.90	0.54
1:C:4:DG:C2'	1:C:5:DG:H5''	2.31	0.54
2:B:43:LYS:CG	2:B:44:THR:HG23	2.36	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:5:DG:C2'	1:D:6:DA:H5'	2.37	0.54
1:D:10:DA:H1'	1:D:11:DT:H5''	1.90	0.53
2:A:84:LYS:HD2	4:A:532:HOH:O	2.07	0.53
2:B:61:LYS:HD2	2:B:114:THR:OG1	2.09	0.53
2:A:57:VAL:CG2	2:A:114:THR:HB	2.39	0.53
2:A:143:MET:O	2:A:178:GLY:HA2	2.09	0.53
2:B:43:LYS:CG	2:B:44:THR:N	2.72	0.52
2:A:124:MET:O	2:A:128:LEU:HG	2.10	0.52
2:B:131:LEU:HB3	2:B:172:GLN:NE2	2.25	0.52
1:C:6:DA:H1'	1:C:7:DT:H5''	1.91	0.52
2:A:134:GLY:HA2	4:A:548:HOH:O	2.10	0.51
2:B:20:LYS:CD	2:B:20:LYS:H	2.23	0.51
2:A:169:THR:HG21	2:A:174:PHE:CG	2.46	0.51
2:A:65:TYR:CD1	2:A:76:ARG:HD3	2.46	0.50
1:D:5:DG:H2'	4:A:551:HOH:O	2.11	0.50
1:C:6:DA:C2'	1:C:7:DT:H5'	2.41	0.50
2:A:144:PRO:HG2	2:A:157:THR:O	2.12	0.50
1:C:12:DA:C2	1:D:1:DT:N3	2.80	0.49
2:B:19:ASP:O	2:B:23:GLN:HG3	2.12	0.49
2:B:195:SER:HA	2:B:198:MET:HE2	1.94	0.49
2:B:38:TRP:O	2:B:39:PRO:O	2.29	0.49
2:A:57:VAL:HG23	2:A:60:ILE:HD12	1.95	0.49
2:A:37:ILE:HG13	2:A:63:LEU:CD1	2.43	0.49
2:A:29:VAL:HG21	2:A:110:MET:HE1	1.96	0.48
2:B:39:PRO:O	2:B:41:THR:N	2.46	0.48
2:A:102:ASN:CG	2:A:103:SER:N	2.64	0.48
2:A:26:TYR:CE2	2:A:30:LYS:HD2	2.48	0.48
1:D:4:DG:H4'	4:A:504:HOH:O	2.14	0.48
2:A:169:THR:HG21	2:A:174:PHE:CD2	2.48	0.48
2:B:21:LEU:HD22	2:B:99:PHE:HB3	1.95	0.47
2:A:94:ASP:HB2	2:A:111:GLU:HA	1.95	0.47
1:C:12:DA:C6	1:D:1:DT:C4	3.02	0.47
2:B:46:THR:HB	2:B:188:VAL:HG11	1.95	0.47
1:D:3:DT:H71	4:D:33:HOH:O	2.15	0.47
2:A:76:ARG:HD2	2:A:94:ASP:O	2.14	0.47
2:A:172:GLN:HG3	2:A:173:PRO:HD2	1.96	0.47
2:A:82:ILE:HD11	2:A:87:LYS:HE3	1.96	0.46
1:C:8:DC:C5	2:B:154:ASP:HB3	2.50	0.46
2:B:38:TRP:CD1	2:B:39:PRO:HD3	2.51	0.46
2:B:189:PRO:HD3	4:B:312:HOH:O	2.16	0.46
1:D:7:DT:C5'	2:A:196:ASP:HB2	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:98:GLU:HB3	2:A:105:LEU:HD11	1.97	0.45
1:D:5:DG:H5'	4:A:551:HOH:O	2.16	0.45
1:D:10:DA:H1'	1:D:11:DT:C5'	2.46	0.45
2:A:115:GLY:HA2	4:A:506:HOH:O	2.17	0.45
2:A:24:GLN:O	2:A:28:GLU:HG3	2.17	0.45
2:A:58:VAL:HB	2:A:59:PRO:HD3	1.99	0.45
2:A:117:ILE:HG12	2:A:161:GLU:HG2	1.98	0.44
2:B:38:TRP:CD1	2:B:39:PRO:CD	3.00	0.44
2:B:43:LYS:HG2	2:B:44:THR:H	1.81	0.44
2:B:3:VAL:HG13	2:B:177:ILE:HG23	1.99	0.44
1:C:3:DT:H6	4:A:622:HOH:O	2.00	0.44
2:A:101:GLU:O	2:A:102:ASN:ND2	2.50	0.44
1:D:8:DC:C5	2:A:154:ASP:HB3	2.52	0.44
2:B:143:MET:O	2:B:178:GLY:HA2	2.18	0.44
2:A:161:GLU:HA	2:B:89:LYS:HE3	2.00	0.44
2:A:110:MET:HG3	2:A:111:GLU:N	2.33	0.43
2:B:38:TRP:CG	2:B:39:PRO:CD	3.01	0.43
2:A:119:SER:HA	2:B:118:SER:HB2	1.99	0.43
2:B:17:SER:HA	4:B:323:HOH:O	2.18	0.43
2:B:33:ILE:HD11	2:B:179:PHE:CD2	2.53	0.43
1:C:4:DG:H1'	1:C:5:DG:O4'	2.19	0.42
2:A:101:GLU:HG3	2:A:102:ASN:HD22	1.85	0.42
1:C:6:DA:N6	4:C:26:HOH:O	2.52	0.42
2:A:110:MET:CG	2:A:111:GLU:N	2.83	0.42
2:A:48:ASN:HA	2:A:185:ASN:O	2.19	0.42
1:C:6:DA:C1'	1:C:7:DT:H5'	2.47	0.42
1:C:11:DT:H1'	1:C:12:DA:H5'	2.01	0.42
2:A:127:LEU:HD13	2:A:139:ALA:HB1	2.01	0.42
2:B:107:ARG:HD3	4:B:298:HOH:O	2.19	0.42
2:B:38:TRP:CB	2:B:39:PRO:CD	2.98	0.41
4:A:573:HOH:O	2:B:146:LYS:HE3	2.19	0.41
1:C:7:DT:O4	2:B:116:ASN:ND2	2.48	0.41
2:B:105:LEU:HD23	4:B:298:HOH:O	2.20	0.41
2:B:117:ILE:HG12	2:B:161:GLU:HG2	2.03	0.41
2:B:19:ASP:HA	2:B:20:LYS:NZ	2.36	0.41
2:B:22:ILE:HD11	2:B:106:LYS:HD2	2.03	0.41
2:A:132:LYS:HB3	2:A:132:LYS:HE2	1.84	0.40
2:A:16:LEU:HD23	2:A:22:ILE:HG22	2.03	0.40
1:D:11:DT:H6	1:D:11:DT:H2'	1.77	0.40
2:A:81:ASP:O	2:A:85:LEU:HB2	2.21	0.40
2:A:167:GLU:C	2:A:169:THR:H	2.25	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:8:ILE:O	2:A:13:LYS:HE3	2.22	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	
2	A	205/213 (96%)	194 (95%)	11 (5%)	0	100	100
2	B	208/213 (98%)	198 (95%)	7 (3%)	3 (1%)	14	6
All	All	413/426 (97%)	392 (95%)	18 (4%)	3 (1%)	26	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	38	TRP
2	B	39	PRO
2	B	40	ALA

### 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	
2	A	188/194 (97%)	179 (95%)	9 (5%)	31	26
2	B	191/194 (98%)	182 (95%)	9 (5%)	32	27

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	379/388 (98%)	361 (95%)	18 (5%)	32	27

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	61	LYS
2	A	95	VAL
2	A	110	MET
2	A	170	GLU
2	A	172	GLN
2	A	174	PHE
2	A	200	LYS
2	A	205	LYS
2	A	207	LYS
2	B	20	LYS
2	B	47	ILE
2	B	101	GLU
2	B	111	GLU
2	B	112	PHE
2	B	132	LYS
2	B	174	PHE
2	B	198	MET
2	B	210	VAL

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

Mol	Chain	Res	Type
2	A	125	ASN
2	A	172	GLN
2	B	121	HIS
2	B	172	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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are 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

### 6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers

EDS was not executed - this section will therefore be empty.