Full wwPDB X-ray Structure Validation Report (i)

Jun 19, 2024 – 12:58 PM EDT

PDB ID	:	4OKY
Title	:	Crystal structure of PvuRts1I, a 5-hydroxymethylcytosine DNA restriction endonuclease
Authors	:	Wang, C.L.; Shao, C.; Zang, J.Y.
Deposited on	:	2014-01-23
Resolution	:	2.90 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 (i) 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 (i)) were used in the production of this report:

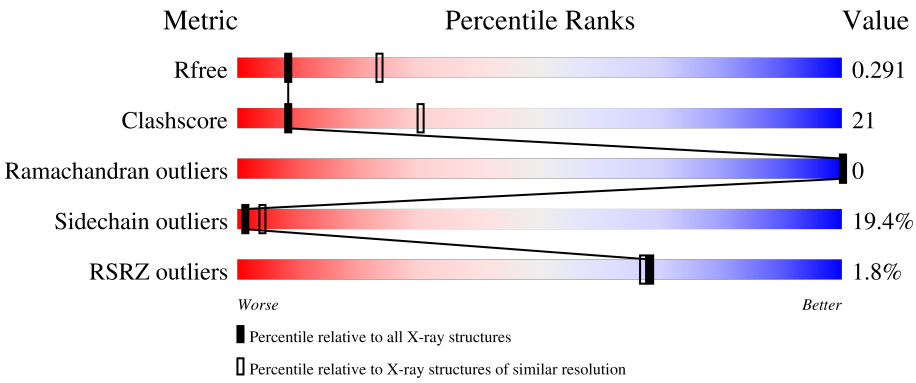
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 <=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	293	2% 55% 33% 8% •

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2184 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 Restriction endonuclease PvuRts1 I.

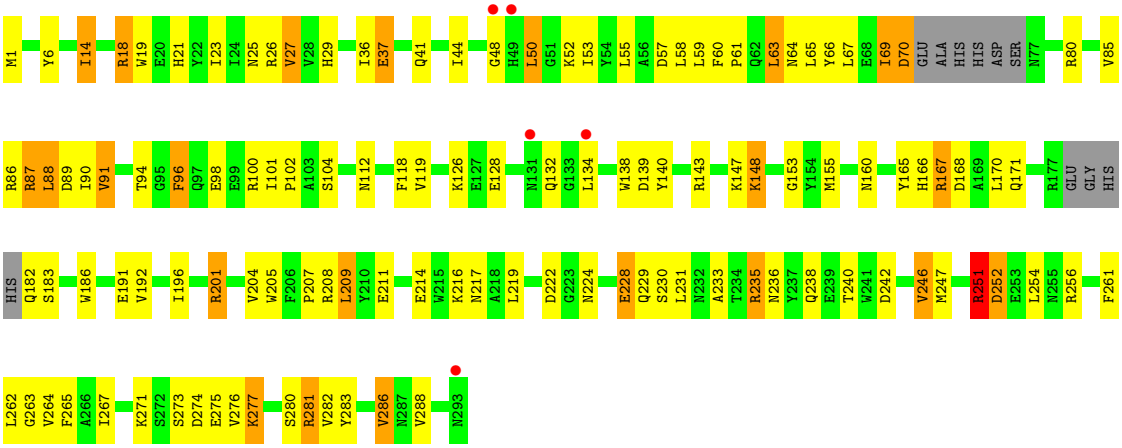
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	C	283	2184	1386	380	411	3	4	0	0	0

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: Restriction endonuclease PvuRts1 I

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	160.21Å 160.21Å 45.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.71 – 2.90 50.66 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.6 (56.71-2.90) 97.7 (50.66-2.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.84 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.252 , 0.287 0.263 , 0.291	Depositor DCC
R_{free} test set	662 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	61.6	Xtriage
Anisotropy	0.228	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 62.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	2184	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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	0.55	1/2225 (0.0%)	0.83	6/3021 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	160	ASN	N-CA	5.10	1.56	1.46

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	208	ARG	CB-CA-C	7.63	125.66	110.40
1	C	208	ARG	N-CA-C	-6.77	92.73	111.00
1	C	160	ASN	CB-CA-C	6.64	123.68	110.40
1	C	251	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	C	209	LEU	N-CA-CB	5.98	122.37	110.40
1	C	286	VAL	CB-CA-C	-5.16	101.61	111.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	2184	0	2027	88	0
All	All	2184	0	2027	88	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 21.

All (88) 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:166:HIS:N	1:C:167:ARG:HH12	1.64	0.95
1:C:167:ARG:HH11	1:C:167:ARG:H	1.11	0.95
1:C:23:ILE:O	1:C:27:VAL:HG12	1.70	0.90
1:C:167:ARG:O	1:C:171:GLN:HG3	1.76	0.85
1:C:166:HIS:N	1:C:167:ARG:NH1	2.26	0.83
1:C:155:MSE:SE	1:C:247:MSE:HE1	2.31	0.80
1:C:273:SER:O	1:C:277:LYS:HE3	1.82	0.79
1:C:166:HIS:H	1:C:167:ARG:HH12	1.28	0.79
1:C:170:LEU:CD2	1:C:204:VAL:HG21	2.16	0.76
1:C:64:ASN:HA	1:C:96:PHE:CE1	2.21	0.75
1:C:167:ARG:HH11	1:C:167:ARG:N	1.87	0.71
1:C:251:ARG:HG2	1:C:251:ARG:HH11	1.55	0.69
1:C:170:LEU:HD23	1:C:204:VAL:HG21	1.76	0.67
1:C:230:SER:OG	1:C:236:ASN:ND2	2.28	0.66
1:C:18:ARG:HH11	1:C:18:ARG:HG3	1.61	0.65
1:C:60:PHE:O	1:C:63:LEU:O	2.16	0.64
1:C:165:TYR:HB3	1:C:167:ARG:NH1	2.13	0.64
1:C:48:GLY:O	1:C:50:LEU:HG	1.99	0.63
1:C:170:LEU:CD1	1:C:182:GLN:CB	2.76	0.63
1:C:1:MSE:HE3	1:C:6:TYR:HB2	1.79	0.63
1:C:66:TYR:OH	1:C:98:GLU:HB2	1.98	0.63
1:C:64:ASN:HB3	1:C:96:PHE:HA	1.81	0.63
1:C:66:TYR:CZ	1:C:98:GLU:HB2	2.33	0.62
1:C:217:ASN:OD1	1:C:228:GLU:HB2	1.99	0.62
1:C:251:ARG:HH11	1:C:251:ARG:CG	2.12	0.62
1:C:70:ASP:OD1	1:C:70:ASP:N	2.30	0.61
1:C:143:ARG:O	1:C:148:LYS:NZ	2.30	0.59
1:C:166:HIS:HB2	1:C:167:ARG:NH1	2.19	0.58
1:C:14:ILE:O	1:C:21:HIS:HB2	2.03	0.57
1:C:170:LEU:HD11	1:C:182:GLN:CB	2.36	0.56
1:C:233:ALA:HA	1:C:236:ASN:ND2	2.21	0.55
1:C:85:VAL:HA	1:C:90:ILE:HD11	1.89	0.55
1:C:41:GLN:OE1	1:C:80:ARG:NH1	2.41	0.54
1:C:252:ASP:HB2	1:C:256:ARG:H	1.71	0.54
1:C:282:VAL:HG12	1:C:283:TYR:CD2	2.43	0.54
1:C:214:GLU:HB2	1:C:235:ARG:HH11	1.72	0.54
1:C:19:TRP:O	1:C:23:ILE:HG13	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:ARG:CD	1:C:112:ASN:HB2	2.37	0.53
1:C:273:SER:OG	1:C:274:ASP:N	2.42	0.53
1:C:37:GLU:OE2	1:C:61:PRO:HG2	2.09	0.53
1:C:252:ASP:HB3	1:C:254:LEU:H	1.74	0.52
1:C:153:GLY:C	1:C:288:VAL:HG13	2.29	0.52
1:C:165:TYR:HB3	1:C:167:ARG:CZ	2.40	0.51
1:C:55:LEU:C	1:C:55:LEU:HD13	2.32	0.51
1:C:59:LEU:HD12	1:C:66:TYR:HB3	1.91	0.51
1:C:90:ILE:HG22	1:C:91:VAL:N	2.26	0.50
1:C:37:GLU:HG2	1:C:138:TRP:CE2	2.47	0.50
1:C:87:ARG:NH1	1:C:89:ASP:OD2	2.42	0.50
1:C:25:ASN:O	1:C:29:HIS:HD2	1.95	0.49
1:C:140:TYR:O	1:C:143:ARG:HB2	2.12	0.49
1:C:236:ASN:HD22	1:C:236:ASN:N	2.10	0.49
1:C:64:ASN:HA	1:C:96:PHE:CD1	2.48	0.48
1:C:166:HIS:CA	1:C:167:ARG:NH1	2.76	0.48
1:C:229:GLN:HB2	1:C:276:VAL:HG22	1.96	0.48
1:C:25:ASN:O	1:C:29:HIS:CD2	2.67	0.48
1:C:50:LEU:HD13	1:C:52:LYS:HB2	1.96	0.48
1:C:224:ASN:OD1	1:C:281:ARG:NH1	2.47	0.47
1:C:246:VAL:HG22	1:C:262:LEU:HB2	1.96	0.47
1:C:18:ARG:HG3	1:C:18:ARG:NH1	2.30	0.47
1:C:261:PHE:CZ	1:C:286:VAL:HG23	2.50	0.47
1:C:201:ARG:N	1:C:242:ASP:OD2	2.44	0.47
1:C:205:TRP:C	1:C:207:PRO:HD3	2.34	0.47
1:C:87:ARG:HB3	1:C:251:ARG:HB2	1.98	0.46
1:C:201:ARG:NH1	1:C:240:THR:OG1	2.48	0.46
1:C:26:ARG:NE	1:C:112:ASN:HB2	2.31	0.46
1:C:67:LEU:HD13	1:C:118:PHE:CD2	2.50	0.46
1:C:165:TYR:O	1:C:168:ASP:HB2	2.15	0.45
1:C:87:ARG:O	1:C:88:LEU:C	2.55	0.45
1:C:211:GLU:HA	1:C:216:LYS:HA	1.99	0.45
1:C:27:VAL:HG23	1:C:119:VAL:HG22	1.98	0.44
1:C:138:TRP:HD1	1:C:139:ASP:N	2.15	0.44
1:C:170:LEU:HD21	1:C:204:VAL:HG21	1.99	0.44
1:C:58:LEU:HD23	1:C:58:LEU:HA	1.92	0.43
1:C:36:ILE:HD11	1:C:126:LYS:HD2	2.00	0.43
1:C:166:HIS:CB	1:C:167:ARG:NH1	2.80	0.42
1:C:222:ASP:OD1	1:C:224:ASN:ND2	2.52	0.42
1:C:132:GLN:HB2	1:C:134:LEU:HD12	2.01	0.42
1:C:57:ASP:OD2	1:C:69:ILE:HG12	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:138:TRP:CD1	1:C:139:ASP:N	2.88	0.41
1:C:261:PHE:CZ	1:C:263:GLY:HA2	2.55	0.41
1:C:246:VAL:HG13	1:C:265:PHE:CD1	2.55	0.41
1:C:170:LEU:HD22	1:C:186:TRP:CG	2.56	0.41
1:C:90:ILE:CG2	1:C:91:VAL:N	2.83	0.41
1:C:118:PHE:CD1	1:C:118:PHE:C	2.93	0.41
1:C:264:VAL:CG2	1:C:286:VAL:HG22	2.50	0.41
1:C:101:ILE:HA	1:C:102:PRO:HD3	1.97	0.41
1:C:205:TRP:O	1:C:207:PRO:HD2	2.21	0.40
1:C:267:ILE:HG23	1:C:267:ILE:O	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
1	C	277/293 (94%)	261 (94%)	16 (6%)	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
1	C	216/254 (85%)	174 (81%)	42 (19%)	1 4

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	14	ILE
1	C	18	ARG
1	C	27	VAL
1	C	37	GLU
1	C	44	ILE
1	C	50	LEU
1	C	53	ILE
1	C	63	LEU
1	C	65	LEU
1	C	69	ILE
1	C	70	ASP
1	C	86	ARG
1	C	87	ARG
1	C	88	LEU
1	C	91	VAL
1	C	94	THR
1	C	96	PHE
1	C	100	ARG
1	C	104	SER
1	C	128	GLU
1	C	147	LYS
1	C	148	LYS
1	C	167	ARG
1	C	183	SER
1	C	191	GLU
1	C	192	VAL
1	C	196	ILE
1	C	201	ARG
1	C	209	LEU
1	C	219	LEU
1	C	228	GLU
1	C	231	LEU
1	C	235	ARG
1	C	238	GLN
1	C	246	VAL
1	C	251	ARG
1	C	252	ASP
1	C	271	LYS
1	C	275	GLU
1	C	277	LYS
1	C	280	SER
1	C	281	ARG

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

Mol	Chain	Res	Type
1	C	62	GLN
1	C	112	ASN
1	C	229	GLN
1	C	236	ASN
1	C	238	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)

There are no ligands in this entry.

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, 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	C	279/293 (95%)	0.12	5 (1%) 68 67	26, 58, 88, 110	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	48	GLY	7.3
1	C	293	ASN	4.3
1	C	49	HIS	4.1
1	C	134	LEU	2.1
1	C	131	ASN	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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.