



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 2ODI / pdb\_00002odi  
Title : Restriction Endonuclease BCNI-Cognate DNA Substrate Complex  
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Deposited on : 2006-12-22  
Resolution : 1.45 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

<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>.

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

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

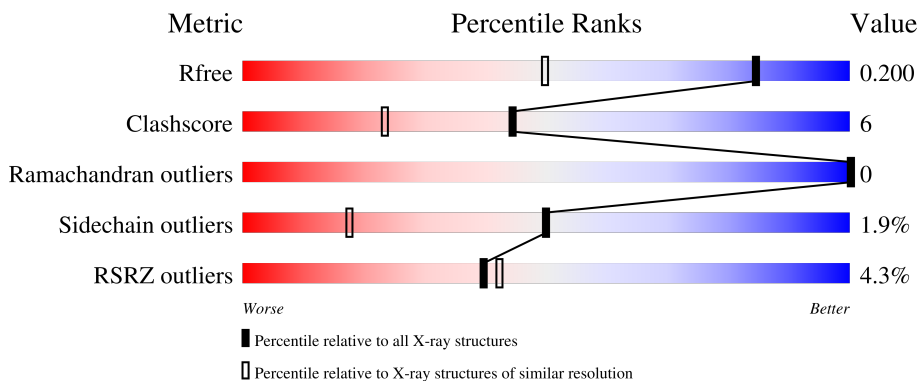
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.45 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.



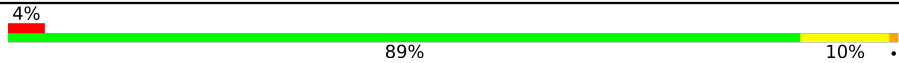
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1756 (1.46-1.46)
Clashscore	190562	1795 (1.46-1.46)
Ramachandran outliers	187476	1776 (1.46-1.46)
Sidechain outliers	187428	1776 (1.46-1.46)
RSRZ outliers	180081	1756 (1.46-1.46)

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  $\geq 3$ , 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	C	11	 82% 18%
1	E	11	 73% 27%
2	D	11	 9% 100%
2	F	11	 9% 82% 18%
3	A	238	 5% 90% 9%

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Mol	Chain	Length	Quality of chain
3	B	238	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a small red segment on the left labeled '4%', a large green segment in the middle labeled '89%', and a small yellow segment on the right labeled '10%'. A small black dot is visible at the far right end of the bar.</p>

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5583 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 5'-D(\*AP\*AP\*CP\*CP\*CP\*GP\*GP\*AP\*GP\*AP\*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	C	11	Total 223	C 106	N 47	O 60	P 10	0	0	0
1	E	11	Total 223	C 106	N 47	O 60	P 10	0	0	0

- Molecule 2 is a DNA chain called 5'-D(\*CP\*TP\*CP\*CP\*GP\*GP\*GP\*TP\*TP\*GP\*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	D	11	Total 222	C 107	N 37	O 68	P 10	0	0	0
2	F	11	Total 222	C 107	N 37	O 68	P 10	0	0	0

- Molecule 3 is a protein called R.BcnI.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
3	A	238	Total 2042	C 1296	N 361	O 375	S 1	Se 9	3	16	0
3	B	238	Total 1996	C 1271	N 346	O 367	S 1	Se 11	0	10	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP Q8RNV8
A	28	MSE	MET	modified residue	UNP Q8RNV8
A	64	MSE	MET	modified residue	UNP Q8RNV8
A	106	MSE	MET	modified residue	UNP Q8RNV8
A	181	MSE	MET	modified residue	UNP Q8RNV8
A	199	MSE	MET	modified residue	UNP Q8RNV8
B	1	MSE	MET	modified residue	UNP Q8RNV8
B	28	MSE	MET	modified residue	UNP Q8RNV8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	64	MSE	MET	modified residue	UNP Q8RNV8
B	106	MSE	MET	modified residue	UNP Q8RNV8
B	181	MSE	MET	modified residue	UNP Q8RNV8
B	199	MSE	MET	modified residue	UNP Q8RNV8

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0
5	B	1	Total Cl 1 1	0	0


- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	37	Total O 37 37	0	1
6	D	39	Total O 39 39	0	0
6	E	32	Total O 32 32	0	0
6	F	45	Total O 45 45	0	1
6	A	261	Total O 261 261	0	4
6	B	236	Total O 236 236	0	1

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

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.

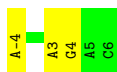
- Molecule 1: 5'-D(\*AP\*AP\*CP\*CP\*CP\*GP\*GP\*AP\*GP\*AP\*C)-3'

Chain C:  82% 18%



- Molecule 1: 5'-D(\*AP\*AP\*CP\*CP\*CP\*GP\*GP\*AP\*GP\*AP\*C)-3'

Chain E:  73% 27%




- Molecule 2: 5'-D(\*CP\*TP\*CP\*CP\*GP\*GP\*GP\*TP\*TP\*GP\*T)-3'

Chain D:  9% 100%



- Molecule 2: 5'-D(\*CP\*TP\*CP\*CP\*GP\*GP\*GP\*TP\*TP\*GP\*T)-3'

Chain F:  9% 82% 18%




- Molecule 3: R.BcnI

Chain A:  5% 90% 9%



- Molecule 3: R.BcnI

Chain B:  4% 89% 10%



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.55Å 87.87Å 79.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.38 – 1.45 28.38 – 1.45	Depositor EDS
% Data completeness (in resolution range)	97.5 (28.38-1.45) 97.5 (28.38-1.45)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.34 (at 1.45Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.174 , 0.199 0.177 , 0.200	Depositor DCC
$R_{free}$ test set	4977 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	11.3	Xtrriage
Anisotropy	0.055	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 35.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.005 for l,-k,h	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5583	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	13.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 56.41 % 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 2.7467e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: CA, CL

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	0.30	0/251	0.56	0/385
1	E	0.29	0/251	0.57	0/385
2	D	0.30	0/247	0.74	0/380
2	F	0.29	0/247	0.76	0/380
3	A	0.45	0/2077	0.65	0/2781
3	B	0.47	0/2031	0.67	0/2720
All	All	0.43	0/5104	0.66	0/7031

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	C	223	0	123	2	0
1	E	223	0	123	3	0
2	D	222	0	127	0	0
2	F	222	0	127	1	0
3	A	2042	0	2094	29	0
3	B	1996	0	2048	23	0
4	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	261	0	0	10	0
6	B	236	0	0	3	0
6	C	37	0	0	1	0
6	D	39	0	0	0	0
6	E	32	0	0	1	0
6	F	45	0	0	0	0
All	All	5583	0	4642	58	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 6.

The worst 5 of 58 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:A:152[B]:LYS:HD2	6:A:829:HOH:O	1.40	1.17
3:A:114:ILE:HD12	3:A:199:MSE:HE3	1.27	1.07
6:C:10:HOH:O	3:A:64[A]:MSE:HG2	1.54	1.06
6:E:81:HOH:O	3:B:64[B]:MSE:HG2	1.53	1.05
3:A:152[B]:LYS:HG2	6:A:845:HOH:O	1.56	1.03

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	
3	A	252/238 (106%)	240 (95%)	12 (5%)	0	100	100
3	B	245/238 (103%)	238 (97%)	7 (3%)	0	100	100
All	All	497/476 (104%)	478 (96%)	19 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	
3	A	230/209 (110%)	223 (97%)	7 (3%)	36	7
3	B	225/209 (108%)	222 (99%)	3 (1%)	61	31
All	All	455/418 (109%)	445 (98%)	10 (2%)	50	13

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	B	1[A]	MSE
3	B	1[B]	MSE
3	B	150	LEU
3	A	127	ASN
3	A	130	HIS

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

Mol	Chain	Res	Type
3	A	84	GLN
3	A	127	ASN
3	A	135	ASN
3	B	135	ASN

### 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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 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 [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	C	11/11 (100%)	0.04	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	8, 9, 16, 17	0
1	E	11/11 (100%)	-0.03	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	8, 8, 16, 16	0
2	D	11/11 (100%)	0.27	1 (9%) <span style="border: 1px solid red; padding: 2px;">15</span> <span style="border: 1px solid red; padding: 2px;">14</span>	8, 10, 17, 21	0
2	F	11/11 (100%)	0.02	1 (9%) <span style="border: 1px solid red; padding: 2px;">15</span> <span style="border: 1px solid red; padding: 2px;">14</span>	8, 9, 14, 19	0
3	A	232/238 (97%)	0.27	11 (4%) <span style="border: 1px solid red; padding: 2px;">36</span> <span style="border: 1px solid red; padding: 2px;">38</span>	5, 11, 19, 28	29 (12%)
3	B	232/238 (97%)	0.27	9 (3%) <span style="border: 1px solid red; padding: 2px;">43</span> <span style="border: 1px solid red; padding: 2px;">46</span>	6, 12, 18, 24	17 (7%)
All	All	508/520 (97%)	0.25	22 (4%) <span style="border: 1px solid red; padding: 2px;">40</span> <span style="border: 1px solid red; padding: 2px;">42</span>	5, 12, 19, 28	46 (9%)

The worst 5 of 22 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	51	ILE	5.9
3	A	52	THR	4.8
3	B	69	ALA	4.1
3	A	131	ALA	3.7
3	B	71[A]	SER	3.5

### 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 oligosaccharides 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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
4	CA	A	703	1/1	0.96	0.26	19,19,19,19	1
5	CL	A	705	1/1	0.99	0.12	18,18,18,18	0
5	CL	B	704	1/1	0.99	0.09	15,15,15,15	0
4	CA	A	701	1/1	1.00	0.07	8,8,8,8	0
4	CA	B	702	1/1	1.00	0.08	7,7,7,7	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.