



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 4DK9 / pdb_00004dk9
Title : Crystal Structure of MBD4 Catalytic Domain Bound to Abasic DNA
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Deposited on : 2012-02-03
Resolution : 2.76 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

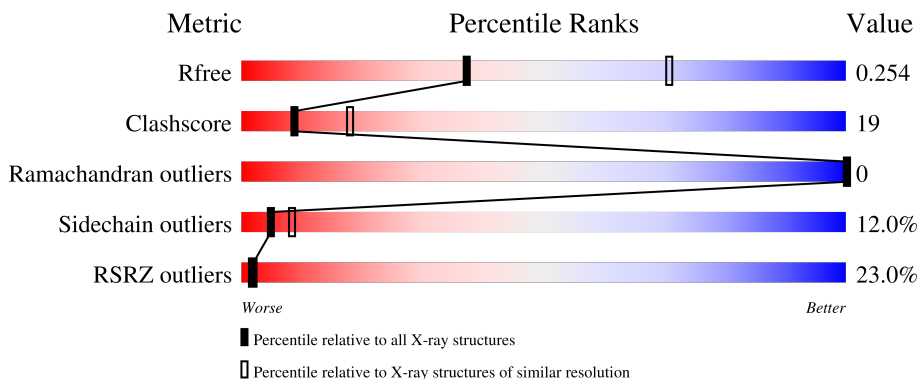
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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 ≥ 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	B	11	27% 27% 64% 9%
2	C	11	73% 18% 9%
3	A	157	22% 59% 25% 6% 11%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 1601 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*GP*AP*CP*GP*TP*GP*GP*AP*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	11	Total	C	N	O	P	0	0	0
			227	108	48	61	10			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	11	Total	C	N	O	P	0	0	0
			210	102	32	66	10			

- Molecule 3 is a protein called Methyl-CpG-binding domain protein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	140	Total	C	N	O	S	0	0	0
			1162	771	189	200	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	424	GLY	-	expression tag	UNP O95243
A	425	SER	-	expression tag	UNP O95243

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	K	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	O	0	0
			1	1		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	41.00Å 55.16Å 122.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	61.10 – 2.76 61.10 – 2.76	Depositor EDS
% Data completeness (in resolution range)	99.8 (61.10-2.76) 99.9 (61.10-2.76)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.234 , 0.259 0.232 , 0.254	Depositor DCC
R_{free} test set	350 reflections (4.62%)	wwPDB-VP
Wilson B-factor (Å ²)	56.6	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 70.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	1601	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.89% 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: K, 3DR

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	B	0.43	0/256	0.86	1/394 (0.3%)
2	C	0.39	0/220	1.08	0/335
3	A	1.08	2/1205 (0.2%)	1.16	9/1647 (0.5%)
All	All	0.94	2/1681 (0.1%)	1.11	10/2376 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	475	ILE	CA-CB	5.87	1.57	1.54
3	A	454	HIS	CG-ND1	-5.64	1.32	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	504	LYS	N-CA-C	-9.87	95.27	109.48
3	A	509	TYR	N-CA-C	6.35	120.23	112.23
3	A	491	ALA	N-CA-C	6.17	118.00	111.28
1	B	2	DA	C2'-C3'-O3'	-6.03	102.46	111.50
3	A	504	LYS	C-N-CD	-5.83	107.78	120.60
3	A	440	PRO	CA-C-N	-5.82	114.60	120.31
3	A	440	PRO	C-N-CA	-5.82	114.60	120.31
3	A	504	LYS	CA-C-N	5.53	140.27	127.00
3	A	504	LYS	C-N-CA	5.53	140.27	127.00
3	A	572	GLU	N-CA-C	5.15	117.63	111.71

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	B	227	0	124	21	0
2	C	210	0	124	4	0
3	A	1162	0	1092	35	0
4	A	1	0	0	0	0
5	B	1	0	0	0	0
All	All	1601	0	1340	54	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 (54) 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:5:DC:H2''	1:B:6:DG:OP1	1.43	1.12
2:C:21:DC:H2''	2:C:22:DT:H5'	1.18	1.12
3:A:438:TRP:HH2	3:A:547:ILE:O	1.44	0.98
1:B:1:DA:H1'	1:B:2:DA:OP1	1.64	0.97
1:B:3:DG:H4'	1:B:3:DG:OP1	1.62	0.95
3:A:447:LEU:HD23	3:A:447:LEU:N	1.88	0.88
3:A:469:THR:HG21	3:A:512:ARG:HH22	1.45	0.82
3:A:447:LEU:HD23	3:A:447:LEU:H	1.43	0.81
3:A:438:TRP:CH2	3:A:547:ILE:O	2.31	0.81
2:C:21:DC:C2'	2:C:22:DT:H5'	2.09	0.80
1:B:3:DG:OP1	1:B:3:DG:C4'	2.29	0.80
3:A:551:ASN:C	3:A:551:ASN:HD22	1.93	0.75
3:A:508:LEU:O	3:A:512:ARG:HB2	1.90	0.71
1:B:1:DA:C1'	1:B:2:DA:OP1	2.39	0.70
1:B:2:DA:H2''	1:B:3:DG:C8	2.26	0.69
1:B:5:DC:C2'	1:B:6:DG:OP1	2.30	0.68
1:B:6:DG:H4'	3:A:509:TYR:CE2	2.28	0.68
2:C:21:DC:H2''	2:C:22:DT:C5'	2.10	0.67
1:B:3:DG:H5''	1:B:3:DG:H8	1.60	0.66
3:A:490:VAL:O	3:A:493:THR:HB	1.96	0.66
3:A:469:THR:CG2	3:A:512:ARG:HH12	2.13	0.61
3:A:447:LEU:N	3:A:447:LEU:CD2	2.62	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:DG:C8	1:B:3:DG:H5''	2.39	0.56
1:B:8:DG:H5'	3:A:505:PRO:HB2	1.89	0.54
3:A:469:THR:HG21	3:A:512:ARG:NH2	2.18	0.53
3:A:442:ARG:HA	3:A:450:GLU:OE1	2.09	0.52
1:B:6:DG:H4'	3:A:509:TYR:HE2	1.72	0.52
1:B:1:DA:H4'	1:B:2:DA:OP2	2.08	0.52
3:A:551:ASN:C	3:A:551:ASN:ND2	2.63	0.52
1:B:3:DG:H2''	1:B:4:DA:O5'	2.11	0.50
1:B:2:DA:OP1	1:B:2:DA:O4'	2.29	0.50
3:A:455:ASP:OD1	3:A:458:LYS:HE3	2.11	0.50
3:A:443:SER:N	3:A:450:GLU:OE1	2.43	0.50
3:A:526:LYS:O	3:A:528:TRP:HD1	1.97	0.47
3:A:447:LEU:HD11	3:A:566:TYR:HB2	1.96	0.47
3:A:455:ASP:CG	3:A:458:LYS:HG3	2.40	0.46
1:B:6:DG:H2''	1:B:7:DT:C7	2.46	0.46
3:A:469:THR:HG21	3:A:512:ARG:HH12	1.79	0.45
3:A:463:THR:HB	3:A:540:TYR:CE2	2.52	0.45
1:B:6:DG:OP2	3:A:510:ASP:HB3	2.18	0.44
3:A:567:HIS:O	3:A:568:ASP:C	2.60	0.44
1:B:5:DC:C2'	3:A:508:LEU:HD23	2.48	0.44
3:A:445:PHE:HB2	3:A:447:LEU:HD21	1.99	0.43
3:A:515:THR:HG23	3:A:535:HIS:CD2	2.53	0.43
3:A:507:GLY:O	3:A:508:LEU:C	2.61	0.43
3:A:562:LYS:HE2	3:A:562:LYS:HB2	1.83	0.43
3:A:532:ILE:HG23	3:A:533:GLU:N	2.34	0.42
1:B:2:DA:H2''	1:B:3:DG:H8	1.79	0.42
1:B:3:DG:H2'	1:B:4:DA:C8	2.54	0.42
3:A:459:LEU:O	3:A:463:THR:HG23	2.19	0.42
1:B:3:DG:C2'	1:B:4:DA:C8	3.03	0.42
2:C:18:3DR:H3'	3:A:466:LEU:O	2.20	0.42
3:A:553:TRP:O	3:A:567:HIS:HE1	2.04	0.41
3:A:485:TYR:HE2	3:A:494:ALA:HB2	1.86	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	
3	A	138/157 (88%)	131 (95%)	7 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	
3	A	117/143 (82%)	103 (88%)	14 (12%)	5	8

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	439	THR
3	A	442	ARG
3	A	447	LEU
3	A	451	THR
3	A	466	LEU
3	A	468	ARG
3	A	469	THR
3	A	489	GLU
3	A	493	THR
3	A	501	GLU
3	A	512	ARG
3	A	550	VAL
3	A	562	LYS
3	A	572	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	446	ASN
3	A	535	HIS
3	A	551	ASN
3	A	557	HIS
3	A	573	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	3DR	C	18	2	8,11,12	0.54	0	7,14,17	1.56	2 (28%)

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	3DR	C	18	2	-	0/3/15/16	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	18	3DR	O4'-C4'-C3'	2.82	107.88	103.73
2	C	18	3DR	C1'-C2'-C3'	2.10	105.51	103.26

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	18	3DR	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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

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	B	11/11 (100%)	1.56	3 (27%) 1 1	106, 119, 185, 200	0
2	C	10/11 (90%)	0.55	0 100 100	94, 105, 111, 138	0
3	A	140/157 (89%)	1.37	34 (24%) 2 1	74, 96, 129, 150	0
All	All	161/179 (89%)	1.33	37 (22%) 2 2	74, 98, 135, 200	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	577	LEU	6.9
3	A	575	GLU	5.0
3	A	506	LEU	4.2
3	A	509	TYR	4.1
3	A	446	ASN	3.8
3	A	504	LYS	3.7
3	A	488	ALA	3.6
3	A	438	TRP	3.5
3	A	469	THR	3.3
1	B	1	DA	3.2
3	A	576	LYS	3.1
3	A	505	PRO	3.0
3	A	439	THR	3.0
3	A	465	PHE	2.9
3	A	559	GLU	2.9
3	A	497	ARG	2.8
3	A	468	ARG	2.8
3	A	511	LEU	2.7
3	A	451	THR	2.5
3	A	467	ASN	2.5
3	A	484	LYS	2.5
1	B	6	DG	2.5
3	A	442	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
3	A	495	ASP	2.4
3	A	474	ALA	2.4
3	A	494	ALA	2.4
3	A	567	HIS	2.3
3	A	487	SER	2.3
3	A	510	ASP	2.3
3	A	496	TRP	2.3
3	A	503	LEU	2.2
3	A	489	GLU	2.2
3	A	447	LEU	2.2
1	B	5	DC	2.1
3	A	493	THR	2.1
3	A	472	LYS	2.1
3	A	508	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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, 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($\text{\AA}^2$)	Q<0.9
2	3DR	C	18	11/12	0.98	0.08	88,91,96,98	0

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, 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($\text{\AA}^2$)	Q<0.9
4	K	A	601	1/1	0.92	0.22	119,119,119,119	0

6.5 Other polymers [i](#)

There are no such residues in this entry.